Proceedings of the 2019 miniKanren and Relational Programming Workshop

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Accessibility
Proceedings of the 2019 miniKanren and Relational Programming Workshop

William E. Byrd
and
Nada Amin

TR-02-19

Computer Science Group
Harvard University
Cambridge, Massachusetts
Preface

This report aggregates the papers presented at the first miniKanren and Relational Programming Workshop, hosted on August 2nd, 2019 in Berlin, Germany and co-located with the twenty-second International Conference on Functional Programming.

The miniKanren and Relational Programming Workshop is a new workshop for the miniKanren family of relational (pure constraint logic programming) languages: miniKanren, microKanren, core.logic, OCanren, Guanxi, etc. The workshop solicits papers and talks on the design, implementation, and application of miniKanren-like languages. A major goal of the workshop is to bring together researchers, implementors, and users from the miniKanren community, and to share expertise and techniques for relational programming. Another goal for the workshop is to push the state of the art of relational programming—for example, by developing new techniques for writing interpreters, type inferencers, theorem provers, abstract interpreters, CAD tools, and other interesting programs as relations, which are capable of being “run backwards,” performing synthesis, etc.

6 papers were submitted to the workshop, and each submission was reviewed by two to three members of the program committee. After deliberation, all submissions were accepted to the workshop.

In addition to the six full papers presented

- William E. Byrd gave a morning tutorial on miniKanren,
- Daniel P. Friedman and William E. Byrd gave a closing Q&A with audience.

Thanks to all presenters, participants, and members of the program committee.

William E. Byrd & Nada Amin

Program Committee

Claire Alvis, Sparkfund
Nada Amin, Harvard University (Program Chair)
Tom Gilray, University of Alabama at Birmingham
Jason Hemann, Northeastern University
Eric Holk, Google
Kanae Tsushima, National Institute of Informatics
William E. Byrd, University of Alabama at Birmingham (General Chair)
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Towards a miniKanren with fair search strategies

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We describe fairness levels in disjunction and conjunction implementations. Specifically, a disjunction implementation can be fair, almost-fair, or unfair. And a conjunction implementation can be fair or unfair. We compare the fairness level of four search strategies: the standard miniKanren interleaving depth-first search, the balanced interleaving depth-first search, the fair depth-first search, and the standard breadth-first search. The two non-standard depth-first searches are new. And we present a new, more efficient and shorter implementation of the standard breadth-first search. Using quantitative evaluation, we argue that each depth-first search is a competitive alternative to the standard one, and that our improved breadth-first search implementation is more efficient than the current one.

1 INTRODUCTION

miniKanren is a family of relational programming languages. Friedman et al. [3, 4] introduce miniKanren and its implementation in The Reasoned Schemer and The Reasoned Schemer, 2nd Ed (TRS2). Hemann et al. [5] describe microKanren, a minimal core of miniKanren comprised of only 54 LOC, miniKanren has been implemented in many other languages, including multiple ones using the same language (e.g. OCaml[7]). As demonstrated in Byrd et al. [2], miniKanren can be used to naturally express difficult computations, such as using an interpreter to perform example-based program synthesis, or using a proof checker as a theorem prover. The papers, talks, and tutorials on miniKanren.org present many other unusual problems, and their solutions in miniKanren.

A subtlety arises when a cond* contains many clauses: not every clause has an equal chance of contributing to the result. As an example, consider the following relation repeat^0 and its invocation.

```scheme
(defrel (repeat^0 x out)
  (cond* 
    ((≡ (,x) out))
    ((fresh (res)
      (≡ (,x ,res) out)
      (repeat^0 x res))))))

> (run 4 q
  (repeat^0 '* q))
'((*) (* *) (* * *) (* * * *))
```

Next, consider the following disjunction of invoking repeat^0 with four different letters.

---

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miniKanren.org/workshop/2019/8-ART1
cond$^c$ intuitively relates its clauses with logical or. And thus an unsuspicious beginner would expect each letter to contribute equally to the result, as follows.

$$
((a) (b) (c) (d) \\
(a a) (b b) (c c) (d d) \\
(a a a) (b b b) (c c c) (d d d))
$$

The cond$^c$ in TRS2, however, generates a less expected result.

$$
((a) (a a) (b) (a a a) \\
(a a a a) (b b) \\
(a a a a a a) (d))
$$

The miniKanren in TRS2 implements interleaving DFS (DFS$^i$), the cause of this unexpected result. With this search strategy, each cond$^c$ clause takes half of its received computational resources and passes the other half to its following clauses, except for the last clause that takes all resources it receives. In the example above, the $a$ clause takes half of all resources. And the $b$ clause takes a quarter. Thus $c$ and $d$ barely contribute to the result.

DFS$^i$ is sometimes powerful for an expert. By carefully organizing the order of cond$^c$ clauses, a miniKanren program can explore more “interesting” clauses than those uninteresting ones, and thus use computational resources efficiently.

DFS$^i$ is not always the best choice. For instance, it might be less desirable for novice miniKanren users—understanding implementation details and fiddling with clause order is not their first priority. There is another reason that miniKanren could use more search strategies than just DFS$^i$. In many applications, there does not exist one order that serves all purposes. For example, a relational dependent type checker contains clauses for constructors that build data and clauses for eliminators that use data. When the type checker is generating simple and shallow programs, the clauses for constructors had better be at the top of the cond$^c$ expression. When performing proof searches for complicated programs, the clauses for eliminators had better be at the top of the cond$^c$ expression. With DFS$^i$, these two uses cannot be efficient at the same time. In fact, to make one use efficient, the other one must be more sluggish. Boskin et al. [1] propose and implement a means to eliminate or re-order disjunctive clauses to accommodate varying search needs such as these.

The specification that gives every clause in the same cond$^c$ equal “search priority” is fair disj. And search strategies with almost-fair disj give every clause similar priority. Fair conj, a related concept, is more subtle. We cover it in the next section.

Our research compares four search strategies with different features in fairness (Table 1). To summarize our contributions, we

- propose and implement **balanced interleaving depth-first search (DFS$^b_i$)**
- propose and implement **fair depth-first search (DFS$^f$)**
- implement in a new way the standard breath-first search. We refer to BFS$_{ser}$ as the original implementation by Seres et al. [9] and BFS$_{imp}$ as our new one. When we use BFS without subscripts, we mean both BFS$_{ser}$
and BFS$imp$. We formally prove that the two implementations are semantically equivalent, however, BFS$imp$ runs faster in all benchmarks and is shorter.

Source code of implementations, examples, benchmarks, and formal proofs are available at the following URL:
https://github.com/LuKC1024/Towards-a-miniKanren-with-fair-search-strategies

<table>
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<tr>
<th>Search Strategies</th>
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<td>BFS</td>
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Table 1. Fairness of all search strategies

2 SEARCH STRATEGIES AND FAIRNESS
In this section, we define fairness levels in disjunction and conjunction implementations. Specifically, a disjunction implementation can be fair, almost-fair, or unfair. And a conjunction implementation can be fair or unfair. Fairness, intuitively, measures how evenly a search strategy allocates computational resource to “sibling” spaces.

Before going further into fairness, we give a short review of the terms: state, space, and goal. A state is a collection of constraints. (Here, we restrict constraints to unification constraints.) Every answer corresponds to a state. A space is a collection of states. And a goal is a function from a state to a space. Every state in the output space includes the input state and possibly more constraints.

Now we elaborate fairness by running more queries about repeat$^*$. We never use run$^*$ here because fairness is more interesting when we ask for a bounded number of answers. It is perfectly fine, however, to use run$^*$ with any search strategy.

2.1 Fair disj
Given the following program, it is natural to expect lists of each letter to constitute 1/4 in the answer list. DFS$si$, TRS2’s search strategy, however, results in many more lists of a than lists of other letters. And some letters (e.g. c and d) are rarely seen. The more clauses, the worse the situation.

```plaintext
;; DFS$si$ (unfair disj)
> (run 12 q
  (cond
    ((repeat o 'a q))
    ((repeat o 'b q))
    ((repeat o 'c q))
    ((repeat o 'd q)))
  '((a) (a a) (b) (a a a)
    (a a a a) (b b)
    (a a a a a) (c)
    (a a a a a a) (b b b)
    (a a a a a a a) (d))
```
Under the hood, the cond here allocates computational resources to four trivially different spaces. The unfair disj in DFSi allocates many more resources to the first space. On the contrary, fair disj would allocate resources evenly to each space.

;; DFS (fair disj)
> (run 12 q
  (cond^'
    ((repeat° 'a q))
    ((repeat° 'b q))
    ((repeat° 'c q))
    ((repeat° 'd q))))
'(a a a) (b b b) (c c c) (d d d)

;; BFS (fair disj)
> (run 12 q
  (cond^'
    ((repeat° 'a q))
    ((repeat° 'b q))
    ((repeat° 'c q))
    ((repeat° 'd q))))
'(a a a) (b b b) (c c c) (d d d)

Running the same program again with almost-fair disj (e.g. DFSbi) gives a similar result, where b and c are swapped. Almost-fair, however, is not completely fair, as shown by the following example.

;; DFSbi (almost-fair disj)
> (run 16 q
  (cond^'
    ((repeat° 'a q))
    ((repeat° 'b q))
    ((repeat° 'c q))
    ((repeat° 'd q))
    ((repeat° 'e q))))
'(a a a a) (b b b b) (c c c c) (d d d d)

DFSbi is fair only when the number of goals is a power of 2, otherwise, it allocates some goals with twice as many resources as the others. In the above example, where the cond has five clauses, DFSbi allocates more resources to the clauses of a, b, and c.

We end this subsection with precise definitions of all levels of disj fairness. Our definition of fair disj is slightly more general than the one in Seres et al. [9], which is only for binary disjunction. We generalize it to a multi-arity one.

**Definition 2.1 (fair disj).** A disj is fair if and only if it allocates computational resources evenly to spaces produced by goals in the same disjunction (i.e., clauses in the same cond).

**Definition 2.2 (almost-fair disj).** A disj is almost-fair if and only if it allocates computational resources so evenly to spaces produced by goals in the same disjunction that the maximal ratio of resources is bounded by a constant.

**Definition 2.3 (unfair disj).** A disj is unfair if and only if it is not almost-fair.
2.2 Fair conj

Given the following program, it is natural to expect lists of each letter to constitute 1/4 in the answer list. Search strategies with unfair conj: DFS$_i$, DFS$_bi$, and DFS$_f$, however, results in many more lists of a than lists of other letters. And some letters are rarely seen. Here again, as the number of clauses grows, the situation worsens.

Although some strategies have a different level of fairness in disj, they have the same behavior when there is no call to a relational definition in cond clauses, including this case.

Under the hood, the cond and the call to repeat are connected by conj. The cond goal outputs a space including four trivially different states. Applying the next conjunctive goal, (repeat o x q), produces four trivially different spaces. In the examples above, all search strategies allocate more computational resources to the space of a. On the contrary, fair conj would allocate resources evenly to each space. For example,

A more interesting situation is when the first conjunct produces an unbounded number of states. Consider the following example: a naive specification of fair conj might require search strategies to produce all sorts of singleton lists, but there would not be any lists of length two or longer, which makes the strategies incomplete. A search strategy is complete if and only if "every correct answer would be discovered after some finite time" [9], otherwise, it is incomplete. In the context of miniKanren, a search strategy is complete means that every correct answer has a position in large enough answer lists.
Naively fair conj

\[
\text{run 6 q}
\]

\[
\text{fresh } (xs)
\]

\[
\text{cond'}
\]

\[
((\text{repeat}^n \ 'a \ xs))
\]

\[
((\text{repeat}^n \ 'b \ xs)))
\]

\[
(\text{repeat}^n \ xs \ q)))
\]

\[
'(((a)) \ ((b))
\]

\[
(((a \ a)) \ ((b \ b))
\]

\[
(((a \ a \ a)) \ ((b \ b \ b)))
\]

Our solution requires a search strategy with fair conj to organize states in buckets in spaces, where each bucket is a finite collection of states and every space contains possibly infinite buckets, and to allocate resources evenly among spaces derived from states in the same bucket. It is up to a search strategy designer to decide by what criteria to put states in the same bucket, and how to allocate resources among spaces related to different buckets.

BFS puts states of the same cost in the same bucket, and allocates resources carefully among spaces related to different buckets such that it produces answers in increasing order of cost. The cost of an answer is its depth in the search tree (i.e., the number of calls to relational definitions required to find the answer) [9]. In the above examples, the cost of each answer is equal to their lengths because we need to apply \(\text{repeat}^n\) \(n\) times to find an answer of length \(n\). In the following example, every answer is a list of a list of symbols, where inner lists in the same outer list are identical. Here the cost of each answer is equal to the length of its inner lists plus the length of its outer list. For example, the cost of \(((a) \ (a))\) is \(1 + 2 = 3\).

Naively fair conj

\[
\text{run 12 q}
\]

\[
\text{fresh } (xs)
\]

\[
\text{cond'}
\]

\[
((\text{repeat}^n \ 'a \ xs))
\]

\[
((\text{repeat}^n \ 'b \ xs)))
\]

\[
(\text{repeat}^n \ xs \ q)))
\]

\[
'(((a)) \ ((b))
\]

\[
(((a \ a)) \ ((b \ b))
\]

\[
(((a \ a \ a)) \ ((b \ b \ b)))
\]

We end this subsection with precise definitions of all levels of conj fairness.

**Definition 2.4 (Fair Conj).** A conj is fair if and only if it allocates computational resources evenly to spaces produced from states in the same bucket. A bucket is a finite collection of states. And search strategies with fair conj should represent spaces with possibly unbounded collections of buckets.

**Definition 2.5 (Unfair Conj).** A conj is unfair if and only if it is not fair.
3 INTERLEAVING DEPTH-FIRST SEARCH

In this section, we review the implementation of DFS₁. We focus on parts that are relevant to other strategies. TRS2, chapter 10 and the appendix, “Connecting the wires”, provide a comprehensive description of the miniKanren implementation but limited to unification constraints (≡). Fig. 1 and Fig. 2 show parts that are later compared with other search strategies. We follow some conventions to name variables: ss name states; gs (possibly with subscript) name goals; variables ending with ∞ name spaces. Fig. 1 shows the implementation of disj. The first function, disj₂, implements binary disjunction. It applies the two disjunctive goals to the input state s and composes the two resulting spaces with append∞. The following syntax definitions say disj is right-associative. Fig. 2 shows the implementation of conj. The first function, conj₂, implements binary conjunction. It applies the first goal to the input state, then applies the second goal to states in the resulting space. The helper function append-map∞ applies its input goal to states in its input space and composes the resulting spaces. It reuses append∞ for space composition. The following syntax definitions say conj is also right-associative.
4 BALANCED INTERLEAVING DEPTH-FIRST SEARCH

DFSₙₚ has almost-fair disj and unfair conj. Its implementation differs from DFSᵢ’s in the disj macro. When there are one or more disjunctive goals, the new disj builds a balanced binary tree whose leaves are the goals and whose nodes are disj₂s, hence the name of this search strategy. In contrast, the disj in DFSᵢ constructs the binary tree in a particularly unbalanced form. We list the new disj with its helper in Fig. 3. The new helper, disj+, takes two additional ‘arguments’. They accumulate goals to be put in the left and right subtrees. The first clause handles the case where there is only one goal. In this case, the tree is the goal itself. When there are more goals, we partition the list of goals into two sublists of roughly equal lengths and recur on the two sublists. We move goals to the accumulators in the last clause. As we are moving two goals each time, there are two base cases: (1) no goal remains; (2) one goal remains. We handle these two new base cases in the second clause and the third clause, respectively.

5 FAIR DEPTH-FIRST SEARCH

DFSᵢ has fair disj and unfair conj. Its implementation differs from DFSᵢ’s in disj₂ (Fig. 4). The new disj₂ calls a new and fair version of append∞. append∞fair immediately calls its helper, loop, with the first argument, s?,

---

Fig. 2. implementation of DFSᵢ (Part II)
(define-syntax disj
  (syntax-rules ()
    ((disj) fail)
    ((disj ... stop recursion when one of the input spaces is obviously finite. The third clause, as we mentioned above, is just for

initialize to #t, which indicates that we haven’t swapped s\^\infty and t\^\infty. The swapping happens at the third cond clause in loop, where s? is updated accordingly. The first two cond clauses essentially copy the cars and stop recursion when one of the input spaces is obviously finite. The third clause, as we mentioned above, is just for
swapping. When the fourth and last clause runs, we know that both $s^\infty$ and $t^\infty$ end with thunks, and that we have swapped them. In this case, we construct a new thunk. The new thunk swaps back the two spaces in the recursive call to loop. This is unnecessary for fairness—we do it to produce answers in a more readable order.

6 BREADTH-FIRST SEARCH

BFS has both fair disj and fair conj. Our first BFS implementation (Fig. 5) serves as a "stepping-stone" toward BFS\textsubscript{imp}. It is so similar to DFS\textsubscript{f} (not DFS\textsubscript{i}) that we only need to apply two changes: (1) rename append-map\textsubscript{\infty} to append-map\textsubscript{\infty}f and (2) replace append\textsubscript{\infty} with append\textsubscript{\infty}f in append-map\textsubscript{\infty}f's body.

This implementation can be improved in two ways. First, as mentioned in subsection 2.2, BFS puts answers in buckets and answers of the same cost are in the same bucket. In the above implementation, however, it is not obvious how we manage cost information—the cars of a space have cost 0 (i.e., they are all in the same bucket), and every thunk indicates an increment in cost. It is even more subtle that append\textsubscript{\infty}f and the append-map\textsubscript{\infty}f respects the cost information. Second, append\textsubscript{\infty}f is extravagant in memory usage. It makes $O(n + m)$ new cons cells every time, where $n$ and $m$ are the sizes of the first buckets of two input spaces. DFS\textsubscript{f} is also space extravagant.

In the following paragraphs, we first describe BFS\textsubscript{imp} implementation that manages cost information in a more clear and concise way and is less extravagant in memory usage. Then we compare BFS\textsubscript{imp} with BFS\textsubscript{ser}.

We simplify the cost information by changing the Space type, modifying related function definitions, and introducing a few more functions. The new type of Space is a pair whose car is a list of answers (the bucket), and whose cdr is either #f or a thunk returning a space. The #f here means the space is obviously finite, just like empty list in other implementations. We list functions related to the pure subset in Fig. 6. The first three functions are space constructors. none makes an empty space; unit makes a space from one answer; and step makes a space from a thunk. The remaining functions are as before. Luckily, the change in append\textsubscript{\infty}f also fixes the miserable space extravagance—the use of append helps us to reuse the first bucket of $t^\infty$. Functions implementing impure features are in Fig. 7. The first function, elim, takes a space $s^\infty$ and two continuations $fk$ and $sk$. When $s^\infty$ contains no answers, it calls $fk$. Otherwise, it calls $sk$ with the first answer and the rest of the space. This function is similar to an eliminator of spaces, hence the name. The remaining functions are as before.

Kiselyov et al. [6] have demonstrated that a MonadPlus hides in implementations of logic programming systems. BFS\textsubscript{imp} is not an exception: append-map\textsubscript{\infty}f is like bind, but takes arguments in reversed order; none, unit, and append\textsubscript{\infty}f correspond to mzero, unit, and mplus, respectively.
Towards a miniKanren with fair search strategies

#| → Space |#
(define (none)
  (() . #f))

#| State → Space |#
(define (unit s)
  ((,s) . #f))

#| (→ Space) → Space |#
(define (step f)
  (() . ,f))

#| Space × Space → Space |#
(define (append\_fair s t)
  (cons (append (car s) (car t))
       (let ((t\_1 (cdr s)) (t\_2 (cdr t)))
          (cond
           (not t\_1) t\_2
           (not t\_2) t\_1
           (else (lambda () (append\_fair (t\_1) (t\_2))))))))

#| Goal × Space → Space |#
(define (append\_map\_fair g s)
  (foldr
   (lambda (s t)
    (append\_fair (g s) t))
   (let ((f (cdr s))
         (step (and f (lambda () (append\_map\_fair g (f))))))
    (car s))))

#| Maybe Nat × Space → [State] |#
(define (take\_n s)
  (let loop ((n n) (vs (car s)))
    (cond
     ((and n (zero? n)) '())
     ((pair? vs)
      (cons (car vs)
        (loop (and n (sub1 n)) (cdr vs))))
     (else (let ((f (cdr s))
                  (if f (take\_n n (f)) '()))))))

Fig. 6. New and changed functions in BFS\_imp that implements pure features
Kuang-Chen Lu, Weixi Ma, and Daniel P. Friedman

\[ \text{Space} \times (\text{State} \times \text{Space} \to \text{Space}) \times (\to \text{Space}) \to \text{Space} \]

(\textit{define} (elim s\textsuperscript{∞} fk sk)
  (let ((ss (car s\textsuperscript{∞}))(f (cdr s\textsuperscript{∞})))
    (cond
     ((pair? ss) (sk (car ss) (cons (cdr ss) f)))
     (f (step (lambda () (elim (f) fk sk))))
     (else (fk)))
  ))

\[ \text{Goal} \times \text{Goal} \times \text{Goal} \to \text{Goal} \]

(\textit{define} (ifte g\textsubscript{1} g\textsubscript{2} g\textsubscript{3})
  (lambda (s)
    (elim (g\textsubscript{1} s)
      (lambda () (g\textsubscript{3} s))
      (lambda (s\textsubscript{0} s\textsuperscript{∞})
        (append map\textsubscript{fair} g\textsubscript{2}
          (append\textsuperscript{∞} (unit s\textsubscript{0}) s\textsuperscript{∞}))))
  ))

(\textit{define} (once g)
  (lambda (s)
    (elim (g s)
      (lambda () (none))
      (lambda (s\textsubscript{0} s\textsuperscript{∞}) (unit s\textsubscript{0}))))

Fig. 7. New and changed functions in BFS\textsubscript{imp} that implement impure features

Now we compare the pure subset of BFS\textsubscript{imp} with BFS\textsubscript{ser}. We focus on the pure subset because BFS\textsubscript{ser} is designed for a pure relational programming system. We prove in Coq that these two search strategies are semantically equivalent, since the result of \textit{(run n ? g)} is the same either way. (See the GitHub repository for the formal proofs.) To compare efficiency, we translate BFS\textsubscript{ser}’s Haskell code into Racket. (See the GitHub repository for the translated code.) The translation is direct due to the similarity of the two relational programming systems. The translated code is longer than BFS\textsubscript{imp}. And it runs slower in all benchmarks. Details about differences in efficiency are in section 7.

7 QUANTITATIVE EVALUATION

In this section, we compare the efficiency of the search strategies. A concise description is in Table 2. A hyphen means “running out of 500 MB memory”. The first two benchmarks are from TRS2. \textit{revers\textsuperscript{a}} is from Rozplokhas and Boulytchev [8]. The next two benchmarks about generating quines are based on a similar test case in Byrd et al. [2]. We modify the relational interpreters because we don’t have disequality constraints (e.g. \textit{absent\textsuperscript{a}}). The sibling benchmarks differ in the \textit{cond\textsuperscript{a}} clause order of their relational interpreters. The last two benchmarks are about synthesizing expressions that evaluate to \textit{’(I love you)}. They are also based on a similar test case in Byrd et al. [2]. Again, we modify the relational interpreters for the same reason. And the sibling benchmarks
Table 2. The results of a quantitative evaluation: running times of benchmarks in milliseconds

<table>
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<th>DFS\textsubscript{i}</th>
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differ in the \textit{cond} clause order of their relational interpreters. The first one has elimination rules (i.e. application, \textit{car}, and \textit{cdr}) at the end, while the other has them at the beginning. We conjecture that DFS\textsubscript{i} would perform badly in the second case because elimination rules complicate the problem when synthesizing (i.e., our evaluation supports our conjecture.)

In general, variants of DFS usually performs better than BFS. The reason might be that BFS tends to remember more states at the same time. Among the three variants of DFS, which all have unfair \textit{conj}, DFS\textsubscript{f} is most resistant to clause permutation in quines and ‘(I love you)s, followed by DFS\textsubscript{bi} then DFS\textsubscript{i}. Thus, we consider DFS\textsubscript{bi} and DFS\textsubscript{f} competitive alternatives to DFS\textsubscript{i}. Among the two implementations of BFS, BFS\textsubscript{imp} constantly performs as well or better.

8 RELATED RESEARCH

In this section, we describe related research. Yang [10] points out that a disjunct complex would be ‘fair’ if it were a full and balanced tree. Seres et al. [9] describe BFS. We present another implementation. Our implementation is semantically equivalent to theirs. But, ours is shorter and performs better in comparison with a straightforward translation of their Haskell code. Rozplokhas and Boulytchev [8] address the non-commutativity of conjunction, while our work about \texttt{disj} fairness addresses the non-commutativity of disjunction.

9 CONCLUSION

We analyze the definitions of fairness. Implementation of \texttt{disj} can be fair, almost-fair, or unfair, depending on how evenly it allocates computational resources to spaces related to disjunctive goals. Implementation of \texttt{conj}
can be fair or unfair, depending on how evenly it allocates computational resources to spaces related to states in the same bucket. Our definition of fair conj, unlike the one by Seres et al. [9], is orthogonal with completeness.

We devise two new search strategies (i.e., DFS$_{bi}$ and DFS$_{f}$) and devise a new implementation of BFS, BFS$_{imp}$. These strategies have different features in fairness: DFS$_{bi}$ has an almost-fair disj and unfair conj. DFS$_{f}$ has fair disj and unfair conj. BFS has both fair disj and fair conj. No search strategy here combines unfair disj and fair conj. This is because we haven’t seen a case where these kinds of search strategies would be interesting.

Our quantitative evaluation shows that DFS$_{bi}$ and DFS$_{f}$ are competitive alternatives to DFS$_{i}$, the current miniKanren search strategy, and that BFS$_{imp}$ is more practical than BFS$_{ser}$.

We prove formally that BFS$_{imp}$ is semantically equivalent to BFS$_{ser}$. But, BFS$_{imp}$ is shorter and performs better in comparison with a straightforward translation of their Haskell code.

Although there are very few benchmarks, this is preliminary work where we are making a point that certain levels of fairness come without cost in some cases, and that each of the search strategies: DFS$_{i}$, DFS$_{bi}$, DFS$_{f}$, and BFS, can co-exist inside one’s head. Constructing a miniKanren with all levels of fairness is future work.

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REFERENCES

First-order miniKanren representation: Great for tooling and search

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We present a first-order implementation of miniKanren that makes it easy to build a miniKanren debugger, and allows any other processes (including a human or a neural network) to guide the search. Typical miniKanren implementations use procedures to represent data structures like goals and streams. Instead, our implementation uses Racket structs, which are transparent, decomposable, manipulable, and not coupled with a particular search strategy. We obtain this first-order implementation by carefully applying defunctionalization rules to a higher-order implementation, deriving two compatible versions with the same search behavior and comparable performance. Decoupling the search in the first-order implementation makes it possible to analyze, transform, and optimize a miniKanren program, even while that program is running. We use a “human guided” search as a miniKanren debugger, and to demonstrate the breadth of supported search strategies. The flexibility in how we interpret goals and streams opens up possibilities for new tools, and we hope to inspire the community to build better miniKanren tooling.

CCS Concepts:
- Software and its engineering → Functional languages; Constraint and logic languages; Constraints.

Additional Key Words and Phrases: miniKanren, microKanren, logic programming, relational programming, Scheme, Racket, program synthesis

1 INTRODUCTION

While miniKanren has been applied successfully to at least seven different programming problems [1], understanding how miniKanren arrives at answers is challenging for all but the simplest examples. Well-hidden typos and inefficient ordering of conjuncts can be difficult to detect. Why is it that after more than a decade of work, miniKanren still lacks developer tools, like a basic debugger?

A seemingly different challenge is supporting different miniKanren search strategies. Different strategies may be better-suited to particular miniKanren programs, so flexibility in choosing search strategies can make programs run faster. Unfortunately, in Byrd et al. [1], changing the search strategy required significant modifications to a miniKanren implementation, as well as a tedious, ad hoc reimplementation of a relational interpreter to work around the default search behavior.

These two problems—building a debugger, and implementing various search strategies—can seem very different on the surface. However, they are actually symptoms of one design decision common to most miniKanren implementations: a higher-order representation of programs coupled to a particular search strategy.

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Most miniKanren implementations represent an executing relational program using goals and streams. In a higher-order implementation, goals and streams are represented using procedures. Procedures encapsulate computational behavior, defining the search strategy for satisfying constraints. Unfortunately, procedures can’t be modified or decomposed. Printing one yields:

```lisp
> (fresh (x y)
    (conde ((== 10 x) (== 20 y))
           ((== 30 x) (== 40 y))))
#<procedure:...>
```

Our proposal is to use a first-order program representation, where goals and streams are data structures not coupled to any search strategy. These data structures are decomposable, inspectable and manipulable:

```lisp
> (fresh (x y)
    (conde ((== 10 x) (== 20 y))
           ((== 30 x) (== 40 y))))
#s(disj
    #s(conj #s(== 10 #s(var x 1)) #s(== 20 #s(var y 2)))
    #s(conj #s(== 30 #s(var x 1)) #s(== 40 #s(var y 2))))
```

The rest of the paper is organized as follows: we begin in Section 2 with an example of how a miniKanren stepper could be useful. Section 3 describes the requirements of a miniKanren implementation to support such a stepper. In Section 4, we implement a first-order version of miniKanren alongside a higher-order version with the same search behavior. In Section 5, we explain the program transformations that make the stepper output easier to understand. Section 6 describes applications of the first-order representation, such as building the stepper, using a neural guided search strategy, and using dynamic search strategies, as well as providing ideas for other miniKanren tools that have been difficult to implement in the past. We hope to inspire you to build these tools.

2 APPENDOH THE DEFICIENT: A MINIKANREN EMERGENCY

We begin this paper with a miniKanren emergency that we want you—dear readers—to help us resolve. Nub Let, a hapless but earnest beginner miniKanren programmer, has accidentally replaced our beloved “appendo the Magnificent” relation with appendoh the deficient.

As might be expected of such a dubious relation, appendoh has been failing left and right. Even worse, Nub Let forgot to email the code for appendoh as part of the bug report. Fortunately, Nub Let included a transcript of an interactive stepper session, enabled by using a first-order representation of miniKanren code.

Will you help us save this paper from the embarrassment of including a buggy five-line program?

Recall that a correct implementation of appendo looks like this:

```lisp
(define-relation (appendo l s ls)
    (conde
        ((== '() l) (= s ls))
        ((fresh (a d res)
            (= ‘,(a ,d) l)
            (= ‘,(a ,res) ls)
            (appendo d s res)))))
```

The output of interactively stepping through an appendoh query is shown in Figure 1. The execution of the query is broken down into steps. Each step of evaluation corresponds to expanding one branch of a logical disjunction.
Stepping through query: (query (x y) (appendoh x y '(1 2 3)))

Initial miniKanren query

Current Depth: 0  Number of Choices: 1

| Choice 1: |
| x = #s(var x 15) |
| y = #s(var y 16) |
| Constraints: |
| * (appendoh #s(var x 15) #s(var y 16) (1 2 3)) |

[h]elp, [u]ndo, or choice number> 1

Current Depth: 1  Number of Choices: 2

| Choice 1: |
| x = () |
| y = (1 2 3) |
| No constraints |

| Choice 2: |
| x = (1 . #s(var b 18)) |
| y = #s(var y 16) |
| Constraints: |
| * (appendoh #s(var b 18) #s(var y 16) (1 2 3)) |

[h]elp, [u]ndo, or choice number> 2

Current Depth: 2  Number of Choices: 2

| Choice 1: |
| x = (1) |
| y = (1 2 3) |
| No constraints |

| Choice 2: |
| x = (1 1 . #s(var b 21)) |
| y = #s(var y 16) |
| Constraints: |
| * (appendoh #s(var b 21) #s(var y 16) (1 2 3)) |

[h]elp, [u]ndo, or choice number>

Fig. 1. Interactive stepper output for the buggy appendoh
Let us explain what you see in Figure 1: at each step, the user picks a disjunct to expand from a list of numbered choices. In the appendoh case, the disjuncts correspond to the conde branches in its definition, each of which constrains the query variables in some way. For each choice, the stepper shows the values the query variables will have if we take that choice. The stepper also shows zero or more constraints associated with a choice. These constraints correspond to goals that must be satisfied if we take the choice.

To prevent an explosion of choices, the stepper only displays the new choices that are descendants of the choice we just made in the previous step. We retain a history of user choices to provide the option of revisiting previous choices by backtracking. Not all choices will have descendants: some may fail entirely, and some successful choices may not have descendants.

Based on the stepper output in Figure 1 alone, can you deduce what the error is in the implementation of appendoh? The stepper output of the correctly specified appendo is in Appendix A.

It turns out that Nub Let hid the appendoh implementation in Appendix B. Even without the code, a keen reader may have noticed that in all the recursive calls to appendoh shown in the interactive stepper (look at the constraints), the value of the last argument never changes from (1 2 3). This suggests that the bug is somewhere in the recursive call portion of the appendoh code, and that perhaps the last argument is incorrect.

3 DESIDERATA

The appendoh example is simple and has a linear search space. Even so, we hope that this example convinces you that a miniKanren stepper can be a useful debugging tool. In particular, being able to choose which disjuncts to explore is even more helpful for complex miniKanren programs with many disjuncts. Only some of the disjuncts might be buggy, so bugs would only manifest in queries that explore that disjunct. Being able to choose the order of execution allows the programmer to determine which portion of the search space is of interest.

To build a stepper like the one in Section 2, our miniKanren implementation must satisfy two requirements:

1. **Streams should be decomposable, and thus inspectable.** We need some other data structure that, unlike procedures, can be programmatically transformed into a human-readable, textual representation.

2. **Small-step execution should be possible.** We need to be able to take one small step to expand the chosen disjunct. Fine-grained control is important. The idea is not to search the disjunct exhaustively, but to see just a little further in order to inform our next decision. We may even end up deciding that continuing search of this disjunct will be a waste of effort, and prefer to revisit an earlier choice.

In short, we need the search state—the data representation of the stream—to be decoupled from the search behavior—the search strategy that miniKanren uses. Unfortunately, typical miniKanren implementations use a higher-order representation where these are hard to tease apart. Streams are represented as procedures that bundle the search state (the data) with the search behavior (a single interpretation of that data). To satisfy our requirements, we need a representation where the search state is accessible, so we can use it in at least two different ways: as information to be examined by the user, and as a stream to be stepped. If we instead use a first-order representation of the search state, we can view the stream as syntax, writing a different interpreter to satisfy each requirement.

If both functionalities are present, then the interactive stepper is straightforward to implement. The stepper extracts the list of disjuncts, presents these disjuncts as choices to the user, displays the query variable values and constraints associated with each choice, reads user input, takes a step to expand the chosen disjunct, and repeats until the user is satisfied.

---

1 Technically, it is possible to use procedures implemented in an object-oriented style, with a method that prints a human-readable textual representation of its internal state. The issue with this approach is that supporting a new behavior requires implementing a new method. To avoid this issue, we could implement a method that provides a complete representation as a data structure suitable for any purpose. But at that point, we might as well just represent the stream as that data structure, and not use a procedure at all.
4 IMPLEMENTATIONS OF MINIKANREN

In this section, we build two miniKanren implementations, one of which uses a first-order representation of the search state to decouple it from the search behavior. That miniKanren implementation satisfies the two requirements from Section 3: that streams should be decomposable, and small-step executions should be possible.

Our plan is to build miniKanren on top of a version of μKanren. We start by implementing a higher-order μKanren to use as a familiar reference point. Our implementation is inspired by that of Hemann and Friedman [2], but modified for improved performance. We encourage interested readers to review Hemann and Friedman [2] for a detailed presentation of the μKanren language.

To obtain first-order μKanren, we will defunctionalize [4] the higher-order version. In particular, the first-order μKanren uses Racket’s struct, which can be decomposed using match expressions (Section 4.1.1). This version of μKanren also has a small step interpreter step that performs a small, finite expansion of a stream (Section 4.3.1). For comparability, both μKanren versions implement the same search behavior.

The first-order and higher-order implementations share a portion of code, presented in Section 4.1. The remaining portions of both implementations are shown side by side in Figure 2 on page 8. We discuss the higher-order portion in Section 4.2, and compare with the first-order portion in Section 4.3. We recover a version of miniKanren in Section 4.4, compatible with both first-order and higher-order μKanren.

4.1 Common Implementation

The common portion defines logic variables, constraint states, unification, and reification:

4.1.1 Logic Variables. Logic variables\(^2\) are defined using struct, which has the benefit of avoiding type collision with vectors. This allows support for vectors as terms, though for simplicity we do not include this support.

The declaration (struct NAME (FIELD...) #:prefab) defines a new record type, automatically generating a constructor named NAME and field accessors named NAME-FIELD for each FIELD. The match form can pattern match against values of this new type. Additionally, using #:prefab gives values of this type the printable notation #s(NAME FIELD ...).

A logic variable contains the name it was given in the program, to help with debugging, and a non-negative integer index used to identify it. Unlike the implementation of Hemann and Friedman [2], a fresh logic variable is constructed with an index allocated from a shared, mutable counter, using var/fresh. This simplifies the role of states, which would otherwise carry a functional counter [2]. Additionally, without this simplification, a first-order implementation would be burdened with managing the functional counter to resolve unbound logic variables\(^3\).

\(\text{;; Logic variables}\
\begin{align*}
\text{(struct var (name index) #:prefab)}
\text{(define (var=? x1 x2)}
\text{(= (var-index x1) (var-index x2)))}
\text{(define initial-var (var #f 0))}
\text{(define var/fresh)}
\text{(let ((index 0))}
\text{\quad (lambda (name) (set! index (+ 1 index))}
\text{\quad (var name index)))}
\end{align*}
\)

\(^2\)Recall that a logic variable is a term that acts as a placeholder for an unknown value.

\(^3\)Because we would like to implement many interpreters, this burden would multiply, because each interpreter would have to include support for resolving unbound logic variables.
4.1.2 States. States containing constraints are defined using struct and track equality information using a substitution. Other kinds of constraints could be supported, but for simplicity we will only consider equality constraints in this implementation.

Substitutions are implemented as in Hemann and Friedman [2], though we choose to enforce acyclic term structure using the occurs? check when extending a substitution via extend-sub. Our choice is consistent with typical miniKanren implementations.

```scheme
;; States
(define empty-sub '())
(define (walk t sub)
  (let ((xt (and (var? t) (assf (lambda (x) (var=? t x)) sub))))
    (if xt (walk (cdr xt) sub) t)))
(define (occurs? x t sub)
  (cond ((pair? t) (or (occurs? x (walk (car t) sub) sub)
                        (occurs? x (walk (cdr t) sub) sub)))
        ((var? t) (var=? x t))
        (else #f)))
(define (extend-sub x t sub)
  (and (not (occurs? x t sub)) '((,x . ,t) . ,sub)))
(struct state (sub) #:prefab)
(define empty-state (state empty-sub))
```

4.1.3 Streams and Unification. Unification enforces equality between two terms and returns a substitution. It is defined as unify, and is implemented as in Hemann and Friedman [2].

```scheme
;; Unification
(define (unify/sub u v sub)
  (let ((u (walk u sub)) (v (walk v sub)))
    (cond ((and (var? u) (var? v) (var=? u v)) sub)
          ((var? u) (extend-sub-u v sub))
          ((var? v) (extend-sub-v u sub))
          ((and (pair? u) (pair? v))
           (let ((sub (unify/sub (car u) (car v) sub)))
            (and sub (unify/sub (cdr u) (cdr v) sub))))
          (else (and (eqv? u v) sub))))
(define (unify u v st)
  (let ((sub (unify/sub u v (state-sub st))))
    (and sub (cons (state sub) #f))))
```

4.1.4 Reification. Reification normalizes the presentation of terms containing logic variables for easier reading and comparison. We implement reify as in Hemann and Friedman [2], except using a local, mutable counter to simplify allocation of normalized variable identifiers.

```scheme
;; Reification
(define (walk* tm sub)
  (let ((tm (walk tm sub)))
    (if (pair? tm)
4.2 Higher-order $\mu$Kanren

Goals, immature streams, and stream maturation are implemented differently in higher- vs first-order $\mu$Kanren. The remaining implementation of higher-order $\mu$Kanren is shown on the left-hand-side of Figure 2 on page 8.

4.2.1 Goals. Goals include binary disjunctions and conjunctions, equality constraints, and calls to user-defined relations.

Binary disjunctions and conjunctions are constructed by disj and conj, which are defined as in Hemann and Friedman [2], except that our implementation introduces explicit pauses.

Equality constraints are constructed with ==, which uses unify.

A call to a user-defined relation applies the relation to argument terms. Each call is represented as a thunk, wrapped with the constructor relate. A user-defined relation is expressed as a Racket procedure, such that forcing a call’s thunk corresponds to applying the procedure to the call’s arguments. (Higher-order $\mu$Kanren places no restrictions on what a Racket procedure implementing a relation may do, but Section 4.3.2 will describe a few restrictions when defunctionalizing relation definitions in first-order $\mu$Kanren.) A use of relate is also passed first-order metadata describing the call, but this information is discarded by our higher-order implementation.

4.2.2 Streams. Streams define a search strategy for finding answers, which are states that satisfy all of a program’s constraints. For improved performance$^4$, we approximate the behavior of the interleaving search used in Byrd et al. [1].

A mature stream is either the empty stream #f, or a pair of an answer and a remaining stream. An immature stream is a thunk that has suspended the search before its completion. Forcing this thunk causes the search to continue. Streams will typically be constructed with the help of mplus, bind, and pause.

In our first-order implementation, we represent suspension more explicitly than in Hemann and Friedman [2]. A pause produces an immature stream by introducing a thunk that delays transition of a goal to a stream. When forced, the thunk applies the goal to a state.

The stream constructor mplus combines two streams. To improve answer diversity, our implementation of mplus interleaves more often to prevent highly productive branches from monopolizing search resources.

$^4$Compared with the search described by Hemann and Friedman [2], this implementation can synthesize quines about twice as fast, and twines and thrines nearly three times as fast, using a simple relational interpreter.
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interleave each time an answer is popped off of a mature stream, rather than popping all available answers before interleaving. To make this possible, bind needs to introduce a pause every time an answer is discovered.

Our implementation of mplus also delays forcing a child stream until it is ready to be examined. This small detail can significantly reduce wasted work and memory usage.

4.2.3 Stream maturation. Stream maturation is implemented by mature. In order to extract an answer from a stream, mature will continue forcing execution of an immature stream until it is mature.

4.3 First-order μKanren

We defunctionalize the higher-order representation of goals and streams to obtain the first-order implementation of μKanren shown in Figure 2. This new version of μKanren represents goal and streams as decomposable data structures, and interprets streams using step, with the help of a goal interpreter start.

The goal constructors disj, conj, relate, and ==, along with the stream constructors bind, mplus, and pause, are now defined using struct to allow decomposition by pattern matching via the keyword match. Mature streams are represented the same way as in higher-order μKanren, as either the empty stream #f, or a pair of an answer and a remaining stream.

We add two new procedures, step and start, which are small-step interpreters for streams and goals respectively. These interpreters work together to implement the same interleaving search strategy defined by the higher-order representation. More specifically, step interprets an immature stream by making a finite amount of progress in searching for an answer. It interprets a mature stream by simply returning the stream. The interpreter start interprets a goal in the context of a state, producing a new stream.

4.3.1 Defunctionalizing goals and streams. Each higher-order goal and stream definition maps to a pattern matching clause in the corresponding interpreter. We have aligned the higher-order and first-order implementation listings in Figure 2 to make this correspondence easy to see: for each line in the higher-order definition on the left, scan along the same line to see the first-order version on the right.

We need to be careful when translating from higher-order to first-order. Failing to apply step at the right time may impede search progress, and change the order in which answers are found. Applying step too soon may lead to an infinite loop when a recursive relation is involved. These issues may not manifest immediately, so when they show up, it is unclear whether they are due to a particular query, the definition of a relation, or a problem with the search implementation. To avoid causing these issues when defunctionalizing the higher-order representation, we must match its order of evaluation. We follow these rewrite rules to defunctionalize:

**Rule:** Top-level higher-order constructors become match clauses

- Every top-level higher-order goal constructor becomes a struct pattern clause of the match expression in start.
- Every top-level higher-order stream constructor becomes a struct pattern clause of the match expression in step.
- In each case, the pattern decomposes the struct, naming its components the same way the higher-order constructors name their parameters. The clause bodies will then be defunctionalized.

**Rule:** Delaying goal/stream interpretation

- Anywhere that lambda is used in the higher-order implementation to delay construction of a goal or stream, we simply apply the corresponding struct constructor.
- **Justification:** though the higher-order constructors are procedures which immediately execute their interpretation when applied, the enclosing lambda leaves them inert until forced. Likewise, newly-constructed struct data remains inert until interpreted by step or start.
Rule: Interpreting a goal/stream

- Anywhere that the higher-order implementation does not delay construction of a goal, we apply start to the result of the corresponding struct constructor.
- Except in the case of pause, anywhere that the higher-order implementation does not delay construction of a stream, we apply step to the result of the corresponding struct constructor.
- Anywhere the higher-order implementation uses pause, we simply apply the pause struct constructor.
- Justification: We apply start and step in this way because the higher-order constructors also immediately execute their interpretation. But pause is an exception because it does not immediately execute its interpretation. Instead, it delays interpretation by producing an immature stream (a thunk).

Rule: Forcing immature streams

- Anywhere that the higher-order implementation forces progress of an immature stream, which corresponds to applying a thunk, we apply step to that stream to also force progress. Look at the definition of mature in Figure 2 for a simple example of this.

4.3.2 Defunctionalizing user-defined relations. As in the higher-order implementation, we represent calls to user-defined relations using thunks wrapped with the constructor relate. Forcing such a thunk applies a Racket procedure to the call’s arguments. Unlike the higher-order implementation, uses of first-order relate will retain the call description metadata passed alongside the thunk. While an entirely first-order representation of user-defined relations is possible, we prefer retaining this use of thunks because it allows us to leverage Racket’s implementation of lexical scope and procedure abstraction, rather than reimplement this ourselves.

Despite continuing to use thunks to represent calls to relations, it is still possible to perform arbitrary manipulation of relation definitions if we make some assumptions about what such definitions may do. Specifically, we will assume the use of Racket is limited to constructing µKanren goals. When these goals include terms, the terms may be expressed using constants, references to lexical variables, pair construction, or fresh logic variable construction.

Given our assumptions, we can obtain a first-order representation of a user-defined relation by applying it to fresh logic variables to represent the parameters. Any goals in the result are those specified in the body of the relation. Any terms in those goals are also either those specified, or they are logic variables. If a logic variable is one of those used to represent a parameter of the relation, then the relation definition must reference that parameter at that location. All other logic variables were created fresh. Additionally, since each use of relate wraps metadata describing the call, we can always reconstruct a complete first-order description of a relation, even if it is recursive.

4.4 Recovering miniKanren

In Appendix C we implement miniKanren on top of µKanren, using syntax-rules definitions similar to those in Hemann and Friedman [2], with a notable addition: we define query, which describes the initial stream built by a corresponding run. Since we are interested in manipulating streams in more ways than just extracting answers, we need to separate stream construction from answer extraction. To extract answers from a stream, we provide stream-take such that:

\[(\text{run } N \text{ body } \ldots) \implies (\text{stream-take } N \text{ (query body } \ldots))\]

We will make use of the metadata wrapped by relate when visualizing program representation in Section 5. The metadata is a list containing the relation procedure (to allow unambiguous identification), the name of that procedure, and all of the argument terms.

---

Hemann and Friedman [2] leverages Scheme in the same way.
4.5 Performance

When comparing the higher-order and first-order implementations, we notice a small performance difference when generating twines and thrines using a simple relational interpreter, and when running relational arithmetic benchmarks. On these tests, the first-order implementation is roughly 5 percent slower and spends about 10 percent more time in the garbage collector. Chez Scheme runs a Scheme version of our benchmarks about 25 percent faster than Racket, but the first-order vs. higher-order performance ratio remains about the same.

5 PROGRAM TRANSFORMATIONS

With the first-order miniKanren implementation given in Section 4, streams are decomposable (and therefore printable), and small-step executions are possible.

You might wonder whether we could stop here. After all, we can apply the procedure step to progress through a query, and print the structure of the stream at each step. Appendix D shows the output of stepping through the query (query (xs) (appendo xs xs ’(a a))) using interleaving search.

Unfortunately, this presentation is intolerably verbose, even for such a simple query. It is difficult to extract useful information about actual available choices. In addition, many useless steps are taken to expand goals that are obviously failing. In order to obtain a simpler program representation that is actually human-readable, we need to **extract the useful information about each disjunct**, and **remove obviously failing disjuncts**.

This section recovers the useful trace information that we saw in Figure 1 by introducing stream transformations. Section 5.1 describes how to prune obviously failing disjuncts, so they need not be considered by a human. Section 5.2 describes how to transform the stream into disjunctive normal form (DNF), to provide a flat tree of disjuncts to choose from.

5.1 Pruning obvious failures

Some streams and goals in Appendix D are obviously failing. Examples include streams of the form (bind #f _), (mplus #f #f), goals of the form (== A B), where A and B are unequal terms, and other streams and goals that depend on failing components. We would prefer to eliminate these immediately so we can focus on meaningful parts of the search state.

We can define a **pruning** transformation that uses equality information to identify constraints that will definitely fail. Since streams and goals depending on these constraints may fail to produce any results, we will prune them accordingly, producing a simplified stream.

Before taking a look at the implementation, it is important to consider that a transformation may change the observed behavior of our program. We should decide what sorts of changes we will tolerate. Viewing a miniKanren program as a search for answers that satisfy a query, we feel the following concessions are acceptable:

1. Answers may be reordered.
2. The contents of a state may be rearranged as long as its constraints express the same meaning.
3. The number of steps required to find an answer may change, but will remain finite.
4. Unproductive, infinite streams may become finite.

With that in mind, let us take a look at the implementation. The interpreters `prune/stream` and `prune/goal` partially evaluate `==` constraints, using the resulting state to propagate equality information. Any failing stream or goal will cause its parent `bind` and `conj` to also fail. A `mplus` or `disj` containing a failing child will be replaced by the remaining child.

```scheme
(define (prune/stream s)
  (match s
    ((mplus s1 s2) (match (prune/stream s1)
                         (match (prune/stream s2)
                           ((con s3) (match (prune/stream s)))))
     (match s
               (match (prune/goal s)
                      ((mplus g1 g2) (match (prune/goal g1)
                                             (match (prune/goal g2)
                                                    ((equal?) (prune/goal s)))))
               (match (prune/goal s)
                      ((disj g1 g2) (match (prune/goal g1)
                                            (match (prune/goal g2)
                                                   ((equal?) (prune/goal s)))))))))
```


Since pruning performs obvious clean up work without spending steps to do so, pruning often leads to a trace containing fewer steps.
5.2 Disjunctive Normal Form

In Figure 1, we actually do one more thing—we provide the user with a flat list of choices. We can obtain those choices by first transforming the stream into disjunctive normal form (DNF).

The general DNF transformation rewrites goals and streams in the following way:

\[
\begin{align*}
\text{(conj (disj A B) C)} & \implies \text{(disj (conj A C) (conj B C))} \\
\text{(conj C (disj A B))} & \implies \text{(disj (conj C A) (conj C B))} \\
\text{(pause st (disj A B))} & \implies \text{(mplus (pause st A) (pause st B))} \\
\text{(bind (mplus A B) C)} & \implies \text{(mplus (bind A C) (bind B C))} \\
\text{(bind C (disj A B))} & \implies \text{(mplus (bind C A) (bind C B))}
\end{align*}
\]

This transformation has the effect of lifting all disjunctions out of conjunctions.

\[
\begin{align*}
\text{(define (dnf/stream s))} & \\
\text{(define (push-pause st g))} & \\
\begin{align*}
\text{(match g} & \\
\text{((disj g1 g2) (mplus (push-pause st g1) (push-pause st g2)))} \\
\text{(g (pause st g)))}
\end{align*} \\
\text{(match s} & \\
\begin{align*}
\text{((bind s g}) & \\
\text{(let loop1 ((s (dnf/stream s)) (g (dnf/goal g)))} \\
\text{(define (loop2 s g}) & \\
\begin{align*}
\text{(match g} & \\
\text{((disj ga gb) (mplus (loop2 s ga) (loop2 s gb))}) & \\
\text{(g (bind s g)))}
\end{align*} \\
\text{(match s} & \\
\begin{align*}
\text{((mplus sa sb) (mplus (loop1 sa g) (loop1 sb g))}) & \\
\text{('st . ,s) } & \\
\text{(mplus (push-pause st g) (loop1 s g))} & \\
\text{(s (loop2 s g)))))
\end{align*} \\
\text{((pause st g) (push-pause st (dnf/goal g))} & \\
\text{(mplus s1 s2) (mplus (dnf/stream s1) (dnf/stream s2))} & \\
\text{('st . ,s) } & \\
\text{((st .,(dnf/stream s)))} & \\
\text{(s s))}
\end{align*}
\]
\]

\[
\begin{align*}
\text{(define (dnf/goal g))} & \\
\text{(match g} & \\
\begin{align*}
\text{((conj g1 g2} & \\
\text{(let loop1 ((g1 (dnf/goal g1)) (g2 (dnf/goal g2)))} \\
\text{(define (loop2 g1 g2)
\end{align*}
\]

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Aside from preparing a stream for choice extraction, DNF has the added benefit of allowing equality information to propagate more deeply, making pruning more effective. Consider the following goal:

\[(\text{conj} (\text{disj} (== x 5) A) B)\]

In this situation, information flow is impeded. Though we constrain \(x\) to be equal to 5, we cannot propagate this information into the goal \(B\), because it is possible that goal \(A\) may not express the same constraint.

We can allow this information to flow more freely by lifting the disjunction out of the conjunction like so:

\[(\text{disj} (\text{conj} (== x 5) B) (\text{conj} A B))\]

Because each child of the disjunction now has its own copy of the goal \(B\), it is safe to propagate the equality information into the affected copy of \(B\).

Another interesting property of DNF is that if we apply the full pruning transformation on a DNF stream, all remaining \(==\) constraints will be trivial in the sense that they are guaranteed to succeed, and will not provide new information. This is because all available equality information will have already been gathered in a pause state. These trivial constraints may be safely removed.

One downside of DNF transformation is that program representations can become larger due to copying children of conjunctions. In the worst case, the increase is exponential in the size of the original program. This potential increase in size may be mitigated by an increased likelihood of prunable, obvious failures, thanks to the improved flow of equality information.

Once we have a pruned DNF stream, we are finally in a position to provide choices of the form shown in Figure 1. In section 6.1 we will explain how we extract this list of choices from such a stream.

6 APPLICATIONS

6.1 The miniKanren stepper

In this section, we implement \texttt{explore/stream}, the miniKanren stepper described in Figure 1. The implementation of \texttt{explore/stream} is shown below. The main stepping loop follows the outline given at the end of Section 3. We describe each facet of this loop in chronological order.

The first thing the \texttt{explore/stream} loop does is use the procedure \texttt{stream->choices} to extract a list of choices from a stream. To prepare for extraction, we transform the stream into disjunctive normal form (DNF) using \texttt{dnf/stream}, and perform pruning with \texttt{prune/stream}. Recall that these procedures were implemented in Section 5. With a pruned DNF stream, we are guaranteed to find all disjunctions at the top (after any available answers), as a \texttt{mplus} tree, with leaves of the form (pause \texttt{STATE GOAL}). Each of these leaves represents a choice, and all choices can be extracted while traversing the tree.

Once we have our list of choices, we need to present each choice in a digestible form. We handle this with \texttt{print-choice}. The \texttt{GOAL} in each choice is either a trivial \(==\) constraint (i.e., both terms are guaranteed to be equal), a relate call, or a \texttt{conj} tree whose leaves are goals of this form. Traversing this \texttt{conj} tree, as in \texttt{goal->constraints}, extracts these goals as a list of the choice’s constraints. As mentioned in Section 5.2, pruning
in DNF has the emergent property that any remaining == constraints are guaranteed to be trivial, so we may discard them, allowing us to focus on the relate calls. Finally, pruning in DNF also guarantees that, of the variables mentioned in STATE, only those reachable from query variables are referenced in non-trivial constraints. Therefore, we can discard STATE, keeping only the query variable values.

With the interesting part of the implementation out of the way, all we have left to do is read user input and act accordingly. When the user makes a choice, we use step to expand the choice, providing the resulting stream as input to the next iteration of \texttt{explore/stream}.

```scheme
(define (explore/stream qvars s)
  (define margin "\ |\")
  (define (pp prefix v) (pprint/margin margin prefix v))
  (define (pp/qvars vs)
    (define qv-prefixes (map qv-prefix qvars)))
  (if (and qv-prefixes
            (for-each (lambda (prefix v) (pp prefix v)) qv-prefixes vs)
            (for-each (lambda (v) (pp "\ *\ " v)) vs)))
    (define (print-choice s)
      (match s
        ((pause st g)
         (pp/qvars (walked-term initial-var st))
         (define cxs (walked-term (goal->constraints st g) st))
         (unless (null? cxs)
           (displayln (string-append margin "Constraints:"))
           (for-each (lambda (v) (pp "\ *\ " v)) cxs))
         (when (null? cxs)
           (displayln (string-append margin "No constraints")))))))
    (let loop ((s (stream->choices s)) (undo '()))
      (define previous-choice
        (and (pair? undo)
          (let* ((i.s (car undo)) (i (car i.s)) (s (cdr i.s)))
            (list-ref (dropf s state?) (- i 1)))))
      (define results (takef s state?))
      (define choices (dropf s state?))
      (display "\n====================================================================")
      (displayln "====================================================================")
      (unless (= (length results) 0)
        (printf "Number of results: ~a\n" (length results))
        (for-each (lambda (st)
                    (pp/qvars (walked-term initial-var st))
                    (newline)) results)
      (when (and previous-choice (null? results))
        (printf "Previous Choice:\n")
        (print-choice previous-choice)
        (newline))
  (let ((i.s (car undo)) (i (car i.s)) (s (cdr i.s)))
    (list-ref (dropf s state?) (- i 1))))
```

(0)
6.2 Search Strategies

A first-order miniKanren with the small-step interpreter \textit{step} gives us fine-grained control over the search strategy. Even though we begin Section 4.2 by reimplementing the biased-interleaving search, we can use whatever search strategy we like. In this section, we describe a few different search strategies to illustrate the potential.

6.2.1 Human-guided search. One way of interpreting what we did in Section 6.1 is that we \textit{hijacked} the search strategy. That is, the miniKanren stepper is actually an implementation of a human-guided search strategy! Since human behavior is as unpredictable as any program can be, the miniKanren stepper shows that any kind of search strategy is possible.

6.2.2 Neural-guided search. Another idea is to replace the human-guided search strategy with an artificial human, namely a neural network. In Zhang et al. [6] a neural network learns to guide the search of first-order miniKanren performing a Programming by Example (PBE) task. In such a task, miniKanren is used to synthesize a procedure specified using input/output pairs.
For example, if we have a relational interpreter (eval-expo program environment output), then an example PBE query with two input-output pairs is:

\[
\text{(query (p)}
\begin{align*}
&\text{(eval-expo p (\’(a) (\’(a a a)))} \\
&\text{(eval-expo p (\’(b) ((b b b)))}
\end{align*}
\]

The neural network evaluates each choice (disjunct) independently. It takes as input all the constraints corresponding to each choice, and chooses the highest-scoring choice to expand in each step.

There are several deviations from Section 6.1 required for this setup. First, unlike in Figure 1, the neural network can choose to expand any choice in each step, not just descendants of the previous step. So, the neural network would not need the "undo" functionality at all. This revised setup means that the neural network does not need to remember any past information, and can treat each step independently.

Another deviation is including a powerful strategy appropriate to PBE problems, which specializes the expansion behavior of eval-expo constraints. In a PBE problem, each input-output example imposes constraints on the same program to be synthesized, using eval-expo. However, a typical miniKanren search will attempt to satisfy each eval-expo constraint completely before moving on to the next. This is unfortunate because only information from one input-output example is considered at any given time, leading to deep exploration of impossible synthesis choices that could have been ruled out immediately. Instead, we implement an expansion step that will simultaneously expand all eval-expo constraints that share the same program argument. Combined with DNF transformation and pruning, simultaneous expansion leads to immediate feedback being provided by all input-output examples for each synthesis choice made. Aside from refuting bad choices more quickly, this simultaneous expansion allows the neural network to observe related constraints all at once.

The network in Zhang et al. [6] is trained on generated PBE problems to which we know the answer, so that the correct choice is known for each step. During training, the neural network guides the search by making choices, and incorrect choices are penalized. We refer interested readers to Zhang et al. [6] for more details about the model architecture and training.

6.2.3 Dynamic Goal Reordering. Though we have mostly discussed transformations while developing the interactive stepper, there are other ways to improve a search strategy. One source of possibilities is Byrd et al. [1], which makes use of a relational interpreter that performs various forms of conjunction reordering based on heuristics specific to the interpreter definition. For instance, goals are reordered based on determinism annotations on conde expressions, where at most one clause of a conde may succeed if we know the value of a particular subset of logic variables. The conde expressions recognized as deterministic are expanded before any others. This reordering would be more natural to implement with a decoupled representation, where it is easier to compose with transformations and other strategies.

6.3 More possibilities

Beyond improvements to the search, a first-order representation is a natural medium for program analysis and compilation. For instance, though we have not explored these ideas, we believe it should be possible to implement garbage collection of constraints and substitutions, and just-in-time compilation, possibly benefiting from mode analysis.

7 RELATED WORK

There have been previous unpublished implementations of miniKanren search using a first-order implementation. For example, Michael Ballantyne has implemented a small-step miniKanren interpreter that uses a first-order
representation. Ballantyne has also implemented a big-step interpreter with a first-order representation of the search tree, in order to implement backjumping search. Both of these implementations decouple the program representation from the search.

To our knowledge, this paper is the first first-order representation of a miniKanren search that preserves the exact search order of the standard higher-order search representation, or that provides rules for a defunctionalization that preserves this search order.

We are happy to report that in this same workshop, two other papers also make use of first-order representations. We hope this is the start of a new trend. Lozov et al. [3] performs supercompilation (as conjunctive partial deduction) using a first-order representation of goals that is similar to our representation of the search state. Rozplokhas et al. [5] also uses a similar first-order goal representation to formalize semantics for miniKanren.

8 CONCLUSION

Both higher-order and first-order representations of miniKanren have their advantages. For both pedagogical reasons and for the beauty and abstractness of the implementation, higher-order representations have been dominant in the miniKanren literature. On multiple occasions, however, implementors have had to switch to first-order representations in order to inspect or control the execution of miniKanren code.

Currently it is a burden to switch representations, not just because of the effort of defunctionalization, but because any changes to the search order may break existing tests, and make it difficult to compare benchmark results.

The rules for defunctionalization presented in Section 4.3, which preserve the exact search order, allow us to have the best of both worlds. We can have a higher-order implementation for pedagogical purposes, for example, and switch to a first-order implementation when it is time to debug or trace the code. And we can be confident that our existing tests, applications, and benchmark programs will not need to change. Of course, while we are living in the first-order world, we can easily change the search, or use an external process to guide the search.

We hope that there will be two positive outcomes to making it easier to switch between higher-order and first-order implementations of miniKanren. First, the wider availability of first-order implementations should make it much easier to produce the tools that the miniKanren ecosystem desperately needs, such as tracers, steppers, and debuggers. The rules in Section 4.3 imply that a programmer can debug a program in a first-order implementation of miniKanren, and be sure the program would show equivalent behavior in a higher-order miniKanren. Secondly, by decoupling the search from the program representation, it should be easier to experiment with novel search techniques, such as driving the search from an external process, or mixing different searches within a single miniKanren program. This should make it easier and faster to explore program synthesis and other advanced relational programming topics.

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First-order miniKanren representation: Great for tooling and search • 2:19


A STEPPER OUTPUT FOR APPENDO

This appendix shows the interactive stepper output for the query (query (x y) (appendo x y '(1 2 3))) with a correctly implemented appendo shown in Section 2. Compare this output with that of a buggy appendo implementation in Figure 1

Stepping through query: (query (x y) (appendo x y '(1 2 3)))

Current Depth: 0 Number of Choices: 1

| Choice 1:
| x = #s(var x 1)
| y = #s(var y 2)
| Constraints:
| * (appendo #s(var x 1) #s(var y 2) (1 2 3))

[h]elp, [u]ndo, or choice number> 1

Current Depth: 1 Number of Choices: 2

| Choice 1:
| x = ()
| y = (1 2 3)
| No constraints

| Choice 2:
| x = (1 . #s(var b 4))
| y = #s(var y 2)
| Constraints:
| * (appendo #s(var b 4) #s(var y 2) (2 3))

[h]elp, [u]ndo, or choice number> 2

Current Depth: 2 Number of Choices: 2

| Choice 1:
| x = (1)
| y = (2 3)
| No constraints

| Choice 2:
| x = (1 2 . #s(var b 7))
| y = #s(var y 2)
| Constraints:
| * (appendo #s(var b 7) #s(var y 2) (3))

[h]elp, [u]ndo, or choice number>
B IMPLEMENTATION OF APPENDOH
This appendix shows the implementation of the buggy appendoh that generated the stepper output in Figure 1.

(define-relation (appendoh l s ls)
  (conde
    ((== '() l) (== s ls))
    ((fresh (a d res)
      (== '(',a ,d) l)
      (== ',(a ,res) ls)
      (appendoh d s ls)))))

C RECOVERING MINIKANREN
We show an implementation miniKanren on top of µKanren. This version of miniKanren is compatible with both the higher-order µKanren implemented in Section 4.2, and the first-order µKanren implemented in Section 4.3.

(define-syntax define-relation
  (syntax-rules ()
    ((_ (name param ...) g ...)
      (define (name param ...)
        (relate (lambda () (fresh () g ...)) '(,name name ,param ...))))))

;; Low-level goals
(define succeed (== #t #t))
(define fail (== #f #t))
(define-syntax conj*
  (syntax-rules ()
    (_ succeed)
    (_ g) g
    (_ gs ... g-final) (conj (conj* gs ...) g-final))))
(define-syntax disj*
  (syntax-rules ()
    (_ fail)
    (_ g) g
    (_ g0 gs ...) (disj g0 (disj* gs ...))))

;; High level goals
(define-syntax fresh
  (syntax-rules ()
    (_ (x ...) g0 gs ...)
    (let ((x (var/fresh 'x)) ...) (conj* g0 gs ...))))

(define-syntax conde
  (syntax-rules ()
    (_ (g gs ...) (h hs ...) ...)
    (disj* (conj* g gs ...) (conj* h hs ...) ...))))

;; Queries
(define-syntax query
  (syntax-rules ()
    (_ (x ...) g0 gs ...)}
(let ((goal (fresh (x ...) (== (list x ...) initial-var) g0 gs ...)))
  (pause empty-state goal)))))

(define (stream-take n s)
  (if (eqv? 0 n) '()
    (let ((s (mature s)))
      (if (pair? s)
          (cons (car s) (stream-take (and n (- n 1)) (cdr s))
                '())))))

(define-syntax run
  (syntax-rules ()
  ((_ n body ...) (map reify/initial-var (stream-take n (query body ...))))))

(define-syntax run*
  (syntax-rules ()
  ((_ body ...) (run #f body ...))))

D RAW APPENDO TRACE

This appendix shows how the search representation changes as we step through the query (query (xs) (appendo xs xs ' (a a))) using interleaving search. We need to transform this representation to recover the easier-to-read textual representations in Figure 1 and Appendix A.

Step 0:
(pause
  (state ()))
(conj
  (== (#s(var xs 1)) #s(var #f 0))
  (relate (appendo #s(var xs 1) #s(var xs 1) (a a)))))

Step 1:
(mplus
  (bind #f (relate (appendo #s(var xs 1) #s(var xs 1) (a a))))
  (pause
    (state (((== #s(var #f 0) (#s(var xs 1))))
      (disj
        (conj (== #s(var xs 1) ()) (== #s(var xs 1) (a a)))
        (conj
          (== (#s(var a 2) . #s(var b 3)) #s(var xs 1))
          (== (#s(var a 2) . #s(var c 4)) (a a)))
          (relate (appendo #s(var b 3) #s(var xs 1) #s(var c 4)))))))

Step 2:
  (pause
    (state (((== #s(var #f 0) (#s(var xs 1))))
      (disj
        (conj (== #s(var xs 1) ()) (== #s(var xs 1) (a a)))
        (conj
          (== (#s(var a 2) . #s(var b 3)) #s(var xs 1))
          (== (#s(var a 2) . #s(var c 4)) (a a)))
          (relate (appendo #s(var b 3) #s(var xs 1) #s(var c 4)))))))
First-order miniKanren representation: Great for tooling and search

\[
(= (\#(var a 2) . \#(var b 3)) \#(var xs 1))
\]
\[
(= (\#(var a 2) . \#(var c 4)) (a a))
\]
\[
(relate (appendo \#(var b 3) \#(var xs 1) \#(var c 4))))
\]

Step 3:
\[
(mplus
\]
\[
(pause
\]
\[
(state ((= #s(var #f 0) (#s(var xs 1))))
\]
\[
(conj
\]
\[
(conj
\]
\[
(= (\#(var a 2) . \#(var b 3)) \#(var xs 1))
\]
\[
(= (\#(var a 2) . \#(var c 4)) (a a))
\]
\[
(relate (appendo \#(var b 3) \#(var xs 1) \#(var c 4))))
\]
\[
(bind \#f ((= \#(var xs 1) (a a))))
\]
Step 4:
\[
(mplus
\]
\[
(bind \#f ((= \#(var xs 1) (a a)))
\]
\[
(mplus
\]
\[
(bind (mplus (bind \#f ((= (\#(var a 2) . \#(var c 4)) (a a))) \#f)
\]
\[
(relate (appendo \#(var b 3) \#(var xs 1) \#(var c 4))))
\]
\[
(pause
\]
\[
(state
\]
\[
((= \#(var c 4) (a))
\]
\[
(= \#(var a 2) a)
\]
\[
(= \#(var xs 1) (a . \#(var b 3)))
\]
\[
(= \#(var #f 0) ((a . \#(var b 3))))))
\]
\[
(disj
\]
\[
(conj ((= \#(var b 3) ()) (a . \#(var b 3)) (a))
\]
\[
(conj
\]
\[
(conj
\]
\[
(= (\#(var a 5) . \#(var b 6)) \#(var b 3))
\]
\[
(= (\#(var a 5) . \#(var c 7)) (a))
\]
\[
(relate (appendo \#(var b 6) (a . \#(var b 3)) \#(var c 7))))))
\]
Step 5:
\[
(mplus
\]
\[
(bind
\]
\[
(mplus (bind \#f ((= (\#(var a 2) . \#(var c 4)) (a a))) \#f)
\]
\[
(relate (appendo \#(var b 3) \#(var xs 1) \#(var c 4))))
\]
\[
(pause
\]
\[
(state
\]
\[
((= \#(var c 4) (a))
\]
\[
(= \#(var a 2) a)
\]
\[
(= \#(var xs 1) (a . \#(var b 3)))
\]
(== #s(var #f 0) ((a . #s(var b 3)))))

(disj
(conj (== #s(var b 3) ()) (== (a . #s(var b 3)) (a)))
(conj
(conj
(== (#s(var a 5) . #s(var b 6)) #s(var b 3))
(== (#s(var a 5) . #s(var c 7)) (a)))
(relate (appendo #s(var b 6) (a . #s(var b 3)) #s(var c 7)))))

Step 6:
(pause
(state
((== #s(var c 4) (a))
(== #s(var a 2) a)
(== #s(var xs 1) (a . #s(var b 3)))
(== #s(var #f 0) ((a . #s(var b 3)))))
(disj
(conj (== #s(var b 3) ()) (== (a . #s(var b 3)) (a)))
(conj
(conj
(== (#s(var a 5) . #s(var b 6)) #s(var b 3))
(== (#s(var a 5) . #s(var c 7)) (a)))
(relate (appendo #s(var b 6) (a . #s(var b 3)) #s(var c 7)))))

Step 7:
Answer: (state ((== #s(var #f 0) ((a)))))
Remaining stream:
(mplus
(pause
(state
((== #s(var c 4) (a))
(== #s(var a 2) a)
(== #s(var xs 1) (a . #s(var b 3)))
(== #s(var #f 0) ((a . #s(var b 3)))))
(conj
(conj
(== (#s(var a 5) . #s(var b 6)) #s(var b 3))
(== (#s(var a 5) . #s(var c 7)) (a)))
(relate (appendo #s(var b 6) (a . #s(var b 3)) #s(var c 7)))))
(mplus (bind #f (== #s(var xs 1) #s(var c 4))) #f))

E PRUNING
We show the code used to eliminate obviously failing goals from our stream.

(define (prune/stream s)
  (match s
First-order miniKanren representation: Great for tooling and search

((mplus s1 s2) (match (prune/stream s1)
  (#f (prune/stream s2))
  (s1 (match (prune/stream s2)
   (#f s1)
   (s2 (mplus s1 s2)))))

(bind s g) (match (prune/stream s)
  (#f #f)
  (‘(,st . #f) (prune/goal st g))
  ((pause st g1)
    (match (prune/goal st g)
      (#f #f)
      (‘(pause st g1) (pause st (conj g1 g)))))
  (s (match (prune/goal empty-state g)
    (#f #f)
    ((pause _ _) (bind s g)))))

((pause st g) (prune/goal st g))

‘(‘(,st . ,s) ‘(‘,st . ,,(prune/stream s)))

(s s))

(define (prune/goal st g)
  (define (prune/term t) (walk* t (state-sub st)))
  (match g
    ((disj g1 g2) (match (prune/goal st g1)
      (#f (prune/goal st g2))
      ((pause st1 g1)
        (match (prune/goal st g2)
          (#f (pause st1 g1))
          ((pause _ g2) (pause st (disj g1 g2))))))
    ((conj g1 g2) (match (prune/goal st g1)
      (#f #f)
      ((pause st g1) (match (prune/goal st g2)
        (#f #f)
        ((pause st g2) (pause st (conj g1 g2)))))
      ((relate thunk d) (pause st (relate thunk (prune/term d))))
    )))

(defun (prune/goal empty-state g)
  (define (prune/goal empty-state g)
    )

F TRANSFORMATION TO DISJUNCTIVE NORMAL FORM

We show the code used to transform our streams into Disjunctive Normal Form (DNF).

(define (dnf/stream s)
(define (push-pause st g)
  (match g
    ((disj g1 g2) (mplus (push-pause st g1) (push-pause st g2)))
    (g (pause st g)))
  (match s
    (bind s g)
    (let loop1 ((s (dnf/stream s)) (g (dnf/goal g)))
      (define (loop2 s g)
        (match g
          ((disj ga gb) (mplus (loop2 s ga) (loop2 s gb)))
          (g (bind s g)))
        (match s
          ((mplus sa sb) (mplus (loop1 sa g) (loop1 sb g)))
          (s (loop2 s g)))
          ((pause st g) (push-pause st (dnf/goal g)))
          ((mplus s1 s2) (mplus (dnf/stream s1) (dnf/stream s2)))
          (',' (dnf/stream s))
          (s s)))))

(define (dnf/goal g)
  (match g
    ((conj g1 g2)
      (let loop1 ((g1 (dnf/goal g1)) (g2 (dnf/goal g2)))
        (define (loop2 g1 g2)
          (match g2
            ((disj g2a g2b) (mplus (loop2 g1 g2a) (loop2 g1 g2b)))
            (g2 (conj g1 g2)))
          (match g1
            ((disj g1a g1b) (disj (loop1 g1a g2) (loop1 g1b g2)))
            (g1 (loop2 g1 g2)))
            ((disj g1 g2) (disj (dnf/goal g1) (dnf/goal g2)))
            (g g)))
Relational Interpreters for Search Problems∗

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We address the problem of constructing a solver for a certain search problem from its solution verifier. The main idea behind the approach we advocate is to consider a verifier as an interpreter which takes a data structure to search in as a program and a candidate solution as this program’s input. As a result the interpreter returns “true” if the candidate solution satisfies all constraints and “false” otherwise. Being implemented in a relational language, a verifier becomes capable of finding a solution as well. We apply two techniques to make this scenario realistic: relational conversion and supercompilation. Relational conversion makes it possible to convert a first-order functional program into relational form, while supercompilation (in the form of conjunctive partial deduction (CPD)) — to optimize out redundant computations. We demonstrate our approach on a number of examples using a prototype tool for OCANREN — an implementation of miniKanren for OCAML, — and discuss the results of evaluation.

CCS Concepts: • Software and its engineering → Constraint and logic languages; Source code generation;

Additional Key Words and Phrases: relational programming, relational interpreters, search problems

1 INTRODUCTION

Verifying a solution for a problem is much easier than finding one — this common wisdom can be confirmed by anyone who used both to learn and to teach. This observation can be justified by its theoretical applications, thus being more than informal knowledge. For example, let us have a language $L$. If there is a predicate $V_L$ such that $\forall \omega : \omega \in L \iff \exists p_\omega : V_L(\omega, p_\omega)$ (with $p_\omega$ being of size, polynomial on $\omega$) and we can recognize $V_L$ in a polynomial time, then we call $L$ to be in the class $NP$ [Garey and Johnson 1990]. Here $p_\omega$ plays role of a justification (or proof) for the fact $\omega \in L$. For example, if $L$ is a language of all hamiltonian graphs, then $V_L$ is a predicate which takes a graph $\omega$ and some path $p_\omega$ and verifies that $p_\omega$ is indeed a hamiltonian path in $\omega$. The implementation of the predicate $V_L$, however, tells us very little about the search procedure which would calculate $p_\omega$ as a function of $\omega$. For the whole class of $NP$-complete problems no polynomial search procedures are known, and their existence at all is a long-standing problem in the complexity theory.

There is, however, a whole research area of relational interpreters, in which a very close problem is addressed. Given a language $L$, its interpreter is a function $eval_L$ which takes a program $p^L$ in the language $L$ and an input $i$ and calculates some output such that

$eval_L(p^L, i) = \llbracket p^L \rrbracket_L(i)$

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miniKanren.org/workshop/2019/8-ART3
where $\llbracket \bullet \rrbracket_L$ is the semantics of the language $L$. In these terms, a verification predicate $V_L$ can be considered as an interpreter which takes a program $\omega$, its input $p_\omega$, and returns true or false. A relational interpreter is an interpreter which is implemented not as a function $\text{eval}_L$, which calculates the output from a program and its input, but as a relation $\text{eval}_L^\omega$, which connects a program with its input and output. This alone would not have much sense, but if we allow the arguments of $\text{eval}_L^\omega$ to contain variables we can consider relational interpreter as a generic search procedure which determines the values for these variables making the relation hold. Thus, with relational interpreter it is possible not only to calculate the output from an input, but also to run a program in an opposite “direction”, or to synthesize a program from an input-output pair, etc. In other words, relational verification predicate is capable (in theory) to both verify a solution and search for it.

Implementing relational interpreters amounts to writing it in a relational language. In principle, any conventional language for logic programming (Prolog [Clocksin and Mellish 2003], Mercury [Somogyi et al. 1996], etc.) would make the job. However, the abundance of extra-logical features and the incompleteness of default search strategy put a number of obstacles on the way. There is, however, a language specifically designed for pure relational programming, and, in a narrow sense, for implementing relational interpreters — miniKanren [Friedman et al. 2005]. Relational interpreters, implemented in miniKanren, demonstrate all their expected potential: they can synthesize programs by example, search for errors in partially defined programs [Byrd et al. 2017], produce self-evaluated programs [Byrd et al. 2012], etc. However, all these results are obtained for a family of closely related Scheme-like languages and require a careful implementation and even some ad-hoc optimizations in the relational engine.

From a theoretical standpoint a single relational interpreter for a Turing-complete language is sufficient: indeed, any other interpreter can be turned into a relational one just by implementing it in a language, for which relational interpreter already exists. However, the overhead of additional interpretation level can easily make this solution impractical. The standard way to tackle the problem is partial evaluation or specialization [Jones et al. 1993]. A specializer $\text{spec}_M$ for a language $M$ for any program $p^M$ in this language and its partial input $i$ returns some program which, being applied to the residual input $x$, works exactly as the original program on both $i$ and $x$:

$$\forall x : \llbracket \text{spec}_M(p^M, i) \rrbracket_M(x) = \llbracket p^M \rrbracket_M(i, x).$$

If we apply a specializer to an interpreter and a source program, we obtain what is called the first Futamura projection [Futamura 1971]:

$$\forall i : \llbracket \text{spec}_M(\text{eval}_L^M, p^L) \rrbracket_M(i) = \llbracket \text{eval}_L^M \rrbracket_M(p^L, i).$$

Here we added an upper index $M$ to $\text{eval}_L$ to indicate that we consider it as a program in the language $M$. In other words, the first Futamura projection specializes an interpreter for a concrete program, delivering the implementation of this program in the language of interpreter implementation. An important property of a specializer is Jones-optimality [Jones et al. 1993], which holds when it is capable to completely eliminate the interpretation overhead in the first Futamura projection. In our case $M = \text{miniKanren}$, from which we can conclude that in order to eliminate the interpretation overhead we need a Jones-optimal specializer for miniKanren. Although implementing a Jones-optimal specializer is not an easy task even for simple functional languages, there is a Jones-optimal specializer for a logical language [Leuschel et al. 2004], but not for miniKanren.

The contribution of this paper is as follows:

- We demonstrate the applicability of relational programming and, in particular, relational interpreters for the task of turning verifiers into solvers.
- To obtain a relational verifier from a functional specification we apply relational conversion [Byrd 2009; Lozov et al. 2018] — a technique which for a first-order functional program directly constructs its relational counterpart. Thus, we introduce a number of new relational interpreters for concrete search problems.
Relational Interpreters for Search Problems

- We employ supercompilation in the form of conjunctive partial deduction (CPD) [De Schreye et al. 1999] to eliminate the redundancy of a generic search algorithm caused by partial knowledge of its input.
- We give a number of examples and perform an evaluation of various solutions for the approach we address.

Both relational conversion and conjunctive partial deduction are done in an automatic manner. The only thing one needs to specify is the known arguments or the execution direction of a relation.

As concrete implementation of miniKanren we use OCANREN [Kosarev and Boulytchev 2016] — its embedding in OCAML; we use OCAML to write functional verifiers; our prototype implementation of conjunctive partial deduction is written in HASKELL.

The paper is organized as follows. In Section 2 we give a complete example of solving a concrete problem — searching for a path in a graph, — with relational verifier. Section 3 recalls the cornerstones of relational programming in miniKanren and the relational conversion technique. In Section 4 we describe how conjunctive partial deduction was adapted for relational programming. Section 5 presents the evaluation results for concrete solvers built using the technique in question. The final section concludes.

2 SEARCHING FOR PATHS IN A GRAPH WITH A RELATIONAL VERIFIER

In this section we demonstrate how to solve a concrete problem of searching for paths in a directed graph with a relational verifier. A directed graph is a tuple \((N, E, start, end)\), where \(N\) is a finite set of nodes, \(E\) is a finite set of edges, functions \(start, end : E \rightarrow N\) return a start and an end nodes for a given edge respectively. A path in a directed graph is a sequence:

\[ \langle n_0, e_0, n_1, e_1, \ldots, n_k, e_k, n_{k+1} \rangle \]

such that

\[ \forall i \in \{0 \ldots k\} : n_i = start(e_i) \text{ and } n_{i+1} = end(e_i). \]

The problem of searching for paths in a graph is to find a set \(\{p \mid p \text{ is a path in } g\}\), where \(g\) is a graph. There are many concrete algorithms which search for paths in a graph. Implementing any of them involves determining in which way to traverse the graph, how to ensure one does not get stuck exploring a cycle in the graph (a cycle is a path in the graph of form \(\langle n_0, e_0, \ldots, n_k, e_k, n_0 \rangle\)), how to ensure one path is not processed multiple times, and so on. A much easier task is to implement a simple verifier, which checks if a sequence is indeed a path in a graph, and generate the path searching routine from it by the relational conversion.

Below is the implementation of the verifier "isPath". This function takes as an input a list of nodes "ns" and a graph "g". We represent the graph as a list of edges, stipulating there are no parallel edges. Each edge \(e\) is represented as a pair of nodes \((n, m)\), where \(n = start(e)\), \(m = end(e)\). Given \(ns = [n_0, \ldots, n_{k+1}]\) and a graph \(g = [e_0, \ldots, e_j]\), the function returns true, if \(\exists 0 \ldots k\) such that \(\langle n_0, e_0, n_1, e_1, \ldots, e_k, n_{k+1} \rangle\) is a path in \(g\).

```haskell
let rec isPath ns g =
  match ns with
  | x1 :: x2 :: xs -> elem (x1, x2) g && isPath (x2 :: xs) g
  | [] -> true

let rec elem e g =
  match g with
  | [] -> false
  | x :: xs -> if eq e x then true else elem e xs
```

The function "elem" checks if an edge "e" exists in the graph "g". We omit the definition of equality check for edges "eq", since it is trivial to implement and is not relevant for the example.
We stipulate that a path must include at least two nodes, since searching for shorter paths is trivial. Line 3 of the “isPath” definition checks that the first two nodes of the list form an edge of the graph. Then it checks that what is left after deleting the first node from the list is still a path in the graph. Line 4 may come off a little counterintuitive, since it states that a path which includes a single arbitrary node is in the input graph. However we only execute this branch by a recursive call of “isPath”, which only happens after we have already ensured with the call to the “elem” function that the said node is in the graph.

The relational conversion of the verifier function “isPath” generates a relation “isPath” defined for a path “ns”, a graph “g” and a boolean value “res”, which is true if “ns” is a path in the graph “g” and false otherwise. The function “elem” is transformed into a relation “elem” defined for an edge “e”, a graph “g” and a boolean value “res”, which is true if “e” is an edge in the graph “g” and false otherwise. The result of the relational conversion of the functions “isPath” and “elem” is presented below.

5 let rec elem e g res = conde [
6   (g ≡ nil () ∧ res ≡ ↑false);
7   (fresh (x xs resEq) (g ≡ x % xs) ∧
8      (eq e x resEq) ∧
9      (conde [resEq ≡ ↑true ∧ res ≡ ↑true);
10     (conde [false ∧ elem e xs res]))))]
11
12 let rec isPath ns g res = conde [
13   (fresh (el) (ns ≡ el % nil ()) ∧
14      (res ≡ ↑true));
15   (fresh (x1 x2 xs resElem resIsPath) (ns ≡ x1 % (x2 % xs)) ∧
16      (elem (pair x1 x2) g resElem) ∧
17      (isPath e (x2 % xs) g resIsPath) ∧
18      (conde [false ∧ res ≡ ↑false);
19      (conde [true ∧ res ≡ resIsPath]))])
20
21 Here we use the syntax of OCANREN. A new relation is defined as a recursive function with the keywords “let rec”. The body of the relation is a goal created with the following goal constructors.

- Disjunction \( g_1 \lor g_2 \), where \( g_1, g_2 \) — some goals. The two goals are evaluated independently and their results are combined.
- Disjunction of goal list \( \text{conde} \) \( [g_1; \ldots; g_n] \), where \( g_1; \ldots; g_n \) — some goals.
- Conjunction \( g_1 \land g_2 \), where \( g_1, g_2 \) — some goals. The goal \( g_2 \) is evaluated only if the evaluation of \( g_1 \) succeeded; the evaluation of \( g_2 \) uses the results of \( g_1 \).
- Syntactic unification \( t_1 \equiv t_2 \), where \( t_1, t_2 \) — some terms. Unification is a basic goal constructor. If \( t_1 \) and \( t_2 \) can be unified, the goal is considered successful and failed otherwise.
- Relation call \( r^n t_1 \ldots t_n \) where \( r^n \) is a name of some \( n \)-ary relation, and \( t_i \) are terms.
- To introduce fresh variables into scope, one should use \( \text{fresh} (\overline{x}) \) \( g \), where \( \overline{x} \) is a list of variable names.
Besides goal constructors we use some syntactic sugar for values and lists. \( \uparrow \) is used to transform a value into a logic value. The empty list is represented as \( \text{”nil ()”} \), and to construct a new list from a value \( \text{”h”} \) and a list \( \text{”t”} \) we use \( \text{”h \% t”} \). A tuple of \( \text{”x”} \) and \( \text{”y”} \) is created with \( \text{”pair x y”} \).

Regrettably, this relational interpreter suffers from poor performance. Query \( \text{”isPath” q <graph> \text{”true”} } \) for path searching takes more than ten minutes even for graphs of 5 nodes. This is somewhat expected, considering that the relational conversion generates a relation which can be used for many different queries, which is excessive when any particular query is in question. This is, of course, not a desirable behaviour. Fortunately, further transformation of the relation can improve the performance.

For example, if we consider a query \( \text{”isPath” q <graph> \text{”true”} } \), we can simplify lines 18 through 24 of its definition. First, we notice that, having \( \text{”res”} \) be equal to \( \text{”\uparrow true”} \), we can safely remove the disjunct in line 23, after what the whole \( \text{”conde”} \) becomes unnecessary and can be removed. After moving the unifications for \( \text{”resElem”} \) and \( \text{”resIsPath”} \) to the top level, we get the following equivalent definition of the \( \text{”isPath”} \) relation.

The specialized version of the relation is much more performant than the original one. Before, searching paths of length 5 took more than 10 minutes while the specialized version finds paths of length 10 in the graph with 100 edges in a few seconds.

This transformation can be performed automatically with conjunctive partial deduction. The result of partially deducting the \( \text{”isPath” p q \text{”true”} } \), where \( \text{”p”} \) and \( \text{”q”} \) are fresh variables is about 40 lines of code long and it has the same performance as the manually transformed relation. We omit the transformed program because of the space concerns, but it can be found in the repository\(^1\).

3 RELATIONAL CONVERSION

In this section we describe how the relational conversion in the form of \texttt{unnesting} [Byrd 2009] is done. Unnesting constructs a relational program by a first-order functional program.

First, a new variable for every subexpression is introduced with the \texttt{let}-expression. Then, all pattern matching and if-expressions are translated into disjunctions, in which unifications are generated for the patterns. Free variables are introduced into scope with the \texttt{fresh}. Every \( n \)-ary function becomes \( (n + 1) \)-ary relation with the last argument being unified with \( \text{”\uparrow true”} \), so further specialization is still possible.

\begin{verbatim}
let rec isPath n s r = conde [ (fresh (el)) (ns ≡ el % nil ()) \& (res ≡ ↑true)).
(fresh (x1 x2 xs resElem resIsPath) (resElem ≡ ↑true) \& (resIsPath ≡ ↑true) \&
(ns ≡ x1 % (x2 % xs)) \& (elem (pair x1 x2) g resElem) \&
(isPath (x2 % xs) g resIsPath))]
\end{verbatim}

The example of unnesting is shown in Fig. 1. The input functional program is presented in Fig. 1a. The result of introducing fresh variables for subexpressions is in Fig. 1b. The relational program before the conjuncts are reordered is shown in Fig. 1c and the result of the unnesting is presented in Fig. 1d.

\(^1\)https://github.com/Lozov-Petr/miniKanren-2019-Relational-Interpreters-for-Search-Problems
let rec append a b = match a with |
  | []    → b
  | x :: xs → x :: append xs b

let rec append \(o\) a b c = (a \(\equiv\) []) \(\wedge\) b \(\equiv\) c \(\lor\)
(fresh (x xs q) (a \(\equiv\) x :: xs) \(\wedge\)
  (append \(o\) xs b q) \(\wedge\)
  (c \(\equiv\) x :: q))

Fig. 1. Example of unnesting

Note, that the unnesting has limitations: it does not support higher-order functions and partial application. A more general method of translation which does not impose the same limitations was developed [Lozov et al. 2018]. Unfortunately, it uses higher-order relations which are not currently supported in conjunctive partial deduction, so we use unnesting.

The forward execution of the relation mimics the execution of the function from which it was constructed by relational conversion. This makes forward execution quite efficient, to the detriment of the execution in the backwards direction. The unnesting can be modified to improve the performance of backward execution. Let us consider the conversion of a functional conjunction "\(f \_1 \_x \_1 \&\& f \_2 \_x \_2\)".

\[
\lambda \text{res} \rightarrow \\
fresh (p) (\ (f \_1 \_x \_1 \ p) \wedge \\
\conde [ \\
(p \equiv \upfalse \wedge \res \equiv \upfalse); \\
(p \equiv \uptrue \wedge f \_2 \_x \_2 \res)])
\]

Mimicking the function evaluation, the forward execution of this code first computes "\(f \_1 \_x \_1\)". If it fails, then the result "\res" is unified with "\false", otherwise the second conjunct "\(f \_2 \_x \_2\)" is executed and its result is unified with the result. This strategy is not efficient in the backward direction, when we know what "\res" is. The following relation is much more performant when executed in the backward direction:

\[
\lambda \text{res} \rightarrow \\
\conde [ \\
(res \equiv \upfalse \wedge f \_1 \_x \_1 \upfalse); \\
(f \_1 \_x \_1 \uptrue \wedge f \_2 \_x \_2 \res)]
\]

In particular, if "\res \equiv \uptrue", both conversions execute "\(f \_2 \_x \_2 \res\)" but when the first conversion computes "\(f \_1 \_x \_1 \ p\)" with fresh "\p", the second executes "\(f \_1 \_x \_1 \uptrue\)". Using the second conversion is enough to significantly
increase the performance in the backward direction. For example, the path search takes several minutes if the first conversion strategy is used, whereas it finishes in less than a second in the second case.

Choosing the second conversion strategy comes with a price for the forward execution. Instead of executing “$f_1 x_1 p$”, where “p” is fresh, the second strategy executes both “$f_1 x_1 \uparrow \text{false}$” and “$f_1 x_1 \uparrow \text{true}$”. In the worst case scenario, when the execution of “$f_1$” does not depend on the last argument, it doubles the number of executions of “$f_1$”.

To sum up, by choosing different strategies of the relational conversion we can achieve significant performance improvement. There is no single right way of doing the conversion which improves the performance of the execution in every possible direction. Choosing a strategy per each relation and each direction manually is not feasible, but it can be achieved with a fully-automatic program transformation, such as conjunctive partial deduction.

4 CONJUNCTIVE PARTIAL DEDUCTION

Specialization [Jones et al. 1993] is a natural way to tackle the problem of redundant computations when a part of the input is known. A fully-automatic specialization technique developed in the domain of logic programming is called partial deduction [Komorowski 1982; Lloyd and Shepherdson 1991]. It is related to the supercompilation of functional languages [Glück and Sørensen 1994; Turchin 1986]. The particular flavour of the partial deduction we are interested in is called conjunctive partial deduction [De Schreye et al. 1999]. As opposed to the partial deduction, conjunctive partial deduction handles conjunctions of atoms, thus being able to perform such optimizations as tupling [Hu et al. 1997] and deforestation [Wadler 1988]. Below we demonstrate by example the features of conjunctive partial deduction.

Deforestation is a program transformation which gets rid of intermediate data structures. The following example demonstrates deforestation. Consider a goal “$\text{append}^o xs ys ts \land \text{append}^o ts zs rs$” (note the shared “ts”), where “$\text{append}^o x y xy$” describes concatenation, “$\text{nil} ()$” constructs the empty list, and “$h \% t$” constructs a new list from the value “h” and another list “t” (similarly to “cons” in Scheme and “::” in OCaml).

$$\text{let rec append}^o x y xy = \text{conde} \left[ \begin{array}{l} (x \equiv \text{nil} ()) \land xy \equiv y) \land \end{array} \right.$$ 

$$(\text{fresh} (h t ty) (\begin{array}{l} (x \equiv h \% t) \land \end{array} \begin{array}{l} (xy \equiv h \% ty) \land \end{array} \begin{array}{l} (\text{append}^o t y ty))) \right]$$

This goal concatenates three lists: “xs”, “ys”, “zs”, constructing an intermediate list “ts”. During the execution of this goal, elements of the list “xs” are examined twice: first when “ts” is constructed, and then when the result “rs” is constructed. What is worse, “ts” is only constructed to be immediately deconstructed. Deforestation gets rid of “ts” in this example.

A better program would be such that does not construct “ts” at all. Such a program be generated from the original definition by conjunctive partial deduction and is shown below:

$$\text{let rec doubleAppend}^o xs ys zs rs = \text{conde} \left[ \begin{array}{l} (xs \equiv \text{nil} ()) \land \text{append}^o ys zs rs) \land \end{array} \right.$$ 

$$(\text{fresh} (h t ts) (\begin{array}{l} (xs \equiv h \% t) \land \end{array} \begin{array}{l} (rs \equiv h \% ts) \land \end{array} \begin{array}{l} (\text{doubleAppend}^o t ys zs ts)))) \right]$$
Conjunctive partial deduction is also capable of *tupling*. This transformation makes sure that the same data structure is traversed once even if computing several results. The following example demonstrates such a case.

The goal \( \text{maxLength} \ o \ xs \ m \ l \) computes both the maximum value of the list \( \text{xs} \) and its length. The elements of the list are Peano numbers with \( \text{zero}() \) as the zero and \( \text{succ} \) as the successor function. The third argument \( b \) of the relation \( \text{le}^o \ x \ y \ b \) is \( \uparrow \text{true} \) if \( x \) is less or equal than \( y \), and \( \uparrow \text{false} \) otherwise. The relation \( \text{gt}^o \ x \ y \ b \) is similar to \( \text{le}^o \ x \ y \ b \), but it checks for \( x \) to be greater than \( y \).

```plaintext
let maxLength^o xs m l = max^o xs m ∧ length^o xs l

let rec length^o xs l = conde [
  (xs ≡ nil () ∧ l ≡ zero ());
  (fresh (h t m) (xs ≡ h % t ∧ l ≡ succ m ∧ length^o t m))]

let max^o xs m = max^1o xs (zero ()) m

let rec max^1o xs n m = conde [
  (xs ≡ nil () ∧ m ≡ n);
  (fresh (h t) (xs ≡ h % t) ∧
   (conde [
    (le^o h n ↑true ∧ max^1o t n m);
    (gt^o h n ↑true ∧ max^1o t h m)]))]

let rec le^o x y b = conde [
  (x ≡ zero () ∧ b ≡ ↑true);
  (fresh (x1) (x ≡ succ x1 ∧ y ≡ zero () ∧ b ≡ ↑false));
  (fresh (x1 y1) (x ≡ succ x1 ∧ y ≡ succ y1 ∧ le^o x1 y1 b))]

let rec gt^o x y b = conde [
  (x ≡ zero () ∧ b ≡ ↑false);
  (fresh (x1) (x ≡ succ x1 ∧ y ≡ zero () ∧ b ≡ ↑false));
  (fresh (x1 y1) (x ≡ succ x1 ∧ y ≡ succ y1 ∧ gt^o x1 y1 b))]
```

Execution of the goal \( \text{maxLength}^o \ xs \ m \ l \) leads to \( \text{xs} \) being traversed twice. There is a way to rewrite the program so that \( \text{xs} \) is traversed once, but this requires fusing together the definitions of \( \text{length}^o \) and \( \text{max}^o \), which either restricts code reuse, or leads to code duplication. A better way is to only fuse the definitions when it is needed, and do it automatically by employing tupling.

The desirable implementation of the \( \text{maxLength}^o \ xs \ m \ l \) relation is the following (the definitions of \( \text{gt}^o \) and \( \text{le}^o \) are left out for brevity). It can be achieved with conjunctive partial deduction as well:
let maxLength^o xs m l = maxLength^o 1 xs m (zero ( ) ) l

let rec maxLength^o xs m n l = conde [  
  (xs ≡ nil ( ) ∧ m ≡ n ∧ l ≡ zero ());  
  (fresh (h t l_1))  
    (xs ≡ h % t) ∧  
    (l ≡ succ l_1) ∧  
    (conde [  
      (le^o h n ∧ maxLength^o t m n l);  
      (gt^o h n ∧ maxLength^o t m h l)])]]

4.1 CPD for Prolog-like languages

Initially, conjunctive partial deduction was developed for Prolog-like languages. Conjunctive partial deduction partially evaluates goals, which are conjunctions of atoms, using two levels of control: local and global [Glück et al. 1996]. The global control determines which atoms are to be partially deduced. The local control — what the definitions for the atoms selected at the global control shall be. Both local and global control construct tree structures which represent the input program.

Local control constructs finite SLD-trees for conjunctions of atoms. The construction is guided with an unfold operator: it selects a literal from the leaf of the partially constructed SLD-tree and adds its resolvents as children at each step. Since, in general, SLD-trees are infinite, a decision to stop unfolding should be made at some point. There are several techniques for doing this, the most promising of them combine determinacy and either some well-founded or well-quasi order, such as homeomorphic embedding, or other measures.

Global control determines the set of the conjunctions for which partial SLD-trees are built. The important goal of the global control is to ensure termination. The termination is achieved with the abstraction. If there is a goal which is embedded into the current goal, it points to the possibility of nontermination. The embedding tells that there is a certain similarity between the two goals, and if a current goal keeps being processed, then their similar subpart will appear again and again, causing nontermination. Whenever the embedding goals are detected, the current goal is abstracted to remove the common subgoal from consideration.

When the partial deduction is done, the only thing left is to construct the residual program. The clauses are generated from a partial SLD-tree, one tree per conjunction at the global level. A conjunction is uniquely renamed to give a name for the predicate being defined. All free variables of the root of the tree become arguments of the predicate. For each non-failing path in the SLD-tree a clause is generated: a substitution collected along the path is substituted into the head of the clause, and the body is generated from what is in the leaf.

4.2 CPD for miniKanren

In this section we describe how we adapted conjunctive partial deduction for miniKanren. We describe the particular unfolding and generalization strategies as well as discuss how the conjunctive partial deduction had to be modified as a response to the differences between PROLOG and miniKanren.

4.2.1 Local Control. Goals in miniKanren are different from those in Prolog-like languages: besides conjunction, disjunction and relation calls, there are explicit unification and introduction of fresh variables. We normalize the input goal so that it was a disjunction of conjunctions of relation calls. To do so, we first pop all the fresh variables to the top level ("fresh (x) (p (x) ∧ fresh (y) (q(x) ∨ r(y, x)))" becomes "fresh (x y) (p(x) ∨ q(x) ∧ r(y, x))"). Then we transform the goal to be a disjunction of conjunctions of relation calls or unifications. All unifications in each conjunction are evaluated to some substitution (or the
conjunct is discarded, if some unification fails). The normalization allows us to only consider conjunctions of relation calls while doing conjunctive partial deduction.

The local control constructs the following tree structure which represents the goal:

```plaintext
type local_tree =
  Fail
  | Success of subst
  | Leaf of goal list * subst
  | Disj of local_tree list
  | Conj of local_tree * goal list
```

Leaf nodes can be either “Fail”, “Success” or “Leaf”. The “Fail” node is created whenever the evaluation of the current goal fails. When the current goal evaluates to some substitution, we create the “Success” node with this substitution. The last leaf node is called “Leaf”, it corresponds to some partially evaluated goal. This type of node contains a substitution which has been computed up to this point, and a residual goal. The goal in this type of node is then examined at the global level.

“Disj” node corresponds to a disjunction in a goal: its children are the local control trees constructed for all disjuncts. The last type of nodes is a “Conj” node. It is a transient node, which keeps track of a conjunction being unfolded.

In general, unfolding replaces some of the relation calls with their bodies and partially evaluates them. The particular unfolding strategy we adhere to is the following. At each step only one relation call is replaced with its body: the leftmost selectable relation call. The selectable relation call is the one which does not embed any of its predecessors — goals which were unfolded in order to get the current goal. Embedding here is the modification of the homeomorphic embedding defined for the conjunctions of goals in conjunctive partial deduction literature [De Schreye et al. 1999]. Since using pure embedding to control unfolding leads to hideously big programs, we also allow only one non-deterministic unfold.

4.2.2 Global Control. The conjunctions in the “Leaf” nodes are processed at the global level. This step is responsible for the termination of the transformation. Generally speaking, the danger for nontermination arises whenever we encounter a subgoal which we have encountered before: processing the same thing will lead to itself over and over again. To break the vicious circle, one needs to stop unfolding the encountered subgoal, this is what abstraction serves for.

The simplest case here is when we come upon the goal which is equal up to variable renaming to any other goal at the global level. When this happens, we stop exploring the goal. This is called variant check in the literature, and it is done both at the global and the local control levels.

The more complicated case is when a subpart of the goal repeats. This case we test with the modification of the homeomorphic embedding relation (strict homeomorphic embedding), initially developed for conjunctions. A conjunction \( \overline{A} \) is considered embedded into a conjunction \( \overline{B} \) when there is an ordered subconjunction within \( \overline{A} \), each conjunct of which is embedded into the corresponding conjunct of \( \overline{B} \):

\[
\overline{A} = A_0 \land A_1 \land \cdots \land A_n \preceq B_0 \land B_1 \land \cdots \land B_m = \overline{B}, \text{ if } \exists \{i_0 \ldots i_m \mid \forall j. i_j < i_{j+1}\} : \forall j \in \{0 \ldots m\}. A_{i_j} \preceq B_j
\]

A single conjunct is embedded into another \( (A_i \preceq B_j) \) when the following relation holds and \( A_i \) is not a strict instance of the second one \( B_j \):

\[
X \preceq Y, \text{ where } X \text{ and } Y \text{ are variables}
\]

\[
f(x_0, x_1, \ldots, x_n) \preceq f(y_0, y_1, \ldots, y_n) \iff \forall i \in \{0 \ldots n\}. x_i \preceq y_i
\]

\[
f \preceq g(y_0, y_1, \ldots, y_m) \iff \exists i \in \{0 \ldots m\}. f \preceq y_i
\]
This check determines two major causes of the growth within the conjunctions. The conjunction can grow in some argument of a relation call or the number of conjunctions itself can grow. To mitigate the first source of the growth, the bigger conjunction can be replaced with a most specific generalization of the two conjunctions. Otherwise we need to split the embedded subconjunction from the rest and start processing them separately. This process called abstraction removes the subconjunctions which cause potential nontermination, and what is left should indeed be processed further.

4.2.3 Residualization. After the transformation is finished, a residual program is constructed from the global control tree. A relation definition is generated for each conjunction at the global level (this is done with the renaming step of the original conjunctive partial deduction). First, a unique name is given for each conjunction. Then free variables of the conjunction are collected to become the arguments of the relation: the constructors and constants are omitted (for example \( f \ x \ (\text{succ} \ y) \land g \ (\text{zero} \ () \ z) \) becomes \( f_G \ x \ y \ z \)). The body of the definition is generated from the local control tree which corresponds to the conjunction under consideration. The body is formed as a disjunction of conjunctions for the non-failure nodes of the local control tree. A computed substitution is transformed into a conjunction of unifications. Suitable definitions are chosen for a goal in a leaf, and the conjunction of their applications is generated. As a final step we perform redundant argument filtering as described in [Leuschel and Sørensen 1996], and introduce fresh variables where necessary.

5 EVALUATION

In this section we present an evaluation of the proposed approach. We have implemented several relational interpreters for different search problems which can be found in the repository mentioned before. Some of the simpler interpreters demonstrate good performance for different directions on their own and for them CPD transformation is not needed. Thus, we will focus on two search problems which are more complex: searching for a path in a graph and searching for a unifier [Baader and Snyder 2001] of two terms. For each problem we compare four programs.

1. The solver generated by the unnesting alone.
2. The solver generated by the unnesting strategy aimed at backward execution.
3. The solver generated by the unnesting and then specialized by conjunctive partial deduction for the backward direction.
4. The interpretation of the functional verifier with the relational interpreter implemented in Scheme [Byrd et al. 2017].

First, let us compare the performance of the solvers for the path searching problem. The implementation of the functional verifier for this problem is described in Section 2. We ran the search on a graph with 20 nodes and 30 edges, consequentially searching for paths of the length 5, 7, 9, 11, 13, and 15. We averaged the execution times over 10 runs of the same query. We limited the execution time by 300 seconds, and if the execution time of some query exceeded the timeout, we put “>300s” in the result table and did not request the execution of queries for longer paths. The results are presented in Table 1.

We can conclude that the execution time increases with the length of the path to search, which is expected, since with the length of the path the number of the subpaths to be explored is increasing as well. The solver generated by the unnesting alone and the interpretation with the relational interpreter demonstrate poor performance. The first one is due to its inherently inefficient execution in backward direction, while the second suffers from the interpretation overhead. Both the unnesting aimed at the backward execution and the solver automatically transformed with conjunctive partial deduction show good performance. Conjunctive partial deduction performs more thorough specialization, thus producing more efficient program.

Now let us consider the problem of finding a unifier of two terms which have free variables. A term is either a variable \( (X, Y, \ldots) \) or some constructor applied to terms \( (\text{nil}, \text{cons}(H, T), \ldots) \). A substitution
Table 1. Searching for paths in the graph

<table>
<thead>
<tr>
<th>Path length</th>
<th>5</th>
<th>7</th>
<th>9</th>
<th>11</th>
<th>13</th>
<th>15</th>
</tr>
</thead>
<tbody>
<tr>
<td>Only conversion</td>
<td>0.01s</td>
<td>1.39s</td>
<td>82.13s</td>
<td>&gt;300s</td>
<td>—</td>
<td>—</td>
</tr>
<tr>
<td>Backward oriented conversion</td>
<td>0.01s</td>
<td>0.37s</td>
<td>2.68s</td>
<td>2.91s</td>
<td>4.88s</td>
<td>10.63s</td>
</tr>
<tr>
<td>Conversion and CPD</td>
<td>0.01s</td>
<td>0.06s</td>
<td>0.34s</td>
<td>2.66s</td>
<td>3.65s</td>
<td>6.22s</td>
</tr>
<tr>
<td>Scheme interpreter</td>
<td>0.80s</td>
<td>8.22s</td>
<td>88.14s</td>
<td>191.44s</td>
<td>&gt;300s</td>
<td>—</td>
</tr>
</tbody>
</table>

Table 2. Searching for a unifier of two terms

<table>
<thead>
<tr>
<th>Terms</th>
<th>f(X, a)</th>
<th>f(a % b % nil, c % d % nil, L)</th>
<th>f(X, X, g(Z, t))</th>
<th>f(a, X)</th>
<th>f(X % XS, YS, X % ZS)</th>
<th>f(g(p, L), Y, Y)</th>
</tr>
</thead>
<tbody>
<tr>
<td>Only conversion</td>
<td>0.01s</td>
<td>&gt;300s</td>
<td>&gt;300s</td>
<td>0.01s</td>
<td>0.11s</td>
<td>2.26s</td>
</tr>
<tr>
<td>Backward oriented conversion</td>
<td>0.01s</td>
<td>0.06s</td>
<td>0.34s</td>
<td>0.01s</td>
<td>0.07s</td>
<td>0.90s</td>
</tr>
<tr>
<td>Conversion and CPD</td>
<td>0.04s</td>
<td>5.15s</td>
<td>&gt;300s</td>
<td>0.04s</td>
<td>5.15s</td>
<td>&gt;300s</td>
</tr>
</tbody>
</table>

maps a variable to a term. A unifier of two terms \( t \) and \( s \) is a substitution \( \sigma \) which equalizes them: \( t\sigma = s\sigma \) by simultaneously substituting the variables for their images. For example, a unifier of the terms \( \text{cons}(42, X) \) and \( \text{cons}(Y, \text{cons}(Y, \text{nil})) \) is a substitution \( \{ X \mapsto \text{cons}(42, \text{nil}), Y \mapsto 42 \} \).

We implemented a functional verifier which checks if a substitution equalizes two input terms. We represent a variable name as a unique Peano number. A substitution is represented as a list of terms, in which the index of the term is equal to the variable name to which the term is bound, so the substitution \( \{ X \mapsto \text{cons}(42, \text{nil}), Y \mapsto 42 \} \) is represented as a list "\[ \text{cons}(42, \text{nil}); 42] \". The verifier returns true if the input terms can be unified with the candidate substitution and false otherwise.

As in the previous problem, we compare four solvers generated for the verifier described. With each solver, we search for a unifier of two terms and compare the execution times. The time comparison is presented in Table 2. The first two rows of each column contain two terms being unified. We use uppercase letters from the end of the alphabet (\( X, Y, \ldots \)) to denote variables, lowercase letters from the beginning of the alphabet (\( a, b, \ldots \)) to denote constants (constructors with zero arguments), identifiers which start from the lowercase letter (\( f, g, \ldots \)) to denote constructors.

Note, we compute a unifier for two terms, but not necessarily the most general unifier. We can implement the most general unification in \textsc{miniKanren}, but achieving the comparable performance using relational verifiers requires additional check that the unifier is indeed the most general. We are currently working on the implementation of such relational verifier.

Here four solvers compare to each other similarly to the previous problem: unnesting demonstrates the worst execution time, relational interpretation in Scheme is a little better, while unnesting aimed at backward execution and conjunctive partial deduction significantly improve the performance.

There exist pairs of terms, for which either of the solvers fails to compute a unifier under 300 seconds. The example of such terms is "\( f(A, B, C, A, B, C, D) \)" and "\( f(B, C, D, x(R, S), x(a, T), x(Q, b), x(a, b)) \)". This is caused by how general and declarative the verifier is: there is nothing in it to restrict the search space. We can modify the verifier with the additional check to ensure that there are no bound variables in the candidate unifier. This modification restricts the search space when there are many variables in the input terms. But it also changes the semantics of the initial verifier and, as a consequence, the solvers: only idempotent unifiers can be found.
To sum up, we demonstrated by two examples that it is possible to create problem solvers from verifiers by using relational conversion and conjunctive partial deduction. Currently conjunctive partial deduction improves the performance the most, as compared to interpreting verifiers with Scheme relational interpreter or doing relational conversion which is solely aimed at backward or forward execution.

6 Conclusion and Future Work

We have presented a way to construct a solver for a search problem from its solution verifier by first doing a relational conversion and then specializing the relation according to the desired execution direction by means of conjunctive partial deduction.

There are a few directions for future work.

Even if we generate a relation optimized for the particular direction, executing it with miniKanren still carries some overhead of interpretation. We believe that the best performance can be achieved by generating a functional program from the relation optimized for the particular direction. This way we can avoid interpretation overhead but still get the benefits of the approach. The second direction is to explore other specialization techniques besides conjunctive partial deduction which are better suited for miniKanren programs.

References


Constructive Negation for MiniKanren

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We present an extension of MINIKANREN with the negation operator based on the method of constructive negation. The idea of this method is to constructively build a stream of answers for the negated goal by collecting and negating individual answers to the positive version of the goal. As we demonstrate on the series of examples constructive negation suits to pure logical nature of MINIKANREN: the relations involving the negation operator still can be “run” in various directions.

CCS Concepts: Software and its engineering → Functional languages; Constraint and logic languages;

Additional Key Words and Phrases: relational programming, constructive negation, OCanren

1 INTRODUCTION

MINIKANREN [Byrd 2010; Friedman et al. 2005] is a minimalistic domain-specific language which brings features of logic programming into a host language. Typical MINIKANREN implementation introduces only a few operators: conjunction, disjunction, unification (which can be seen as equality constraint) and fresh variable introduction (existential quantification). Although this basis constitutes a Turing-complete language, in practice there are some cases when the availability of negative reasoning is desirable.

For example, consider the following problem. Suppose we want to write a program which removes the first occurrence of a given element from a list. In order to do that in MINIKANREN we have to first define a ternary relation remove which binds the desired element, original list, and the same list after the deletion. Substituting the first argument of the relation with some element e, the second argument with some list xs and the third argument with a free variable q gives us a goal which is a proof search procedure. When passed to the run function, the goal will produce a lazy stream of answers. Each answer is represented by substitution which binds free variables to some terms. In the case of remove, we would expect a single answer, which binds q to the list, equal to xs, except that the first occurrence of e is removed.

The code on Listing 1 demonstrates a possible implementation of remove. It consists of three disjuncts, which represent three different cases. First, if the original list is empty, then the resulting list should also be empty. If the original list is not empty and its head is equal to the given element, then the resulting list should be equal to the tail. Finally, in the case when the head is not equal to the given element, the resulting list should be equal to the original, from the tail of which the occurrence of the element e is deleted.

Given the definition of remove from Listing 1, the following query run (remove 2 [1;2;3] q) will wrongly return two answers: q = [1;3] and q = [1;2;3]. The redundant (and, indeed, incorrect) answer q = [1;2;3] arose because of the third disjunct from remove definition, which always succeeds and thus generates a copy of the original list. We can prevent this behavior using disequality constraint, yet another primitive, which some MINIKANREN implementations provide. Adding constraint x ≠ e to the third case makes all disjuncts mutually exclusive.

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miniKanren.org/workshop/2019/8-ART4
disjoint. With the fixed definition of remove (Listing 2), the query given above returns the single answer \( q = [1;3] \) as expected.

```ocaml
let remove e xs ys = (xs ≡ [] ∧ ys ≡ []) ∨ fresh (x xs') (x ≡ e ∧ xs ≡ x::xs' ∧ ys ≡ xs') ∨ fresh (x xs' ys') (xs ≡ x::xs' ∧ ys ≡ x::ys' ∧ remove e xs' ys')
```

Listing 1. A flawed definition of remove relation

```ocaml
let remove e xs ys = (xs ≡ [] ∧ ys ≡ []) ∨ fresh (x xs') (x ≡ e ∧ xs ≡ x::xs' ∧ ys ≡ xs') ∨ fresh (x xs' ys') (xs ≡ x::xs' ∧ ys ≡ x::ys' ∧ remove e xs' ys')
```

Listing 2. The correct definition of the remove relation

```ocaml
let remove p xs ys = (xs ≡ [] ∧ ys ≡ []) ∨ fresh (x xs') (p x ∧ xs ≡ x::xs' ∧ ys ≡ xs') ∨ fresh (x xs' ys') (x ≠ e ∧ xs ≡ x::xs' ∧ ys ≡ x::ys' ∧ remove e xs' ys')
```

Listing 3. A generalized version of the remove relation

A disequality constraint represents a very limited form of negation, which is often not sufficient. Imagine that we want to generalize remove, so that instead of taking an element, it takes a predicate \( p \) and removes the first element of the list which satisfies the predicate (Listing 3). In order to do that and avoid similar pitfalls as in our first attempt to define remove, we need to ensure that the third disjunct succeeds only when the predicate \( p \) fails on the head element of \( xs \). Thus we need a general negation operator. Unfortunately, none of the existing MiniKanren implementations provide this feature.

In the world of Prolog, negative reasoning is usually implemented using so-called negation as failure approach. Under this rule, a goal \( \neg p \) succeeds whenever \( p \) fails, and it fails whenever \( p \) succeeds. Unfortunately, negation as failure is unsound if the negated goal \( p \) contains free variables. For example, the query `run (q ≡ 0) ∧ \neg(q ≡ 1)` succeeds, although `run \neg(q ≡ 0) ∧ (q ≡ 0)` fails. The failure in the last case is due to the fact that at the time of negation execution the variable \( q \) is free, thus \( q ≡ 1 \) succeeds, and conversely \( \neg(q ≡ 1) \) fails. This example demonstrates that negation as failure goes against the philosophy of MiniKanren, which positions itself as a pure relational language without non-logical features.

In this work we present an implementation of the negation operator based on the method of constructive negation. It overcomes several shortcomings of negation as failure and provides a more expressive form of negative reasoning than disequality constraints. The idea of this method is to constructively build a stream of answers for the negated goal. It is achieved by first collecting all answers to the positive version of the goal and then negating them. The constructive negation approach while being an improvement over negation as failure still has its own drawbacks. For example, applying the negation operator to the goal that has infinitely many answers results in a non-terminating computation.

Although constructive negation was first proposed as an extension of Prolog, to the best of our knowledge our work is the first attempt to adapt it for MiniKanren. We have developed a prototype implementation for OCANREN [Kosarev and Boulytchev 2018], a MiniKanren dialect embedded in OCAML. Like the rest of
MiniKanren, our version of constructive negation is developed in a pure functional style using persistent data structures. This makes it different from the earlier implementations for Prolog.

Our paper is structured as follows. In Section 2 we give more examples of constructive negation usage. Section 3 describes our implementation. Next, Section 4 presents the evaluation of the negation on a series of examples. In Section 5 we discuss the limitations of our approach and directions for future work. Section 6 reviews related works. The final Section 7 concludes.

2 MOTIVATING EXAMPLES

In this section we give further motivation for adding the negation into relational programming. We present several examples of how the negation can be used.

2.1 Relational If-then-else

In the Prolog one can simulate the conditional if-then-else operator using so-called soft cut [Naish 1995]. The behavior of the soft cut \( c \rightarrow^* t ; e \) can be described as follows:

- if the goal \( c \) succeeds (i.e. produces at least one answer) then the result of \( c \rightarrow^* t ; e \) is equivalent to \( c \land t \);
- if the goal \( c \) fails (i.e. produces no answers at all) then the result of \( c \rightarrow^* t ; e \) is equivalent to \( e \).

The soft cut is an example of a non-relational feature. Such features usually do not compose well in the sense that they are sensitive to the order in which they appear in a program. For example, consider the goal \((c \rightarrow^* t ; e) \land g\), and assume that \( c \land g \) always fails regardless of the order of conjuncts. Then if \( c \) succeeds the result of the above goal will be equivalent to \( c \land t \land g \) and thus it will fail. Suppose we reorder the conjuncts as follows: \( g \land (c \rightarrow^* t ; e) \). Now the goal \( c \), when executed after \( g \), certainly fails, and thus the result of the whole goal will be equivalent to \( g \land e \) which does not fails necessarily. One can see that the simple reordering of subgoals in the program can lead to the different results.

With the help of constructive negation if-then-else can be simply expressed as follows:

\[
\text{let } \text{ifte } c \ t \ e = (c \land t) \lor (\neg c \land e)
\]

The behavior of if-then-else defined in this way subsumes the behavior of the soft cut. That is, every answer to the query \( \text{run } (c \rightarrow^* t ; e) \) is also an answer to the query \( \text{run } (\text{ifte } c \ t \ e) \). Moreover, ifte is not sensitive to the order of subgoals in the program.

2.2 Classical Implication and Universal Quantification

With the negation added to the language, one can easily express other connectives of the classical first-order logic, namely the implication and universal quantification\(^1\), using the well-known equivalences.

\[
\text{let } (\Rightarrow) : \text{goal } \rightarrow \text{goal } \rightarrow \text{goal } = \left( g_1 \rightarrow (\neg g_1 \lor (g_1 \land g_2)) \right)
\]

\[
\text{let } \forall : (\forall \ a \rightarrow \text{goal}) \rightarrow \text{goal } = \left( g \rightarrow \neg \text{fresh} \ (x) \ (\neg g \ x) \right)
\]

However, we should make a few remarks here. It is well known that the search implemented in conventional MiniKanren is complete, meaning that every answer to an arbitrary query will be found eventually. In the presence of constructive negation (and thus implication and universal quantification defined through negation)

\(^1\) It is easy to see that \( c \Rightarrow t \) is equivalent to ifte \( c \ t \ succ \)
the search becomes incomplete as we will later see. Moreover, constructive negation is computationally heavy and thus the double usage of it, as in the definition of \texttt{forall}, can be inefficient in some cases.

Despite all this trouble we have found that the above definitions are still useful. Some of the previous MiniKanren implementations introduced \textit{eigen} variables, adopted from \texttt{\lambda}Prolog [Miller and Nadathur 2012]. Eigen variables act as a universally quantified variables. Yet, to the best of our knowledge, there is no sound implementation of eigen variables with the support of disequality constraints. We observed that our implementation of universal quantifications through double negation works nicely with disequalities (we give some examples in the Section 4).

2.3 \textbf{Graph Unreachability Problem}

One of the classical examples of negation application in logic programming is a problem of checking whether one node of the graph is unreachable from the another [Przymusinski 1989]. The code on Listing 4 defines binary relation \texttt{edge}, which binds pairs of nodes in graph, connected by some edge, and binary relation \texttt{reachable}, which is nothing more than a transitive closure of the \texttt{edge} relation. Then the relation \texttt{unreachable} is simply negation of \texttt{reachable}.

\begin{verbatim}
let edge x y =
  (x, y) \equiv ('a', 'b') \lor
  (x, y) \equiv ('b', 'a') \lor
  (x, y) \equiv ('b', 'c') \lor
  (x, y) \equiv ('c', 'd')

let reachable x y =
  x \equiv y \lor
  fresh(z) (edge x z \land reachable z y)

let unreachable x y =
  \neg(reachable x y)

Listing 4. Unreachability in a graph
\end{verbatim}

Given this definition the query \texttt{run unreachable 'c' 'a'} will succeed. A knowledgeable reader might notice that constructive negation is not necessary in this case because negation as failure will deliver the same result. But the query \texttt{run unreachable 'c' q} will fail under negation as failure because of the free variable \texttt{q} which will appear under the negation. However constructive negation will succeed delivering the constraint \texttt{q \neq 'd'}.

2.4 \textbf{Unreachability in Labeled Transition Systems}

One can consider a special kind of graphs — \textit{labeled transition systems} [Baier and Katoen 2008]. Labeled transition system is defined by a set of states \( S \), a set of labels \( L \) and a ternary transition relation \( R \subseteq S \times L \times S \). By existential quantification over labels one can then obtain a binary relation. Taking its transitive closure gives the reachability relation. The negation of the reachability relation can be used to check that some state \( s' \) is not reachable from the initial state \( s \). The Listing 5 shows an encoding of an abstract labeled transition system in OCANREN.

Labeled transition systems are often used to describe the behavior of imperative languages. Although the naive encoding of transition relation in OCANREN with simple enumeration of reachable states is often not tractable for
checking reachability (or unreachability) in huge state spaces arising in practical imperative programs, it still can be used, for example, for the task of prototyping the semantics of such languages.

```ml
module type LTS = sig
  type state
  type label
  val transition : state -> label -> state -> goal
end

module LTSExt (T : LTS) = struct
  let reachable : T.state -> T.state -> goal =
    fun s s'' ->
      s ≡ s''
    ∨
    fresh (l s') ( T.transition s l s' ∧
                      reachable s' s''
     )

  let unreachable : T.state -> T.state -> goal =
    fun s s' ->
      ¬(reachable s s')
end
```

Listing 5. Unreachability in a labeled transition system

3 IMPLEMENTATION

In this section we present our implementation of constructive negation. We start with the general ideas behind the method (Section 3.1). We describe how the constructive negation behaves in concrete examples, starting from trivial ones and moving to more sophisticated. During this presentation, we will observe, that in order to implement constructive negation, we need a solver for universally quantified disequality constraints. We will show that such solver can be implemented on top of existing MiniKanren disequality solver with just a few modifications (Section 3.2). In the Section 3.3, we describe how the OCanren search can be extended to support constructive negation. We will also discuss how negation interacts with recursion and present the notion of stratification (Section 3.4).

3.1 General Ideas

Constructive negation is based on the following idea: given a goal \( \neg g \), one can construct an answer for this goal by collecting all answers to its positive version \( g \) and then taking their complementation. In order to do that, a notion of “negation” of an answer is needed. Since each answer can be matched to some logical formula [Przymusinski 1989; Stuckey 1991], a “negation” of an answer corresponds to the logical negation of this formula.
Example 1. Consider the goal \( \neg(q \equiv 1 \land r \equiv 2) \). Its positive version \( q \equiv 1 \land (r \equiv 2) \) has single answer, a substitution \( \{ q \mapsto 1, r \mapsto 2 \} \) which corresponds to the formula \( q = 1 \land r = 2 \). By negating this formula we obtain \( q \neq 1 \lor r \neq 2 \). This formula still can be represent by a single substitution \( \{ q \mapsto 1, r \mapsto 2 \} \). However, we now treat this substitution differently, as a disequality constraint.

Some MiniKanren implementations (including OCAnren) already have the support for disequality constraints. A programmer can use them with the help of \( \neq \) primitive, as we have seen in the remove example (Listing 2). Usually, the support for disequalities is implemented as follows.

- A current state is maintained during the search. The state consists of a substitution, which represents positive information, and a disequality constraint store, which represents negative information. A constraint store can be implemented simply as a list of substitutions, or as a more efficient data structure.
- Each time a subgoal of the form \( t \neq u \) is encountered during the search, its satisfiability in current substitution is checked. If it is satisfiable, then disequality is added to the constraint store.
- Each time a subgoal of the form \( t = u \) is encountered during the search, its satisfiability in current substitution is refined by the result of unification of terms \( t \) and \( u \). Then the satisfiability of disequality constraints is rechecked in the refined substitution.

Various optimizations can be applied to the scheme above. For example, there is no need to recheck every disequality in the store after each unification. We will not discuss these optimizations here, as they are irrelevant to our goals.

Unfortunately, as the next example illustrates, disequalities presented above are not sufficient to implement constructive negation.

Example 2. Consider a goal \( \neg(\text{fresh} \ (x) \ (q \equiv (x, x))) \), which states that \( q \) should not be equal to some pair of identical terms. The subgoal \( \text{fresh} \ (x) \ (q \equiv (x, x)) \) succeeds, delivering the substitution \( \{ q \mapsto (x, x) \} \). Because the variable \( x \) occurs under fresh, the corresponding formula is existentially quantified: \( \exists x. q = (x, x) \).

By the negation of this formula we obtain \( \forall x. q \neq (x, x) \). This formula differs from disequality formula from example 1 as it contains universally quantified variable \( x \).

Thereby, in order to support the negation of goals, containing fresh, we need to extend disequality constraint solver, so it can check the satisfiability of universally quantified disequality constraints in the form \( \forall x. t \neq u \) \(^2\) [Chan 1988; Stuckey 1991]. Later, in Section 3.2 we will show how it can be done, for now let us assume we have such a solver.

We took care about fresh under negation. It led us to a more complicated representation of the state. During the search we maintain a pair of a current substitution and a universally quantified disequality constraint store. But now an interesting question arises: is this representation closed under negation? If we perform negation one more time, will we obtain a finite number of states in a similar form?

Lucky for us, it is the case. In order to verify it, let us consider a logical formula which corresponds to the representation of the state:

\[
\exists \Xi. \left( \bigwedge_i (v_i = t_i) \land \bigwedge_j \forall y_j. \bigvee_k (w_{jk} \neq u_{jk}) \right) \tag{1}
\]

Here \( v_i \) and \( w_{jk} \) denote some variables, \( t_i \) and \( u_{jk} \) denote some terms. Existentially quantified variables \( \Xi \) correspond to the variables occurred under fresh. The left conjunct corresponds to the substitution, the right conjunct corresponds to the constraint store. The constraint store itself is represented as a conjunction of individual universally quantified disequalities. As we have seen in example 1, each disequality corresponds to a

\(^2\) \( \Xi \) notation denotes a vector of variables, \( t \) and \( u \) are terms that may or may not contain variables from \( \Xi \)
disjunction of individual inequalities over variables\(^3\). Besides existential variables \(x\) and universal variables \(y_j\), there are free variables \(z\) that may occur in the formula (such as variables \(q\) and \(r\) in the examples 1, 2).

By logical negation of the above formula, we get the following:

\[
\forall x_i \left( \bigwedge_{i} (u_i \neq t_i) \lor \bigwedge_{j} \exists y_i \bigwedge_k (w_{jk} = u_{jk}) \right) 
\]

(2)

The left disjunct, which corresponded to the substitution in the original formula, now became a disequality constraint. Looking at the right disjunct, we can see that each disequality has transformed into a substitution. However, there is one subtlety here. Variables \(w_{jk}\) may be universally quantified, that is \(w_{jk} \in \overline{x}\) for some \(j, k\). Moreover, the terms \(u_{jk}\) may contain universally quantified variables as well. \(\text{Vars}(w_{jk}) \subseteq \overline{x}\) for some \(j, k\). In the section 3.2 we will show, that each disjunct \(\exists y_i \bigwedge_k (w_{jk} = u_{jk})\) is either unsatisfiable or could be rewritten as \(\exists y_i \bigwedge_k (w_{jk}^* = u_{jk}^*)\), such that universally quantified variables do not occur among variables \(w_{jk}^*\) or in terms \(u_{jk}^*\). [Liu et al. 1999].

Taking this into account, we can rewrite the formula 2 as follow:

\[
(\bigwedge_{j} \exists y_i \bigwedge_k (w_{jk}^* = u_{jk}^*)) \lor \forall x_i (u_i \neq t_i) 
\]

(3)

In the obtained formula each disjunct corresponds to one state in the form, similar to the given in formula 1. In the left disjunct, each sub-disjunct corresponds to a substitution with an empty disequality constraint, the right disjunct corresponds to the single universally quantified disequality constraint with an empty substitution. Thereby, the proposed representation of states is closed under negation.

One can perform further manipulations on the formula 3. Given that equivalence \(a \lor \neg b = a \land b \lor \neg b\) holds in classical logic, we can rewrite formula 3 in the following way:

\[
(\exists x_i \bigwedge_{i} (u_i = t_i) \land \bigwedge_{j} \exists y_i \bigwedge_k (w_{jk}^* = u_{jk}^*)) \lor \forall x_i (u_i \neq t_i) 
\]

(4)

The latter transformation, while vacuous from the logical point of view, could improve the performance of the search in practice. It follows from the fact, that the subpart \(\exists x_i \bigwedge_{i} (u_i = t_i)\) of the formula extends each substitution with additional mappings, thus delivering more positive information. If the negation constitutes a subpart of some larger goal, this positive information could lead to the earlier failure during the search.

Finally let us consider the negation in general case. A goal can be matched to a logical formula in the following way. Each answer to the goal corresponds to the state which itself corresponds to a logical formula in the form similar to one in formula 1. Goal can have multiple answers (even an infinite number). In the corresponding logical formula these answers will be connected by the disjunction:

\[
\bigwedge_{n} \left( \exists x_{ni} \bigwedge_{i} (u_{ni} = t_{ni}) \land \bigwedge_{j} \exists y_{nj} \bigwedge_k (w_{njk} \neq u_{njk}) \bigg) \right) 
\]

(5)

The negation of this formula after an application of the transformations described above will become:

\[
\bigwedge_{n} \left( \exists x_{ni} \bigwedge_{i} (u_{ni} = t_{ni}) \land \bigwedge_{j} \exists y_{nj} \bigwedge_k (w_{njk}^* = u_{njk}^*) \lor \bigwedge_{l} (u_{ni} \neq t_{ni}) \right) 
\]

(6)

\(^3\)We will give further explanation in section 3.2
In this way, we have obtained the result of the negation of the goal as a conjunction of the negation of individual answers to the positive version of the goal. However, one of the pitfalls of this construction is that the process will not terminate if the positive version of the goal has an infinite number of answers. Thus, in the general case, it makes the OCanren search incomplete for the goals involving negation.

3.2 Constraint Solver, Formally

In this section we will formally define the satisfiability of universally quantified disequality constraints and quantified equalities, mentioned in section 3.1. We will also present a simple decision procedure for satisfiability checking. To do that we need a rather standard notions of terms, substitutions, unifiers, etc. For the sake of completeness we give these definitions here. We also mention a standard unification algorithm as a decision procedure for equality constraints. This view of unification bridges the gap between the conventional and constraint logic programming.

Let us start with definitions of terms and substitutions.

Definition 1. Given an infinite set of variables $V$ and a finite set of constructor symbols $C = \{ C^i_n \}_i$, each with associated arity $n_i$, the set of terms $T$ is inductively defined as follow:

- $\forall v \in V, v \in T$ — every variable is a term;
- $\forall c_k \in C, \forall t_1 \ldots t_k \in T\cdot c_k(t_1, \ldots, t_k) \in T$ — every application of $k$-ary constructor symbol to $k$ terms is a term.

From now on we will assume that terms are untyped (unordered) and that there exists an infinitely many constructors of any arity.

Definition 2. Two terms $t$ and $u$ are syntactically equal, denoted as $t = u$, iff either

- $t = u$ and $s = v$ for some variable $v$;
- $t = c_k(t_1, \ldots, t_k), s = c_k(s_1, \ldots, s_k)$ and $\forall i \in \{1..k\}. t_i = s_i$.

Definition 3. A substitution $\sigma$ is a function from variables to terms: $\sigma : V \rightarrow T$, s.t. $\sigma(x) \neq x$ only for a finite number of variables. Every substitution $\sigma$ can be represented as a finite list of pairs $\{ x_1 \mapsto t_1, \ldots, x_n \mapsto t_n \}$. By $\text{dom}(\sigma)$ we denote the set $\{ x_1, \ldots, x_n \}$ and by $\text{codom}(\sigma)$ we denote the set $\{ t_1, \ldots, t_n \}$. We denote empty substitution as $\bot$. We also extend the set of substitutions defined above with the one additional element $\perp$.

Definition 4. A substitution can be applied to a term. The result of an application of $\sigma$ ($\sigma \neq \perp$) to $t$, written as $t\sigma$, is a term defined in the following way:

- $x\sigma \triangleq \sigma(x)$;
- $c_k(t_1, \ldots, t_k)\sigma \triangleq c_k(t_1\sigma, \ldots, t_k\sigma)$.

The result of the application of $\perp$ to any term is undefined.

Lemma 1. Given two substitutions $\sigma$ and $\theta$, if $\forall v \in V. \sigma(v) = \theta(v)$ then $\forall t. t\sigma = u\theta$.

Proof. Can be proved by the induction on $t$. $\square$

We are now ready to define the the satisfiability of equality constraint.

Definition 5. Equality constraint $t \equiv u$ is

- satisfiable if $\exists \sigma. t\sigma = u\sigma$; such $\sigma$ is called a unifier of $t$ and $u$;
- unsatisfiable otherwise.

Next, we show that the standard unification algorithm can be seen as a decision procedure for checking satisfiability of equality constrains. Before that we need to introduce several definitions.
Definition 6. A term $t$ is subsumed by the term $u$, denoted as $t \sqsubseteq u$, iff $\exists \sigma. \ t = \sigma u$. If $t \sqsubseteq u$ we will also say that $t$ is a more specific term than $u$, or $u$ is a more general term than $t$.

Definition 7. A substitution $\sigma$ is subsumed by a substitution $\tau$, denoted as $\sigma \sqsubseteq \tau$, iff $\forall t. \ t\sigma \sqsubseteq t\tau$. If $\sigma \sqsubseteq \tau$ we also say that $\sigma$ is a more specific substitution than $\tau$, or $\tau$ is a more general substitution than $\tau$.

Definition 8. Given terms $t$ and $u$ their unifier $\sigma$ is called the most general unifier, iff for every other unifier $\tau \sigma \neq u\sigma$. □

Theorem 1. Given terms $t$ and $u$ they are either not unifiable (meaning that $\forall \sigma. \ t\sigma \neq u\sigma$), or there exists their most general unifier.

Proof. Proof of this statement can be found in [Robinson et al. 1965]. Proof of the termination and correctness of the unification algorithm, used by the most MiniKanren implementations, can be found in [Kumar and Norrish 2010]. □

From now on we will denote the most general unifier of two terms $t$ and $u$ as $\text{mgu}(t, u)$. In case of $t$ is not unifiable with $u$, we assume that $\text{mgu}(t, u) = \bot$.

Remark 1. Note we can associate with an equality constraint $t \equiv u$ a logical first-order formula $t = u$. Additionally, we can associate with each substitution a logical first-order formula by the following rules:

- empty substitution $\top$ is associated with the truth constant $\top$;
- $\bot$ is associated with the falsity constant $\bot$;
- $\{x_1 \mapsto t_1, \ldots, x_n \mapsto t_n\}$ is associated with the formula $x_1 = t_1 \land \cdots \land x_n = t_n$.

Now we can have yet another view on unification. We can say that giving the problem of deciding satisfiability of the formula $t = u$, a unification algorithm reduces it to checking satisfiability of a simpler formula, which corresponds to a substitution. Such a formula is either trivially unsatisfiable (in the case of $\bot$) or trivially satisfiable.

Later on we will need the notion of idempotent substitution and idempotent unifier.

Definition 9. Substitution $\sigma$ is idempotent iff $\forall t. \ t\sigma\sigma = t\sigma$

Lemma 2. If two terms are unifiable, there exists their idempotent unifier.

Proof. For the proof of this statement (for the case of unification algorithm, used in MiniKanren), we refer an interested reader to [Kumar and Norrish 2010]. □

We are ready to move on to disequality constraints. We start with the regular (not quantified) disequations.

Definition 10. A disequality constraint $t \not\equiv u$ is

- satisfiable if $\exists \sigma. \ t\sigma \neq u\sigma$;
- unsatisfiable otherwise.

Next lemma gives us a simple decision procedure for checking satisfiability of disequations.

Lemma 3. Disequality constraint $t \not\equiv u$ is

- satisfiable if $\text{mgu}(t, u) \neq \top$;
- unsatisfiable otherwise.

Proof. Let $\theta = \text{mgu}(t, u)$. Let us first show that if $\theta = \top$ then disequality is unsatisfiable. By the definition of $\top$ we have $t\theta = t$ and $u\theta = u$, by the definition of unifier $t\theta = u\theta$, and thus $t = u$. From that, it is easy to show that $\forall \sigma. \ t\sigma = u\sigma$, which means that the disequality is unsatisfiable according to the definition 10. If $\theta \neq \top$ then there exists a substitution $\sigma$, s.t. $\theta \sqsubseteq \sigma$ (e.g. $\sigma = \top$). Since $\theta$ is most general unifier, and $\sigma$ is more general that $\theta$, then $\sigma$ is not a unifier, and thus $t\sigma \neq u\sigma$. □
Given lemma 3, the satisfiability of constraint \( t \neq u \) can be checked easily. One need to compute \( \text{mgu}(t, u) \) and if it is not an empty substitution, then constraint is satisfiable.

**Remark 2.** Given disequality constraint \( t \neq u \), a substitution \( \text{mgu}(t, u) \), can be matched to the logical formula in the following way:

- the empty substitution \( \top \) is associated with falsity constant \( \bot \);
- \( \bot \) is associated with the truth constant \( \top \);
- \( \{x_1 \mapsto t_1, \ldots, x_n \mapsto t_n\} \) is associated with the formula \( x_1 \neq t_1 \lor \cdots \lor x_n \neq t_n \).

The following definition introduces universally quantified disequalities.

**Definition 11.** Universally quantified disequality constraint \( \forall x. t \neq u \) is

- satisfiable iff \( \exists \sigma. \forall \tau, \text{dom}(\tau) \subseteq x. t\sigma \neq u\sigma \)
- unsatisfiable iff \( \forall \sigma. \exists \tau, \text{dom}(\tau) \subseteq x. t\sigma = u\tau \)

For the decision procedure of this type of disequalities, we need one auxiliary lemma.

**Lemma 4.** Given terms \( t \) and \( u \) consider \( \theta = \text{mgu}(t, u) \). Assume without the loss of generality that \( x \not\subseteq \text{codom}(\theta) \) (if \( v \mapsto x \in \theta \) for some \( x \in x \) consider \( \hat{\theta} \) such that it is equal to \( \theta \) except that instead of mapping \( v \) to \( x \) it maps \( x \) to \( v \)). Universally quantified disequality constraint \( \forall x. t \neq u \) is

- satisfiable if \( \theta = \bot \) or \( \text{dom}(\theta) \not\subseteq x \)
- unsatisfiable if \( \text{dom}(\theta) \subseteq x \)

*Proof.* First, let us show that \( \theta = \bot \) or \( \text{dom}(\theta) \not\subseteq x \) implies satisfiability. We need to show that \( \exists \sigma. \forall \tau, \text{dom}(\tau) \subseteq x. t\sigma \neq u\sigma \). Take \( \sigma = \top \). Thus \( t\sigma = t\tau \) and \( u\sigma = u\tau \). It is left to show that \( \forall \tau, \text{dom}(\tau) \subseteq x. t\tau \neq u\tau \). If \( \theta = \bot \) then \( t \) and \( u \) are not unifiable, which implies the above statement. Otherwise, consider some \( \tau \) such that \( \text{dom}(\tau) \subseteq x \). If \( t\tau = u\tau \) then \( \tau \) is a unifier of \( t \) and \( u \). Thus \( \tau \not\subseteq \theta \). If we will show that \( \text{dom}(\theta) \subseteq \text{dom}(\tau) \subseteq x \) we will get a contradiction with the our assumptions and therefore \( t\tau \neq u\tau \). Indeed, consider \( v \in \text{dom}(\theta) \). If \( v \not\in \text{dom}(\tau) \) consider two cases:

- \( v \mapsto w \in \theta \) for some \( w \in V \). From the assumptions follows that \( w \not\in x \) and thus \( w \not\in \text{dom}(\tau) \). Consider the term \( f(v, w) \) for some binary constructor \( f \). It is easy to see that \( f(v, w)\tau = f(v, w)\theta \). Thus it should be that \( v \in \text{dom}(\tau) \).
- \( v \mapsto s \in \theta \) for some term \( s \notin V \). Then, trivially \( v\tau = v\theta \). Thus it should be that \( v \in \text{dom}(\tau) \).

Finally, let us show that if \( \text{dom}(\theta) \subseteq x \) then \( \exists \sigma. \exists \tau, \text{dom}(\tau) \subseteq x. t\sigma = u\tau \). Indeed, consider some \( \sigma \). Take \( \tau = \theta \). Then \( t\tau = u\tau \) and thus \( t\sigma = u\sigma \).

By the above lemma, if the substitution \( \text{mgu}(t, u) \) maps only universally quantified variables, then the disequality is unsatisfiable and satisfiable otherwise.

Finally, it is left to show how to check satisfiability of the quantified equalities of the form \( \forall x. \exists y. t \equiv u \). As we will see soon, if the constraint of this form is satisfiable, then the logical formula, corresponding to the constraint, \( \forall x. \exists y. t = u \) is equivalent to the formula \( \exists y. \wedge_i v_i = t_i \) such that \( v_i \in V \) and \( v_i \not\in x \) and \( \text{Vars}(t_i) \cap x = \emptyset \) for all \( t \) [Liu et al. 1999].

**Definition 12.** Quantified equality constraint of the form \( \forall x. \exists y. t \equiv u \) is

- satisfiable iff \( \exists \phi. \forall \tau, \text{dom}(\phi) \subseteq x. \exists \sigma. \text{dom}(\phi) \subseteq y. t\sigma = u\phi \)
- unsatisfiable iff \( \forall \phi. \exists \tau, \text{dom}(\phi) \subseteq x. \forall \sigma. \text{dom}(\phi) \subseteq y. t\sigma \neq u\phi \)
Lemma 5. If terms \( t \) and \( u \) are not unifiable then the constraint is unsatisfiable. Otherwise let \( \theta \) be an idempotent unifier of \( t \) and \( u \) (Lemma 2 states that if terms are unifiable there exists their idempotent unifier). Let \( \theta_y \triangleq \{ y \mapsto t \mid y \mapsto t \in \theta \land y \notin \overline{y} \} \). Let \( \hat{\theta} = \{ v \mapsto t \mid v \mapsto t \in \theta \land v \notin \overline{y} \} \). Then the quantified equality constraint of the form \( \forall x \in \overline{x}. t \equiv u \) is

- satisfiable if \( \text{dom}(\hat{\theta}) \cap \overline{x} = \emptyset \) and \( \forall p \in \text{codom}(\hat{\theta}). \text{Vars}(p) \cap \overline{x} = \emptyset \)
- unsatisfiable if \( \text{dom}(\hat{\theta}) \cap \overline{x} \neq \emptyset \) or \( \exists p \in \text{codom}(\hat{\theta}). \text{Vars}(p) \cap \overline{x} \neq \emptyset \)

Proof. First, it is obvious that if \( t \) and \( u \) are not unifiable then the constraint is unsatisfiable. Next, let us prove the statement involving satisfiability. We need to show that if the given condition is met, then \( \forall \sigma. \text{dom}(\tau) \subseteq \overline{x}. \exists \phi, \text{dom}(\phi) \subseteq \overline{y}, t\phi \sigma = u\phi \sigma \). Take \( \sigma = \hat{\theta} \). Given an arbitrary \( \tau \) such that \( \text{dom}(\tau) \subseteq \overline{x} \) take \( \phi \) to be equal to \( \theta_y \). To complete the proof we will need an auxiliary statement.

- \( \forall s. s \hat{\theta} = s\hat{\theta} \) where \( \hat{\theta} = \{ x \mapsto t \mid x \mapsto t \in \tau \} \).

Proof. By the Lemma 1 it is sufficient to show that \( \forall \nu \in V. \nu \hat{\theta} = \nu\hat{\theta} \). Consider the cases:
- \( \nu \in \overline{x} \). Then \( \nu \hat{\theta} = \tau(\nu)\hat{\theta} \) since \( \text{dom}(\hat{\theta}) \cap \overline{x} = \emptyset \). Then \( \nu \hat{\theta} = \nu \hat{\theta} \) and by the construction \( \nu \hat{\theta} = \nu \hat{\theta} \).
- \( \nu \notin \overline{x} \). Then \( \nu \hat{\theta} = \hat{\theta}(\nu) \). Since \( \forall p \in \text{codom}(\hat{\theta}). \text{Vars}(p) \cap \overline{x} = \emptyset \), \( \nu \hat{\theta} = \nu \hat{\theta} \) and trivially \( \nu \hat{\theta} = \hat{\theta}(\nu) \).

By this statement \( t\theta_y \tau \hat{\theta} = t\theta_y \hat{\theta} \) and \( u\theta_y \tau \hat{\theta} = u\theta_y \hat{\theta} \). Because \( \hat{\theta} \) is idempotent \( t\theta = t\theta \hat{\theta} \) and \( u\theta = u\theta \hat{\theta} \).

Finally, since \( \hat{\theta} \) is a unifier \( t\theta = u\theta \).

It is left to prove the statement involving unsatisfiability. In fact, we will prove more general statement.

- Let \( \tilde{t} \) and \( \tilde{u} \) be two arbitrary unifiable terms, let \( \hat{\theta} \) be their unifier. Then \( \forall \overline{x} \subseteq \text{dom}(\hat{\theta}) \cup \bigcup_{p \in \text{codom}(\hat{\theta})} \text{Vars}(p). \overline{x} \neq \emptyset. \forall \sigma. \exists \tau. \text{dom}(\tau) \subseteq \overline{x}. t\tau \sigma \neq \tilde{u}\tau \sigma \)

Proof. By the induction on \( t \):
- \( t = \nu \) for some \( \nu \in V \). Consider the cases for \( u \):
  * \( \tilde{u} = w \) for some \( w \in V \). Then \( \hat{\theta} = \{ \nu \mapsto w \} \) and either \( \nu \in \overline{x} \) or \( w \in \overline{x} \).
    Let the former be true (the other case is similar). Given some arbitrary \( \sigma \) take \( \tau \triangleq \{ \nu \mapsto z \mid z \in V \setminus (\text{dom}(\sigma) \cup \bigcup_{p \in \text{codom}(\sigma)} \text{Vars}(p)) \} \). Then \( \nu \tau \sigma = z \) and \( \nu \tau \sigma \neq z \).
  * \( \tilde{u} = g(\tilde{u}_1, \ldots, \tilde{u}_m) \) for some constructor \( g \). Then \( \hat{\theta} = \{ \nu \mapsto u \} \).
    If \( \nu \in \overline{x} \) then pick some constructor \( f \neq g \) (because we assume there exists an infinite number of constructor symbols, we can always do it).
    Take \( \tau \triangleq \{ \nu \mapsto f(z_1, \ldots, z_n) \mid z_i \in V \} \). Then \( \nu \tau \sigma = \nu \tau \sigma = g(\tilde{u}_1, \ldots, \tilde{u}_m) \), and thus these terms are not equal. If \( \nu \notin \overline{x} \) then take some \( x \in \overline{x} \).
    For some \( i \) it should be that \( x = \text{Vars}(\tilde{u}_i) \).
    Given \( \sigma \) consider \( \sigma(\nu) \), pick some \( s \) such that \( \sigma(\nu) \neq g(\tilde{u}_1, \ldots, \tilde{u}_m) \) (it can be done by the induction on \( \sigma(\nu) \)). Then \( \tau \triangleq \{ x \mapsto s \} \).
- \( t = f(t_1, \ldots, t_n) \). Consider the cases for \( u \):
  * \( \tilde{u} = w \) for some \( w \in V \). Then the proof proceeds in the same way as in the previous case.
  * \( \tilde{u} = f(\tilde{u}_1, \ldots, \tilde{u}_n) \) (\( \tilde{u} \) cannot be equal to some constructor \( g \neq f \) by our assumption of unifiability of terms). Then by our assumption \( t_1 \equiv \tilde{u}_1 \land \cdots \land t_n \equiv \tilde{u}_n \).
    For some \( i \) it should be the case that \( \overline{x} \subseteq \text{dom}(\tilde{t}_i) \cup \bigcup_{p \in \text{codom}(\tilde{t}_i)} \text{Vars}(p) \) where \( \tilde{t}_i \triangleq \text{mg}(t_i, \tilde{u}_i) \).
    By the induction for an arbitrary \( \sigma \) there exists \( \tau \) such that \( t_1 \tau \sigma \neq \tilde{u}_i \tau \sigma \) and thus \( t \tau \sigma \neq \tilde{u} \tau \sigma \).

\[\square\]
Given this lemma, we compute \( \text{mgu}(t, u) \) in order to check the constraint \( \forall x \exists y. t \equiv u \). By \( \text{mgu}(t, u) \) we can construct an idempotent substitution \( \theta \) that is also a unifier of \( t \) and \( u \). We take \( \hat{\theta} \) — a part of \( \theta \) that does not bind existentially quantified variables \( \overline{y} \). Then we check if \( \hat{\theta} \) binds variables from \( \overline{x} \), or some term from codomain of \( \hat{\theta} \) contains variables from \( \overline{x} \). If it does then the constraint is unsatisfiable, because in this case we can always pick an assignment for \( \overline{x} \) that will make the terms not unifiable. Otherwise it is satisfiable.

### 3.3 Extending the Search

In this section we describe how OCanren search interacts with negation. Also we finally present the code of negation operator itself.

In OCanren (as in any typical MiniKanren implementation) the search is implemented on top of backtracking lazy stream monad [Kiselyov et al. 2005]. During the search the current state is maintained. The state contains accumulated constraints plus some supplementary information stored in the environment (for example, the identifier of last allocated variable). A goal is simply a function which takes a state and returns a lazy stream of states. All logical primitives, such as individual constraints, conjunction, disjunction, and fresh variable introduction, can be implemented based on this representation of goals (an interested reader may refer to [Hemann and Friedman [n. d.]; Kiselyov et al. 2005]).

Now we can define the negation operator \(~\) (see Listing 6). Let us describe it in details.

```plaintext
let (~) g st = let sts' = g st in
      let cexs = Stream.map (diff st) sts' in
      let sub ss cex = let ss' = negate cex in
                         Stream.bind ss (fun s ->
                         Stream.bind ss' (fun s' ->
                                             Stream.unit (merge s s'))
                         ))
      in Stream.fold sub (Stream.unit st) cexs
```

Listing 6. Implementation of the negation operator

Negation operator is a function which takes a goal and returns a negated goal. Because a goal is itself a function taking a state, \(~\) takes two arguments: the goal \( g \) and the state \( st \) (line 1).

The first step of constructive negation is to run the positive version of the goal, as code in line 2 does. We run it in the current state \( st \) and thus the call \( g \ st \) returns a stream of refined states \( st \) \( st ' \). Each state from this stream will contain the constraints from the original state \( st \) as its subpart. However, we need to negate only constraints originated from \( g \) solely. Thus, on the line 3 we map every state from the stream \( st \) \( st ' \) to its difference with respect to the original substitution \( st \). In order to compute difference of two states \( st \) and \( st ' \) (Listing 7), given that \( st \) is more general that \( st ' \) we need to compute difference of their substitutions and disequality constraints stores. The difference of substitution \( s ' \) with respect to \( s \) (Listing 8) is just a substitution containing all mappings from \( s ' \) which are not simultaneously in \( s \). The difference of constraint store \( c ' \) with respect to \( c \) is a constraint store...
containing all disequalities that are in \( c' \) but not in \( c \) (Listing 8). As long as persistent data structures are used to implement substitutions and constraint stores, the \texttt{diff} can be computed\(^4\).

Line 4 defines auxiliary function \texttt{sub} which takes two arguments: a stream of states \( ss \) and some state \( cex \), and returns another stream. The purpose of this function is to "subtract" \( cex \) from every element of \( ss \). It is done as follows. First, the state \( cex \) is negated as described in section 3.1 (line 5). As we have seen, as a result of the negation of a single state the stream of states (the disjunction of formulas) can be obtained. Thus the result of the call \texttt{negate cex} is the stream of states \( ss' \). For every combination of some state \( s \) from the given stream \( ss \) (line 6) and some state \( s' \) from the stream \( ss' \) representing the result of negation (line 7) we compute their conjunction (line 8). The conjunction of two states is computed by the function \texttt{merge} (Listings 7, 8).

Finally, on the line 11 \texttt{fold} is called on the stream \( cexs \), which is a stream of answers for the positive version of the goal \( g \), with function \texttt{sub} defined above and the initial accumulator \texttt{Stream.unit st}. The function \texttt{Stream.fold} is implemented as a regular left fold over a possibly infinite list. Intuitively with folding over stream \( cexs \) we “subtract” from the original state \( st \) every answer obtained from the goal \( g \).

\(^4\) In the Listing 8 we present a simple representation of the constraint store as a list of substitutions. In the actual implementation, we use more sophisticated representation, that also provides the \texttt{diff} function. The simpler version presented here gives some intuition on how to implement \texttt{diff} for the constraint stores.
module State = struct
  type t = Env.t * Subst.t * CStore.t
...

let diff (e, s, cs) (e′, s′, cs′) =
  let e = Env.diff e e′ in
  let s = Subst.diff s s′ in
  let cs = CStore.diff cs cs′ in
  (e, s, cs)

let merge (e, s, cs) (e′, s′, cs′) =
  let e = Env.merge e e′ in
  match Subst.merge s s′ with
  | None → None
  | Some s →
    let cs = CStore.merge cs cs′ in
    match CStore.recheck s cs with
    | None → None
    | Some cs → (e, s, cs)
end

Listing 7. Implementation of the auxiliary functions

module VarMap = Map.Make(Var)

module Subst = struct
  (* Substitution is a mapping from variables to terms *)
  type t = Term.t VarMap.t
...

  let diff s s′ =
    VarMap.fold (fun v t a →
      if not (VarMap.mem v s) then
        VarMap.add v t a
      else a
    ) s′ VarMap.empty

  let merge s s′ =
    VarMap.fold (fun v t →
      function
      | None → None
      | Some a → unify a v t
    ) s′ (Some s)
end

module CStore = struct
  (* Constraint store is a list of substitutions *)
  type t = Subst.t list
...

  (* cs′ must be obtained from cs by
   * the addition of new constraints
   *)
  let diff cs cs′ =
    if cs′ = cs then []
    else
      match cs′ with
      | _ :: cs′ → diff cs cs′
      (* cs′ = [] implies cs = [] *)
      | _ → assert false

  let merge cs cs′ =
    List.append cs′ cs
end

Listing 8. Implementation of the auxiliary functions
3.4 Stratification

A negation, when combined with recursion, might become a source of confusion for a programmer. Consider the program in Listing 9. It encodes a two-players game. The positions in the game are given as single-character strings ‘a’, ‘b’, ‘c’ and ‘d’. A binary relation move encodes the game field as the set of possible moves. An unary relation winning determines the set of winning positions, meaning that if the first player starts from some winning position, by making “good” moves the player has an opportunity to win the game. According to the definition of winning, the position is winning if there exists a move from this position to some non-winning position. Clearly, every position with no moves from it is losing.

Given the suchlike definition of winning, there is no doubt, that the position ‘d’ is losing position, and thus the goal winning ‘d’ should fail, which, in turn, means that winning ‘c’ should succeed. However, whether the goal winning ‘a’ (or winning ‘b’) should fail or succeed is not clear.

```prolog
let move x y =
    (x, y) ≡ (‘a’, ‘b’) ∨
    (x, y) ≡ (‘b’, ‘a’) ∨
    (x, y) ≡ (‘b’, ‘c’) ∨
    (x, y) ≡ (‘c’, ‘d’)

let winning x =
    fresh (y) (move x y ∧ ¬(winning y))
```

Listing 9. Encoding of two-players game

The problem with the semantics of program in Listing 9, originates from the interaction of negation and recursion. Definition of the relation winning refers to itself under negation. Logic programs that have this property are called non-stratified [Przymusinski 1989]. Vice versa, programs that do not have loops over negation, are called stratified.

Our current implementation handles only stratified programs. We leave the task of supporting non-stratified programs as a direction for future work.

4 EVALUATION

In this section, we present an evaluation of implemented constructive negation on a series of examples.

4.1 If-then-else

Using relational if-then-else operator, presented in section 2.1, we have implemented several higher-order relations over lists, namely find (Listing 10), remove (Listing 11) and filter (Listing 12). These relations are almost identical (syntactically) to their functional implementations. We have tested that these relations can be run in various directions and produce the expected results. For example, the goal filter p q q with the predicate p equal to

```prolog
fun l → fresh (x) (l ≡ [x])
```

stating that the given list should be a singleton list, starts to generate all singleton lists. Vice versa, the goal filter p q [] with that same p generates all lists, constrained to be not a singleton list.

Note, this implementation differs from the one in Section 1, but it is easy to see that these two are semantically equivalent.
Listings 13-16 give more concrete examples of queries to these relations. In the listing the syntax `run n q g` means running a goal `g` with the free variable `q` taking the first `n` answers ("*" denotes all answers). After the sign `≈` the result of the query is given. The result `fail` means that the query has failed. The result `succ `{ ... `}` means that the query has succeeded delivering `n` answers. Each answer represents a set of constraint on free variables. Constraints are of two forms: equality constraints, e.g. `q = (1, _, _0)`, or disequality constraints, e.g. `q ≠ (1, _, _0)`. The terms of the form `_i` in the answer denote some universally quantified variables.

**Listing 10.** A definition of `find` relation

```scheme
let find p e xs =
  fresh (x xs' ys') (xs ≡ x::xs' ∧
         ifte (p x)
               (e ≡ x)
         (find p e xs'))
```

**Listing 11.** A definition of `remove` relation

```scheme
let remove p xs ys =
  (xs ≡ [] ∧ ys ≡ []) ∨
  fresh (x xs' ys') (xs ≡ x::xs' ∧
                   ifte (p x)
                         (ys ≡ x::ys' ∧
                          (ys ≡ x::ys' ∧
                           remove p xs' ys'))
                     filter p xs' ys')
```

**Listing 12.** A definition of `filter` relation

```scheme
let filter p xs ys =
  (xs ≡ [] ∧ ys ≡ []) ∨
  fresh (x xs' ys') (xs ≡ x::xs' ∧
                   (ifte (p x)
                         (ys ≡ x::ys')
                       ∧
                       filter p xs' ys'))
```

**Listing 13.** Definition of the predicate `p`

```scheme
let p l = fresh (x) (l ≡ [x])
```

**Listing 14.** Example of queries to `find` relation

```scheme
run 3 q (fresh (e) find p e q)
≈ succ {
    { q = [-0] :: [-1] }
    { q = -0 :: [1] :: [-2];
      -0 ≠ [-3] }
    { q = -0 :: [-1] :: [-2] :: [-3];
      -0 ≠ [-4]; -1 ≠ [-5] }
}
```

**Listing 15.** Example of queries to `remove` relation

```scheme
run 3 q (fresh (e) remove p q []
≈ succ {
    { q = [ ] }
    { q = [-0]; -0 ≠ [-1] }
    { q = [-0]; -1;
      -0 ≠ [-2]; -1 ≠ [-3] }
}
```

```scheme
let filter p xs ys =
  (xs ≡ [] ∧ ys ≡ []) ∨
  fresh (x xs' ys') (xs ≡ x::xs' ∧
                   (ifte (p x)
                         (ys ≡ x::ys')
                       ∧
                       filter p xs' ys'))
```

```scheme
let remove p xs ys =
  (xs ≡ [] ∧ ys ≡ []) ∨
  fresh (x xs' ys') (xs ≡ x::xs' ∧
                   ifte (p x)
                         (ys ≡ x::ys')
                     (ys ≡ x::ys') ∧
                     remove p xs' ys')
```

```scheme
let find p e xs =
  (xs ≡ [] ∧ ys ≡ []) ∨
  fresh (x xs' ys') (xs ≡ x::xs' ∧
                     ifte (p x)
                         (e ≡ x)
                     (find p e xs'))
```

```scheme
let p l = fresh (x) (l ≡ [x])
```

```scheme
run 3 q (fresh (e) find p e q)
≈ succ {
    { q = [-0] :: [-1] }
    { q = -0 :: [1] :: [-2];
      -0 ≠ [-3] }
    { q = -0 :: [-1] :: [-2] :: [-3];
      -0 ≠ [-4]; -1 ≠ [-5] }
}
```

```scheme
run 3 q (fresh (e) remove p q []
≈ succ {
    { q = [ ] }
    { q = [-0]; -0 ≠ [-1] }
    { q = [-0]; -1;
      -0 ≠ [-2]; -1 ≠ [-3] }
}
```
run 3 q (filter p q q)
\leadsto\text{succ} \{
    \{ q = [ ] \}
    \{ q = [\cdot_0] \}
    \{ q = [\cdot_0; \cdot_1] \}
\}

run 3 q (filter p q [1])
\leadsto\text{succ} \{
    \{ q = [[[1]]] \}
    \{ q = [\cdot_0; [1]]: \cdot_0 \neq [\cdot_1] \}
    \{ q = [[1]; \cdot_0]: \cdot_0 \neq [\cdot_1] \}
\}

run 3 q (filter p q [ ])
\leadsto\text{succ} \{
    \{ q = [ ] \}
    \{ q = [\cdot_0]; \cdot_0 \neq [\cdot_1] \}
    \{ q = [\cdot_0; \cdot_1]; \cdot_0 \neq [\cdot_2]; \cdot_1 \neq [\cdot_3] \}
\}

Listing 16. Example of queries to \text{filter}

4.2 Universal quantification
In the Section 2.2 we presented the \textit{forall} goal constructor which is implemented through the double negation. We have observed, that although \textit{forall} \textit{g} does not terminate when the goal \textit{g} \textit{x} has an infinite number of answers (assuming \textit{x} is a fresh variable), it does terminate in the case when \textit{g} \textit{x} has a finite number of answers. The behavior of \textit{forall} in this case is sound even in the presence of disequality constraints or nested quantifiers. The Table 1 gives some concrete examples. The left column contains the tested goals\textsuperscript{6} and the right column gives the obtained results. For the results we use the same notation as in the previous section.

5 LIMITATIONS AND FUTURE WORK
In this section we discuss the limitations of constructive negation in general and our implementation in particular. Also we consider possible directions for future work.

5.1 Type Constraints
Although the program written in \textsc{OCAnren} typechecks statically (thus, for example, preventing the user from unifying two terms of distinct types), at runtime the type information is erased. In the presence of even regular disequality constraints it can lead to the incorrect results. As an example, consider the following program:

\textsuperscript{6} We typeset the goals in terms of first-order logic syntax instead of \textsc{OCAnren} syntax for brevity and clarity.
\[
\begin{array}{|l|}
\hline
\forall x. x = q & \text{fail} \\
\forall x. \exists y. x = y & \text{succ } \{[q = _\cdot 0]\} \\
\forall x. \exists y. x = y \land y = q & \text{fail} \\
\forall x. q = (1,x) & \text{fail} \\
\forall x. \exists y. y = (1,x) & \text{succ } \{[q = _\cdot 0]\} \\
\forall x. \exists y. x = (1,y) & \text{fail} \\
\forall x. x \neq q & \text{fail} \\
\forall x. \exists y. x \neq y & \text{succ } \{[q = _\cdot 0]\} \\
\forall x. \exists y. x \neq y \land y = q & \text{fail} \\
\forall x. q \neq (1,x) & \text{succ } \{[q \neq (1, _\cdot 0)]\} \\
(\exists x. q = (1,x)) \land (\forall x. q \neq (1,x)) & \text{fail} \\
\forall x. (x,x) \neq (0,1) & \text{succ } \{[q = _\cdot 0]\} \\
\forall x. (x,x) \neq (1,1) & \text{fail} \\
\forall x. (x,x) \neq (q,1) & \text{succ } \{[q \neq 1]\} \\
\exists a. b. q = (a,b) \land \forall x. (x,x) \neq (a,b) & \text{succ } \{[q = (_\cdot 0, _\cdot 1); _\cdot 0 \neq _\cdot 1]\} \\
\hline
\end{array}
\]

Table 1. forall evaluation

**type** bool = true | false

**let** g =

fresh (x y z : bool) (  
    (x ≠ y)  
    (y ≠ z)  
    (z ≠ x)  
)

The goal g states that there exists at least three different non-equal terms of type bool, which, as we know, is not true. Yet the query run g will succeed.

In order to prevent unsoundness in cases like this, type information in the form of type constraints should be somehow attached to the variables at runtime. The satisfiability of type constraints then should be rechecked each time when the new disequality is added to some variable. An extension of OCanren with type constraints is a direction for future work.
5.2 Non-stratified Programs
As we have already discussed in the section 3.4 our current implementation handles only stratified logic programs. One of the possible extensions is to support non-stratified programs, such as one given in Listing 9, with respect to well-founded and/or stable model semantics (see section 6 for the details).

5.3 Negation of Goals With an Infinite Number of Answers
Consider the following program:

\[
\text{let } \text{zeros } l =
\begin{align*}
1 & \equiv [0] \\
\lor \\
\text{fresh } (l') ( \\
(1 \equiv 0 :: l') \\
(zeros l')
\end{align*}
\]

The unary relation zeros defines lists consisting of zeros. Now, intuitively, the query run \( \neg(\text{zeros } q) \) should enumerate all lists that are not built out of zeros only. Yet this query will fail to deliver even a single answer. Why? Consider its operational behavior. First the positive version of the goal, that is \( \text{zeros } q \), should be executed. Then all answers to this goal should be collected and complemented. However, there is an infinite number of answers to \( \text{zeros } q \) and thus this process will never terminate.

It is a significant drawback of constructive negation that the negation of the goal cannot be computed if the goal has an infinite number of answers. This limitation cannot be avoided in general, however in some cases it is possible to narrow the number of answers to some subgoal by the reordering of surrounding subgoals. For example, the query run \( \neg(\text{zeros } q) \land (q \equiv [1]) \) can be executed in finite time by the reordering of conjuncts. It seems that the best strategy is to delay negative subgoals as long as possible, but we do not have a formal proof of that.

6 RELATED WORKS
There are two directions of work in the process of incorporating negative reasoning in the logic programming: the first considers the semantics of negation, and the second is focused mainly on implementation aspects.

The first attempt to give a semantics for negation in logic programming was done by Clark [Chan 1988; Clark 1978] with his completion semantics. It was then realized, that Clark’s semantics has various drawbacks [Van Gelder et al. 1991].

Przymusinski [Przymusinski 1989] has studied the semantics of stratified logic programs. He introduced the notion of \textit{perfect model semantics} for such programs. Stratified logic programs have a variety of good properties, including the property that each stratified program has a unique minimal model.

In an attempt to extend the semantics of negation to non-stratified programs the \textit{well-founded semantics} was proposed [Van Gelder et al. 1991]. However, this semantics is three-valued, meaning that for some queries it can return answer unknown. For example, given the relation \textit{winning} (section 3.4, listing 9), queries \textit{winning} ‘\( a \)’ and \textit{winning} ‘\( b \)’ would return unknown.

An alternative approach is \textit{stable model semantics} [Gelfond and Lifschitz 1988]. Under this semantics, non-stratified logic program can have several stable models. Program, that defines \textit{winning}, has two stable models, in one of these models goal \textit{winning} ‘\( a \)’ succeeds and \textit{winning} ‘\( b \)’ fails, in the other \textit{winning} ‘\( a \)’ fails and \textit{winning} ‘\( b \)’ succeeds. Logic programming under stable model semantics is also known under the name answer set programming (ASP).
The works [Dovier et al. 2000; Stuckey 1991] are theoretical studies of constructive negation in the context of constraint logic programming. They give a necessary and sufficient condition for the constraint structures that are compatible with constructive negation. Namely, the constraint structure should be admissible closed.

From an implementation side, Chen et al. [Chen et al. 1995; Chen and Warren 1996] developed a Prolog system based on SLG resolution, which is sound with respect to well-founded semantics. However, they used negation as failure with delaying of non-ground negative subgoals. [Liu et al. 1999] is an extension of this system with the support of the constructive negation. Works [Álvez et al. 2004; Barták 1998] implement a constraint logic programming systems with the support of constructive negation. Yet, as with our implementation, the constructive negation in these systems supports only equality and disequality constraints over first-order terms. We are not aware of any practical implementation that is parametric over arbitrary admissible closed constraint structures.

Many tools were developed to compute stable models of logic programs, among them are [Gebser et al. 2007; Giunchiglia et al. 2006]. These systems usually require to perform grounding of logic program. The problem of finding stable models of ground logic program then is encoded as propositional formula and solved by some SAT solver. Unfortunately, some logic programs do not have finite grounding, but even if a program has it, grounding may cause an exponential blow-up. Recently, a goal-directed system for computing stable models was developed [Arias et al. 2018; Marple et al. 2012, 2017]. To the best of our knowledge, it is the only ASP system, that does not require grounding. The key components of this system are the usage of tabling, constructive negation, coinductive logic programming, and non-monotonic reasoning check. It is an interesting and challenging task to extend MiniKanren with the support of stable model semantics in the spirit of this line of work.

7 CONCLUSION
We have presented an implementation of constructive negation for relational programming language OCanren, a dialect of MiniKanren. Unlike the negation as failure, constructive negation is consistent with the pure logical nature of MiniKanren. As we have demonstrated the negative reasoning increases the expressive power of relational language by allowing to compose more relations in a natural and intuitive form.

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Constructive Negation for MiniKanren


Certified Semantics for miniKanren∗

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We present two formal semantics for the core miniKanren. First, we give denotational variant which corresponds to the minimal Herbrand model for definite logic programs. Second, we present operational semantics which models interleaving, and prove its soundness and completeness w.r.t. denotational semantics. Our development is supported by formal Coq specification, thus making it certified.

CCS Concepts: • Theory of computation → Constraint and logic programming; Denotational semantics; Operational semantics;

Additional Key Words and Phrases: Relational programming, denotational semantics, operational semantics, certified programming

1 INTRODUCTION

The introductory book on miniKanren [Friedman et al. 2005] describes the language by means of an evolving set of examples. In the series of follow-up papers [Alvis et al. 2011; Byrd and Friedman 2007; Hemann and Friedman 2013, 2015; Hemann et al. 2016; Swords and Friedman 2013] various extensions of the language were presented with their semantics explained in terms of Scheme implementation. We argue that this style of semantic definition is fragile and not self-evident since it requires the knowledge of semantics of concrete implementation language. In addition the justification of important properties of relational programs (for example, refutational completeness [Byrd 2009]) becomes cumbersome. In the area of programming languages research a formal definition for the semantics of language of interest is a de-facto standard, and in our opinion in its current state miniKanren deviates from this standard.

There were some previous attempts to define a formal semantics for miniKanren. Lozov et al. [2017] present a variant of nondeterministic operational semantics, and Rozplokhas and Boulytchev [2018] use another variant of finite-set semantics. None of them was capable of reflecting the distinctive property of miniKanren search — interleaving [Kiselyov et al. 2005], thus deviating from the conventional understanding of the language.

In this paper we present a formal semantics for core miniKanren and prove some its basic properties. First, we define denotational semantics similar to the least Herbrand model for definite logic programs [Lloyd 1984]; then we describe operational semantics with interleaving in terms of labeled transition system. Finally, we prove the soundness and completeness of the operational semantics w.r.t the denotational one. We support our development

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$C = \{ C_{ki} \}$ constructors with arities
$T_X = X \cup \{ C_{ki}(t_1, \ldots, t_{k_i}) \mid t_j \in T_X \}$ terms over the set of variables $X$
$D = T_\emptyset$ ground terms
$X = \{ x, y, z, \ldots \}$ syntactic variables
$\mathcal{A} = \{ \alpha, \beta, \gamma, \ldots \}$ semantic variables
$\mathcal{R} = \{ R_{ki} \}$ relational symbols with arities
$\mathcal{G} = T_X \equiv T_X$ unification
$\mathcal{G} \land \mathcal{G}$ conjunction
$\mathcal{G} \lor \mathcal{G}$ disjunction
$\textbf{fresh } X \cdot \mathcal{G}$ fresh variable introduction
$R_{ki}(t_1, \ldots, t_{k_i}), t_j \in T_X$ relational symbol invocation
$S = \{ R_{ki} = \lambda x_1 \ldots x_{k_i} . g_i ; \}$ specification

Fig. 1. The syntax of the source language

\[
\mathcal{FV}(x) = \{ x \} \\
\mathcal{FV}(C_{ki}(t_1, \ldots, t_{k_i})) = \bigcup \mathcal{FV}(t_i) \\
\mathcal{FV}(t_1 \equiv t_2) = \mathcal{FV}(t_1) \cup \mathcal{FV}(t_2) \\
\mathcal{FV}(g_1 \land g_2) = \mathcal{FV}(g_1) \cup \mathcal{FV}(g_2) \\
\mathcal{FV}(g_1 \lor g_2) = \mathcal{FV}(g_1) \cup \mathcal{FV}(g_2) \\
\mathcal{FV}(\textbf{fresh } x . g) = \mathcal{FV}(g) \setminus \{ x \} \\
\mathcal{FV}(R_{ki}(t_1, \ldots, t_{k_i})) = \bigcup \mathcal{FV}(t_i)
\]

Fig. 2. Free variables in terms and goals

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five types of goals: unification of terms, conjunction and disjunction of goals, fresh variable introduction and invocation of some relational definition. Thus, unification is used as a constraint, and multiple constraints can be combined using conjunction, disjunction and recursion. For the sake of brevity we abbreviate immediately nested “fresh” constructs into the one, writing “fresh $x$ . fresh $y$ . . .” instead of “fresh $x$ . fresh $y$ . . . $g$”. The final syntactic category is specification $S$. It consists of a set of relational definitions and a top-level goal. A top-level goal represents a search procedure which returns a stream of substitutions for the free variables of the goal. The definition for set of free variables for both terms and goals is given on Figure 2; as “fresh” is the sole binding construct the definition is rather trivial. The language we defined is first-order, as goals can not be passed as parameters, returned or constructed at runtime.

We now informally describe how relational search works. As we said, a goal represents a search procedure. This procedure takes a state as input and returns a stream of states; a state (among other information) contains a substitution which maps semantic variables into terms over semantic variables. Then five types of scenarios are possible (dependent on the type of the goal):

- Unification “$t_1 \equiv t_2$” unifies terms $t_1$ and $t_2$ in the context of the substitution in the current state. If terms are unifiable, then their MGU is integrated into the substitution, and one-element stream is returned; otherwise the result is an empty stream.
- Conjunction “$g_1 \land g_2$” applies $g_1$ to the current state and then applies $g_2$ to the each element of the result, concatenating the streams.
- Disjunction “$g_1 \lor g_2$” applies both its goals to the current state independently and then concatenates the results.
- Fresh construct “fresh $x$ . $g$” allocates a new semantic variable $\alpha$, substitutes all free occurrences of $x$ in $g$ with $\alpha$, and runs the goal.
- Invocation “$R_{ki}(t_1, \ldots, t_k)$” finds a definition for relational symbol $R_{ki} = \lambda x_1 \ldots x_k . g_i$, substitutes all free occurrences of formal parameter $x_j$ in $g_i$ with term $t_j$ (for all $j$) and runs the goal in the current state.

We stipulate, that the top-level goal is preceded by an implicit “fresh” construct, which binds all its free variables, and the final substitutions for these variables constitute the result of the goal evaluation.

Conjunction and disjunction form a monadic [Wadler 1995] interface with conjunction playing role of “bind” and disjunction — of “mplus”. In this description we swept a lot of important details under the carpet — for example, in actual implementations the components of disjunction are not evaluated in isolation, but both disjuncts are being evaluated incrementally with the control passing from one disjunct to another (interleaving); instead streams the implementation can be based on “ferns” [Byrd et al. [n. d.]] to defer divergent computations, etc.

As an example consider the following specification:

```latex
append0 = \lambda x y xy .
\quad ((x \equiv \text{Nil}) \land (xy \equiv y)) \lor
\quad \text{fresh} h t ty .
\quad (x \equiv \text{Cons} \ (h, t)) \land
\quad (xy \equiv \text{Cons} \ (h, ty)) \land
\quad (\text{append0} \ y \ t \ ty)
);
revers0 = \lambda x y .
\quad ((x \equiv \text{Nil}) \land (y \equiv \text{Nil})) \lor
\quad \text{fresh} h t t' .
\quad (x \equiv \text{Cons} \ (h, t)) \land
```

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Here we defined two relational symbols — “appendo” and “reverso”, — and specified a top-level goal “reverso x x’. The symbol “appendo” defines a relational concatenation of lists — it takes three arguments and performs a case analysis on the first one. If the first one is an empty list (“Nil”), then the second and the third arguments are unified. Otherwise the first argument is deconstructed into a head “h” and a tail “t”, and the tail is concatenated with the second argument using a recursive call to “appendo” and additional variable “ty”, which represents the concatenation of “t” and “y”. Finally, we unify “Cons (h, ty)” with “xy” to form a final constraint. Similarly, “reverso” defines relational list reversing. The top-level goal represents a search procedure for all lists “x”, which are stable under reversing, i.e. represent palindromes. Running it results in an infinite stream of substitutions:

\[ \begin{align*}
\alpha \mapsto & \text{Nil} \\
\alpha \mapsto & \text{Cons (} \beta_0, \text{Nil) } \\
\alpha \mapsto & \text{Cons (} \beta_0, \text{Cons (} \beta_0, \text{Nil) ) } \\
\alpha \mapsto & \text{Cons (} \beta_0, \text{Cons (} \beta_1, \text{Cons (} \beta_0, \text{Nil) ) ) } \\
\vdots
\end{align*} \]

where “\(\alpha\)” — a semantic variable, corresponding to “x”, “\(\beta_i\)” — free semantics variables.

The notions above can be formalized in Coq in a natural way using inductive data types. We have made a few non-essential simplifications and modifications for the sake of convenience. Specifically, we restrict the arities of constructors to be either zero or two:

\textbf{Inductive term : Set :=}
| Var : var \to \text{term} \\
| Cst : con_name \to \text{term} \\
| Con : con_name \to \text{term} \to \text{term} \to \text{term}. \\

Here “\text{var}” and “\text{con_name}” — types representing variables and constructor names, whose definitions we omitted for the sake of brevity. Similarly, we restrict relations to always have exactly one argument:

\textbf{Definition rel : Set := \text{term} \to \text{goal}.}

These restrictions do not make the language less expressive in any way since we can represent a sequence of terms as a list using constructors Nil and Cons.

We also introduce one additional type of goals — failure — for deliberately unsuccessful computation (empty stream). As a result, the definition of goals looks as follows:

\textbf{Inductive goal : Set :=}
| Fail : \text{goal} \\
| Unify : \text{term} \to \text{term} \to \text{goal} \\
| Disj : \text{goal} \to \text{goal} \to \text{goal} \\
| Conj : \text{goal} \to \text{goal} \to \text{goal} \\
| Fresh : (\text{var} \to \text{goal}) \to \text{goal} \\
| Invoke : rel_name \to \text{term} \to \text{goal}. \

\[
\begin{align*}
  x[t/x] & = t \\
y[t/x] & = y \quad (y \neq x) \\
C_{ki}^{k_i}(t_1, \ldots, t_{k_i})[t/x] & = C_{ki}^{k_i}(t_1[t/x], \ldots, t_{k_i}[t/x]) \\
(t_1 \equiv t_2)[t/x] & = t_1[t/x] \equiv t_2[t/x] \\
(g_1 \land g_2)[t/x] & = g_1[t/x] \land g_2[t/x] \\
(g_1 \lor g_2)[t/x] & = g_1[t/x] \lor g_2[t/x] \\
(fresh \ x \cdot g)[t/x] & = fresh \ x \cdot g \\
(fresh \ y \cdot g)[t/x] & = fresh \ y \cdot (g[t/x]) \quad (y \neq x) \\
(R_{ki}^{k_i}(t_1, \ldots, t_{k_i})[t/x] & = R_{ki}^{k_i}(t_1[t/x], \ldots, t_{k_i}[t/x])
\end{align*}
\]

Fig. 3. Substitutions for terms and goals

Note that in our formalization we use the higher-order abstract syntax for variable binding [Pfenning and Elliott 1988]. We preferred it to the first-order syntax because it gives us the ability to use substitution and inductive principle provided by Coq. However, we still need to carefully ensure some expected properties on the structure of syntax trees. For example, we should require that the definitions of relations do not contain unbound variables:

**Definition** 
`closed_goal_in_context (c : list var) (g : goal) : Prop :=`
\[\forall \ n, \ is_fv_of_goal \ n \ g \rightarrow \mathsf{In} \ n \ c.\]

**Definition** 
`closed_rel (r : rel) : Prop :=`
\[\forall \ (arg : term), \ closed_goal_in_context (fv_term arg) (r \ arg).\]

**Definition** 
`def : Set := {r : rel | closed_rel r}.`

In the snippet above “def” corresponds to a set of relational symbol definitions in a specification.

We set an arbitrary environment (a map from relational symbol to the definition of relation) to use further throughout the formalization. Failure goals allow us to define it as a total function:

**Definition** 
`env : Set := rel_name \rightarrow \mathsf{def}.`

**Variable** 
`Prog : env.`

3 DENOTATIONAL SEMANTICS

In this section we present a denotational semantics for the language we defined above. We use a simple set-theoretic approach which can be considered as an analogy to the least Herbrand model for definite logic programs [Lloyd 1984]. Strictly speaking, instead of developing it from scratch we could have just described the conversion of specifications into definite logic form and took their least Herbrand model. However, in that case we would still need to define the least Herbrand model semantics for definite logic programs in a certified way. In addition, while for this concrete language the conversion to definite logic form is trivial, it may become less trivial for its extensions (with, for examples, nominal constructs [Byrd and Friedman 2007]) which we plan to do in future.

To motivate further development, we first consider the following example. Let us have the following goal:
x ≡ Cons (y, z)

There are three free variables, and solving the goal delivers us the following single answer:

α ↦→ Cons (β, γ)

where semantic variables α, β and γ correspond to the syntactic ones “x”, “y”, “z”. The goal does not put any constraints on “y” and “z”, so there are no bindings for “β” and “γ” in the answer. This answer can be seen as the following ternary relation over the set of all ground terms:

\{ (Cons (β, y), β, y) \mid β, y \in D \} ⊂ D^3

The order of “dimensions” is important, since each dimension corresponds to a certain free variable. Our main idea is to represent this relation as a set of total functions

f : A ↦→ D

from semantic variables to ground terms. We call these functions representing functions. Thus, we may reformulate the same relation as

\{ (f(α), f(β), f(γ)) \mid f ∈ [α ≡ Cons (β, γ)] \}

where we use conventional semantic brackets “[•]” to denote the semantics. For the top-level goal we need to substitute its free syntactic variables with distinct semantic ones, calculate the semantics, and build the explicit representation for the relation as shown above. The relation, obviously, does not depend on concrete choice of semantic variables, but depends on the order in which the values of representing functions are tupled. This order can be conventionalized, which gives us a completely deterministic semantics.

Now we implement this idea. First, for a representing function

f : A → D

we introduce its homomorphic extension

\[ \bar{f} : \mathcal{T}_A \rightarrow D \]

which maps terms to terms:

\[ \bar{f}(α) = f(α) \]
\[ \bar{f}(C_i^{k_i}(t_1, \ldots, t_{k_i})) = C_i^{k_i}(f(t_1), \ldots, f(t_{k_i})) \]

Let us have two terms \( t_1, t_2 \in \mathcal{T}_A \). If there is a unifier for \( t_1 \) and \( t_2 \) then, clearly, there is a substitution \( θ \) which turns both \( t_1 \) and \( t_2 \) into the same ground term (we do not require \( θ \) to be the most general). Thus, \( θ \) maps (some) ground variables into ground terms, and its application to \( t_{i(2)} \) is exactly \( \bar{θ}(t_{i(2)}) \). This reasoning can be performed in the opposite direction: a unification \( t_1 ≡ t_2 \) defines the set of all representing functions \( f \) for which \( \bar{f}(t_1) = \bar{f}(t_2) \).

Then, the semantic function for goals is parameterized over environments which prescribe semantic functions to relational symbols:

\[ Γ : \mathcal{R} → (\mathcal{T}_A^* \rightarrow 2^{A→D}) \]

An environment associates with relational symbol a function which takes a string of terms (the arguments of the relation) and returns a set of representing functions. The signature for semantic brackets for goals is as follows:
It maps a goal into the set of representing functions w.r.t. an environment $\Gamma$.

We formulate the following important completeness condition for the semantics of a goal $g$:

$$\forall \alpha \not\in \text{FV}(g) \forall d \in D \forall f \in \llbracket g \rrbracket \ \exists f' \in \llbracket g \rrbracket : f'(\alpha) = d \land \forall \beta \neq \alpha : f'(\beta) = f(\beta)$$

In other words, representing functions for a goal $g$ restrict only the values of free variables of $g$ and do not introduce any "hidden" correlations. This condition guarantees that our semantics is complete in the sense that it does not introduce artificial restrictions for the relation it defines. It can be proven that the semantics of goals always satisfy this condition.

We remind conventional notions of pointwise modification of a function

$$f[x \leftarrow v](z) = \begin{cases} f(z), & z \neq x \\ v, & z = x \end{cases}$$

and substitution of a free variable with a term in terms and goals (see Figure 3).

For a representing function $f : A \to D$ and a semantic variable $\alpha$ we define the following generalization operation:

$$f \uparrow \alpha = \{ f \mid \alpha \leftarrow d \mid d \in D \}$$

Informally, this operation generalizes a representing function into a set of representing functions in such a way that the values of these functions for a given variable cover the whole $D$. We extend the generalization operation for sets of representing functions $\mathfrak{F} \subseteq A \to D$:

$$\mathfrak{F} \uparrow \alpha = \bigcup_{f \in \mathfrak{F}} (f \uparrow \alpha)$$

Now we are ready to specify the semantics for goals (see Figure 4). We’ve already given the motivation for the semantics of unification: the condition $\overline{f}(t_1) = \overline{f}(t_2)$ gives us the set of all (otherwise unrestricted) representing functions which “equate” terms $t_1$ and $t_2$. Set union and intersection provide a conventional interpretation for disjunction and conjunction of goals, and the semantics of relational invocation reduces to the application of corresponding function from the environment. The only interesting case is “fresh $x . g$”. First, we take an arbitrary semantic variable $\alpha$, not free in $g$, and substitute $x$ with $\alpha$. Then we calculate the semantics of $g[\alpha/x]$. The interesting part is the next step: as $x$ can not be free in “fresh $x . g$”, we need to generalize the result over $\alpha$ since in our model the semantics of a goal specifies a relation over its free variables. We introduce some non-determinism, by choosing arbitrary $\alpha$, but it can be proven by structural induction, that with different choices of free variable, semantics of a goal won’t change. Consider the following example:

fresh $y . (\alpha \equiv y) \land (y \equiv \text{Zero})$

As there is no invocations involved, we can safely omit the environment. Then:
\[ [t_1 \equiv t_2]_\Gamma = \{ f : A \rightarrow D \mid f (t_1) = f (t_2) \} \text{ [UNIFY]} \]
\[ [g_1 \land g_2]_\Gamma = [g_1]_\Gamma \land [g_2]_\Gamma \text{ [AND]} \]
\[ [g_1 \lor g_2]_\Gamma = [g_1]_\Gamma \lor [g_2]_\Gamma \text{ [OR]} \]
\[ \left[ \text{fresh } x \cdot g \right]_\Gamma = (\left[ g(\alpha/x) \right]_\Gamma) \uparrow \alpha, \alpha \not\in FV(g) \text{ [FRESH]} \]
\[ [R(t_1, \ldots, t_k)]_\Gamma = (\Gamma R) t_1 \ldots t_k \text{ [INVOKED]} \]

Fig. 4. Denotational semantics of goals

\[
\begin{align*}
\left\{ f \mid f (\bar{a}) = \bar{f}(\bar{a}) \right\} \uparrow \bar{\alpha} & = (\text{by } \text{FRESH}) \\
\left\{ f \mid f (\alpha) \equiv \bar{f}(\bar{\alpha}) \right\} \uparrow \bar{\alpha} & = (\text{by } \text{FRESH}) \\
\left\{ f \mid f (\alpha) \equiv \bar{f}(\bar{\alpha}) \right\} \uparrow \bar{\alpha} & = (\text{by } \text{CONJ}) \\
\left\{ f \mid f (\alpha) \equiv \bar{f}(\bar{\alpha}) \right\} \uparrow \bar{\alpha} & = (\text{by } \text{DISJ}) \\
\left\{ f \mid f (\alpha) = \text{Zero} \right\} \uparrow \bar{\alpha} & = (\text{by } \text{THE TOTALITY OF REPRESENTING FUNCTIONS})
\end{align*}
\]

In the end we’ve got the set of representing functions, each of which restricts only the value of free variable \( \alpha \).

The final component is the semantics of specifications. Given a specification

\[ \{ R_i = \lambda x_1^i \ldots x_{k_i}^i . g_i ; i = 1 \ldots n \} g \]

we have to construct a correct environment \( \Gamma_0 \) and then take the semantics of the top-level goal:

\[ [\{ R_i = \lambda x_1^i \ldots x_{k_i}^i . g_i ; i = 1 \ldots n \} g]_\Gamma = [g]_{\Gamma_0} \]

As the set of definitions can be mutually recursive we apply the fixed point approach. We consider the following function

\[ F : (\mathcal{T}_A^+ \rightarrow 2^{A \rightarrow D})^n \rightarrow (\mathcal{T}_A^+ \rightarrow 2^{A \rightarrow D})^n \]

which represents a semantic for the set of definitions abstracted over themselves. The definition of this function is rather standard:

\[ F(p_1, \ldots, p_n) = (t_1^i \ldots t_{k_i}^i \mapsto [g^i[t_1^i/x_1^i, \ldots, t_{k_i}^i/x_{k_i}^i]]_\Gamma, \ldots, t_n^i \ldots t_{k_n}^i \mapsto [g^n[t_1^n/x_1^n, \ldots, t_{k_n}^n/x_{k_n}^n]]_\Gamma) \]

where \( \Gamma R_i = p_i \)

Here \( p_i \) is a semantic function for \( i \)-th definition; we build an environment \( \Gamma \) which associates each relational symbol \( R_i \) with \( p_i \) and construct a \( n \)-dimensional vector-function, where \( i \)-th component corresponds to a function which calculates the semantics of \( i \)-th relational definition application to terms w.r.t. the environment \( \Gamma \). Finally, we take the least fixed point of \( F \) and define the top-level environment as follows:

\[ \Gamma_0 R_i = (fix F)[i] \]
where “[i]” denotes the i-th component of a vector-function.
The least fixed point exists by Knaster-Tarski [Tarski 1955] theorem — the set \((T^* \rightarrow 2^{A \rightarrow D})^n\) forms a complete lattice, and \(T\) is monotonic.
To formalize denotational semantics in Coq we can define representing functions simply as Coq functions:

**Definition**\(^\star\) \texttt{repr\_fun} : \texttt{Set} := \texttt{var} \rightarrow \texttt{ground\_term}.

We define the semantics via inductive proposition “\texttt{in\_denotational\_sem\_goal}” such that

\[
\forall g,f : \texttt{in\_denotational\_sem\_goal} \; g \; f \iff f \in \llbracket g \rrbracket T \Gamma.
\]

The definition is as follows:

**Inductive** \texttt{in\_denotational\_sem\_goal} : \texttt{goal} \rightarrow \texttt{repr\_fun} \rightarrow \texttt{Prop} :=
| \texttt{dsgUnify} | \forall f t1 t2, \texttt{apply\_repr\_fun} f t1 = \texttt{apply\_repr\_fun} f t2 \rightarrow \texttt{in\_denotational\_sem\_goal} \texttt{(Unify t1 t2)} f |
| \texttt{dsgDisjL} | \forall f g1 g2, \texttt{in\_denotational\_sem\_goal} g1 f \rightarrow \texttt{in\_denotational\_sem\_goal} \texttt{(Disj g1 g2)} f |
| \texttt{dsgDisjR} | \forall f g1 g2, \texttt{in\_denotational\_sem\_goal} g2 f \rightarrow \texttt{in\_denotational\_sem\_goal} \texttt{(Disj g1 g2)} f |
| \texttt{dsgConj} | \forall f g1 g2, \texttt{in\_denotational\_sem\_goal} g1 f \rightarrow \texttt{in\_denotational\_sem\_goal} g2 f \rightarrow \texttt{in\_denotational\_sem\_goal} \texttt{(Conj g1 g2)} f |
| \texttt{dsgFresh} | \forall f fn a fg, \texttt{is\_fv\_of\_goal} a (\texttt{Fresh} fg) \rightarrow \texttt{in\_denotational\_sem\_goal} \texttt{(fg a)} fn \rightarrow (\forall x, x <> a \rightarrow \texttt{fn} x = f x) \rightarrow \texttt{in\_denotational\_sem\_goal} \texttt{(Fresh} fg) f |
| \texttt{dsgInvoke} | \forall r t f, \texttt{in\_denotational\_sem\_goal} \texttt{(proj1\_sig} (\texttt{Prog} r) t) f \rightarrow \texttt{in\_denotational\_sem\_goal} \texttt{(Invoke} r t) f. |

Here we refer to a fixpoint “\texttt{apply\_repr\_fun}” which calculates the extension “\(\vec{\cdot}\)” for a representing function, and inductive proposition “\texttt{is\_fv\_of\_goal}” which encodes the set of free variables for a goal.

Recall that the environment “\texttt{Prog}” maps every relational symbol to the definition of relation, which is a pair of a function from terms to goals and a proof that it has no unbound variables. So in the last case “(\texttt{proj1\_sig} (\texttt{Prog} r) t)” simply takes the body of the corresponding relation; thus “\texttt{Prog}” in Coq specification plays role of a global environment \(\Gamma\).

It is interesting that in Coq implementation we do not need to refer to Tarski-Knaster theorem explicitly since the least fixpoint semantic is implicitly provided by inductive definitions.
4 OPERATIONAL SEMANTICS

In this section we describe operational semantics of miniKanren, which corresponds to the known implementations with interleaving search. The semantics will be given in the form of labeled transition system (LTS). From now on we assume the set of semantic variables to be linearly ordered ($\mathcal{A} = \{\alpha_1, \alpha_2, \ldots\}$).

We introduce the notion of substitution

$$\sigma : \mathcal{A} \rightarrow \mathcal{T}$$

as a (partial) mapping from semantic variables to terms over the set of semantic variables. We denote $\Sigma$ the set of all substitutions, $\text{Dom}(\sigma)$ — the domain for a substitution $\sigma$, $\text{VRan}(\sigma) = \bigcup_{\alpha \in \text{Dom}(\sigma)} \text{FV}(\sigma(\alpha))$ — its range (the set of all free variables in the image).

The states in the transition system have the following shape

$$S = \mathcal{G} \times \Sigma \times \mathbb{N} | S \oplus S | S \otimes \mathcal{G}$$

As we will see later, an evaluation of a goal is separated into elementary steps, and these steps are performed interchangeably for different subgoals. Thus, a state has a tree-like structure with intermediate nodes corresponding to partially-evaluated conjunctions (“\$\otimes\$”) or disjunctions (“\$\oplus\$”). A leaf in the form $(g, \sigma, n)$ determines a goal in a context, where $g$ — a goal, $\sigma$ — a substitution accumulated so far, and $n$ — a natural number, which corresponds to a number of semantic variables used to this point. For a conjunction node its right child is always a goal since it cannot be evaluated unless some result is provided by the left conjunct.

We also need extended states

$$\bar{S} = \diamond | S$$

where $\diamond$ symbolizes the end of evaluation, and the following well-formedness condition:

**Definition 1.** Well-formedness condition for extended states:

- $\diamond$ is well-formed;
- $(g, \sigma, n)$ is well-formed iff $\text{FV}(g) \cup \text{Dom}(\sigma) \cup \text{VRan}(\sigma) \subset \{\alpha_1, \ldots, \alpha_n\}$;
- $s_1 \oplus s_2$ is well-formed iff $s_1$ and $s_2$ well-formed;
- $s \otimes g$ is well-formed iff $s$ is well-formed and for all leaf triplets $(\_\_\_, n)$ in $s$ $\text{FV}(g) \subseteq \{\alpha_1, \ldots, \alpha_n\}$.

Informally the well-formedness restricts the set of states to those in which all goals use only allocated variables. Finally, we define the set of labels:

$$L = \diamond | \Sigma \times \mathbb{N}$$

The label “$\diamond$” is used to mark those steps which do not provide an answer; otherwise a transition is labeled by a pair of a substitution and a number of allocated variables. The substitution is one of the answers, and the number is threaded through the derivation to keep track of allocated variables; we ignore it in further explanations.

The transition rules are shown on Figure 5. The first two rules specify the semantics of unification. If two terms are not unifiable under the current substitution $\sigma$ then the evaluation stops with no answer; otherwise it stops with the answer equal to the most general unifier.

The next two rules describe the steps performed when disjunction (conjunction) is encountered on the top level of the current goal. For disjunction it schedules both goals (using “$\oplus$”) for evaluating in the same context as the parent state, for conjunction — schedules the left goal and postpones the right one (using “$\otimes$”).

The rule for “fresh” substitutes bound syntactic variable with a newly allocated semantic one and proceeds with the goal; no answer provided at this step.
\[\langle t_1 \equiv t_2, \sigma, n \rangle \xrightarrow{\diamond} mgu(t_1, t_2, \sigma) \quad \text{[UNIFYFAIL]}\]

\[\langle t_1 \equiv t_2, \sigma, n \rangle \xrightarrow{mgu(t_1, t_2, \sigma, n)} \diamond \quad \text{[UNIFYSUCCESS]}\]

\[\langle g_1 \lor g_2, \sigma, n \rangle \xrightarrow{\diamond} (g_1, \sigma, n) \oplus (g_2, \sigma, n) \quad \text{[DISJ]}\]

\[\langle g_1 \land g_2, \sigma, n \rangle \xrightarrow{\diamond} (g_1, \sigma, n) \otimes g_2 \quad \text{[CONJ]}\]

\[\langle \text{fresh } x . g, \sigma, n \rangle \xrightarrow{\diamond} (g[\alpha_{n+1}/x], \sigma, n + 1) \quad \text{[FRESH]}\]

\[R^{k_i}_i = \lambda x_1 \ldots x_{k_i} \cdot g \quad \text{[INVOK]}\]

\[\langle R^{k_i}_i (t_1, \ldots, t_{k_i}), \sigma, n \rangle \xrightarrow{\diamond} \left\langle g[t_1/x_1 \ldots t_{k_i}/x_{k_i}], \sigma, n \right\rangle \quad \text{[DISJSTOP]}\]

\[s_1 \xrightarrow{\diamond} \quad \text{[DISJSTOPANS]}\]

\[s_1 \xrightarrow{r} \diamond \quad \text{[CONJSTOP]}\]

\[s \xrightarrow{\diamond} \quad \text{[CONJSTOPANS]}\]

\[s \xrightarrow{(\sigma, n)} \quad \text{[DISJSTOPANS]}\]

\[s_1 \xrightarrow{\diamond} s'_1 \quad \text{[DISJSTEP]}\]

\[s_1 \xrightarrow{r} s'_1 \quad \text{[DISJSTEPANS]}\]

\[s \xrightarrow{(\sigma, n)} s' \quad \text{[CONJSTEP]}\]

\[s \xrightarrow{(\sigma, n)} (s' \otimes g) \quad \text{[CONJSTEPANS]}\]

\[s \xrightarrow{(\sigma, n)} (s' \otimes g) \quad \text{[CONJSTEPANS]}\]

Fig. 5. Operational semantics of interleaving search

The rule for relation invocation finds a corresponding definition, substitutes its formal parameters with the actual ones, and proceeds with the body.

The rest of the rules specify the steps performed during the evaluation of two remaining types of the states — conjunction and disjunction. In all cases the left state is evaluated first. If its evaluation stops with a result then the right state (or goal) is scheduled for evaluation, and the label is propagated. If there is no result then the
conjunction evaluation stops with no result (ConjStop) as well while the disjunction evaluation proceeds with the right state (DisjStop).

The last four rules describe interleaving, which occurs when the evaluation of the left state suspends with some residual state (with or without an answer). In the case of disjunction the answer (if any) is propagated, and the constituents of the disjunction are swapped (DisjStep, DisjStepAns). In case of conjunction, if the evaluation step in the left conjunct did not provide any answer, the evaluation is continued in the same order since there is still no information to proceed with the evaluation of the right conjunct (ConjStep); if there is some answer, then the disjunction of the right conjunct in the context of the answer and the remaining conjunction is scheduled for evaluation (ConjStepAns).

The introduced transition system is completely deterministic. There was, however, some freedom in choosing the order of evaluation for conjunction and disjunction states. For example, instead of evaluating the left substate first we could choose to evaluate the right one, etc. In each concrete case we would end up with a different (but still deterministic) system which would prescribe different semantics to a concrete goal. This choice reflects the inherent non-deterministic nature of search in relational (and, more generally, logical) programming. However, as long as deterministic search procedures are sound and complete, we can consider them "equivalent".

A derivation sequence for a certain state determines a trace—a finite or infinite sequence of answers. We may define a set of finite or infinite sequences $X^{\omega}$ over an alphabet $X$ as a set of functions from natural numbers into a lifted set $X_{\perp} = X \cup \{\perp\}$:

$$X^{\omega} = \{\omega : \mathbb{N} \rightarrow X_{\perp} \mid \forall n \in \mathbb{N}, \omega(n) = \perp \Rightarrow \omega(n + 1) = \perp\}$$

Informally speaking, we represent a sequence as a function which maps positions (treated as natural numbers) into the elements of the sequence. We use "$\perp$" to specify that there is no element at given position, and we stipulate, that there are no "holes" in this representation: if there is no element at given position then there are no elements at greater positions as well.

For this representation we may define the empty sequence $\varepsilon$ and operations of prepending a sequence $\omega$ with an element $a$ and taking a suffix of a sequence $\omega$ from a position $n$ as follows:

$$\varepsilon = i \mapsto \perp$$

$$a\omega = i \mapsto \begin{cases} a, & i = 0 \\ \omega(i - 1), & \text{otherwise} \end{cases}$$

$$\omega[n:] = i \mapsto \omega(n + i)$$

For a given state $s$ a trace $\mathcal{T}_r \in L^{\omega}$ is a sequence of labels, defined as follows simultaneously with the sequence of states $\{s_i\}$:

$$s_0 = s$$

$$\mathcal{T}_r(n) = a, \quad s_{n+1} = s' \quad \text{if} \quad s_n \neq \diamond, \quad s_n \xrightarrow{a} x'$$

$$\mathcal{T}_r(n) = \perp, \quad s_{n+1} = \diamond \quad \text{if} \quad s_n = \diamond$$

The trace corresponds to the stream of answers in the reference MiniKanren implementations.

To formalize the operational part in Coq we first need to define all preliminary notions from unification theory [Baader and Snyder 2001] which our semantics uses.

\footnote{There still can be differences in observable behavior of concrete goals under different sound and complete search strategies: a goal can be refutationally complete [Byrd 2009] under one strategy and non-complete under another.}
In particular, we need to implement the notion of the most general unifier (MGU). As is well-known [McBride 2003] all standard recursive algorithms for calculating MGU are not decreasing on argument terms, so we can't define it as a simple recursive function in Coq due to the termination check. There is no such obstacle when we define MGU as a proposition:

\textbf{Inductive} \textsc{Mgu} : \text{term} \to \text{term} \to \text{option subst} \to \text{Set} := ... \\
However, we still need to use a well-founded induction to prove the existence of the most general unifier and its defining properties:

\textbf{Lemma} \textsc{Mgu}_\textsc{ex} : \forall t1 t2, \{ r & \textsc{Mgu} t1 t2 r \}.

\textbf{Definition} \textsc{unifier} (s : subst) (t1 t2 : term) : \text{Prop} := \text{apply_subst} s t1 = \text{apply_subst} s t2.

\textbf{Lemma} \textsc{Mgu}_\textsc{unifies}:
\forall t1 t2 s, \textsc{Mgu} t1 t2 (\text{Some} s) \to \textsc{unifier} s t1 t2.

\textbf{Definition} \textsc{more}_\textsc{general} (m s : subst) : \text{Prop} := \exists (s’ : subst), \forall (t : \text{term}), \text{apply_subst} s t = \text{apply_subst} s’ (\text{apply_subst} m t).

\textbf{Lemma} \textsc{Mgu}_\textsc{most}_\textsc{general}:
\forall (t1 t2 : \text{term}), (m : subst), \textsc{Mgu} t1 t2 (\text{Some} m) \to
\forall (s : subst), \textsc{unifier} s t1 t2 \to \textsc{more}_\textsc{general} m s.

\textbf{Lemma} \textsc{Mgu}_\textsc{non}_\textsc{unifiable}:
\forall (t1 t2 : \text{term}),
\textsc{Mgu} t1 t2 \text{None} \to \exists s, \neg (\text{unifier} s t1 t2).

For this well-founded induction we use the number of free variables in argument terms as a well-founded order on pairs of terms:

\textbf{Definition} \textsc{terms} := \text{term} \times \text{term}.

\textbf{Definition} \textsc{fvOrder} (t : \text{terms}) := \text{length} (\text{union} (\text{fv_term} (\text{fst} t)) (\text{fv_term} (\text{snd} t))).

\textbf{Definition} \textsc{fvOrderRel} (t p : \text{terms}) := \textsc{fvOrder} t < \textsc{fvOrder} p.

\textbf{Lemma} \textsc{fvOrder_wf} : \text{well}_\text{fed} \textsc{fvOrderRel}.

After this preliminary work, the described transition relation can be encoded naturally as an inductively defined proposition (here "state'" stands for an extended state):

\textbf{Inductive} \textsc{eval_step} : \text{state} \to \text{label} \to \text{state'} \to \text{Set} := ...

We state the fact that our system is deterministic through existence and uniqueness of a transition for every state:

\textbf{Lemma} \textsc{eval_step_ex} : \forall (st : \text{state}), \{ l : \text{label} & \{ st' : \text{state'} & \textsc{eval_step} st l st' \} \}.
Lemma eval_step_unique:
\[
\forall (st : \text{state}) (l1 l2 : \text{label}) (st'1 st'2 : \text{state}'),
\quad \text{eval_step } st \text{ l1 } st'1 \rightarrow \text{eval_step } st \text{ l2 } st'2 \rightarrow l1 = l2 \land st'1 = st'2.
\]

To work with (possibly) infinite sequences we use the standard approach in Coq — coinductively defined streams:

Context \{ A : \text{Set} \}.

CoInductive stream : \text{Set} :=
| Nil : stream
| Cons : A → stream → stream.

Although the definition of the datatype is coinductive some of its properties we are working with make sense only when defined inductively:

Inductive in_stream : A → stream → Prop :=
| inHead : ∀ x t, in_stream x (Cons x t)
| inTail : ∀ x h t, in_stream x t → in_stream x (Cons h t).

Inductive finite : stream → Prop :=
| fNil : finite Nil
| fCons : ∀ h t, finite t → finite (Cons h t).

Then we define a trace coinductively as a stream of labels in transition steps and prove that there exists a unique trace from any extended state:

Definition trace : \text{Set} := @stream label.

CoInductive op_sem : state' → trace → \text{Set} :=
| osStop : op_sem Stop Nil
| osState : ∀ st l st' t, eval_step st l st' →
\quad \text{op_sem } st' t →
\quad \text{op_sem } (\text{State } st) (Cons l t).

Lemma op_sem_ex(st' : state') : {t : trace & op_sem st' t}.

Lemma op_sem_unique :
\[
\forall st' t1 t2, \text{op_sem } st' t1 \rightarrow \text{op_sem } st' t2 \rightarrow \text{equal_streams } t1 t2.
\]

Note, for the equality of streams we need to define a new coinductive proposition instead of using the standard syntactic equality in order for coinductive proofs to work [Chlipala 2013].

One thing we can prove using operational semantics is the interleaving properties of disjunction. Specifically, we can prove that a trace for a disjunction is a one-by-one interleaving of streams for its disjuncts:

CoInductive interleave : stream → stream → stream → Prop :=
| interNil : ∀ s s', \text{equal_streams } s s' → interleave Nil s s'
| interCons : ∀ h t s rs, interleave s t rs → interleave (Cons h t) s (Cons h rs).
Lemma sum_op_sem: ∀ st1 st2 t1 t2 t, op_sem (State st1) t1 →
   op_sem (State st2) t2 →
   op_sem (State (Sum st1 st2)) t →
   interleave t1 t2 t.

This allows us to prove the expected properties of interleaving in a more general setting of arbitrary streams:

- the elements of the interleaved stream are exactly those of two interleaved streams;
- the interleaved stream is finite iff both interleaving streams are finite.

The corresponding Coq lemmas are as follows:

Lemma interleave_in: ∀ s1 s2 s, interleave s1 s2 s →
   ∀ x, in_stream x s ↔ in_stream x s1 ∨ in_stream x s2.

Lemma interleave_finite: ∀ s1 s2 s, interleave s1 s2 s →
   (finite s ↔ finite s1 ∧ finite s2).

5 SEMANTICS EQUIVALENCE

Now when we defined two different kinds of semantics for miniKanren we can relate them and show that the results given by these two semantics are the same for any specification. This will actually say something important about the search in the language: since operational semantics describes precisely the behavior of the search and denotational semantics ignores the search and describes what we should get from mathematical point of view, by proving their equivalence we establish completeness of the search which means that the search will get all answers satisfying the described specification and only those.

But first, we need to relate the answers produced by these two semantics as they have different forms: a trace of substitutions (along with numbers of allocated variables) for operational and a set of representing functions for denotational. We can notice that the notion of representing function is close to substitution, with only two differences:

- representing function is total;
- terms in the domain of representing function are ground.

Therefore we can easily extend (perhaps ambiguously) any substitution to a representing function by composing it with an arbitrary representing function and that will preserve all variable dependencies in the substitution. So we can define a set of representing functions corresponding to substitution as follows:

\[[\sigma] = \{ f \circ \sigma \mid f : \mathcal{A} \mapsto \mathcal{D} \}\]

And denotational analog of an operational semantics (a set of representing functions corresponding to answers in the trace) for given extended state \(s\) is then defined as a union of sets for all substitution in the trace:

\[[s]_{op} = \sqcup_{(\sigma, n) \in T_r} [\sigma]\]

This allows us to state theorems relating two semantics.

Theorem 1 (Operational semantics soundness). For any specification \(\{\ldots\} g\), for which the indices of all free variables in \(g\) are limited by some number \(n\)

\[[\langle g, e, n \rangle]_{op} \subset [[\ldots] g].\]
It can be proven by nested induction, but first, we need to generalize the statement so that the inductive hypothesis would be strong enough for the inductive step. To do so, we define denotational semantics not only for goals but for arbitrarily extended states. Note that this definition does not need to have any intuitive interpretation, it is introduced only for proof to go smoothly. The definition of the denotational semantics for extended states is on Figure 6. The generalized version of the theorem uses it:

**Lemma 1 (Generalized soundness).** For any top-level environment $\Gamma_0$ acquired from some specification, for any well-formed (w.r.t. that specification) extended state $s$

$$[s]_{op} \subseteq [s]_{\Gamma_0}.$$  

It can be proven by induction on the number of steps in which a given answer (more accurately, the substitution that contains it) occurs in the trace. The induction step is proven by structural induction on the extended state $s$.

It would be tempting to formulate the completeness of operational semantics as the inverse inclusion, but it does not hold in such generality. The reason for this is that denotational semantics encodes only dependencies between the free variables of a goal, which is reflected by the completeness condition, while operational semantics may also contain dependencies between semantic variables allocated in "fresh". Therefore we formulate the completeness with representing functions restricted on the semantic variables allocated in the beginning (which includes all free variables of a goal). This does not compromise our promise to prove the completeness of the search as miniKanren provides the result as substitutions only for queried variables, which are allocated in the beginning.

**Theorem 2 (Operational semantics completeness).** For any specification $\{ \ldots \} g$, for which the indices of all free variables in $g$ are limited by some number $n$

$$\{ f | \{ a_1, \ldots, a_n \} | f \in [\{ \ldots \} g]_{op} \} \subseteq \{ f | \{ a_1, \ldots, a_n \} | f \in [\langle g, \sigma, n \rangle]_{\Gamma_0} \}.$$  

Similarly to the soundness, this can be proven by nested induction, but the generalization is required. This time it is enough to generalize it from goals to states of the shape $[\langle g, \sigma, n \rangle]_{\Gamma_0}$. We also need to introduce one more auxiliary semantics — bounded denotational semantics:

$$[\bullet]^l : \mathcal{G} \to 2^{\mathcal{A} \to D}$$  

Instead of always unfolding the definition of a relation for invocation goal, it does so only given number of times. So for a given set of relational definitions $\{ R_{ki}^l = \lambda x_1^i \ldots x_{k_i}^i . g_i ; \}$ the definition of bounded denotational semantics is exactly the same as in usual denotational semantics, except that for the invocation case:

$$[\langle R_{ki}^l (t_1, \ldots, t_{k_i}) \rangle]^{l+1} = [\langle g_i[t_1/x_1^i, \ldots, t_{k_i}/x_{k_i}^i] \rangle]^l$$

It is convenient to define bounded semantics for level zero as an empty set:
Bounded denotational semantics is an approximation of a usual denotational semantics and it is clear that any answer in usual denotational semantics will also be in bounded denotational semantics for some level:

**Lemma 2.** \([g]_0 \subseteq \bigcup_l [g]^l\)

Formally it can be proven using the definition of the least fixed point from Tarski-Knaster theorem: the set on the right-hand side is a closed set.

Now the generalized version of the completeness theorem is as follows:

**Lemma 3 (Generalized completeness).** For any set of relational definitions, for any level \(l\), for any well-formed \((w.r.t.\ that\ set\ of\ definitions)\ state \(\langle g, \sigma, n \rangle\),

\[
\{f|_{\alpha_1, \ldots, \alpha_n} | f \in [g] \cap [\sigma] \} \subset \{f|_{\alpha_1, \ldots, \alpha_n} | f \in [(g, \sigma, n)]\}_{op}\}
\]

It is proven by induction on the level \(l\). The induction step is proven by structural induction on the goal \(g\).

The proofs of both theorems are certified in Coq, although the proofs for a number of (obvious) technical facts about representing functions and computation of the most general unifier as well as some properties of denotational semantics, proven informally in Section 3, are admitted for now. For completeness we can not just use the induction on proposition \(\text{in\_denotational\_sem\_goal}\), as it would be natural to expect, because the inductive principle it provides is not flexible enough. So we need to define bounded denotational semantics in our formalization too and perform induction on the level explicitly:

**Inductive** \(\text{in\_denotational\_sem\_lev\_goal} : \text{nat} \rightarrow \text{goal} \rightarrow \text{repr\_fun} \rightarrow \text{Prop} := \)

\(
\ldots
| \text{dslgInvoke} : \forall l \ t \ f, \hspace{1cm} \\
\text{in\_denotational\_sem\_lev\_goal} \ l \ (\text{proj1\_sig} (\text{Prog} \ t)) \ f \rightarrow \hspace{1cm} \\
\text{in\_denotational\_sem\_lev\_goal} \ (\text{S} \ l) \ (\text{Invoke} \ t \ f).
\)

The lemma relating bounded and unbounded denotational semantics is translated into Coq:

**Lemma** \(\text{in\_denotational\_sem\_some\_lev}: \forall (g : \text{goal}) \ (f : \text{repr\_fun}), \hspace{1cm} \hspace{1cm} \\
\text{in\_denotational\_sem\_goal} \ g \ f \rightarrow \hspace{1cm} \hspace{1cm} \\
\exists l, \text{in\_denotational\_sem\_lev\_goal} \ l \ g \ f.
\)

The statements of the theorems are as follows:

**Theorem** \(\text{search\_correctness}: \forall (g : \text{goal}) \ (k : \text{nat}) \ (f : \text{repr\_fun}) \ (t : \text{trace}), \hspace{1cm} \hspace{1cm} \\
\text{closed\_goal\_in\_context} \ (\text{first\_nats} \ k) \ g \rightarrow \hspace{1cm} \hspace{1cm} \\
\text{op\_sem} \ (\text{State} (\text{Leaf} \ g \ \text{empty\_subst} \ k)) \ t \rightarrow \hspace{1cm} \hspace{1cm} \\
\text{in\_denotational\_analog} \ t \ f \rightarrow \hspace{1cm} \hspace{1cm} \\
\text{in\_denotational\_sem\_goal} \ g \ f.
\)

**Theorem** \(\text{search\_completeness}: \forall (g : \text{goal}) \ (k : \text{nat}) \ (f : \text{repr\_fun}) \ (t : \text{trace}), \hspace{1cm} \hspace{1cm} \\
\text{closed\_goal\_in\_context} \ (\text{first\_nats} \ k) \ g \rightarrow \hspace{1cm} \hspace{1cm} \\
\text{op\_sem} \ (\text{State} (\text{Leaf} \ g \ \text{empty\_subst} \ k)) \ t \rightarrow \hspace{1cm} \hspace{1cm} \\
\text{in\_denotational\_sem\_goal} \ g \ f \rightarrow \hspace{1cm} \hspace{1cm} \\
\exists (f' : \text{repr\_fun}), \ (\text{in\_denotational\_analog} \ t \ f') \land \hspace{1cm} \hspace{1cm} \\
\forall (x : \text{var}), \ \text{In} x (\text{first\_nats} \ k) \rightarrow f x = f' x.
\)
One important immediate corollary of these theorems is the correctness of certain program transformations. Since the results obtained by the search on a specification are exactly the results from the mathematical model of this specification, after the transformations of relations that do not change their mathematical meaning the search will obtain the same results. Note that this way we guarantee only the stability of results as the set of ground terms, the other aspects of program behavior, such as termination, may be affected. This allows us to safely (to a certain extent) apply such natural transformations as:

- changing the order of constituents in conjunction or disjunction;
- swapping conjunction and disjunction using distributivity;
- moving fresh variable introduction.

and even transform relational definitions to some kinds of normal form (like all fresh variables introduction on the top level with the conjunctive normal form inside), which may be convenient, for example, for metacomputation.

6 CONCLUSION AND FUTURE WORK

In this paper we presented a formal semantics for core miniKanren and proved some its basic properties, which are believed to hold in existing implementations. We consider our work as an initial setup for a future development of miniKanren semantics. The language we considered here lacks many important features, which are already introduced and employed in many implementations. Integrating these extensions — in the first hand, disequality constraints, — into the semantics looks a natural direction for future work. We also are going to address the problems of proving some properties of relational programs (equivalence, refutational completeness, etc.).

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Relational Processing for Fun and Diversity
Simulating a CPU relationally with miniKanren

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Defining a central processing unit relationally using miniKanren is proposed as a new approach for realizing assembly code diversification. Software diversity has long been championed as a means of protecting digital ecosystems from widespread failures due to cyberattacks and faults, but is often difficult to achieve in practice. Using relational programming to simulate a processor allows large-scale automatic synthesis of assembly-level code. Early experiments with the technique indicate that such synthesis might lead to better automation of code diversification by breaking the synthesis problem into manageable chunks. An early prototype is presented, with some sample synthesis tasks and discussion of possible future applications.

CCS Concepts:
• Software and its engineering → Search-based software engineering; Software safety; Interpreters; Source code generation;
• Security and privacy → Software security engineering;

Additional Key Words and Phrases: program synthesis, relational programming, miniKanren, software artificial diversity, malware polymorphism

1 INTRODUCTION

Software diversity has long been recognized as valuable for protecting digital ecosystems from widespread failures due to cyberattacks and faults [Cohen 1993]. Higher diversity of software implementations reduces the likelihood that a single, common flaw pervades all its deployments, and therefore that a single attack can compromise all members of the ecosystem or affect them all in the same way. Unfortunately, most software ecosystems today remain highly monocultural—all deployments of a given software product for a given architecture are almost identical, save for minute differences. This homogeneity often allows individual, low-cost cyberattacks to have a devastating impact on large numbers of computer systems. For example, the 2015 Stagefright 2.0 vulnerability left nearly all Android devices susceptible to remote compromise due to almost identical multimedia library implementations being used by nearly all apps on all the devices [Peters 2015].

One reason why software monoculture continues to abound despite its brittleness against attack is the uniformity of most present-day tools for developing software products. Compilers are typically designed to solve an optimization problem that translates a given source program into a single, efficient, equivalently behaved object program. Compiler design goals therefore typically include semantic transparency (behavioral preservation), runtime efficiency, and space efficiency, but not diversity. The source semantics of mainstream imperative languages (e.g., C/C++), the optimization stages of their compilers, and their backend code generation algorithms are all designed to search for a single, good solution to this optimization problem.

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Our research seeks to shift the compiler optimization problem to pursue diversity as a design goal of software development. In particular, a diversifying software development methodology should attempt to yield a maximally dissimilar collection of object code implementations that are all semantically transparent to the original source code and that meet a certain baseline efficiency.

Although a few code diversification strategies have attained widespread deployment in practice, most provide only limited forms of diversity that continue to leave ecosystems vulnerable to exploitation. For example, Address Space Layout Randomization (ASLR) diversifies programs by randomly choosing the base addresses of libraries at load-time. While this does help mitigate some attacks, the resulting diversity is low, leaving the software vulnerable to derandomization attacks [Shacham et al. 2004]. We believe that changing implementations in a more fundamental way, such as modifying how values are computed, will protect against broader classes of attacks.

Historically, however, this type of diversity has been expensive to obtain and difficult to achieve in an automated fashion. Prior work [Lundquist et al. 2016] proposes merging the fields of artificial diversity of software with that of program synthesis, leveraging the natural diversity of search-style problems in order to create a plethora of program implementations. Because they spring from search problems, these implementations could be fundamentally different ways of solving a given computational problem.

One approach to program synthesis makes use of miniKanren1 [Byrd 2009], a family of logic languages typically implemented as an embedded Domain-Specific Language (DSL). Since it is a relational language, miniKanren programmers write specifications that relate values, and users submit queries that yield sets of values within the relation. Since any value can queried by the system, values in computations that are traditionally thought of as “inputs” or “outputs” need not be. Querying for “inputs” that relate to a specified “output” effectively reverses the computation.

Prior work [Byrd et al. 2017, 2012] has proposed realizing program synthesis in miniKanren by implementing the relational specification of an interpreter that relates input code to the output values it produces. Reversing the computation by querying for inputs from given outputs then produces possible code that produces the desired output values.

Program synthesis in miniKanren seems particularly well suited to the goal of code diversity given miniKanren’s natural ability to easily produce multiple answers for a query. The user simply specifies how many answers (at a maximum) the system is to search for, and a list of results is returned.

In the interest of being source code-agnostic with our diversification efforts, we would like to synthesize programs at the assembly level. This potentially has the advantages of breaking the synthesis problem into manageable chunks, as well as allowing existing code (of any origin) to be diversified in a largely automated fashion.

Since we wish to achieve program diversity at the assembly level, we propose writing a relational specification of a central processing unit. This miniKanren program interprets assembly code by relating that code to the states of the processor before and after code execution. A user can then query for assembly code that produces a desired output state.

2 MODES OF OPERATION
We envision the following ways of using an assembly language synthesis system, such as our miniKanren prototype:

2.1 For Diversification
Our main goal is to explore automated diversification of assembly-level code. Potential approaches include the following:

1See http://minikanren.org
Use of code sketches:
A standard approach to program synthesis is that of the program sketch [Solar-Lezama 2008, 2009]: a partial program with holes to be filled with synthesized code. In a logic programming environment, a sketch is a partial program with portions (here instruction mnemonics, instruction arguments, etc.) represented by logic variables. Those variables are then queried using a run command to determine possible values. To control which output is being computed, a sketch must also include a formal specification of what is to be computed. In assembly diversification, these specifications are constraints dictating the possible processor states after a computation runs. (We use sketches in this paper.) Constraints collectively define correctness of synthesized answers. Correct synthesis algorithms yield only answers that are correct with respect to the constraints. For code diversification, constraints must therefore also define what is meant by program equivalence. For example, a user may decide to specify result values for processor status flags, or leave them unconstrained. The latter option broadens the definition of equivalence to include relations on processor states. For example, diversification might be required to preserve register and memory values but be permitted to vary status flag values.

When diversifying existing code, either the program to be diversified already has a formal (mathematical) specification, or (more commonly) there will a priori be no formal notion of correctness. In the ideal former case, we need only convert the specification into miniKanren constraints. In the latter case, quality assurance and testing processes must be applied to check program correctness or equivalence; however, for diversity to be tractable, this assurance process should be (semi-)automated to easily apply it to the multitude of variants generated.

Basic blocks bracketed by control flow instructions:
Generating proper control-flow using program synthesis is known to be a hard problem [Solar-Lezama 2008]. To avoid this, sketches can explicitly specify most or all control-flow transfer instructions [Solar-Lezama 2008]. Instructions in the basic block(s) between control-flow transfers are synthesized in entirety or in part. Program constraints include loop invariants to dictate what is synthesized. (The GCD examples in the following section illustrate this approach.)

Use of effect traces:
In addition to dividing code up into basic blocks, another approach is to generate code based on a known effect trace, synthesizing portions between calls to other functions, system routines, or other side effects not otherwise modeled by our processor relation. The format and construction of arguments to these external routines are determined by constraints provided in the sketch. Since trace equivalence is a common way to define program equivalence, this approach provides an intuitive starting point for solving the equivalence problem mentioned above.

Diversification of existing code:
One of our primary goals for this project is the automated diversification of existing code. To achieve this, processor states could be generated by running existing assembly code forwards in our system. Enhanced by constraints describing known properties of how the code should function, these processor states become the end goals for new code to satisfy.

Gadget-oriented program composition:
A gadget can be defined as any arbitrary string of assembly language instructions, usually one found in pre-existing code. As shown in prior work [Lundquist et al. 2016; Mohan and Hamlen 2012] (and as shown by Return-Oriented Programming in general [Schwartz et al. 2011]), programs can be constructed by stringing chains of gadgets together in an order determined by the synthesis engine as a means of reaching a particular goal. Return-Oriented Programming (ROP) is an exploit technique that repurposes gadgets found in existing benign code to implement attack payloads. While gadget-oriented programming was
originally used for malicious purposes, we speculate that this technique could be adapted for more general synthesis tasks.

The overall theme of the techniques described above is that of breaking the problem of synthesis into small, manageable chunks at a low level. Rather than trying to synthesize an entire program at once, a program is broken down into sub-pieces, which are then synthesized. This avoids traditional problems associated with synthesizing large or complex code fragments, and mitigates state-space explosion.

2.2 For Use by a Compiler

Since programming using processor states and constraint sets is likely to be difficult, we envision this system as a target of higher level tools, such as compilers or interpreters. Replacing the back-end code generation of a compiler with a synthesis system allows for automatic diversification while keeping programming tasks manageable. Compilers are potentially in a unique position to know what constraints must be preserved from higher-level source in the generated assembly. Further, allowing compilers and other code-generation tools to programmatically break the result into known intermediate states allows for smaller and more frequent synthesis tasks, once again keeping with the theme of reducing synthesis problems into smaller sub-problems. Sub-problems can be divided at whatever level is appropriate for the tool based on its knowledge and analysis of the code being generated, or based on what size tasks the synthesis engine is capable of handling efficiently.

We plan to use our system to explore each of the above approaches in future work.

3 IMPLEMENTATION

To explore assembly language synthesis, we have developed a prototype relational assembly interpreter in miniKanren.\(^2\) Our prototype uses a subset of Intel x86 assembly language instructions with the limited set of 32-bit general purpose registers available for that architecture. The implementation uses a modified version of faster-miniKanren\(^3\) written in Racket.

A full scale system must support enough instructions and processor features to synthesize programs that work; but like existing compiler back-ends (which emit only a subset of the available instructions), this does not necessitate supporting the entire instruction set architecture (ISA). Supporting more of the ISA allows for more diversity in the resulting programs, potentially at the cost of synthesis time. However, prior work [Mohan and Hamlen 2012] has found that only a small portion of the architecture is needed to achieve high code diversity.

3.1 The x86\(^{o}\) Relation

Our processor relation \(x86^{o}\) relates assembly code, an input processor state, and an output processor state. Processor state is modeled as a set of association lists, one mapping registers to values and another mapping memory locations to values. The memory mapping is a partial function, only containing those values that have been updated by the program. Addresses read from an uninitialized address return a default value, typically 0.

The interpreter models assembly code as a list of instructions, using a fall-through approach to execute (or synthesize) a basic block. Since we do not yet represent code addresses in our model, execution always necessarily continues to the subsequent instruction in the list. Each instruction is a list containing an instruction mnemonic and the appropriate number of operands for the instruction. Each operand describes a register, an immediate value, or a memory address.


\(^3\) Obtained from https://github.com/greggr/tutorial-relational-interpreters. This version contains modifications by Greg Rosenblatt to speed up the \texttt{eval}\(^{o}\) relational interpreter. The original faster-miniKanren can be obtained from https://github.com/michaelballantyne/faster-miniKanren/blob/master/README.md.
The interpreter calls sub-relations to decode arguments (which relate operands to their values with respect to some processor state) and then interprets the instruction by operating on those values and creating an updated processor state. A fragment of the relation with a few instructions is shown below:

;; The x86 processor relation.
(define (x86 o code rstore new-rstore)
  (conde
    [(≡ '() code) (≡ rstore new-rstore)]
    [(fresh (opcode arglist morecode op1 op2 res rstore1)
      (≡ code ((,opcode ,arglist) ,morecode))
      (conde
        ;; add
        [(≡ opcode 'add)
          (decode-2args o arglist op1 op2 rstore)
          (plus o op1 op2 res)
          (update-arg o arglist res rstore rstore1)
          (x86 o morecode rstore1 new-rstore)]
        . . .
        ;; div
        ;; uses implicit arguments -
        ;; dividend is edx:eax,
        ;; divisor is op1,
        ;; destination (result quotient) is eax,
        ;; (result) remainder is edx
        [(≡ opcode 'div)
          (decode-1arg o arglist op1 rstore)
          (fresh (edx eax n rem rstore2)
            (lookup o 'R_EDX rstore eax)
            (lookup o 'R_EAX rstore edx)
            (append o eax edx n) ; n=dividend: eax=low order bits, edx=high order bits
            (div o n op1 res rem)
            (update o 'R_EAX res rstore rstore1)
            (update o 'R_EDX rem rstore1 rstore2)
            (x86 o morecode rstore2 new-rstore))]
        . . .
        ;; xor
        [(≡ opcode 'xor)
          (decode-2args o arglist op1 op2 rstore)
          (xor o op1 op2 res)
          (update-arg o arglist res rstore rstore1)
          (x86 o morecode rstore1 new-rstore)]
        . . .]}

...
Numeric values in our system are represented as Oleg numerals—little-endian lists of binary digits that encode the base-2 representation of the number. Most arithmetic operations are the definitions found in The Reasoned Schemer [Friedman et al. 2018], with a few of our own added to implement missing operators (e.g., logical operations) needed for the instructions we’ve included. For example, here is our implementation of integer division and remainder:

; integer division with remainder
(define (div^ o a b q r)
  (conde
   [((≡ q ') (≡ r a) (<l^ a b))]
   [(fresh (p)
     (≡p b a)
     (<l^ r b)
     (plus^ p r a)
     (*^ b q p))])))

3.2 Example Queries

3.2.1 Synthesis 1: Generating Some Assembly. Our first example interaction demonstrates a query for generating lots of assembly code quickly. The goal of each synthesized program is to place the result value 1875 (Oleg numeral (1 1 0 0 1 0 1 0 1 1 1)) into register EAX.

We begin by providing the following assembly program sketch:

mov ECX ?_x
mov EDX ?_y
mov EAX ?_p
mov EBX ?_q
mov EDI ?_s
mov ESI ?_t

?_r ECX EDI
?_f EDX EDI
?_d EAX ECX
?_e EBX EDX
?_c EAX EBX

The question marks denote holes to be filled in by the synthesizer, subscripted with the name of the logic variable used to denote the hole. The first six instructions denote mov instructions with a particular destination operand (one for each general-purpose register) and a hole to be filled in for the source operand. The remaining five instructions give concrete destination and source operands but leave the choice of instruction open to be synthesized.

Constraints are then added to further limit what will be synthesized. The first set of constraints prevent any of the final instructions from being additional mov instructions. The second set prevents some source operands from

\footnote{Obtained from https://github.com/miniKanren/CodeFromTheReasonedSchemer2ndEd.}
being the same as some of the other source operands. The third set of constraints disallow source operand holes from containing 1875, which prevents the goal value from being moved directly into a register. Finally, source operands must be positive (non-zero).

We use our processor relation to relate the sketch code, an initial empty processor state, and a final processor state. Constraining this final state by the value in EAX forces our desired goal to be met. We then query for the fully synthesized code (in variable v1) and the resulting processor state (in variable z).

For our queries, we make use of Racket's time operator, which reports (in milliseconds) CPU time, real time, and garbage collection time spent for expression evaluation. The full query and first 2 (of 200) results are as follows:

```scheme
> (time
  (run 200 (v1 z)
    (fresh (d x p e q y c s t f g)
      (≡ v1
        ((mov R_ECX ,x)
         (mov R_EDX ,y)
         (mov R_EAX ,p)
         (mov R_EBX ,q)
         (mov R_EDI ,s)
         (mov R_ESI ,t)
         ,g��X R_ESI)
       ,f��X��DI)
       ,d��X��CX)
       ,e��X��BX��EDX)
       ,c��X��AX��EDX)))

(cpu time: 21109 real time: 21481 gc time: 9204
'(((mov��X��ECX ,0 . ,1))
  (mov��X��EDX ,2 . ,3))
  (mov��X��AX (1))
  (mov��X��BX (0 1 0 0 1 0 1 0 1 1 1))
  (mov��X��DI (0 . ,2))
  (mov��X戝SI (0 . ,3))
  (sub��X戝CX��ESI)
  (sub戝X戝CX戝DI))

(((mov��X戝CX��ESI)
  (mov��X戝CX��AX (0 1 0 0 1 0 1 0 1 1 1))
  (mov��X戝CX戝DI (0 . ,2))
  (mov戝ESI (0 . ,3))
  (sub戝CX戝DI)
  (sub戝CX戝DI)))
```
(add R_EAX R_ECX)
(add R_EBX R_EDX)
(add R_EAX R_EBX))
((R_EAX 1 1 0 0 1 0 1 0 1 1 1) (R_EBX 0 1 0 0 1 0 1 0 1 1 1) (R_ECX) (R_EDX)
(R_EDI _.2 . _.3) (R_ESI _.0 . _.1)))
(=/= ((_.0 0) (_.1 (1 0 0 1 0 1 0 1 1 1)))
((_.0 1) (_.1 (1 0 0 1 0 1 0 1 1 1)))
((_.0 1) (_.1 ()))
((_.0 _.2) (_.1 _.3))
((_.2 0) (_.3 (1 0 0 1 0 1 0 1 1 1)))
((_.2 1) (_.3 (1 0 0 1 0 1 0 1 1 1)))
((_.2 1) (_.3 ()))))))
(((mov R_ECX (_.0 . _.1))
(mov R_EDX (_.2 . _.3))
(mov R_EAX (0 1 0 0 1 0 1 0 1 1 1))
(mov R_EBX (1))
(mov R_EDI (_.2 . _.3))
(mov R_ESI (_.0 . _.1))
(sub R_ECX R_ESI)
(sub R_EDX R_EDI)
(add R_EAX R_ECX)
(add R_EBX R_EDX)
(add R_EAX R_EBX))
((R_EAX 1 1 0 0 1 0 1 0 1 1 1) (R_EBX 1) (R_ECX) (R_EDX) (R_EDI _.2 . _.3) (R_ESI _.0 . _.1)))
(=/= ((_.0 0) (_.1 (1 0 0 1 0 1 0 1 1 1)))
((_.0 1) (_.1 (1 0 0 1 0 1 0 1 1 1)))
((_.0 1) (_.1 ()))
((_.0 _.2) (_.1 _.3))
((_.2 0) (_.3 (1 0 0 1 0 1 0 1 1 1)))
((_.2 1) (_.3 (1 0 0 1 0 1 0 1 1 1)))
((_.2 1) (_.3 ()))))))

3.2.2 Synthesis 2: Reverse ALU. Next we give an example of Angelic Execution—determining which sets of inputs successfully result in a particular output [Bodik et al. 2010; Chandra et al. 2011]. Here the output value 65,535 must be assigned to register EAX after computing the following assembly fragment:

mul EAX EBX
or ECX EDX
add ESI EDI
dec EAX
xor ESI ECX
inc ECX
xor EAX ECX
Each synthesized output contains a set of possible positive input values, one for each of the six general-purpose registers. Each input set successfully results in the desired output.

The full query and first 5 (of 200) results are shown below:

```plaintext
> (time
  (run 200 (v1 v2 v3 v4 v5 v6)
   (fresh (a b c x y z str)
     (pos^ a) (pos^ b) (pos^ c) (pos^ x) (pos^ y) (pos^ z)
     (x86^ ((mov R_EAX , x)
       (mov R_EBX , y)
       (mov R_ECX , z)
       (mov R_EDX , a)
       (mov R_ESI , b)
       (mov R_EDI , c)
       (mul R_EAX R_EBX)
       (or R_ECX R_EDX)
       (add R_ESI R_EDI)
       (dec R_EAX)
       (xor R_ESI R_ECX)
       (inc R_ECX)
       (xor R_EAX R_ECX))
     initial-store
     str)
   (lookup^ "R_EAX str "(1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1))
   (≡ (R_EAX , x) v1)
   (≡ (R_EBX , y) v2)
   (≡ (R_ECX , z) v3)
   (≡ (R_EDX , a) v4)
   (≡ (R_ESI , b) v5)
   (≡ (R_EDI , c) v6)
  )))
```

```
CPU time: 14062 real time: 14129 gc time: 4384
'(((R_EAX (0 1 1 1 1 1 1 1 1 1 1 1 1 1 1)) (R_EBX (1)) (R_ECX (1))
   (R_EDX (1)) (R_ESI (1)) (R_EDI (1)))
[((R_EAX (1)) (R_EBX (0 1 1 1 1 1 1 1 1 1 1 1 1 1 1)) (R_ECX (1))
   (R_EDX (1)) (R_ESI (1)) (R_EDI (1)))
[((R_EAX (0 1 1 1 1 1 1 1 1 1 1 1 1 1 1)) (R_EBX (1)) (R_ECX (1))
   (R_EDX (1)) (R_ESI (1)) (R_EDI (0 _.0 . _.1)))
[((R_EAX (1 1 1 1 1 1 1 1 1 1 1 1 1 1 1)) (R_EBX (0 1)) (R_ECX (1))
   (R_EDX (1)) (R_ESI (1)) (R_EDI (1)))
[((R_EAX (0 1)) (R_EBX (1)) (R_ECX (1))
   (R_EDX (0 0 1 1 1 1 1 1 1 1 1 1 1)) (R_ESI (1)) (R_EDI (1)))
```

...
3.2.3 Specifying Loop Constraints: Euclid’s GCD Algorithm. For our final example, we use a slightly higher-level code fragment to demonstrate how we handle looping or stitching together multiple code blocks. Our prototype currently only handles straight-line code with no jumps. To specify code with loops or more complicated control flow, one can specify how the blocks are to be sequenced using miniKanren.

As our example, we use Euclid’s algorithm for finding the Greatest Common Divisor (GCD) of two integers \(a\) and \(b\). Recall that the algorithm proceeds as follows:

1. Find the remainder \(r\) when dividing \(a\) and \(b\). (Equivalently, find integers \(q\) and \(r\) such that \(a = qb + r\)).
2. If \(r = 0\) then \(b\) already divides both \(a\) and \(b\), and is the largest integer that divides both \(a\) and \(b\); return \(b\).
3. Otherwise, any number that divides \(a\) and \(b\) must also divide \(b\) and \(r\). Repeat from step (1) to find the GCD of \(b\) and \(r\).

To implement this algorithm, first we adopt the convention that the values for \(a\) and \(b\) are held in registers EAX and EBX, respectively. We then write a miniKanren relation gcd_tester to run the loop. Given an assembly code body and input values, the code is run in a processor environment with values in their proper registers. It then checks the remainder result, and either associates the output with the final value (if the remainder is zero) or recursively calls itself on the resulting \(b\) and \(r\) values.

```lisp
(define (gcd_tester code a b gcd)
  (fresh (s s+ stmp)
    (update 'R_EAX a initial-store stmp)
    (update 'R_EBX b stmp s)
    (gcd_block_code code s s+)
    (conde
      [(lookup 'R_EBX s+ '()) ; final remainder is 0
       (lookup 'R_EAX s+ gcd)] ; result is in EAX
      [(fresh (r)
        (lookup 'R_EBX s+ r)
        (≡ r '()) ; r is non-0
        (gcd_tester code b r gcd)))])
)
```

Relation gcd_block_code specifies a loop invariant for our algorithm. It runs our x86 simulator on the given code, assuming the above register convention and specifying mathematical constraints for outputs that a correct run should produce. In this case, the constraints require that after the code runs, the new \(a\) value is the original \(b\) value and the new \(b\) value is the remainder of \(a\) and \(b\).

```lisp
(define (gcd_block_code code s s+)
  (fresh (old_eax old_ebx new_eax new_ebx q r)
    (lookup 'R_EAX s old_eax) ; a
    (lookup 'R_EBX s old_ebx) ; b
    (x86 code s s+)
    (lookup 'R_EAX s+ new_eax)
    (lookup 'R_EBX s+ new_ebx)
    (≡ new_eax old_ebx) ; a ← b
    (div 'old_eax old_ebx q r) ; a = qb + r, r < b
    (≡ new_ebx r)))
```

Using this infrastructure, we now run an algorithm on our processor and verify the answers computed are correct.

; euclid's gcd algorithm
(define euclid '(
  (xor R_EDX R_EDX)
  (div R_EBX)
  (mov R_EAX R_EBX)
  (mov R_EBX R_EDX)))

; test euclid's algorithm and get an answer
> (time (run* (d) (gcd_tester^ euclid (build-num 12) (build-num 9) d)))
  cpu time: 78 real time: 80 gc time: 0
  '((1 1))

> (time (run* (d) (gcd_tester^ euclid (build-num 42) (build-num 30) d)))
  cpu time: 203 real time: 213 gc time: 15
  '((0 1 1))

We see that we get the correct answers of 3 and 6, respectively (in Oleg numeral representation). As with any miniKanren relation, we can query for multiple answers at once; here we simultaneously obtain the gcd of 12 and all values of b up to 12, and verify that the answers are correct:

> (time (run* (b d) (<= o b (build-num 12)) (gcd_tester^ euclid (build-num 12) b d)))
  cpu time: 1329 real time: 1336 gc time: 937
  '((((0 0 1 1) (0 0 1 1))
    ((1) (1)))
  ((0 0 1) (0 1))
  ((1 1) (1 1))
  ((0 0 1) (0 0 1))
  ((0 1 1) (0 1 1))
  ((1 0 1) (1))
  ((1 0 1) (1))
  ((0 1 0 1) (0 1))
  ((0 0 0 1) (0 0 1))
  ((1 0 0 1) (1 1)))

Interestingly, we see three different classes of answers to this query, generated in order: the first six are those for which \( b = d \), namely the factors of 12 (12, 1, 2, 3, 4, and 6). The next group (11 and 5) are those that are relatively prime with 12, for which the only common divisor is 1. The remaining three answers (10, 8, and 9) are those that are not factors of 12, yet still have common factors (2, 4, and 3, respectively) with 12.

3.2.4 Synthesis 3: Synthesizing Euclid’s Algorithm with a Sketch. We now synthesize a similar algorithm by querying for the code. An initial, naïve approach attempts to synthesize the code directly using the relations we have so far:

> (time (run 1 (c) (gcd_tester^ c (build-num 12) (build-num 9) (build-num 3))))
...  
  ; [Fails with out-of-memory error]
The out of memory error occurs because this naïve query is too vague, resulting in a state space explosion. To narrow the search space, we employ a program sketch. Using the intuition that only a division operation and moving some data around should be sufficient, we create a program sketch that only allows \texttt{div} and \texttt{mov} instructions, with a specified program length. We take advantage of relational programming to auto-generate a suitable sketch:

\begin{verbatim}
(define (gen-mov-div-list o depth out)
 (conde
   [(≡ depth '()) (≡ out '())]
   [((≠ depth '()))
    (fresh (n 1)
     (minus o depth (1) n)
     (conde
      [(fresh (op1 op2) (≡ out ((mov ,op1 ,op2) . ,l) ))]
      [(fresh (op) (≡ out ((div ,op) . ,l) ))]))
     (gen-mov-div-list n l)]))

> (time (run 1 (c) (gen-mov-div-list (build-num 4) c)
  (gcd_tester o c (build-num 12) (build-num 9) (build-num 3))))
 cpu time: 361204 real time: 362308 gc time: 84412
 '(((mov R_ECX ()) (div R_EBX) (mov R_EAX R_EBX) (mov R_EBX R_EDX)))
\end{verbatim}

Here we have limited possible sketches to programs of length 4 with only \texttt{div} or \texttt{mov} instructions, and the system is now able to produce an answer. Upon running the test queries below, we see that this new code produces the same correct GCD answers as the original code above.

\begin{verbatim}
(define new-code '(((mov R_ECX ()) (div R_EBX) (mov R_EAX R_EBX) (mov R_EBX R_EDX)))

> (time (run* (d) (gcd_tester o new-code (build-num 12) (build-num 9) d)))
 cpu time: 110 real time: 107 gc time: 46
 '((1 1))

> (time (run* (d) (gcd_tester o new-code (build-num 42) (build-num 30) d)))
 cpu time: 265 real time: 262 gc time: 31
 '((0 1 1))

> (time (run* (b d) (<= o b (build-num 12)) (gcd_tester o new-code (build-num 12) b d)))
 cpu time: 422 real time: 437 gc time: 76
 '(((0 0 1 1) (0 0 1 1))
  ((1) (1))
  ((0 1) (0 1))
  ((1 1) (1 1))
  ((0 0 1) (0 0 1))
  ((0 1 1) (0 1 1))
  ((1 1 0 1) (1))
  ((1 0 1) (1))
  ((0 1 0 1) (0 1))
  ((0 0 0 1) (0 0 1))
  ((0 0 0 1) (1 1)))
\end{verbatim}
4 RELATED WORK

Rosette [Torlak and Bodik 2013] is another DSL embedded in Racket capable of program synthesis and angelic execution. Instead of backtracking search, Rosette uses a satisfiability modulo theories (SMT) solver to find solutions to synthesis and constraint problems. This gives it the potential to be more efficient than miniKanren in finding solutions that are arithmetic in nature. While any sufficiently general program synthesis system should be able to synthesize assembly code in a similar fashion to our approach, we found the Rosette system to be focused on returning a single optimal answer to queries. This makes it more cumbersome to achieve large-scale diversity with Rosette than with miniKanren in our experience.

Minimips\footnote{https://github.com/orchid-hybrid/minimips} is a miniKanren implementation of a MIPS architecture assembler/dissassembler. Its relation converts between MIPS assembly language programs and their binary encodings. Minimips contains a full syntactic description for MIPS instructions, but doesn’t appear to have an interpreter. It therefore has no semantic description of the instructions, or the ability to synthesize code that satisfies a formal specification. We chose to use the x86 architecture rather than a RISC architecture such as MIPS because a CISC architecture naturally allows for more diversity of implementations. As future work we plan to implement a similar relational assembler/disassembler for our x86 system.

Automated generation of assembly code is a common task of compilers and other similar tools. As noted in Section 1, these tools do not have diversity of implementation as a goal. Some systems (e.g., [Hong and Gerber 1993; Pu et al. 1988]) do synthesize assembly code for specific tasks using algorithms unique to the task. However, these systems do not synthesize general-purpose assembly code from arbitrary program constraints, or for the purpose of implementation diversity.

5 CONCLUSION

This paper proposed the implementation of an assembly-level interpreter in miniKanren for the purpose of synthesizing assembly code, motivated by the need for increased software diversity. This allows synthesis problems to be broken up into small, manageable pieces. Such problems can be subdivided using specific sketches, basic blocks, or effect traces; and can be driven by various inputs, including effect traces, processor states obtained from existing code, availability of particular gadgets, or compiler-driven information. Our working prototype implements an assembly interpreter for a small subset of x86, and experiments demonstrate its use for synthesizing code in both straight-line and looping programs. Synthesis examples show the potential to synthesize large numbers of diverse implementations.

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