# DECLARATIVE PROGRAMMING VIA TERM REWRITING 

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## Abstract

I present a new approach to implementing weighted logic programming languages. I first present a bag-relational algebra that is expressive enough to capture the desired denotational semantics, directly representing the recursive conjunctions, disjunctions, and aggregations that are specified by a source program. For the operational semantics, I develop a term-rewriting system that executes a program by simplifying its corresponding algebraic expression.

I have used this approach to create the first complete implementation of the Dyna programming language. A Dyna program consists of rules that define a potentially infinite and cyclic computation graph, which is queried to answer data-dependent questions. Dyna is a unified declarative framework for machine learning and artificial intelligence researchers that supports dynamic programming, constraint logic programming, reactive programming, and object-oriented programming. I have further modernized Dyna to support functional programming with lambda closures and embedded domain-specific languages.

The implementation includes a front-end that translates Dyna programs to bag-relational expressions, a Python API, hundreds of term rewriting rules, and a procedural engine for determining which rewrite rules to apply. The rewrite rules generalize techniques used in constraint logic programming. In practice, our system is usually able to provide simple answers to queries.

Mixing disparate programming paradigms is not without challenges. We had to rethink the classical techniques used to implement logic programming languages. This includes the development of a novel approach for memoization (dynamic programming) that supports partial memoization of fully or partially simplified algebraic expressions, which may contain delayed, unevaluated constraints. Furthermore, real-world Dyna programs require fast and efficient execution. For this
reason, I present a novel approach to just-in-time (JIT) compile sequences of term rewrites using a custom tracing JIT.

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I dedicate this to Angela and our posse of fluffy pets who fill every day with joy.

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## Chapter 1

## Introduction

The promise of usable declarative programming has attracted many to the declarative programming paradigm. Declarative programming is a paradigm where the program will specify what the result of the program should be without specifying how the program should perform the computation [100]. This allows the programmer to focus on the task while leaving maximum flexibility to the implementation of the declarative programming language. As such, the declarative paradigm is successful in a number of different domains. For example: databases with SQL [36, 44], optimization and search problems with SMT/SAT [15, 45], and logic programming languages such as Prolog [35, 38] and Datalog [30, 76].

Our group's research is in the space of Declarative Logic Programming designed for Machine Learning (ML) and Artificial Intelligence (AI) Algorithms. This research is called the Dyna programming language project, which investigates how Dataloginspired programming languages can be used to better encode AI algorithms [1, 59].

This dissertation represents the latest iteration of the Dyna programming language project. The work here is based on the culminated learning from three different implementations of Dyna I created during my Ph.D. This dissertation is the first to fully support the vision of Dyna proposed back in 2011 by Eisner and Filardo [59]. The approach within has the potential to enable a new class of powerful declarative logic programming languages, such as Dyna. Our approach is based on term rewriting on top of a relation algebra we call R-exprs (short for Relational expression). We have figured out how to mix a number of language features previously thought incompatible. This includes: Prolog-style backward
chaining, Datalog-style forward chaining, mixed-chaining [67], memoization, reactive programming, folding \& speculation, constraints satisfaction, functional, weighted terms and aggregation of multiple values for a single term (chapter §2).

Developing a new approach to implement declarative logic programming is not without its challenges. As such, we have investigated ways to make R-exprs practical. This includes novel (and more powerful) approaches for memoization, object-oriented programming, implementation of $\mathbf{R}$-exprs and their rewrite rules, and compilation of sequences of rewrites.

### 1.1 Dissertation Outline

This dissertation focuses on the applicability of term rewriting to the implementation of modern logic/weighted declarative programming languages with applications to ML/AI researchers. As such, I believe there are three different kinds of people who might choose to read this dissertation: (1) AI/ML researchers who might be interested in using the Dyna language implemented in this dissertation, (2) logic programming researchers who are interested in understanding how term rewriting can be used, and (3) term rewriting researchers who are looking for interesting applications of term rewriting.

For those interested in using Dyna, chapter §2 is written as a user manual for Dyna and should be the only chapter that one needs to read to use Dyna. Chapters 5 to 8 introduces our relational algebra, R-exprs, how Dyna programs are converted into a R-exprs and introduces our how our rewriting process is setup. Chapters 9 to 12 build on chapter $\S 8$ to discuss some "more advanced" techniques that I have implemented on top of term rewriting to implement features such as iteration of variable domains, memoization and compilation of multiple steps of rewriting. Chapters 13 to 15 include miscellaneous discussion around Dyna and R-expr rewriting. The dependency graph between chapters is approximately as shown in figure 1-1, and readers should feel free to jump around between chapters.


Figure 1-1. Dependency Graph Between Chapters

### 1.2 Brief History of the Dyna Project

The Dyna project was originally started in 2003 as an umbrella project to develop a programming language for ML researchers [60]. Most algorithms that ML researchers implement can be expressed in a few lines of math. ${ }^{1}$ In the process of researching new algorithms, researchers often have to iterate many times, refining their algorithms. This means that they first revise the mathematical concept of their algorithm and then recode their program to match. The project Dyna aims to reduce the distance between mathematical concepts and executable code.

[^0]This discrepancy between non-executable mathematical notation and executable programs was the central motivation for the Dyna project and led to the development of Dyna 1.0 [60, 61]. Dyna 1.0 extended Datalog [76] ${ }^{2}$ by replacing the boolean semiring used in logic programming to allow the use of any semirings. In other words, this meant that Dyna 1.0 was a notation for dynamic programs. As such, Dyna 1.0 was successfully used in several research papers: Dreyer and Eisner [54], Dreyer et al. [55], Eisner et al. [62], Eisner and Smith [63], Karakos et al. [95], Schafer [117], Smith and Eisner [121, 122, 123], Smith and Smith [124], Smith and Eisner [125, 126, 127, 128], Smith et al. [129, 130].

On the heels of Dyna 1.0 's success, Dyna 2.0 was proposed to rectify many of the limitations of Dyna 1.0 [59]. Dyna 1.0 requires all rules to use the same semiring ${ }^{3}$. Dyna 2.0 removes this restriction. Instead Dyna 2.0's rules define general functions. Dyna 1.0 is a dialect of Datalog, and as such, requires all terms derived using forward chaining to only contain ground terms ${ }^{4}$. This allowed the Dyna 1.0 compiler to generate programs that loop over the entire domain of an expressionmuch like a scan of a database table. Dyna 2.0 has no such restriction. Instead, Dyna 2.0 allows for variables in the program to remain free ${ }^{5}$ as in Prolog. Dyna 2.0 performs unification similar to Prolog where expressions like $a(X)=a(Y)$ unifies $X$ and $Y$ together without knowing the value of $X$ or $Y$. Dyna 2.0 also supports lazy expression allowing for expressions like $X+Y=Z$ to remain "unevaluated". Dyna 2.0 can also eagerly compute and memoize any expression to avoid recomputing the same expression many times. Dyna 2.0 also introduced a prototype-based inheritance mechanism (dynabases), which is useful for building larger programs.

It was shown in Eisner and Filardo [59] that the proposed features of Dyna 2.0

[^1]allowed for many ML algorithms to be concisely expressed. However, the unforeseen consequence of Dyna 2.0's design was that existing mainstream techniques for logic programming no longer worked. This spurred the current research phase to develop the necessary theory and techniques to implement Dyna 2.0: Filardo [66], Filardo and Eisner [67, 68], Francis-Landau et al. [70], Vieira [141], Vieira et al. [142]. Our group's research on Dyna also includes two Ph.D. dissertations prior to this one. The first, by Nathaniel W. Filardo, constructed a theoretical foundation for understanding the well-foundedness of Dyna 2.0 programs [66]. The second dissertation, by Tim Vieira, investigated automated program transformations, which can improve the asymptotic runtime of Dyna 1.0 dynamic programs [141].

This dissertation focuses on the implementation of the Dyna 2.0 language. I accomplish this by developing a novel representation of Dyna programs using R-exprs, a relational algebra, and develop an execution strategy using term rewriting. Much of the work presented in this dissertation is brand new work on Dyna, while other work in this dissertation will adapt the techniques that were previously developed by Eisner, Filardo, and Vieira to the R-expr presentation in this dissertation. Namely, chapter $\S 10$ will introduce how memoization works with R-exprs, which builds on Filardo's dissertation work, and chapter $\S 14$ will discuss how the kinds of program transformations, that are central to Vieira's dissertation work, can be implemented using $\mathbf{R}$-exprs.

From this point forward, the name "Dyna" will only refer to the Dyna 2.0 version of Dyna defined in chapter $\S 2$. The version of Dyna presented in chapter $\S 2$ contains some additional features and slight modifications since the original Dyna 2.0 paper of 2011 [59].

## Chapter 2

## The Dyna Programming Language

In this chapter, I will detail the surface-level syntactic features of Dyna as implemented in this dissertation. ${ }^{6}$ This chapter is intended as a user guide for a developer who is interested in writing Dyna programs. If you (the reader) are only interested in learning how to use Dyna, and do not care about how Dyna works internally, then this chapter should be the only chapter that you need to read.

The syntax for Dyna was originally proposed in Eisner and Filardo [59] and builds on the earlier Dyna papers (section §1.2). I continue to build on this design and extend it in this dissertation. Dyna started with the syntax of logic programming and deviates in a few ways in an attempt to modernize logic programming and make it more palatable for Machine Learning (ML) and Artificial Intelligence (AI) researchers. Primarily, this means having the ability to perform weighted reasoning. For example, a weight can be used to assign a probability to each expression in the program (section §2.2). Furthermore, most ML and AI researchers are primarily writing programs using Python. As such, my Dyna implementation provides an API to be able to easily interact with Python (section §2.3.1).

Our hope for the Dyna project is to enable new approaches to ML and AI to be easily explored, as well as existing approaches to be easily implemented. By this, I mean that we have attempted to incorporate support for commonly

[^2]used programming techniques as well as techniques that are rarely included in programming languages due to their inherent complexity. This includes: invariance to expression ordering-much like a constraint engine-(section §2.4), fixed-point computation and memoization (sections § 2.5 and 2.7)—which is good for handling cyclic reasoning in graphs-, non-ground reasoning (section §2.6), support for domain-specific languages (section §2.10).

As a new take on logic programming, Dyna also includes a number of niceties that have become commonplace among modern ${ }^{7}$ mainstream programming languages. This includes: higher-order functions (section §2.8.1), lambda functions/closures (section §2.8.2), prototype-based inheritance/objects (dynabases) (section §2.9), dependent types (section §2.8.3), and builtin commonly used data structures (such as a hash maps) (section §2.1.1.1).

Admittedly, all of these features have appeared before in other programming languages. What makes Dyna unique and interesting is that it is the first programming language (to our knowledge) that combines all of these features at the same time.

Furthermore, one of the core philosophies of the Dyna research project has been that any program which is "syntactically valid" and "logically consistent" should generally "just work". This means that Dyna tries extremely hard to make programs work even when other logic programming languages would result in an error or non-termination with the same program (chapter §4).

I believe that the combination of features included in Dyna is a sufficiently interesting challenge and demonstrates the main academic contributions of this dissertation (which starts in chapter §5). Additionally, I will discuss in chapter §4 some of the challenges and speculate on the reasons why I believe that Dyna is the only programming language to combine all of these features together at the same time.

[^3]
### 2.1 Dyna's Roots in Logic Programming

Dyna builds on the foundation of Prolog [35, 38] and Datalog [30, 76, 82]-with any Datalog or Prolog program also being a semantically valid ${ }^{8}$ Dyna program. Dyna has a number of syntactic changes compared to Prolog/Datalog in an attempt to modernize logic programming and make it more palatable to developers familiar with scripting languages (such as Python, commonly used in ML applications).

Dyna includes Prolog's and Datalog's notation, which makes it easy to express constraints, don't cares, unions, selections, and joins. For example, we can define a user-defined term ' $a$ ' as the join of two relations ' $b$ ' and ' $c$ '.

1 a(I,K) :- b(I,J), c(J,K).
A logic program is a programming representation of a set or $\mathrm{bag}^{9}$ that contains expressions that return true according to the program. For example, we might say that a program is a function $\mathcal{P}$ that maps expressions, such as $a(2,5)$, to the value true or false depending on whether $a(2,5)$ is contained in the program. For example, we can say that the program defines the term $a(2,5)$ if $\mathcal{P}(a(2,5))$ returns true. Similarly, it is sometimes useful to think of the program as a set/bag containing all terms the program defines as true. In this case, we can write $a(2,5) \in\{t: \mathcal{P}(t)\}$, where the set $\{t: \mathcal{P}(t)\}$ is defined using set-builder notation to contain all terms $t$ where the predicate $\mathcal{P}(t)$ returns true. For convenience, I will blur the distinction between function $\mathcal{P}(\cdot)$ and the set it defines and instead will write $a(2,5) \in \mathcal{P}$.

When working with a logic program, we are not only interested in checking whether a term such as $a(2,5)$ is defined according to the program. Instead, we are often interested in asking the program to fill in a templated expression that contains variables with all assignments to the variables such that the expression returns true under the function $\mathcal{P}$. For example, if we query the program with " $a(X, Y) \in \mathcal{P}$ ", the logic programming system will return a bag of all possible assignments to the variables $X$ and $Y$. E.g. $\{\langle X=2, Y=5\rangle,\langle X=7, Y=12\rangle, \ldots\}$

To see how this actually works, let us work through a small Dyna program that

[^4]has been adapted from commonly used logic programming examples:

```
% a number of facts about parent relationships
% this is parent(ParentName, ChildName).
parent("charles", "james").
parent("elizabeth", "james").
parent("james", "george").
parent("sophia", "george").
parent("sophia", "george").
% rules can combine facts as well as other rules
grandparent(X, Y) :- parent(X, Z), parent(Z, Y).
greatgrandparent(X, Y) :- grandparent(X, Z), parent(Z, Y).
married(X, Y) :- parent(X, Z), parent(Y, Z), X != Y.
related(X,X).
related(X,Y) :- related(X,Z), parent(Z,Y).
related(X,Y) :- parent (X,Z), related(Z,Y).
```

On lines 4 to 7 we define a number of parent-child facts. A fact is a rule without any side conditions (expressions on the right-hand side of :-). A fact represents a term that is true, and can be "looked up" in the program to check whether the expression is true. It is appropriate to think of a fact in a logic program as the same as inserting a tuple into a database table. ${ }^{10}$

Rules, such as those on lines 11 to 16 , define terms in accordance with expressions on the right-hand side of the aggregator :- which return true. For example, the grandparent rule on line 11 combines the parent facts to define terms such as grandparent("charles", "george"). A rule can be seen as similar to performing a join in a database between relations. ${ }^{11}$ The expressions which are used to define rules can reference the all user-defined terms, include all facts, other rule defined terms, and even its own term as shown with related $(\cdot, \cdot)$ on line 14 depending on itself.

To interact with this program, a user of Dyna will make queries against the program using the Dyna runtime, much like an SQL database. A query ends with a question mark '?' and will return the set of bindings to all variables in the expression.

[^5]For example, using the program on lines 4 to 16 , a user could make the following queries:

```
grandparent("charles", X) ? % make a query against the Dyna program
     \\langleX="george"\rangleS % the result is returned as assignments to variables
parent("sophia", X) ?
    {\X="george"\rangle\ % duplicates are combined
```

Note: On line 20 that the duplicate facts from line 7 and 8 is only returned to the user returned once, rather than twice. This differs from what Prolog would do, and the reason for this will be explained in section $\S 2.2$.

### 2.1.1 Structured Terms

While logic programming using only primitive types such as string and int in the previous section is sufficient to construct simple databases, where logic programming really shines is when constructing more complicated structured terms. In Dyna, we denote these structured terms using square brackets []$^{12}$ as: termName[Argument1, Argument2, $\ldots$, ArgumentN]. Note that this syntax for structured terms is different from what has been used for years in Prolog and Datalog. I will discuss why we have chosen to make these changes in section §2.2.1.

To see how structured terms can be used to revise our previous program, consider the following example:

```
% facts represented as structured terms
people(person["james", 50, "charles", "elizabeth"]).
people(person["george", 10, "james", "sophia"]).
% extract information from the structured term
name(person[Name, Age, Father, Mother], Name).
parent(person[Name, Age, Father, Mother], Father).
parent(person[Name, Age, Father, Mother], Mother).
% return the names of all known people
names(Name) :- people(Person), name(Person, Name).
```

[^6]On lines 22 and 23, we define two people using structured terms. We can get the name of a person using a rule like line 26 , which will select the first argument that contains the name of the person. We can further make queries against this program as follows:

```
people(Person), name(Person, "james"), parent(Person, Parent) ?
    Results returned:
        \\langlePerson=person["james",10,"charles","elizabeth"],Parent="charles"\rangle,
            <Person=person["james",10, "charles", "elizabeth"],Parent="elizabeth"\\
```

On line 32, we start with all of the people defined on lines 22 and 23, and then we then filter the people, selecting only those with the name "james", using name(Person, "james"). Finally, the parent rule is used to extract both the father and mother fields from the structured term.

### 2.1.1.1 Builtin Structured Terms

Dyna additionally includes built-in structured terms for convenience. Following Prolog's design, Dyna includes a linked list that can be constructed using square brackets ([]) when there is no term name (e.g. line 36). A vertical bar is used to create a new list where one or more elements have been added to the head of the linked list (e.g. lines 37 and 38). Recursive functions such as list_length (line 41) can be used to scan through a linked list by accessing the first elements of the linked list.

```
list = [1,2,3]. % Construct a list
list_prepend = [4 | list]. % The vertical bar is used to prepend or
list_remove = Rest for list = [Head|Rest]. % return the head and tail
% A recursive predicate defines 〈list, length〉 terms
list_length([], 0).
list_length([Head|Rest], N+1) :- list_length(Rest, N).
```

The equals sign on line 36 shows a major syntactic change from logic programming in that terms have values. Line 36 defines the term list as having the value [1, 2, 3]. The body of a rule is defined by an expression that returns a value. For example, on line 38, the expression 'Rest for list = [Head|Rest]' returns the value assigned
to the variable Rest when the expression 'list = [Head|Rest]' returns true. ${ }^{13}$
Furthermore, inspired by modern scripting languages that make extensive use of associative maps, I have extended Dyna to also include support and syntax for a built-in dictionary type ${ }^{14}$ (line 43). A dictionary can be constructed using curly braces (\{\}). Similarly to linked lists, a vertical bar (line 52) is used to both access a key in the dictionary (line 57) and add/remove keys from the dictionary (line 50).

```
person_dictionary_map = { % a dictionary
    "Name" -> "james",
    "Age" -> 50,
    "Father" -> "charles",
    "Mother" -> "elizabeth",
}.
person_with_address = { % new dictionary with keys added
    "Address" -> "123 North st.",
    | person_dictionary_map % the vertical bar is used to create a new
                % dictionary with new elements added or existing elements
                % removed, just like the linked list on lines 37 and 38
}.
name({Name | Rest}, Name). % The variable name can be used to...
name_alt({ "Name" -> Name | Rest}, Name). % match a key in the dictionary
get_item_by_key({Key -> Value | Rest}, Key, Value).
```

The 'Key -> Value' ${ }^{15}$ pair in the dictionary can use any values in the Herbrand universe. When adding a key to a dictionary, as on line 50 the key must not be contained in the dictionary, otherwise the result of this expression will be null. The reason is that when pulling a key out of the dictionary, as on line 57 , will result in that key not being contained in the dictionary represented by Rest, and we require that all uses of an operator, such as the vertical bar, behave identically regardless of how it is used (in this case to add or remove a key).

[^7]
### 2.2 Weighted Rules

As a Dyna extension to traditional logic programming, Dyna associates a value with each term in the program. As such, we call rules Horn equations rather than Horn clauses as in Prolog and Datalog. A Horn equation returns the value computed from the expression in the equation (which appears on the left-hand side of the aggregator). Returning a value from terms moves Dyna from a logic programming (which only returns true or null (undefined/false)) towards a functional programming language. A single term in Dyna can have values contributed either from a single rule or from a collection of different rules. To handle this, we combine the contributed values using an aggregator. An aggregator is an associative and commutative binary operator written between the head of the rule on the left-hand side and the body of the rule on the right-hand side. For example, the aggregator ' $+=$ ' sum of all contributions, 'max=' selects the maximum contribution, and the aggregator ': -' that we have already seen checks that there is some expression that returns true.

Using the ' $+=$ ' aggregator, we can write a matrix multiplication between two functors ' $b$ ' and ' $c$ ', which represent matrices as follows:
${ }_{60} \mid a(I, K)+=b(I, J)$ * $c(J, K)$.
Here, if we have the term $b(2,8)=3$. and $c(8,5)=7$., then this will define $a(2,5)=21$. One major difference between Dyna and other logic programming languages is that aggregators such as ' $+=$ ' combine all possible values-whereas boolean logical aggregators such as ':-' ${ }^{16}$ can stop once it has found a proof that a term is true. To illustrate this point, let us consider what happens if we add rules $b(2,9)=11$. and $c(9,5)=4$ to our program. In this case, we will have that $a(2,5)=65$. which is the sum of 21 and 44 . Conversely, if this was a logic program as in section $\S 2.1$, then the value of $a(2,5)$ already has been proven true by $b(2,8)$ and $c(8,5)$, and defining new values for the terms $b(2,9)$ and $c(9,5)$ does not change the truthfulness of $a(2,5)$.

In addition to the sum aggregator, Dyna also includes other aggregators, which are commonly found in Al and ML applications. This includes 'min=' and 'max=' which find the min or max value respectively. ' $x=$ ' computes the product of the

[^8]contributions much like ' $+=$ ' computes the sum. ' $I=$ ' and ' $\ell=$ ' compute the boolean logical OR and logical AND between contributions. Dyna also includes a special aggregator ' $:=$ ', which selects the value from the rule defined on the "last line" that returns a non-null value. This allows us to define rules that override the value returned by previously defined rules for a term. ${ }^{17}$ Finally, we have the aggregator ' $=$ ' which ensures there only one contributed value (otherwise it will error), and the aggregator '?=' which is allowed to arbitrarily pick any of the contributed values. ${ }^{18}$ 19

To see how aggregation in logic programming can make expressing AI programs easy, let us consider the task of computing the shortest path in a graph. On lines 62 and 63 , we define a recursive rule to compute the path in a graph. The rule will look for the minimum value associated with each node. We use the keyword arg ${ }^{20}$ to track the argument associated with the minimum value at each node. ${ }^{21}$ This, in essence, tracks backpointers at every node in the graph with \$arg(path(.)) being equivalent to a mathematical expression using argmin.

```
% Horn equation defines how to compute path and tracks back edges
path(start) min= 0 arg [start].
path(X) min= edge(X,Y) + path(Y) arg [X | $arg(path(Y))].
% weighted facts define the edges in a graph
edge("bal", "nyc") = 200.
edge("dc" , "bal") = 20.
```

[^9]```
edge("dc" , "nyc") = 300.
start = "nyc".
% assert }\mp@subsup{}{}{22}\mathrm{ checks the expression is true conditioned on the lines above
assert path("dc") = 220.
assert $arg(path("dc")) = ["dc", "bal", "nyc"].
```


### 2.2.1 Evaluation by Default

One of the major syntactic differences between Dyna and other logic programming languages is that Dyna evaluates an expression in place by default. The reason for this change is that most terms have ${ }^{23}$ a meaningful value, much like how a function returns a value in a functional programming language. Conversely, in logic programming languages such as Prolog or Datalog, terms only "return" the value of true.

To see how this manifests, let us consider the program that computes $\sin (x)^{2}+$ $\cos (x)^{2}$ written in Dyna vs Prolog:

```
% Dyna rule for computing the trig function
trig(X) = sin(X)**2 + cos(X)**2.
% Prolog rule for computing the trig function
trig(X,Result) :- sin(X,S), pow(S,2,SS), cos(X,C), pow(C,2,CC),
    Result $= SS + CC. }\mp@subsup{}{}{24
```

[^10]In Prolog and other logic programming languages, whenever we want to get the value from a term, we have to represent the returned value using an extra argument such as $S$ and $C$. The reason is that terms like $\sin (X, S)$ only "return" true. Hence, it would be meaningless to write an expression like pow $(\sin (X), 2)$ in Prolog with evaluation by default, as it would be equivalent to pow(true, 2). Instead, Prolog chooses to use this syntax to represent structured terms (section §2.1.1), with the expression pow $(\sin (X), 2)$ being interpreted as pow $(\sin [X], 2)$ which is at least potentially meaningful as long as 'pow' is able to do something with the structured term sin[X].

### 2.3 A User's Interaction With Dyna

Dyna's intended audience is ML and AI researchers. This means that Dyna is more focused on expressing mathematical models quickly and getting them to work. Dyna also encourages interactive experimentation, which is useful when debugging a ML model. Dyna also hopes to be a tool which complements existing tooling and infrastructure used by ML researchers-it is unreasonable to expect Dyna to have a thriving ecosystem on its own. As such, Dyna is designed to be used from a "driver" program written in another language. ${ }^{25}$ In this way, Dyna is similar to a database where a program submits queries and updates to run against the Dyna program. As such, Dyna provides a Read-Eval-Print-Loop (REPL) for interacting with the Dyna system, as well as a Python and Java API.

[^11]
### 2.3.1 Python API

Most ML and AI research is currently being done using Python. As such, we expect that the Python API will be the way most users will interact with Dyna. The Java API also provides the same interface as the Python API, so we will omit detailing the Java API here.

Interacting with the Dyna system is similar to interacting with an in-memory database (such as a SQL database). A new Dyna runtime instance is created using the dyna. Dyna() method (line 82). Programs can be loaded into the Dyna runtime using the run method (lines 84 and 92), passing either a string of code or a file containing Dyna code. The defined rules and facts will persist between calls to the run function, just like a database.

```
from dyna import Dyna
dyna_runtime = Dyna() # create a new instance of the Dyna runtime
dyna_runtime.run("""
factorial(N) := factorial(N-1) * N.
factorial(0) := 1.
print factorial(5). % prints 120
""")
# load rules from a file
dyna_runtime.run(open('dyna_program.dyna'))
```

The query method (line 93) is used to return the values that are calculated by the Dyna program. We can return primitive types, such as numbers and strings, as well as more complicated types, such as lists, dictionaries, and dynabases (objects inside of Dyna, which will discussed further in chapter §13):

```
result = dyna_runtime.query("""
factorial(10)? % The first query
factorial(11)? % The second query
factorial_up_to(N) := [factorial(N) | factorial_up_to(N-1)].
factorial_up_to(0) := [].
factorial_up_to(5)? % Third query returns a list
```

assert result [0] == 3628800 the results of queries are returned
assert result[1] == 39916800
assert result[2] $==[120,24,6,2,1]$

Like SQL database APIs, Dyna supports query parameters for passing values into Dyna without having to encode the value as a string. The query parameters are denoted using a dollar sign followed by a number: $\$ 0, \$ 1, \$ 2 \ldots \$ n$. We can pass any value into the Dyna runtime. When a value can be cast into Dyna's Herbrand universe then it is automatically cast and usable inside of Dyna. If a value cannot be cast to Dyna, then it is passed around as an opaque pointer. This allows Dyna to integrate with other Python libraries without having to support everything itself.

```
result = dyna_runtime.query("""
factorial($0)?
""", 5)
assert result[0] == 120
class MyClass: pass
dyna_runtime.run("""
class_reference = $0. % save a reference to the class
""", MyClass()).
result = dyna_runtime.query("""
class_reference?
""")
assert isinstance(result[0], MyClass)
```

Dyna's Python API also supports defining external functions that can be called from the Dyna program. This allows Dyna to leverage existing functions and libraries without requiring all features to be fully reimplemented in Dyna. Unlike with Dyna terms, invoking an externally defined function requires that all arguments have known values. ${ }^{26}$

[^12]```
@dyna_runtime.define_function('my_function')
def my_function(a,b,c):
    return a*3 + b*5 + c*11
dyna_runtime.run("""
assert my_function(1,2,3) == 1*3 + 2*5 + 3*11.
""")
```

As a more useful example of Dyna interacting with external datatypes, we can leverage PyTorch's [111] GPU tensors. Operations such as tensor multiplication can be performed by calling back to external functions:

```
import torch
gpu_tensor = torch.tensor(...).cuda()
@dyna_runtime.define_function('tensor_multiply')
def tensor_multiply(A, B):
    return A @ B
dyna_runtime.run("""
pytorch_tensor = $0.
tensor_squared = tensor_multiply(pytorch_tensor, pytorch_tensor).
""", gpu_tensor)
```

Currently, Dyna does not have built-in support for GPUs, so this approach can at least serve as a "stop gap" for working with GPUs, which are critical for making modern (neural) ML models efficient. In section $\S 16.7$ I will discuss potential future work to add native GPU support to Dyna.

### 2.3.2 Multi-file Programs

Just like other programming languages, Dyna has the ability to write a program using multiple files and import the definitions across files. ${ }^{27}$

Terms in Dyna are scoped to the file in which they were defined. This means that a term $f \circ o(X)$ defined in file a.dyna, will be different from $f \circ o(X)$ in file b. dyna. Terms can be imported from other files using import at the top of the file as follows:

[^13]${ }^{143}$ |:- from "a.dyna" import foo/1.
Dyna also supports importing default terms that were declared as exported. This way, they do not have to be explicitly listed when importing terms into another file.

```
% in file a.dyna
foo(X) = 123.
:- export foo/1.
% in file do_import.dyna
:- import "a.dyna".
assert foo(1) = 123. % use foo defined in file a.dyna
```


### 2.4 Invariance to Expression Order

One of the core principles of Dyna is that the way that a program is written should have as little impact as possible on how the program executes. The Dyna runtime is required to find some solution to the equations/constraints written, but otherwise, it is flexible in how it executes a program. We hope that this approach will make it easier for developers to write Dyna programs as they can focus on the "business logic" ${ }^{28}$ of their ML/AI model while not having to worry about the internal state of the Dyna program.

One way in which Dyna manifests this property is that Dyna is invariant to the order in which expressions and most rules are defined in the program. In this way, Dyna is closer to a constraint satisfaction engine than other logic programming languages, which have a fixed execution order (section §3.1.1). For example, a wellstudied example Prolog program is computing the permutations of a list (lines 151 to 154). Following the mathematical definition of a permutation, given a list $l$ and the permutation $p$ of list $l$, will have that $l$ is also a permutation of the list $p$. E.g. the list $[1,3,2]$ is a permutation of $[1,2,3]$ and vice-versa.

```
deleteone([X|Xs], Xs, X).
deleteone([X|Xs], [X|Ys], Z) :- deleteone(Xs, Ys, Z).
permute([], []).
permute(As, [Z|Bs]) :- deleteone(As, Rs, Z), permute(Rs, Bs).
```

[^14]```
permute([1,2,3], Out)? % works under backchaining (Prolog)
permute(Out, [1,2,3])? % does not work under backchaining
```

This program, when executed under Prolog's fixed execution order, only line 156 will work, while line 157 will cause Prolog to get stuck in an infinite recursion. The reason is that Prolog executes in a fixed left-to-right, top-to-bottom rule ordering. As a result of this fixed order, when Prolog executes line 154, the call to deleteone will be evaluated before the recursive call to permute is evaluated. This means that the variables As and Rs will both be un-ground. This causes deleteone to enumerate the infinite pairs of all lists where the list Rs will contain one less element that is equal to $Z$.

Conversely, when this program is executed under Dyna, we do not commit to execute the rules in any particular order. So, while Dyna runtime is allowed to explore the deleteone function before As and Rs are grounded to specific values, it does not commit to running the deleteone function before the permute function.

### 2.5 Fixed-Point Computations

Dyna is designed as a superset of Prolog and Datalog styles of logic programming. Datalog's execution strategy is notably different from Prolog's. A Prolog program is executed using backward chaining. This means that Prolog starts with the expression that it is trying to prove to be true and performs a depth-first search "backward" looking for other rules and facts that support the expression it is trying to prove. In contrast, Datalog operates via forward chaining. When Datalog starts, it looks at all of the ground facts that are currently asserted by the program. Datalog will then deduce new facts using the rules in the program.

```
158 a :- b. % a is true if b is true
159 b :- a. % b is true if a is true
160 c :- d. % c is true if d is true
1 6 1 ~ a . ~ \% ~ a s s e r t ~ t h a t ~ a ~ i s ~ t r u e ~ r e g a r d l e s s
```

On line 161, we define the fact that the term a is true. Datalog using forwardchaining will use line 159 to deduce that $b$ is also true. Datalog will then stop processing as there are no new rules which can be deduced as true. ${ }^{29}$ In this way, Datalog

[^15]has reached a fixed-point of the program where: $\mathcal{P}_{T}=\operatorname{DeduceAllTrue}\left(\mathcal{P}_{T}\right)$. Here, we can think of the function DeduceAllTrue as taking the set of all terms in the program $\mathcal{P}_{t}$ at time $t$, and computing a new set of terms $\mathcal{P}_{t+1}$ at time $t+1$ which includes the new terms which have also been deduced. Once no additional terms are proven as true, the fixed-point condition is satisfied $\mathcal{P}_{T}=\operatorname{DeduceAllTrue}\left(\mathcal{P}_{T}\right)$.

As mentioned above, Dyna is a superset of Datalog's fixed-point computation. As such, Dyna also includes support for fixed-pointing weighted expressions (whereas Datalog can only fix-point boolean values that are proven true). For example, we can use a fixed point to solve the expression $e=1+e / 2$ by writing lines 162 to 163.

162 e += 1.
163 e += e/2.
Dyna will solve this program by first assigning the value of 1 to the term 1, and then it will iterate line line 163 until it reaches numerical convergence.

The place where having fixed-point semantics is important is in the case of cyclic programs, such as when representing a cycle on a graph. For example, when computing the shortest path in a graph in section §2.2, we were only interested in finding the minimum length to a given node in the graph. If we had a large set of edges, it is likely it would have contained cycles and many alternate paths. Furthermore, the fewest number of steps might not have been the cheapest path in the graph. Hence, the value assigned to a node in the graph might change multiple times during the program's run.

Note: A fixed-point might not exist for a program, or there might exist multiple different fixed-points for a single program. For example, if we define that a term is true if it is false, then the Dyna runtime can cycle forever, proving that ne is true, and then switching back to ne is false:

```
164 ne :- not ne. % ne is true if ne is false
```

Datalog is able to reject programs like this due to stratification [30, 76], which limits the kinds of programs supported by Datalog. Dyna does not limit programs with stratification. Furthermore, Dyna programs are Turning complete (section §15.2) so proving that a general program terminates or cycles is impossible. ${ }^{30}$ Furthermore, we can write a Dyna program which never repeats a state by writing

[^16]a non-terminating program that counts to infinity (lines 165 and 166), or has a negative-weighted-cycle in the shortest path graph (e.g example line 62 ).
165 count_to_infinity += 1.
166 count_to_infinity += count_to_infinity.
As for programs that have multiple fixed-points, Dyna is only looking for an assignment to all terms in the program that are consistent with all of the rules that are defined. ${ }^{31}$ This means that it is possible to define a term that will take on the value true or false, or even a term that can take on any value in the Herbrand universe.
167 true_or_false :- true_or_false.
168 any_value = any_value.
In practice, if either of these terms are queried, they will return the result null or undefined as there is no base case to their recursion. However, this behavior should not be depended on as what value is returned by these kinds of expressions should be treated as purposely underspecified behavior.

### 2.6 Non-ground Reasoning

As mentioned a few times already, Dyna is a superset of both Prolog and Datalog; this means that it supports features of both languages at the same time. One such feature is non-ground reasoning. Non-ground reasoning means that the value assigned to a variable is not known at the time of execution. This technique is most commonly associated with Prolog-style systems and is not supported at all by Datalog. An example of non-ground reasoning would be an expression like:
${ }_{169} \mid \mathrm{a}:-\mathrm{X}>5, \mathrm{X}<3$.
Here, the value of a will be null, as there is no assignment to the variable x , which is both greater than 5 and less than 3 at the same time. This reasoning can also be applied across rules boundaries. For example, the following will still deduce that the term a is false.

[^17]```
170 b(foo[Z]) :- Z > 5.
171 a :- b(foo[Y]), Y < 3.
```

What makes Dyna's approach to non-ground reasoning more powerful than Prolog is that it is able to run non-ground reasoning while simultaneously performing Datalog-style fixed-pointing. To see an example of what this means, let us modify our shortest path program from line 62 to remove the hard-coded starting node and turn it into the all-pairs shortest path program:

```
path(Start,X) min= edge(X,Y) + path(Start,Y).
path(Start,Start) min=0. 32
edge("bal", "nyc") = 200.
edge("nyc", "bal") = 180.
edge("dc" , "bal") = 20.
edge("dc" , "nyc") = 300.
```

Observe this program cannot be executed under Prolog-style backward-chaining or Datalog-style forward-chaining. The reason that Prolog fails to execute this program is that there exists a cycle in the edge graph ("bal" $\rightarrow$ "nyc" $\rightarrow$ "bal"); hence depth-first backwards-chaining would get stuck in a cycle. While Datalog can handle the cycles in the program, it cannot support line 173, as Datalog does not handle non-ground expressions like this.

### 2.7 Memoization, Dynamic Programming, and Reactive Programming

One central feature in Dyna is support for memoization. Memoization essentially is the act of replacing a compute operation in a program with a "recall from memory" operation. For certain kinds of programs, this can have a significant impact on the runtime of a program. For example, the Fibonacci sequence (line 179), will take exponential time without memoization but can run in linear time when memoization is enabled:

```
fib(0) \(+=0\).
fib(1) += 1.
fib(N) += fib(N-1) for \(N>1\); \% semicolon defines two rules at once
```

[^18]```
182
print fib(10). % run in 177 operations without memoization
$memo(fib[N:$ground]) = "unk". % enable memoization for fib where N}\mp@subsup{N}{}{33
    % must be a ground value (such as 0,1,2 etc)
print fib(10). % run in 10 operations with memoization
```

Memoization traditionally requires that the memoized result of a function does not change. However, Dyna's memoization implementation is reactive. Reactive means that Dyna automatically tracks dependencies between anything that a memoized value depends on, and that the memoized value will be recomputed or updated if anything upstream is changed. The value assigned to a term by the Dyna program can change as a result of the system converging towards final values or as the result of the program being modified through the addition of new rules or facts (e.g. line 97 in the Python API example, section §2.3.1).

This means that if some input to a currently memoized value changes, then any dependent memoized values will also be updated. For example, if we modify the definition of fib(3) by defining a new rule, this will cause all of the values of the Fibonacci terms to change.

```
189 print fib(10). % print 55 using the existing memoized value
190 fib(3) += 1. % modify the definition of fib
191 print fib(10). % print }76\mathrm{ using the new modified version of fib
```


### 2.7.1 Prioritization of Updates

In a Dyna program, there are often many terms in the program that are memoized. As a result, there can be many different terms that have pending updates waiting to be processed at any given point in time. However, the Dyna runtime cannot process all updates simultaneously. Additionally, the order in which these updates are processed can significantly impact the program's overall runtime. To handle these cases, we allow the user to specify a prioritization function that controls the order in which updates can be processed. The definition of the prioritization function does not impact the correctness of the program; however, a badly defined prioritization function can cause the program to exhibit much worse runtime-for

[^19]example, a badly defined prioritization function will cause the Fibonacci example (line 179) to run in exponential time instead of linear.

A priority function is defined by defining a \$priority term just like the \$memo show above.

```
$priority(fib[N]) = -N. % prioritize small values first (0,1,2 etc)
% this is the optimal prioritization function
    % for fibonacci and runs in O(n)
$priority(fib[N]) = N. % prioritize large values first (10,9,8 etc)
```

The priority function will map a term to a number that is used as the priority. Higher priorities will be run first. ${ }^{34}$ The priority function is computed only when the pending update is created. ${ }^{35}$ This means that \$priority should be considered a meta-user-definable term used only to control the inner operations of the system and is not like the other Dyna terms in the language, which can be updated reactively.

### 2.7.2 Memoization with non-ground variables

Memoization in a logic programming language, such as Dyna, is not as simple as memoizing the returned value. Instead, terms in logic programming languages support being called in different modes. A mode is the signature which a term supports being "called" with in terms of which variables must be known ground values in the Herbrand universe, and which variables are allowed to be free and have an unknown, yet to be determined, value. We have already seen an example of ground and free variables in the matrix multiplication example repeated here.
${ }_{196} \mid a(I, K)+=b(I, J) * c(J, K)$.
Observe that the only constraints on the variable $J$ are the terms $b(I, J)$ and $c(J, K)$. This means that either ' $b$ ' or ' $c$ ' must allow for the variable $J$ to be free. In which case, that functor will be able to enumerate an upper bound on the domain of the possible values for the variable J. And by extension, any memoized version of that functor will also have to allow for the variable $J$ to be free.

[^20]To declare a policy for memoization, we use the declarations \$free, \$ground and the meta-term \$memo. When a variable is marked as \$ground in a \$memo policy, this means that the value assigned to the variable must be known before we can even check if there are any relevant memoized terms in the memo table. Similarly, \$free specifies that the variable's value is allowed to be unknown and allows us to check for matching terms in the memo table. Unlike with \$ground, when a variable is marked with $\$$ free the memo table will guarantee that all relevant terms will be memoized.

For example, we can write a memo policy for the matrix example on line 196 as follows:
${ }^{197} \mid$ \$memo(b[I:\$ground, J:\$free]) = "null".
Here on line 197, we are specifying that the first variable I must be known before we can check the memo table where we will find all of the relevant terms for different values of J . The $\$$ memo $(\cdot)$ returns the string "null" to indicate what is the default value for the memoized $b(\cdot, \cdot)$ terms. By stating "null", we declare that any term not found in the memo table has the associated value null, and is therefore undefined. Alternately, the \$memo function can return "unk"36 to indicate that the value for a term unknown and must be computed and should be saved into the memo table, or it can return "none" to indicate that there is no value in the memo table for a term and the value should not be saved into the memo table. ${ }^{37}$

### 2.8 Modern PL Constructions

Dyna's syntax was primarily inspired by logic programming languages which date back to the 1970s. Since then, more modern programming languages have been adding syntactic sugar to make common operations easier to write. Dyna has

[^21]followed suit and includes higher-order functions, ${ }^{38}$ lambda functions, and type annotation. ${ }^{39}$

### 2.8.1 Higher-Order Functions

Dyna allows for "function pointers" to be passed around the runtime and for those pointers to be used via indirect calls. This enables Dyna to support higher-order functions. A function pointer in Dyna is represented using the same structured term (section §2.1.1) that we have seen already. When an indirect call is performed, Dyna looks up the user-defined term by name and passes any additional parameters. In Dyna, user-defined terms are scoped to the file in which they were defined. This means that there can potentially be many user-defined terms that share the same name but are otherwise unrelated. To handle this, structured terms track which file they were constructed in via additional hidden metadata. This metadata is only used when an indirect call is performed. A structured term can also be scoped to a dynabase (an object), allowing indirect calls to be performed against terms defined on a dynabase object.

To make an indirect call in a higher-order function, we can simply use a variable in place of the functor's name as on line $198 .{ }^{40}$

```
perform_call(Calling) = Calling(1,2,3).
called(X,Y,Z) = X*Y + Z.
called(X,Y,Z,W) = X*Z + Y+W.
assert perform_call(called[]) = 1*2 + 3.
assert perform_call(called[4]) = 4*2 + 1*3. % curried functions can
    % include extra parameters
```

More usefully, this can allow us to define higher-order functions such as map, which applies a function to all elements in a list line 205.

[^22]```
map(Func, []) = [].
map(Func, [X|Rest]) = [Func(X) | map(Func, Rest)].
assert map('*'[2], [1,2,3]) = [2,4,6].
```


### 2.8.2 Lambda functions

Dyna includes syntax for creating lambda functions. Lambda functions are given a generated name and converted into a standard function like we have been writing already. Given that lambda functions are the same as regular Dyna rules, this means that we also define an aggregator and perform aggregation inside of the lambda function. For example, if we want to create a function that takes the max of two arguments, we can write: ( $(X, Y) \max =X ; Y)$. The semicolon splits between different body expressions, as was done on line 181. The outer parentheses encapsulate the entire expression, with the first inner parentheses denoting additional arguments that can be passed to the lambda expression. If no arguments are included, then the lambda is immediately evaluated in place. For example, we can write $X=5, Y=7,7=$ (max $=X ; Y$ ) where the variables $X$ and $Y$ are captured in a closure, and then the expression is immediately evaluated using the max= aggregator.

Together this can be combined with the higher-order-functions to write the following expressions, which shows how these concepts can be syntactically combined:

209 assert Seven=7, map(((Arg1) max= Seven; Arg1), $[1,5,10])=[7,7,10]$.

### 2.8.3 Type Declarations

Dyna includes support for adding type restrictions onto any expression in the language. Type annotation is denoted using a colon followed by a term. For example, the expression $X$ : int, denotes that the variable $X$ is of type int. The syntax $X$ : int adds the constraint int $(X)$ into the body, where int $(X)$ the same kind of call to the int term that we have seen already. The difference with the colon expression is that instead of returning the value of the $\operatorname{int}(X)$ call, this expression returns the value of $x$. This means that we can easily annotate variables with types throughout the program (e.g. line 213) and can define our own types (e.g. line 210).

```
my_custom_type(foo[X,Y:float,Z:int]) :- Y > Z % define a custom type
my_custom_type(bar[X:string]). % with side conditions
my_function(X: my_custom_type) = .... % check the type matches
```

Furthermore, we can use this approach to construct "templated" types. This is accomplished by passing arguments to the term that appears after the colon. The type-constrained variable is appended as the last argument. For example, a templated list type can be defined as on line 214:

```
list(Type, []).
list(Type, [Head | Rest]) :- Type(Head), list(Type, Rest).
% Check the argument L matches a list of a particular type
list_of_my_type(L : list(my_custom_type[])) = ....
list_of_greater_than_5(L : list(((X) :- X > 5))) = ....
```

The 'Type(Head)' expression turns the list( $\cdot, \cdot$ ) term into a higher order function which indirectly calls the term referenced by the variable Type on line 215

Union types can also be written inline through the use of a vertical bar like $x$ :int|string. Conceptually, this is the same as defining a new term that attempts to match against $\operatorname{int}(X)$ or string( $X$ ) using the :- aggregator. ${ }^{41}$

### 2.8.3.1 Type Checking and Type Errors

In Dyna, "types" are not distinct from "normal" function calls. This is not a significant issue, though, as in Dyna, the Dyna system is allowed to evaluate any expression at any point. This means that the Dyna runtime is allowed to evaluate expressions and function calls at "compile-time" if there is sufficient information to deduce the resulting value of the expression. In other words, if we can statically deduce the types, then it is possible to avoid the runtime overhead of performing type checks. However, in the cases where there is a complicated type, the Dyna runtime is allowed to perform runtime type checks as needed.

A consequence of this design is that Dyna does not have "type errors" that can be

[^23]concisely reported to the user. Instead, Dyna can check and warn about dead code which will never be evaluated under any circumstances. For example, the expression $5<3$ is always returns false. Similarly, the expression int(X), string(X), which places two incompatible type constraints on the variable $x$, hence this expression will always return false and can be reported as a warning. ${ }^{42}$

### 2.9 Object-Oriented Programming (OOP) via Dynabases

When building a larger, more complicated AI and ML system, it is necessary to have some way to make smaller units that can be composed together. In most programming languages, this is done via object-oriented programming (OOP) or via modules (section §2.3.2). Dyna includes a prototype-based approach to objectoriented programming we call dynabases. Dynabases are designed to resemble more typical approaches to OOP, as in a procedural programming language. Dynabases are denoted using curly braces \{\} with Dyna rules inside. The optional keyword new can be used to syntactically distinguish between dictionaries (section §2.1.1.1) and dynabases when needed. For example, we can define a simple Dynabase on line $220 .{ }^{43}$

```
220 e = { count += 1.
    count += 2. }.
assert e.count = 3.
```

We can also modify a dynabase after it has been created:

```
2 2 3 ~ e . c o u n t ~ + = ~ 3 . ~
2 2 4 ~ a s s e r t ~ e . c o u n t ~ = ~ 6 . ~
```

Dynabases can be duplicated as well as extended with additional rules by using the keyword new. Rules defined in a dynabase can reference the dynabase itself using the keyword \$self.

[^24]```
copy_e = new e. % create a copy of a dynabase
extend_e = new e { % extend with additional rules
    % $self}\mp@subsup{}{}{44}\mathrm{ is used to reference fields and functions on the dynabase
    bar(X) = $self.foo * X.
}.
```

The extended and copied dynabases are distinct from their parent and can be modified individually. However, changes to the parent dynabase are visible in the children even after the children have been "created." The reason for this is that Dyna is declarative. Hence the order in which rules are defined does not impact the expressions in the program. For example, we can define a new rule foo on the e dynabase on line 230 . This rule will be visible on all descendants dynabases, and it is even possible that rules defined before foo was defined (line 228), can still reference the definition of foo.

```
e.foo = 123. % define a new field foo on the parent
copy_e.count += 4.
assert copy_e.foo = 123. % the parent changes are visible in children
assert copy_e.count = 10. % existing rules are extended
assert extend_e.bar(2) = 123*2.
assert e.count = 6. % the parent is not impacted by children
```


### 2.9.1 Dynabases vs Procedural Programming OOP

As mentioned above, dynabases are a bit different from classes in a procedural programming language since Dyna is a declarative programming language that can be evaluated out-of-order. Another difference between Dyna's dynabases and procedural programming's OOP is that a dynabase might not actually be "created." To see what I mean by this, consider for a moment that due to aggregation, we can compute the aggregated result of multiple contributions at the same time. In fact, we can even combine an infinite number of expressions. As such, it is possible to write a program where we create an infinite number of dynabases (e.g. lines 236 to 237).

236

```
return_dynabase(X,Y,Z) = { value = ... . }.
result min= return_dynabase(X,Y,Z).value.
```

[^25]We are able to handle programs with an "infinite" number of dynabases as long as we are able to deduce some representation for the value term, which can be used for solving the min= aggregator on line 237.

### 2.10 Embedded Domain Specific Languages

Support for Domain-Specific Languages (DSLs) is a useful feature for programming languages to have as it enables library developers to support a wide variety of different problem domains. Early languages such as LISP were popular among AI researchers in part for this reason. These days, ML researchers are clearly demonstrating that they want access to DSLs despite the fact that Python (the language that is often used for ML) does not easily support macros. This has led libraries such as JAX [23] and PyTorch Script [2], which use Python's reflection to access and manipulate Python ASTs for a function. Given that Dyna is targeting the same kinds of researchers, we believe having the ability to write macros to transform Dyna's AST before it is loaded and also use multiline strings as embedded languages will allow for other languages to be embedded in a Dyna program. ${ }^{45}$

### 2.10.1 String DSLs

A multiline string in Dyna is escaped between the symbol ' $\{$ and \}. The string can contain \{ and \} as long as they are balanced. Furthermore, Dyna-style comments, a percent sign \% followed by a new line, are also stripped from the string. This is done to encourage ' $\}$ to be used to embed DSLs that have similar behavior to the Dyna language. Furthermore, Dyna includes special syntactic handling inspired by languages such as Ruby and Lua for passing a block of \{\} to a function. When \{\} appears after a function call in any form (string, dictionary, or dynabase), it will be passed as the last argument to the function. Together, this allows us to write a program as follows for embedding a context-free grammar (line 238) and a linear program (line 257):

[^26]```
my_grammar = grammar'{
    S -> NP VP
    NP -> Det N
    NP -> NP PP
    VP -> V NP
    VP -> VP PP % inline dyna comments are removed
    PP -> P NP % when converted to a string
    NP -> Papa {} % balanced {} can be included in the string
    N -> caviar
    N -> spoon
    V -> spoon
    V -> ate
    P -> with
    Det -> the
    Det -> a
}.
my_lp = solution[X,Y,Z,Objective]
    for Objective = linear_program({X,Y,Z}) '{
Maximize % embed a linear program using the
        obj: X - 2.3Y + 0.5Z % LP file format
Subject
        c1: X - Y + S <= 10.75
            -Z + 2X - S >= -100
}.
```

For grammar defined on line 238, the entire grammar from lines 238 to 253 is passed to the grammar(•) term as its only argument. The grammar term can create dynabase to represent the grammar. For the linear programming example, the first argument to linear_program( $\cdot, \cdot$ ) is the dictionary $\{" X "->X, " Y "->Y$, "Z"->Z\} and the second argument is the linear program represented as a string. The implementation of linear_program $(\cdot, \cdot)$ is responsible for matching the linear program's variable names with the dictionary's variable names.

### 2.10.2 Macros

Macros have access to the AST of the Dyna program. The Dyna AST is represented using the same structured-term object that is passed around in Dyna programs.

This means that we can use the same pattern-matching tools that we presented in section §2.1.1. Additionally, we have backtick '(), which is used to escape the AST and embed variables used to match against part of the AST. In the following example, on line 266 , we match ${ }_{-}+2$ and replace it with $+3:^{46}$

```
:- macro my_macro/1.
my_macro(X) := X. % no change to AST if it does not match
my_macro(`(`X + 2)) :=`(`X + 3). % change AST
f1 (W) = my_macro(W + 1).
f2(W) = my_macro(W + 2).
assert f1(3) = 4.
assert f2(3) = 6.
```

More practically, macros could be used to implement symbolic auto differentiation (just like JAX [23]), against the AST of a Dyna program. Line 273 shows a conceptual example of how an automatic differentiated neural network could be used in Dyna. ${ }^{47}$

```
network(InputMatrix) = auto_differentiate {
    relu(X) max= X. % define a neural net unit using a dynabase
    relu(X) max= 0.
    layer1(I, J) = InputMatrix.elem(I, J).
    layer2_input(I, J) += $self.layer1(I,K) * $self.weight(K, J).
    layer2(I, J) = $self.relu($self.layer2_input(I,J)).
}.
gradient = network(...).gradient(...). % gradient term added via macro
```

The '.gradient' term is added to the 'network' dynabase by the auto_differentiate macro called on line 273.

[^27]
## Chapter 3

## Related Work

In this chapter, I will cover some related projects and programming languages and how their implementations work. With the Dyna project, it is a bit difficult to make direct comparisons. The reason is that the Dyna language contains combinations of features that are not entirely supported by other languages. A lot of the challenges with Dyna have been a result of how features interact with each other, rather than the addition of any individual features-this will be discussed further in chapter $\S 4$.

Further complicating this chapter, some of the systems/languages that I compare against are not specific implementations but rather a large class of related systemsfor example, I will discuss SQL databases and how they relate to Dyna.

I note this chapter is focused on the implementation of similar declarative and logic programming systems. This chapter is not intended as a tutorial on the features or syntax of other systems. The reason for this is that, except for chapter $\S 2$, this dissertation is focused on implementation and not the Dyna language itself.

I have attempted to write the rest of this dissertation starting at chapter $\S 5$ with few dependencies on this chapter, so this chapter can be skipped for those who are only interested in the academic contributions of this dissertation.

### 3.1 Logic Programming Languages

We start with logic programming as exemplified by languages such as Prolog [9, $35,38,52,101,108,132,144,147]$ and Datalog [30, 76, 119, 140].

Logic programming is commonly associated with the declarative programming paradigm, where the programmer should only have to specify what, rather than how, to perform a computation [100]. Unfortunately, in my opinion, logic programming languages fail to deliver on being "truly" declarative, as I will show when discussing how these languages work.

The way logic programming pursues its declarative goal is by defining a program via logical clauses instead of procedural operations. The usual syntax is similar to the syntax of Dyna that we saw in section §2.1. The only "aggregator" that logic programming languages support is the :- aggregator, which is intended to look like a backward implication symbol ( $\Longleftarrow$ ). The head of the expression, or the part to the left of the :-, is the expression that we are trying to prove by depending on the predicates that appear on the right-hand side of the :--

$$
\begin{array}{l|l}
283 & a(X):-b(X, Y), c(Y) . \\
284 & a(X):-d(X) .
\end{array}
$$

Variables represented placeholder values, just like with Dyna.
The way in which logic programming languages "execute" differs greatly depending on the kind of logic programming language used. That said, one of our goals with Dyna is to unify the two major approaches for executing logic programs. We hope that our execution approach will simultaneously benefit from the advantages both approaches have to offer while minimizing their respective weaknesses.

### 3.1.1 Prolog Language

The oldest and most iconic family of logic programming languages is Prolog. For reference, the first appearances of Prolog happened in 1972, the same year the C programming language was developed [5, 35, 38, 144]. Over the years, there have been countless implementations and extensions of the Prolog language [9, 21, 43, 52, 137, 144, 147, 152]. The diversity in Prolog implementations is not surprising when one considers that a simple Prolog can be implemented in less than 200 lines of code in a procedural programming language [137].

Prolog's approach to execution certainly shows some taint of the limited memory environment in which it was invented. In fact, writing a large Prolog program often requires extensive knowledge of how the Prolog system executes a program.

Prolog searches for the existence of a proof, utilizing a greedy backtrackingbased search method. The way this works is that Prolog searches through declared predicates in a top-to-bottom, left-to-right order, using unification [102] to check if the current assignment to variables is consistent. As long as the current partial assignment is consistent, Prolog continues to expand the program. When Prolog eventually reaches the "end" of unification, it returns the value "true" (typically printed as "yes") to the user as well as the assignment to variables.

To make this description more concrete, consider the program in figure 3-1.

```
a(X) :- X = 0.
a(1). % conceptually equivalent to line 285, instead with 1 instead of 0
a(2).
a(3).
a(4).
a(5).
a(6).
a(7).
a(8).
a(9).
b(A,B,C,D,E,F,G,H,I,J) :- a(A), a(B), a(C), a(D), a(E),
    a(F),a(G),a(H),a(I), a(J).
```

Figure 3-1. Simple Prolog/Datalog Program used to illustrate the differences in language execution. This program defines $a / 1^{48}$ as true for the integer values of 0-9, lines 285 to 294, and b/10 as true for all terms between 0000000000 and 9999999999.

If we query $b(A, B, C, D, E, F, G, H, I, J)$ on line 295, then the Prolog engine returns a lazy stream of bindings to the variables $A, B, C, D, E, F, G, H, I, J$. A binding is the current assignment to the variable. A binding can be either a ground value (such as the number $0,1,2$ etc.), another variable, or a structure (such as in section §2.1.1). The Prolog engine searches for a binding to all variables, which is consistent with the rules of the program. It does this by expanding rules in the standardized left-to-right, top-to-bottom evaluation order. In the case of a/1 and b/10, it starts by expanding $b / 10$, where it encounters the first clause $a(A)$. Then it will unify the

[^28]variable $A$ with the variable $X$ on line 285. At this point, neither the variable $A$ nor $X$ has any ground value assigned to it. The Prolog engine continues to evaluate the program in its left-to-right order and encounters $\mathrm{X}=0$ on line 285. This causes the Prolog engine to assign the value 0 to $X$, which causes the value of $A$ to also be set to 0 as a result of the previous unification between $A$ and $x$. Because there are no more clauses on line 285, the Prolog engine returns control flow back to line 295 where $a(B)$ will then be evaluated, and so on. However, before this return of control flow occurs, the Prolog engine marks a point in its execution, which is used for backtracking.

Once the Prolog engine has reached the end of line 296, it will return the current binding to the caller of $b / 10$. In this case, it was called from the toplevel user query, so the assignment $A=0, B=0, C=0, D=0, E=0, F=0, G=0, H=0, I=0, J=0$ is returned to the user. If we ask the Prolog engine for the next assignment, then it will go to its stack of backtracking locations and pop the most recent location. In this case, it would be where $J$ is assigned the value 0 . The engine then advances to the next disjunctive branch of $\mathrm{a} / 1$, in this case line 286, assigning the value 1 to J. Again, there are no more conditions on line 286 or line 296 which need to be handled, so the new binding is returned to the user, which is now $A=0, B=0, C=0, D=0, E=0, F=0, G=0, H=0, I=0, J=1$. This process will continue until $A=9, B=9, C=9, D=9, E=9, F=9, G=9, H=9, I=9, J=9$, at which point there are no more possible bindings to consider for the variables A-J.

This procedure is conceptually quite simple and has the advantage of only using a limited amount of memory-it only requires storage for the current assignments of the variables and records of where branching decisions were made so it can backtrack. Unfortunately, writing good Prolog programs requires that the programmer think about how the Prolog engine works, and therefore is not entirely "declarative".

For example, suppose that we want to use b/10 to query if there is an integer where the second digit is larger than the first digit, e.g., a number like 0100000000. This can be done using something like line 297:
$297 \mid c:-b(A, B, C, D, E, F, G, H, I, J), A<B$.
Figure 3-2. Checking if there is a number where the second digit is greater than the first digit. For example, something like 0100000000.

Now, line 297 does "work"; however, it is extremely inefficient. The Prolog engine always evaluates left-to-right, which means that it creates a complete assignment to the variables A-J before it checks A < B. This means that Prolog loops through $10^{8}$ combinations before finding an assignment that satisfies $A<B$.

Ideally, we would like to intermix the checking of A < B with the assignment to variables A and B. Thankfully, modern Prolog implementations have a solution to this. Prolog (with a CLP extension, section §3.1.5) allows tracking of delayed constraints, which are constraints (like A < B) that cannot be immediately evaluated. These constraints are conceptually similar to the $A=X$ constraint that was caused by the unification on line 285. However, these constraints are "more powerful" in that instead of depending on manipulation of pointers in internal data structures (see Martelli and Montanari [102]), these constraints are allowed to define their own arbitrary code for handling when a variable's state is modified.

Now, when using delayed constraints, we must still remember that Prolog evaluates left-to-right, as adding the delayed constraint after b/10 does not help:

298 d :- A \#< B, b(A, B, C, D, E, F, G, H, I, J).
Figure 3-3. Delayed Constraint in Prolog. Delayed constraints are annotated using a hash symbol \#. [9, 147]

The way that A \#< B works is that it is saved as a delayed constraint into the constraint store, or program's state (which previously only held variable binding), and will run the \#< code anytime that A or B is "modified". This means that when B is assigned the value 0 , and $A$ already has the value 0 assigned, it will quickly fail the check of A \#< B, causing the Prolog engine to immediately backtrack. ${ }^{49}$

Now, this is great! But there is still some behavior that is a bit "annoying" (and

[^29]also differs from Dyna's approach). If we query for ' $d$ ', the Prolog engine will quickly prove it is true but returns a lazy stream that keeps returning solutions. In other words, rather than finding one solution to prove ' $d$ ', the Prolog engine enumerates all $45 * 10^{8}$ proofs of ' $d$ '. This happens because every time that we ask for the next item in the lazy stream of solutions, Prolog simply backtracks to the previous variable, even if the returned solution (that ' $d$ ' is true) is identical to the previously returned solution. This differs from Dyna, which would only return ' $d$ ' is true once due to the aggregation performed by :- over all of the answers.

To fix this, we can use a cut in Prolog, which prevents the Prolog engine from backtracking over different assignments to variables. A cut in Prolog is annotated with an exclamation mark '!'.

$$
299 \text { e : - A \#< B, b(A, B, C, D, E, F, G,H,I,J), !. }
$$

Figure 3-4. Adding a cut to avoid backtracking through all possible assignments.

Now the 'e' rule will only find one solution and return true one time.
This again shows how Prolog programmers need to be aware of how the Prolog engine works internally to make their programs work well. Conversely, Dyna's is designed to not need operations such as cut. We believe that aggregators such as :- and ?= provide sufficient mechanisms for handling these cases.

### 3.1.1.1 Infinite Relations in Prolog

Prolog treats unification between variables as a first-class built-in operation. This means that expressions such as $X=Y$ are handled without backtracking over possible assignments to $X$ or $Y$. This is achieved by tracking which variables are unified together by essentially updating the internal pointers backing $X$ and $Y$ so that the space in memory reserved for holding the value of $X$ and $Y$ become the same space [102]. This also extends to structural terms. So unification like $X=g(Y, 7)^{50}$ are handled in a similar way.

[^30]This approach lets the Prolog engine work even when there are infinite relations by tracking "constraints" and leaving the exact value of some variable as unknown.

Therefore, the extension of modern Prolog with delayed constraints (as in figure 33) can be seen as an extension of unification from tracking only variable unification via pointer reassignments to allowing arbitrary code to run when a variable's "unification status" is changed.
${ }_{300} \mid f(X, Y):-Y=g(X, 7, V), V \#=X+3$.
Figure 3-5. An example infinite relation in Prolog. $f(X, Y)$ does not require any backtracking to evaluate. It will define an infinite number of terms like $f(1, g(1,7,4)$.

### 3.1.2 Datalog Language

Datalog is the other major approach to logic programming [30, 76] and was the primary inspiration for Dyna 1.0 [60, 61]. A short description of how Datalog works is that it uses the same high-level syntax as Prolog for dynamic programming [18] over a boolean semiring. ${ }^{51}$ Let us break this down a little. Datalog's approach can be summarized as, the Datalog engine stores all true facts. A Datalog fact is a statement that is known to be true and is represented as an term (without any variables). For example, $a(1)$ is a fact, but $a(X)$ is not a fact as it contains the variable X .

Datalog uses the rules in the program and all the stored facts to deduce new facts-storing those as well. A Datalog engine finishes running when the deduced facts are "stabled"-meaning that it has reached a fixed point, just like Dyna section §2.5. Because everything is stored, this creates different kinds of opportunities for how to execute a Datalog program.

First, because everything is stored, this means that even simple programs can cause a Datalog system to be quite inefficient. For example, figure 3-1 would end up storing $10^{10}$ facts in memory to represent $\mathrm{b} / 10$. The simplicity of Datalog means it does not have a mechanism to efficiently represent this program.

Conversely, having a simple "store everything" approach does have some advantages. For example, Datalog systems will often employ a host of efficient join

[^31]techniques as in a database system (e.g. [106]). Datalog can also employ brute force strategies such as "loop over everything" and be guaranteed that these strategies terminate. The reason is that a Datalog program can only represent finite relations. ${ }^{52}$ Being finite is both part of Datalog's design but also follows as a result of Datalog storing all facts, and computers have a finite amount of storage. Hence, there can only be a finite amount of facts stored.

### 3.1.3 Datalog is Breath First, Prolog is Depth First

Another advantage of Datalog's approach is that it avoids Prolog's greedy search behavior. For example, in the following program:

```
301 h :- h.
302 h :- true.
```

Figure 3-6. Datalog program which deduces that ' $h$ ' is true due to line 302.

Datalog easily deduces that ' $h$ ' is true due to line line 302. Once ' $h$ ' is stored in memory, the Datalog system does not get stuck handling line 301 and can easily solve this program. Whereas, Prolog gets stuck on this program. The reason is that Prolog would backtrack through line 301 and reencounter ' $h$ ' endlessly, without ever attempting to run line 302, due to Prolog's top-to-bottom, left-to-right evaluation order.

In this way, we can roughly think of Datalog's approach as breath-first evaluating of logic programs while Prolog is depth-first evaluation. Both of these languages exhibit the relative advantages and disadvantages of these breath vs depth first search: memory overhead, robustness to search order, complexity of internal state (with unified variables in Prolog), etc. In Dyna, we are essentially looking to combine the best of both Datalog and Prolog. We want to avoid cases where a bad execution order causes the system to not terminate or be inefficient. We also want to be able to store deduced facts (like Datalog), while still representing infinite relations (like Prolog).

[^32]
### 3.1.4 Aggregation in Logic Programming

In Dyna, a central feature is aggregation, which appears everywhere throughout a Dyna program and is built into the Dyna syntax (section §2.2). Both Datalog and Prolog "support" aggregation, though their implementations are less flexible than Dyna's aggregation and it is essentially bolted on. The way aggregation has been implemented is by adding "meta predicates" that internally perform the operation of aggregation. In Prolog, this might look something like figure 3-7:

```
j(X,AggResult) :- bagof(InputToAggregator,
                        body_getting_aggregated(X, InputToAggregator),
                        AggList),
    sumlist(AggList,AggResult).
body_getting_aggregated(1,1).
body_getting_aggregated(1,2).
body_getting_aggregated(1,1).
```

Figure 3-7. Aggregation supported by Prolog ${ }^{53}$

Here bagof/3 will collect all of the assignments to InputToAggregator into a list AggList.

Evaluation of bagof/3 requires that all possible assignments to body_getting_aggregated/2 are completed before bagof/3 "returns". This prevents any opportunity to make the aggregator more efficient, as we will see later in section § 6.5 with Dyna.

Note that line 307 and line 309 both assert that 1,1 is true. As a result, the aggregation from Prolog will give a different result than Datalog. In Prolog, it will count the number of times that something has been deduced. In this case with Prolog, we will have 'AggList $=[1,2,1]$ '. However, with Datalog, we get 'AggList $=[1,2]$ ', since the second number 1 was already deduced and not counted a second time.

[^33]
### 3.1.5 Constraint Logic Programming

Constraint Logic Programming (CLP) is a derivative of Prolog-style logic programming. As we already saw in figure 3-3, a delayed constraint is an expression that cannot be immediately evaluated. These constraints are stored in a constraint store, alongside unifications between variables. Constraints are re-evaluated when variables are assigned. Constraints can also interact with other constraints to infer new constraints through propagation [72].

For example, if we have ' $j$ ' defined on line 310:
${ }_{310} \mid \mathrm{j}:-\mathrm{A} \#>5, \mathrm{~B} \#=\mathrm{A}-2, \mathrm{~B} \#<0$.
Figure 3-8. The variable $A$, must be some value greater than 5 , and $B$ must be less than 0 . Additionally, we have that $B=A-2$, hence there is no valid assignment to both $A$ and $B$.

Then there is no possible way to assign $A$ and $B$ such that ' $j$ ' is true. This program can be handled by CLP using interval constraints. First, A \#> 5 will track that A must have a value greater than 5 (the interval $(5, \infty)$ ). Next, the constraint B \#= A - 2 identifies that $B$ is two less than $A$, therefore, it will have an interval of greater than 3 associated with it (the interval $(3, \infty)$ ). Finally, the constraint $B \#<\theta$ will associate an interval of less than 0 , which does not overlap with the greater than 3 interval $((-\infty, 0) \cap(3, \infty)=\emptyset)$. Hence, the CLP engine initiates Prolog-style backtracking, removing constraints and unifications from the constraint store.

To make CLP work, a CLP system will have hundreds of different propagation and simplification rules. These rules are sometimes called Constraint Handling Rules, usually referred to as CHR [72]. Development of new CHRs and ways to implement CHRs has been an area of research over the years [72, 105, 118, 131]. Some of the rewrite rules presented in chapter $\S 6$ are similar to the rules found in CHR [72].

### 3.1.5.1 MiniKanren

MiniKanren [27, 71] was designed as a small implementation of logic programming that can be implemented in less than 200 lines of Scheme. As a result of MiniKan-
ren's small core, it has been the basis for much research on logic programming over the years [28, 91]. MiniKanren's approach to logic programming differs from Prolog in that it does not mutate any global data structures state ${ }^{54}$ but instead returns lazy streams of binding states. Each operation is implemented as a function that modifies the maps from a stream of bindings to a new binding, returning no elements in the event that there is an inconsistency between the variable bindings [26].

The approach taken by miniKanren is similar to our approach in that we are going to avoid mutating global state (like Prolog); however, we do not depend on passing the host's language functions around (like miniKanren) and instead will develop an explicit relational expression representation ( $\mathbf{R}$-exprs in chapter §5). This will allow us to be more flexible with the kinds of mutations and expressions that we can represent and to have more flexibility in picking the order of evaluation.

### 3.1.6 Constraint Satisfaction Programming

Constraint satisfaction-based solving techniques, such as SMT solvers [15, 16, 47, $48,50,107$ ] or answer set programming [64, 79] work by representing a problem as variables and constraints between variables. These systems will often have some ability to handle "function calls" by expanding the function up to some depth. ${ }^{55}$ The representation of a problem is then translated into a mathematical theory that can be solved. For example, the SMT formalism to multiple calls to a modified SAT solver [16]. These approaches have the advantage of residing on the foundation of a sound mathematical theory. It is possible that we could solve Dyna programs the same way. However, we have opted not to take this approach. The reason is that we do not believe that it will work well with the kinds of problems we are interested in. The kinds of Dyna programs we expect will realistically have many more variables than can be realistically solved using these kinds of approaches. ${ }^{56}$ Instead, we have opted for a fixed-point based solving technique, which will iterate until a solution is found (section §2.5).

[^34]
### 3.2 Probabilistic Programming

Dyna is technically not a probabilistic programming language, though it is frequently compared to probabilistic programming languages.

A probabilistic program assigns probabilities to different assignments of the variables [29, 37, 65, 81, 112, 138, 139]. In some ways, we can think of this as a generalization of logic programming, which only assigns true or false with being able to assign factional chances of being true. At a high level, the probability is similar to Dyna's ability to weigh the result from a rule and combine them with an aggregator (section §2.2). However, the difference is that Dyna allows for general weights, which might not be a probability, and probabilistic programming requires that the weights be probabilities. Because probabilistic programming works with a more restricted problem, they will often provide features that are useful for modeling probabilities. For example, probabilistic programming systems have the ability to fit parameters to observed data. ${ }^{57}$

Like Dyna, probabilistic programming systems are usually not complete languages but are either embedded as a DSL or used as a library from a driver program like Dyna (as in section §2.3.1).

### 3.3 Relational Algebra

In chapter $\S 5$, I will go into detail about how we use a relational algebra we call R-exprs to model and implement the Dyna programming language. Chapter $\S 5$ will provide an introduction to relational algebra, so here, I will instead focus on surveying related work that uses a relational algebra.

The most iconic systems that are built on a relational algebra are $\mathrm{SQL}^{58}$ database systems [36, 44]. A SQL database consists of database tables and is queried using SQL queries. Relations in a database are represented by database tables. Tables are combined (intersected or unioned) with other tables, filtered for particular values, and projected to select for a subset of the database table. The result is represented

[^35]as another database relation. The relation can be either returned to the user (as in the case of a query), saved back into the database as a new relation, or further extended and used as a component of a larger query. An example SQL query is shown in figure 3-9:

```
SELECT column_a, column_b, column_c \trianglerightSelect columns using projection
FROM table_1
JOIN table_2 on table_1.column_a = table_2.column_f DIntersect relations
WHERE column_d = 'identifier_1' and table_2.column_e > 11 \trianglerightFilter
```


## (a) SQL

```
result(A,B,C) :- table_1(A,B,C,'identifier_1'), % filter with value
    table_2(E,A), % intersect by reuse of var A
    E > 11. % filter as external constraint
```

(b) Equivalent Dyna

Figure 3-9. Example SQL query

Internally, SQL databases work over finite materialized relations. The relations provide internal APIs to access the underlying data. This can include filtering using a particular key or looping over all tuples in the relation using an iterator [134]. Database systems include query optimizers that rearrange operations in the query to automatically figure out the most efficient way to run queries. ${ }^{59}$

Just like Dyna, SQL supports aggregation and grouping [46, 83, 98]. In both SQL and Dyna, aggregation is represented as reducing a relation into another equivalent relation with the result of aggregation. An example is shown in figure 3-10.

```
SELECT column_a, sum(column_b)
FROM table_1 }\mp@subsup{}{321}{*}|\operatorname{result(A) += B for table_1(A,B,C,D).
GROUP BY column_a
```

(b) Equivalent Dyna

## (a) SQL

Figure 3-10. Example SQL with aggregation and GROUP BY.

[^36]
### 3.4 Term rewriting

Term rewriting [ $10,12,13,17,32-34,51,77,78,93,94,114,143$ ] is a central idea used in this dissertation (chapters 5 and 6). Essentially, given an expression or term, it is rewritten into an semantically equivalent representation. This rewriting process corresponds with execution of the program. For example, if we have the term $2+3$ it can be rewritten as 5 , which is semantically equivalent and represents the evaluation of the plus sign.

The abstraction of term rewriting is too broad to completely cover, so I will instead focus on a few subtopics that are related to the ideas explored in this dissertation.

### 3.4.1 Implementation of Term Rewriting

A term rewriting system is defined using both the structure being rewritten and the rewrite rules performed against the structure. In this dissertation, I will create our own implementation of term rewriting based on the rewrite rules that we defined (chapters 8 and 11). While creating a term rewriting system from scratch is entirely reasonable, there are frameworks such as maude [33, 34] or k-framework [114] that create an implementation from the definition of rewrite rules only. As will be discussed in chapter $\S 8$, the reason that we choose to create our own "rewriting framework" for Dyna, rather than using an existing framework, is that the way in which we apply rewrite rules differs from what existing frameworks are designed to support. The existing frameworks are designed to support procedural languages, where the order of execution is deterministic. Therefore, they only have to match their rewrite rules against the next instruction to execute. Conversely, in Dyna, our execution is non-deterministic, and we implement this by allowing any applicable rewrite rule to match any part of the term.

### 3.4.2 Term Rewriting a Relational Algebra for Logic Programming

Term rewriting on top of a relational algebra for the implementation of logic programming has been experimented with before: Arias et al. [10], Bellia and Occhiuto
[17], Gallego Arias et al. [77, 78]. Bellia and Occhiuto [17] was the first that we are aware of to make the connection in 1993, they created a "variable free" ${ }^{60}$ representation of logic programming they called c-expressions. They also identified that their relational algebra representation can be manipulated with rewrite rules, as we will do in chapter §6. Arias et al. [10], Gallego Arias et al. [77, 78] is an ongoing project which also uses a term rewriting-based formalism to implement logic programming. Like [17], their systems is focused on logic programming and does not support aggregation as we do with Dyna.

### 3.4.3 Functional Logic Programming

Dyna started as an extension of Datalog and has added features of Prolog and eventually added enough features to (in my opinion) be comparable to functional programming. ${ }^{61}$ As such, I think that functional logic programming languages and the version of Dyna in this dissertation share a number of similarities.

The space of functional logic programming has been around for several years, with a journal running from 1995 to 2008 [3] and at least two other programming language research projects in this space: the Curry language (Antoy [6], Antoy et al. [7], Antoy and Hanus [8], Braßel et al. [24], Hanus [84, 85], Hanus et al. [86, 87], Hanus and Prehofer [88], Hanus and Sadre [89]) and Verse language being developed at Epic games (Augustsson et al. [11]). Both Curry and Verse are Haskellesque languages with the addition of logic programming features. The major addition to functional programming is the ability to represent non-deterministic results from functions and assignments to variables. For example, a function can be called when its arguments' values are unknown, and only the return value is known. This is just like logic programming, which allows arguments to functions to be variables with unknown values.

The formalism for these languages is done using term rewriting, just like I do in this dissertation. There are two approaches to handling the non-determinism:

[^37]narrowing and delayed evaluation ${ }^{62}$ (as we see later in this dissertation).
The term Narrowing refers to the idea of representing a program as a set of equations with variables whose value is unknown and then solving those equations. Narrowing does not refer to a specific strategy for solving the equations, but there has been work on developing narrowing strategies for different classes of equations and for more efficient search strategies [6, 7, 85, 88].

### 3.5 Memoization \& Reactive

A long review of what memoization is can be found in section §10.1.
Memoization is the programming technique of saving the result of some computation and reusing it later rather than recomputing it every time [104]. The simplistic view of memoization requires that memoized computation is functional and unchanging so that the stored result does not have to be invalidated later. This can be fixed by making a computation reactive, upstream dependencies for any memoized value are tracked, and whenever there is a change, all downstream dependents are recomputed [14]. (The exact details of how reactive programming is implemented differ greatly between implementations.)

In the context of logic programming, languages such as Datalog (section §3.1.2) are entirely based on memoization, with everything stored instead of "calling a function to perform computation". XSB is an extension of Prolog which adds memoization [135, 145, 151]. XSB works by placing markers on terms that are expanded during backtracking. If a cycle is detected during backtracking, it memoizes that the term where the cycle is detected is memoized as false. If something is later deduced as true later in the computation, it triggers recomputation.

Prior work from the Dyna project has also focused on developing a formalism for memoization and reactive programming: Filardo [66], Filardo and Eisner [67]. This work sought to formalize programs as "computational circuits" where the memoized values are vertices in a graph and edges in the graph track computation dependencies. A vocabulary of different message types that can be passed along the edges was developed that corresponds to different kinds of implementations of reactive programming. For example, there are different options around how

[^38]the recomputation is scheduled and when changes are visible to downstream dependents.

### 3.6 Tracing JIT Compilation

In chapter §12, I will discuss my efforts to compile Dyna to make it run faster. Our compilation method is inspired by traced-based JIT compilation [31, 74, 75] and will also make use of partial evaluation [42, 57, 73, 116, 149]. Trace-based JITting is a very flexible technique for implementing a JIT compiler has been previously used on the first JavaScript JIT compilers [74], the PyPy Python JIT [20], LuaJIT [109], the TorchScript for neural PyTorch models [2], and even Prolog [21]. ${ }^{63}$ Tracing differs from method at a time compilation in that instead of compiling methods that appear in the program, it compiles a sequence of steps performed when executing the program. When there is a conditional branch in the program, the tracing compiler inserts a check that the conditional branches the same way each time. The JIT compiler will insert a stub that is used to resume the trace, and only when the branch is actually taken does it generate the code.

To see how trace-based JIT compilation works, let us work through an example of compiling the program presented in figure 3-11 using tracing. This program starts with ExampleFunction and uses the MyPrint function to indirect to the built-in Print call.

[^39]```
function ExampleFunction(n)
    for i\in[0,n):
        if i<5 :
            MrPrint("hello")
        else if i> 100000000000 :
            MyPrint("never")
        else
            MyPrint("world")
    MyPrint("done")
    return
function MyPrint(x)
    Print("something " + x)
```

Figure 3-11. Example function getting traced

When we generate a trace of figure 3-11 in figure 3-12, observe that we have not generated all of the code from figure 3-11. Instead, we have some lines that are marked "not yet generated". These lines of code contain sufficient metadata and jump statements to get back into the JIT compiler to resume tracing.

```
i\leftarrow0
if i\geqn: \triangleright Condition to check
    not yet generated
if i\geq5: \triangleright Condition to check
    not yet generated
if i> 100000000000:
    not yet generated
Print("something hello") \triangleright The MyPrint function is embedded
i\leftarrowi+1
goto 2 \triangleright Return to top of loop
```

Figure 3-12. First generated version of the function ExampleFunction from figure 311 using tracing.

Observe that the trace has stopped with line 10. The reason is that when tracing, the control flow has jumped back to a location in the code that has already
been generated. Therefore, the tracer compiler will generate the equivalent jump statement to the previously generated statement. Furthermore, observe that the function MyPrint does not appear in the generated output of figure 3-12. The reason is that the trace only contains useful operations, such as incrementing the variable $i$ and calling the built-in Print, but the user's function MyPrint simply gets "absorbed" in by the processes of tracing. ${ }^{64}$

As the program continues to run, it will eventually hit one of the "not yet generated" branches, and it will update jump statements and generate additional code as in figure 3-13.

```
\(i \leftarrow 0\)
if \(i \geq n\) : \(\triangleright\) Condition to check
    not yet generated
if \(i \geq 5: \quad \triangleright\) Updated to branch to newly generated code
    goto 13
if \(i>10000000000\) :
    not yet generated
Print("something hello")
\(i \leftarrow i+1\)
goto 2
\(\triangleright\) Additional compilation to fill in code from line 5
Print("something world")
\(i \leftarrow i+1\)
goto 2
```

Figure 3-13. After the first "not yet generated" branch has been hit and additional code has been added.

This process of running the generated code and replacing the "not yet generated" branches with generated code will continue. Eventually, the program is done and will hit the return statement, as in figure 3-14. Not all parts of the program have been executed, so there are parts of the generated code that contain "not yet

[^40]generated" on some branches.
1: $i \leftarrow 0$
2: if $i \geq n$ :
3: $\quad$ goto 18
if $i \geq 5$ :
goto 13
if $i>10000000000$ :
not yet generated
Print("something hello")
$i \leftarrow i+1$
goto 2
11:
12: $\triangleright$ Additional compilation to fill in code from line 5
PRINT("something world")
$i \leftarrow i+1$
goto 2
16:
17: $\triangleright$ Additional compilation to fill in code from line 3
18: Print("something done")
return $\triangleright$ Generate return to caller code
Figure 3-14. The program has ended as it hits the return statement. Not all branches of the code have been hit, so there can still be un-generated parts of the compiled code.

## Chapter 4

## Challenges in Dyna

Usually the design and implementation of a programming language are done at the same time. This means that difficult-to-implement features and combinations of features are frequently left out of programming languages. Dyna did not have this luxury. Dyna's design was proposed in 2011 by Eisner and Filardo [59] without an implementation. ${ }^{65}$ Furthermore, the long-term vision for the Dyna project is that all programs that are "conceptually sound" (as read by a human, not a computer) and describe the programmer's intent (in a declarative way) should "just work". A more formal way of stating this vision would be to say, "all syntactically correct programs which pass some elementary level of type checking should 'work' and return some 'useful' result to the user when queried".

Now, this vision is an amazing pitch. Nevertheless, it certainly complicates Dyna's implementation. For instance, this vision does not provide us with any insight into the methods to use for Dyna's implementation. However, we can reject many previous approaches to logic programming implementation, as they are incapable of supporting Dyna's vision.

My intention in this chapter is to provide you, the reader, with an understanding of our motivation before I subject you to hundreds of pages about implementation. Note: I am not selling the vision of Dyna-this dissertation is about implementation. I consider the vision of Dyna prior work published in 2011. I encourage anyone interested in more details on this vision to read the original 2011 paper by Eisner

[^41]and Filardo [59]. Additionally, I will state that the implementation presented in this dissertation does not completely deliver on Dyna's vision-there exist syntactically correct and conceptually sound programs that we currently cannot run. However, it is our belief that the techniques presented in this dissertation provide a sufficient foundation for Dyna, such that the vision of Dyna can be approached "in the limit" with additional work and features being added.

### 4.1 Features in Dyna

We start by looking at the features that Dyna supports and how this compares to other declarative frameworks, as shown in table 4-I.

A high-level outline of features supported by Dyna, as compared to other declarative frameworks, is shown in table 4-I. A high-level description for each feature is as follows:

- Finite - All systems are able to define finite relations. By finite, we mean that the set of all declared entries is finite $|\{\cdots\}|<\infty$. For example, the integers between 0 and 10 is a finite set $|\{1,2,3,4,5,6,7,8,9\}|<\infty$. Conversely, an example of a non-finite relation would be all natural numbers as $|\mathbb{N}| \nless \infty$.

Only supporting finite means that simple strategies that enumerate everything are workable. The reason is that brute force strategies are guaranteed to terminate. Because of the termination guarantee, it makes it easier to experiment with different execution orders in finite systems without worrying about secondary non-termination issues.

- Deductive - Deductive means that the system deduces new facts in the language. In logic programming, this is represented as:

322 deduced :- requirement_1, requirement_2.
Logic programming languages, such as Prolog and Datalog, are both deductive, whereas languages such as SQL are generally not considered deductive. The SQL language does not automatically perform inferences for new entries in the database tables, though this can be added manually through the use of database triggers.

| Feature | $\begin{aligned} & \mathrm{SQL} \\ & \S 3.3 \end{aligned}$ | Datalog <br> §3.1.2 | Prolog §3.1.1 | $\begin{aligned} & \text { CLP } \\ & \S 3.1 .5 \end{aligned}$ | Probabilistic Programming §3.2 | $\begin{aligned} & \text { SMT } \\ & \text { §3.1.6 } \end{aligned}$ | Dyna 1.0 [60] | Dyna 2.0 (This dissertation) §2 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| Finite | $\checkmark$ | $\checkmark$ | $\checkmark$ | $\checkmark$ | $\checkmark$ | $\checkmark$ | $\checkmark$ | $\checkmark$ |
| Deductive | $x$ | $\checkmark$ | $\checkmark$ | $\checkmark$ | $\approx$ | $\approx$ | $\checkmark$ | $\checkmark$ |
| Updatable | $\checkmark$ | $\checkmark$ | $\checkmark$ | $\checkmark$ |  | $\checkmark$ | $\checkmark$ | $\checkmark$ |
| Semiring Weighted | $x$ | $x$ | $x$ | $x$ | $\approx$ | $x$ | $\checkmark$ | $\checkmark$ |
| General Weighted | $x$ | $x$ | $x$ | $x$ | $x$ | $x$ | $x$ | $\checkmark$ |
| Aggregation | $\checkmark$ | $\chi_{\text {§3.1.4 }}$ | $\chi_{\text {§3.1.4 }}$ | $x$ | $\checkmark$ | $X_{\text {[50] }}$ | $\checkmark$ | $\checkmark$ |
| Memoization | $\checkmark$ | $\checkmark$ | $\chi_{\text {[135] }}$ | $x$ | $x$ | $x$ | $\checkmark$ | $\checkmark$ |
| Turing Complete | $x$ | $x$ | $\checkmark$ | $\checkmark$ | $x$ | $\approx$ | $\checkmark$ | $\checkmark$ |
| Unconstrained Execution Order | $\checkmark$ | $\checkmark$ | $x$ | $\approx$ | $x$ | $\checkmark$ | $\checkmark$ | $\checkmark$ |
| Constraints | $x$ | $x$ | $x$ | $\checkmark$ | $x$ | $\checkmark$ | $x$ | $\checkmark$ |
| Object Oriented | $x$ | $x$ | $\chi_{[49,146]}$ | $x$ | $x$ | $x$ | $x$ | $\checkmark$ |
| Syntactic Sugar |  | $x$ | $\approx$ |  |  |  | $x$ | $\checkmark$ |

Table 4-I. Different Declarative Programming Languages (Paradigms) compared by feature. The definitions for the features are given on the next page. This table represents what is most commonly associated with a particular programming paradigm. References are included in the table when some further research has been done to add a particular feature.

- Updatable - If the program or data can be changed after the program has started running. Most of the systems in table 4-I provide some kind of REPL where they can be started and maintain state between interactions.
- Weighted - Every expression in the system has a weight. Logic programming languages only have the single weight of true. Some probabilistic programming languages associate a probability with every expression in the system in the case that the language is semiring based (e.g. [65]). Other probabilistic programming languages will randomly sample results from the program and compute the probability for each result depending on how probabilistic factors were used in the execution of the program (e.g. [29, 81, 138]). This allows a probabilistic programming language to represent a result as being $50 \%$

Dyna allows for a value to be associated with every expression. The value is not required to be a numerical value or even a probability.

- Aggregation - Dyna and SQL build aggregation into the language as a central feature. Logic programming with Prolog or Datalog does technically support aggregation as shown in section §3.1.4, but the syntax is generally quite awkward and there are many footguns ${ }^{66}$, in that using the aggregation feature can result in unexpected behavior, as previously described.
- Memoization - If there is a mechanism to avoid performing the same computation multiple times. This is admittedly more of an implementation feature than a language feature. However, the design of the language does (usually) influence the implementation.

Datalog and SQL materialize the result of all intermediate computations and hence are entirely memoized. Prolog generally does not perform memoization at all, but there are some extensions that add memoization (see [135, 145, 151]). The Dyna implementation supports memoization as a central feature, section §2.7 and chapter §10.

- Turing Complete - Many declarative programming frameworks are not Turing-complete, such as SQL and Datalog. The reason for this is that these

[^42]systems instead prioritize termination guarantees, making it easier to rearrange operations/optimize database queries.

Some SMT solvers are also Turing-complete despite the fact that they internally reduce their computation to SAT solving, which is not Turing-complete. The reason is that SMT solvers attempt many reductions, increasing the size without bounds until a solution is found.

- Unconstrained Execution Order - Reordering expressions in a language provides optimization opportunities. SQL is well known for having a query optimizer that attempts to figure out the best order to run operations and perform joins between database tables. Advanced Datalog systems will also reorder operations for better performance. As stated before, Prolog does not allow for automatic reordering of the expressions (section §3.1.1). SMT achieves out-of-order execution and Turing completeness by reducing to a SAT solver, which internally can perform the out-of-order evaluation.
- Constraints - Constraints are logical expressions that can be combined together. SQL, Datalog, and Prolog are generally not considered to have constraints, the logic expressions are evaluated immediately when encountered, though SQL and Datalog do allow query optimizers to reorder expressions before execution starts. In the previous chapter (section §3.1.5), we mostly focused on constraint logic programming (CLP), which is the extension to Prolog that allows for delayed constraints.
- Object Oriented - Most declarative frameworks do not support objectoriented programming. This is a feature that has been added to Dyna (section §2.9). The complication here is that the standard approaches to object-oriented programming are procedural (such as set field on an object, and call function which mutates fields on the object) and do not work well in a declarative paradigm.
- Syntactic Sugar - Dyna aims to be a more modern programming language with the addition of syntactic sugar. For example, we have built-in notation for aggregators, dynabases, lambda functions, etc. The other languages in this table are either Prolog-style or SQL, which do not have this level of syntactic sugar. Admittedly, this does not increase the complexity of implementation
too much as this is entirely handled by the front-end parser, which is scarcely mentioned in this dissertation.


### 4.2 Examples of Difficult Programs

To illustrate some of these features of Dyna and how their combination can make executing programs difficult, let us take a look at a few example programs.

### 4.2.1 All Pairs Shortest Path

The first program is the all-pairs shortest path in a directed graph program, which previously appeared in section §2.6. Each edge of the graph is defined using an edge rule, whose values (weight) is the length of the edge. By construction, we have for any Start and End, such that there exists a path between Start and End, the value of path(Start, End) is the length of the shortest path.

```
path(Start,Start) min= 0. % base case
path(Start,End) min= path(Start,Mid) + edge(Mid,End). % recursive case
edge("baltimore", "washington dc") = 38. % example rules for edges
edge("baltimore", "new york") = 195.
    % many other edge rules omitted
```

This program is not a valid Datalog program and is also likely to not terminate when evaluated under Prolog. The reason why this is not Datalog is that line 323 defines an infinite relation, which is not allowed under Datalog. For example, the distance between path("does not exist", "does not exist") is 0 due to line 323. This line can be modified to work with Datalog by changing it to 'path(Start, Start) min= 0 for city(Start).' in which case 'city(Start)' ensures that the set of starting locations are all valid city names. However, this does change the semantics of the program as path("does not exist", "does not exist") changes from the value of 0 to no longer being defined or null.

A Prolog-style backtracking solver allows Start to be an unknown value and would attempt to answer queries using its depth-first backward chaining strategy. Unfortunately, Prolog's strategy would recurse forever due to line 324 even on the query path("atlantis", "baltimore"). The reason is that path("atlantis",
"baltimore") would recursively depend on path("atlantis", $Y$ ), which would recurse to itself. Prolog essentially gets stuck and never evaluates edge(Mid,End).

This example perfectly demonstrates the limitations of Datalog and Prolog, as the query path("atlantis", "baltimore")? can be easily recognized by a human as the shortest path in a graph between two cities. Even the query path("atlantis", Y)? can be recognized as a single-source shortest path problem and can be solved using Dijkstra's algorithm [53]. A human programmer can also identify that the query path $(X, Y)$ ? is solvable using Bellman-Ford [19, 69] and augmenting the returned pairs of city distance with path(Start,Start) min= 0 .

Our challenge with Dyna is to design a system that is able to handle the fact that Start is a variable while still avoiding the complications of cycles. Furthermore, we would like to recover the efficient strategies of Bellman-Ford or Dijkstra's when evaluating path $(X, Y) .{ }^{67}$

### 4.2.2 "Infinite" Neural Network

Our second example program was written on a whim for our 2017 paper [142] as a natural example Dyna program, and we did not even realize that it is a difficult program to execute until we analyzed this program at a later date. This program (lines 328 to 337), defines a two-dimensional convolutional neural network.

```
\sigma(X) = 1/(1+exp(-X)). % define sigmoid function for all X
in(J) += out(I) * edge(I,J). % vector-matrix product
out(J) += \sigma(in(J)). % activation of node J
out(input[X,Y]) += pixel_brightness(X,Y). % activation for input nodes
loss += (out(J) - target(J))**2. % L2 loss of the predictions
edge(input[X,Y],hidden[X+DX,Y+DY]) = weight_conv(DX,DY). % layer 1
edge(hidden[XX, YY],output[Prop])=weight_output(XX, YY, Prop). % layer 2
weight_conv(DX,DY) := random(*,-1,1) % init with random
    for range(DX, 4,4), range(DY, -4,4).
weight_output(XX, YY, Property) := random(*, -1, 1).
```

[^43]The input to the neural network is pixel_brightness $(X, Y)$ on line 331, and could be defined as a real number between 0 and 1, for example. Lines 328 to 330 define a feed-forward neural network where each neuron in the neural network is defined by a ground value. For example, we can have neurons named input[12,34], hidden $[14,31]$ or even output["kitten"]. The output activation, out (J), of neuron ' $J$ ' is calculated as a sigmoid linear combination of the activations of other neurons ' $I$ '. An interesting facet of this example is that lines 328 to 330 does not define the topology of the neural network. Instead, the topology is entirely determined by edge ( $\mathrm{X}, \mathrm{Y}$ ), which is defined on lines 333 to 337. Specifically, in this example, hidden $(X X, Y Y)$ is activated by the $9 \times 9$ square of neurons centered at input ( $X X, Y Y$ ). Then, for each image property, Prop, all hidden units are pooled to activate the output unit, output (Prop), whose activation represents the degree to which the image is predicted to have that property.

For a given finite image input, this program defines a neural network with finitely many edges. Nevertheless, it is difficult to solve this system of equations using standard strategies. The difficulty arises from line 333. Using a Datalog's forward-chaining strategy, the value for weight_conv( $2,-3$ ) would forward-chain through line 333, which establishes values for infinitely many edges of the form edge(input $(X, Y)$, hidden $(X X, Y Y)$ ) where $X X=X+2$ and $Y Y=Y-3$, despite the fact that only a finite number of edges are used for a given finite image. Conversely, using a Prolog style backwards-chaining with queries like out (output["kitten"]) would lead to an internal query like edge( $I$, hidden $[X X, Y Y]$ ) with free variables $X X$ and YY. That query must return the entire infinite set of input-to-hidden edges, even though for a given finite image, only finitely many of these edges will touch input units that actually have values. This does not work.

To handle this program, we need to represent the infinite relation of edge ( $X, Y$ ) without having to enumerate all of the edges with a lazy iterator or having to materialize all of the edges in a table.

### 4.2.3 "Infinitely" Many Dynabases

This was already mentioned in section §2.9.1, but Dyna's object-oriented programming cannot directly adapt the existing techniques. The reason is that we can easily write a program where there is an infinite number of objects. For example,
lines 338 to 341 defines a program that computes $\operatorname{argmin} x^{2}$.

```
a(X: float) = {
    value = X**2.
}.
best_X = $arg((min= a(X).value arg X)).
```

Line 338 defines a dynabase to represent an input to the function getting optimized. In the "constructor" to the dynabase, it takes the current location where the function is being evaluated. Conceptually, line 338 defines an infinite number of dynabases, as it is defined for all real numbers.

If we have a system that must fully evaluate a subset of the program at a time, then this is impossible to solve. To handle this program, we must represent the entire program symbolically at the same time. In this case, we can recognize the dynabase as a layer of abstraction and can solve the expression $\operatorname{argmin}_{x} x^{2}$ using any number of optimization techniques.

### 4.2.4 "Infinite" Identity Matrix

Another common issue with Dyna programs is that the details that are useful for running the program are not included in the source code. Recall, at the start of this chapter, Dyna's stated goal is that a program that is understandable by humans (not computers) should "just work".

A small example of this is an "infinite" identity matrix:

```
identity_matrix = new matrix \{
    \(\operatorname{elem}(X, Y):=0\).
    \(\operatorname{elem}(X, X):=1\).
\}.
```

Here, we are defining that any element in the matrix such that both of its first and second argument are the same will have a value 1 (line 344), and 0 otherwise (line 343).

This matrix could be used with other matrices. For example, we can define a matrix multiplication on line 353:

```
matrix = new {
    row(X) :- $self.elem(X,_).
    rows += 1 for $self.row(_).
```

```
    col(Y) :- $self.elem(_,Y).
    cols += 1 for $self.col(_).
}.
multiply_matrices(A,B) = new matrix {
    elem(I,K) += A.elem(I,J) * B.elem(J,K).
}.
result = multiply_matrices(identity_matrix, some_other_matrix).
```

This matrix representation is understandable by humans and even has some nice properties. For example, we are not required to use numerical integers as the identifies for the rows and columns. For example, we could have a row identified by the structured term noun["kitten"], and the code continues to work.

However, this code has some difficulties when trying to design execution strategies. For starters, we do not know the number of rows and columns of the identity matrix. In this case, the programmer forgot to inform us that identity matrices are square matrices. As such, if the query 'result. rows?' is attempted, we can not use any information available from the some_other_matrix on line 357 to compute this query. Now, as long as the Dyna program does not need to know the value of 'result.rows', then this is not a problem.

### 4.3 A Common Theme

All of these hard examples that can work share a common theme. There exists a subset of the program that is insufficient to run it. A naive representation-such as Datalog's materialize everything or Prolog lazy bindings to variables-is insufficient when trying to "greedily" evaluate the program, one subset at a time. We need to be able to represent a subset of the program, even when it cannot be evaluated. Once there is enough information, there needs to be some approach to apply the relevant strategies to run the program.

We will develop the necessary techniques in chapters 5 and 6 .

## Chapter 5

## Relational Expressions for Logic programming

In this chapter, I introduce our relational expression notation, which we shorten to $\mathbf{R}$-exprs. This chapter is one of the central contributions of this dissertation and was originally introduced in Francis-Landau et al. [70]. R-exprs is our internal representation, and is used to represent and execute Dyna programs. R-exprs represent both functions-much like how bytecode is the internal representation in procedural programming languages-and bag relations-much like a database table. In essence, we will use $\mathbf{R}$-exprs to represent both code and data (e.g. section §5.1.1).

As we will see throughout this dissertation, the homogeneity of R-exprs allows the Dyna runtime to be flexible, as yet-to-be-evaluated "delayed" code is stored alongside data. This turns out to be necessary to handle a number of tricky programs (chapter §4).

In the remainder of this chapter, I will define the semantic interpretation of R-exprs. In chapter $\S 7$, I will show how Dyna programs can be converted into the R-expr representation. Then in chapters 6 and 8 , I will discuss how $\mathbf{R}$-exprs are used to execute programs using term rewriting. In chapters 10 to 12 , I will address some of the challenges of using a term rewriting system to implement a programming language.

### 5.1 Representing Programs Using Bags

Before discussing our bag-relational algebra in the next section, let us review bag algebras and look at a few high-level examples.

A bag is a generalization of a set with elements in the bag being counted one or more times. Bags are written using $\rceil \cdots \int$. We refer to the number of times an element appears in a bag as multiplicity, and we will denote it using an at-sign @. As an example, a bag containing four, two-element tuples, where only two of which are distinct, looks like:

$$
\begin{equation*}
\left\{\langle 1,2\rangle @ 1,\langle 3,4\rangle @ 3 \int\right. \tag{5.1}
\end{equation*}
$$

where $\langle 1,2\rangle$ is contained once, and $\langle 3,4\rangle$ is contained three times.

### 5.1.1 Bags of Named Tuples

When representing programs, we often have dozens of variables in an expression. Hence, the above notation with values in a tuple marked by their index becomes cumbersome. Instead, from this point forward, we will use bags over named-tuples with bindings to variables. For example:

$$
\begin{equation*}
\eta\langle X=1, Y=2\rangle @ 1,\langle X=3, Y=4\rangle @ 3 \int \tag{5.2}
\end{equation*}
$$

is the same as the expression before, but we have now denoted the first variable as $X$ and the second variable as $Y$.

For our purposes, we say that bags represent relations over all variables, not just the variables explicitly mentioned. Any variable that is not explicitly mentioned is considered to as taking on "any/all" values:

$$
\begin{equation*}
\left.\left.\mathcal{Z} \mathrm{X}=1\rangle @ 1 \int \equiv \mathcal{Z}=1, Y\right\rangle @ 1: Y \in \mathcal{G}\right\} \tag{5.3}
\end{equation*}
$$

At first, this might appear a bit awkward, however, this is useful when we are combining multiple bags together using intersection $冈$ and union $\uplus$. For example, consider the intersection between two bags with different variables:

$$
\begin{equation*}
\text { l } \left.\langle x=1\rangle @ 2 \int 风\right\urcorner\langle Y=4\rangle @ 3 \int \equiv \eta\langle x=1, Y=4\rangle @ 6 \int \tag{5.4}
\end{equation*}
$$

Observe, with bag intersection 冈, the resulting bag is both the intersection of the bags and the product of the multiplicities. Note, we do not have a min multiplicity $\mathbf{R}$-expr. ${ }^{68}$ The reason is that we want to count the number of ways a tuple is contained in the bag. The counts of the number of times a tuple is in the bag is used by the aggregators (section §5.2.2.10).

In example 5.4, the variables were different, so the intersection only appears as a tuple that contains the assignments to both $X$ and $Y$. However, the power of bag intersection is that it checks the consistency between tuples contained in the bag. For example, in example 5.5, we have that the variable $x$ appears in both bags that we are intersecting:

$$
\begin{equation*}
\mathcal{Z}\langle x=1\rangle @ 7,\langle x=3\rangle @ 5 \int \propto \lambda\langle x=3, Y=11\rangle @ 2 \int \equiv \eta\langle x=3, Y=11\rangle @ 10 \int \tag{5.5}
\end{equation*}
$$

The assignment $\langle x=1\rangle$ is incompatible with the second bag, hence it is eliminated by the bag intersection.

Bag union $\uplus$ behaves similarly to bag intersection. Tuples that are compatible with each other have their multiplicities added together. Incompatible tuples behave as independent elements, appearing as their own elements in the resulting bag.

$$
\begin{array}{r}
\tau\langle X=1\rangle @ 1,\langle X=2\rangle @ 2, \int \uplus \eta\langle X=1\rangle @ 3,\langle Y=2\rangle @ 9 \int \\
\equiv \eta\langle X=1\rangle @ 4,\langle X=2\rangle @ 2,\langle Y=2\rangle @ 9 \int \tag{5.6}
\end{array}
$$

### 5.1.2 Representations of Constraints in Bags

To make bags useful for representing computation, we extend our bag notation with a "bag-builder" notation. This bag-builder notation is conceptually similar to set-builder ${ }^{69}$ notation where a set is defined using a boolean predicate. However,

[^44]bags associate a positive multiplicity with each element in the bag instead of the boolean true. This means that our bag-builder notation will return a nonnegative integer which represents the number of times in which a named-tuple is contained in the bag.

As an example, let us define the bag that contains all values assigned to the variable $x$ that are less than the number 5 :

Here, $c$ is the variable that corresponds with the multiplicity of the tuple $\langle x\rangle$ for some value of $x$. The multiplicity is computed using the expression $\begin{cases}1 & \text { if } x<5 \\ 0 & \text { otherwise }\end{cases}$ which returns 1 if it is contained in the bag, and 0 otherwise.

As in section §5.1.1, we can intersect bags to create composite expressions. For example, we can represent bag of integers between $[0,5)$ by intersecting the constraints for $x \in \mathbb{Z}, x \geq 0, x<5$ :

$$
\begin{array}{r}
\eta\langle x\rangle @ c: c=\left\{\begin{array}{ll}
1 & \text { if } x \in \mathbb{Z} \\
0 & \text { otherwise }
\end{array}\right\} \\
冈\left\langle\langle x\rangle @ c: c=\left\{\begin{array}{ll}
1 & \text { if } x \geq 0 \\
0 & \text { otherwise }
\end{array}\right\}\right.  \tag{5.8}\\
冈\left\langle\langle x\rangle @ c: c= \begin{cases}1 & \text { if } x<5 \\
0 & \text { otherwise }\end{cases} \right.
\end{array}
$$

### 5.1.3 A First Step Towards Computation with Bags

Equation (5.8) above has 3 different constraints represented with bag-builder notation. This expression is perfectly acceptable and is something that we are capable of handling and representing in our Dyna implementation. We may even return an expression like this to the user via the REPL in some cases.

However, our goal is to use expressions like example 5.8 to perform computation. This means that we need some way to manipulate expressions like example 5.8, and our manipulations should correspond with computation. The way we accomplish this is to rewrite an expression into a "simpler", semantically equivalent expression. I
will formally define our rewrites in chapter $\S 6$, and what simpler means in chapter §15.

In the case of eq. (5.8), it can be rewritten as a bag which enumerates the possible values of $X$ :

Here, we have rewritten from example 5.8 into example 5.9. In this example, and throughout this dissertation, rewrites are denoted using the right arrow $\rightarrow$. I claim that the rewritten bag in example 5.9 is simpler than example 5.8 , as we can directly read the values $0,1,2,3,4$ (the integers between $[0,5)$ ) from the bag. ${ }^{70}$ For now, an intuitive definition of simpler is that a simpler $\mathbf{R}$-expr requires less computation to get the "final" answer. For example, the value -1 is simpler than $e^{i \pi}$, which requires evaluating the exponential to determine the value of $e^{i \pi}$. I will give a formal definition of simpler in section §15.3.

### 5.1.4 A More Convenient Notation

The notation that I have used so far is admittedly very verbose and tedious to write. As such, we have designed a much more concise notation for $\mathbf{R}$-exprs.

First, observe that all bag expressions are over all variables. Hence, naming variables in the $\langle\cdots\rangle$ tuple is redundant. Second, we will say that all R-exprs "return" a multiplicity. Hence, it is unnecessary to write the if-case expression as in examples 5.7 and 5.8. Third, we are going to represent operations using term names instead of mathematical symbols. This means that an expression like $x \in \mathbb{Z}$ is instead written as int $(X)$. Additionally, we will write $x \geq 0$ as lessthaneq $(0, X)$ and $X<5$ as lessthan $(X, 5)$. Variables, such as $X$, follow the same convention as

[^45]Dyna code, they start with a capital letter and are colored Green in this dissertation. Finally, for intersection and union, we will write * and + in monospaced typewriter font instead of $\otimes$ and $\uplus$.

All together, this will allow us to write example 5.8 as the R-expr:

```
int(X)*lessthaneq(0,X)*lessthan(X,5)
```


### 5.2 Semantics of R-exprs

As noted in section §5.1.3, when rewriting an R-expr to perform "computation", we are rewriting the $\mathbf{R}$-expr into another simpler $\mathbf{R}$-expr which is semantically equivalent. As such, we need to define semantic equivalence for $\mathbf{R}$-exprs.

### 5.2.1 Ground Values

Let us first start by defining the set of ground terms $\mathcal{G}$ built from a ranked set of functors $\mathcal{F}$ and primitive values such as integers, floats, and strings. For example, $\mathcal{G}$ includes 123, 3.14, "hello world" and structured terms like foo[123, "hello", $\operatorname{bar}[4,5,6]$ (as introduced in section §2.1.1).

We further define $\mathcal{M}=\mathbb{N} \cup\{\infty\}$ as the set of multiplicities. Therefore, a bag containing an $n$-arity tuple of ground terms $\left(\mathcal{G}^{n}\right)$, can be represented as a function that map from $\mathcal{G}^{n}$ to the multiplicity $M \in \mathcal{M}$ of that tuple contained in the bag: $\mathcal{G}^{n} \mapsto \mathcal{M}$.

That said, we do not use the integer position in an $n$-arity tuple as it is inconvenient. Instead, we use a map from variable names to ground terms. Let $\tilde{\mathcal{V}}$ be an infinite set of unique variable names. A environment $E(\cdot)$ is, a function that maps from a finite set of variables to ground terms. More specifically, for any $\mathcal{U} \subset \tilde{\mathcal{V}}$, the $\mathcal{U}$-tuples are maps $E(\cdot): \mathcal{U} \mapsto \mathcal{G}$. Using this representation, a bag over $\mathcal{U}$-tuples can be represented by a function $F(\cdot)$ where $F(\cdot):(\mathcal{U} \mapsto \mathcal{G}) \mapsto \mathcal{M} .^{71}$

[^46]For convenience, we define the set of value types as $\mathcal{V}=\tilde{\mathcal{V}} \cup \mathcal{G}$. We extend the $E(\cdot)$ function to $\mathcal{V}$ by mapping any ground value $g \in \mathcal{G}$ to itself. With this, we can think of the $E(\cdot)$ function as a kind of "get value" function. ${ }^{72}$

1. $\llbracket \mathrm{U}=\mathrm{V} \rrbracket_{E}=$ if $E(\mathrm{U})=E(\mathrm{v})$ then 1 else $0, \quad$ where $\mathrm{U}, \mathrm{v} \in \mathcal{V}$
2. $\llbracket \mathrm{T}=\mathrm{f}\left(\mathrm{X}_{1}, \cdots, \mathrm{X}_{n}\right) \rrbracket_{E}=$ if $E(\mathrm{~T})=f\left[E\left(\mathrm{X}_{1}\right), \cdots, E\left(\mathrm{X}_{n}\right)\right]$ then 1 else 0 , where $\mathrm{T}, \mathrm{X}_{1}, \cdots, \mathrm{X}_{n} \in \mathcal{V}$ and $\mathrm{f} \in \mathcal{F}$
3. $\llbracket \mathrm{plus}(\mathrm{I}, \mathrm{J}, \mathrm{K}) \rrbracket_{E}=$ if $E(\mathrm{I})+E(\mathrm{~J})=E(\mathrm{~K})$ then 1 else $0, \quad$ where $\mathrm{I}, \mathrm{J}, \mathrm{K} \in \mathcal{V}$
4. $\llbracket \mathrm{R}+\mathrm{S} \rrbracket_{E}=\llbracket \mathrm{R} \rrbracket_{E}+\llbracket \mathrm{S} \rrbracket_{E}, \quad$ where $\mathrm{R}, \mathrm{s} \in \mathcal{R}$
5. $\llbracket \mathrm{R} * \mathrm{~S} \rrbracket_{E}=\llbracket \mathrm{R} \rrbracket_{E} \cdot \llbracket \mathrm{~S} \rrbracket_{E}, \quad$ where $\mathrm{R}, \mathrm{s} \in \mathcal{R}$
6. $\llbracket \mathrm{M} \rrbracket_{E}=\mathrm{M}, \quad$ where $\mathrm{M} \in \mathcal{M}$
7. $\llbracket \mathrm{if}(\mathrm{Q}, \mathrm{R}, \mathrm{S}) \rrbracket_{E}=$ if $\llbracket \mathrm{Q} \rrbracket_{E}>0$ then $\llbracket \mathrm{R} \rrbracket_{E}$ else $\llbracket \mathrm{S} \rrbracket_{E}, \quad$ where $\mathrm{Q}, \mathrm{R}, \mathrm{S} \in \mathcal{R}$
8. $\llbracket \operatorname{proj}(\mathrm{X}, \mathrm{R}) \rrbracket_{E}=\sum_{x \in \mathcal{G}} \llbracket \mathrm{R} \rrbracket_{E}[\mathrm{X}=x], \quad$ where $\mathrm{X} \in \mathcal{V}, \mathrm{R} \in \mathcal{R}$
9. $\llbracket \mathrm{A}=\operatorname{sum}(\mathrm{X}, \mathrm{R}) \rrbracket_{E}=$ if $E(\mathrm{~A})=\sum_{x \in \mathcal{G}}\left(x * \llbracket \mathrm{R} \rrbracket_{E[\mathrm{X}=x]}\right)$ then 1 else 0 , where $A, x \in \mathcal{V}, R \in \mathcal{R}$.
10. $\llbracket \mathrm{f}\left(\mathrm{T}_{1}, \ldots, \mathrm{~T}_{n}\right) \rrbracket_{E}=\llbracket \mathrm{R}_{\mathrm{f}}\left\{\mathrm{X}_{1} \mapsto \mathrm{~T}_{1}, \ldots, \mathrm{X}_{n} \mapsto \mathrm{~T}_{n}\right\} \rrbracket_{E}$, where $\mathrm{T}_{1}, \ldots, \mathrm{~T}_{n} \in \mathcal{V}$

Figure 5-1. Semantic definitions of all R-expr kinds. The semantics $\llbracket \mathrm{R} \rrbracket_{E}$ of an R-expr $R$ is defined as the multiplicity $(\mathcal{M}=\mathbb{N} \cup\{\infty\})$ of $E$ in the bag defined by R. $E$ is a tuple of named variables and their ground assignments (section §5.1.1). Additional details about the R-exprs' definitions are provided in section §5.2.2.

### 5.2.2 Inductive Definition of R-exprs Semantics

In the remainder of this section, I will give the definitions of $\mathbf{R}$-exprs and their semantics. $\mathbf{R}$-exprs are terms that are defined recursively in terms of sub-R-exprs

[^47]and value types $\mathcal{V}$. As we will see shortly, the semantics of $\mathbf{R}$-exprs is defined as a system of equations where the semantic interpretation each individual $\mathbf{R}$-expr is defined as a multiplicity given environment $E(\cdot)$.

To make this more formal, let $\mathcal{R}$ denote the set of all $\mathbf{R}$-exprs.
Each $\mathbf{R}$-expr $R \in \mathcal{R}$ has a finite set of free variables denoted by $\operatorname{vars}(R)$. A free variable is a variable in an $\mathbf{R}$-expr that is not bound by another $\mathbf{R}$-expr. Note, $\operatorname{vars}(R) \subseteq \tilde{\mathcal{V}}$ as $\operatorname{vars}(R)$ returns a set of named variables, not value types.

Next, for each environment $E$, the denotation function $\llbracket \cdot \rrbracket_{E}: \mathcal{R} \mapsto \mathcal{M}$ interprets an $\mathbf{R}$-expr in the environment given by the named tuple $E$. It defines a multiplicity $\llbracket \mathrm{R} \rrbracket_{E}$ for any $\mathbf{R}$-expr R whose $\operatorname{vars}(\mathrm{R}) \subseteq$ domain $(E) .{ }^{73}$

We will now define and explain $\llbracket \mathrm{R} \rrbracket_{E}$ and vars( R ) for the different kinds of $\mathbf{R}$-exprs in our system, which are also shown in figure 5-1.

### 5.2.2.1 Equality Constraints

First, we define equality constraints between two value types $U, V \in \mathcal{V}$. An equality constraint is true in an environment when the ground value assigned to $U$ and $V$ is equal. This is represented using multiplicity 1 , which indicates that the current environment is consistent with this $\mathbf{R}$-expr and is therefore contained in the bag represented by this $\mathbf{R}$-expr. False is represented with multiplicity $0 .{ }^{74}$

1. $\llbracket \mathrm{U}=\mathrm{V} \rrbracket_{E}=$ if $E(\mathrm{U})=E(\mathrm{v})$ then 1 else $0, \quad$ where $\mathrm{U}, \mathrm{v} \in \mathcal{V}$

Notice that we did not write $\llbracket \cup \rrbracket_{E}=\llbracket \mathrm{V} \rrbracket_{E}$ but $E(\mathrm{U})=E(\mathrm{~V})$-our denotation function $\llbracket \cdot \rrbracket_{E}$ maps $\mathbf{R}$-exprs to multiplicities, and $E(\cdot)$ maps variables and ground terms in $\mathcal{V}$ to ground terms in $\mathcal{G}$.

### 5.2.2.2 Structured Term Equality Constraints

Next, we have equality constraints with a structured term with $T, X_{1}, \cdots, X_{n} \in \mathcal{V}$ and $f \in \mathcal{F}$ being the name of the structured term (previously introduced in section §2.1.1).

[^48]2. $\llbracket \mathrm{T}=\mathrm{f}\left[\mathrm{X}_{1}, \cdots, \mathrm{X}_{n}\right] \rrbracket_{E}=$ if $E(\mathrm{~T})=\mathrm{f}\left[E\left(\mathrm{X}_{1}\right), \cdots, E\left(\mathrm{X}_{n}\right)\right]$ then 1 else 0 ,
where $T, X_{1}, \cdots, X_{n} \in \mathcal{V}$ and $f \in \mathcal{F}$

### 5.2.2.3 Builtin R-exprs Constraints

We also have a number of built-in R-expr kinds. These R-exprs correspond to "lowlevel" operations on primitive ground terms, such as addition between numbers or concatenating strings. Built-ins are defined as 1 when the assignment to their arguments is consistent with their definition. For example, the built-in addition, which henceforth will be referred to as plus for clarity, is defined as follows:
3. $\llbracket \mathrm{plus}(\mathrm{I}, \mathrm{J}, \mathrm{K}) \rrbracket_{E}=$ if $E(\mathrm{I})+E(\mathrm{~J})=E(\mathrm{~K})$ then 1 else $0, \quad$ where $\mathrm{I}, \mathrm{J}, \mathrm{K} \in \mathcal{V}$

Note, in this definition, the + sign corresponds with addition between the values of $E(\mathrm{I})$ and $E(\mathrm{~J})$, which are both in the $\mathcal{G}$ that is the Herbrand universe of values of the Dyna language.

### 5.2.2.4 Constraints

The R-exprs we have so far are constraints. A constraint is any R-expr that has a multiplicity of 0 or 1 in any environment. It is useful to track constraints as they can be duplicated as needed or removed when they are redundant. ${ }^{75}$

### 5.2.2.5 Disjunctions

Disjunctions are the union of two or more R-exprs, and are denoted using +.

$$
\text { 4. } \llbracket \mathrm{R}+\mathrm{S} \rrbracket_{E}=\llbracket \mathrm{R} \rrbracket_{E}+\llbracket \mathrm{S} \rrbracket_{E}, \quad \text { where } \mathrm{R}, \mathrm{~s} \in \mathcal{R}
$$

Disjuncts correspond with the $\uplus$ as we saw before in section §5.1.1. Disjuncts are usually not constraints as their multiplicity is often greater than $1 .{ }^{76}$ This time, the + sign on the right-hand side of this definition is the addition between the

[^49]multiplicities returned by $\llbracket \mathrm{R} \rrbracket_{E}$ and $\llbracket \mathrm{S} \rrbracket_{E}$. The typewriter font plus sign '+’ in the semantic brackets is only used to represent disjunctions, as done here.

The free variables in a disjunction are the union of the free variables in disjunctive sub-R-exprs: $\operatorname{vars}(R+S)=\operatorname{vars}(R) \cup \operatorname{vars}(S)$. This is in contrast to the earlier $\mathbf{R}$-expr kinds where the free variables are simply all of the variables that appeared in the R-expr.

### 5.2.2.6 Conjunctions

Conjunctions are the intersection of two or more R-exprs, and are denoted using *.
5. $\llbracket \mathrm{R} * \mathrm{~S} \rrbracket_{E}=\llbracket \mathrm{R} \rrbracket_{E} \times \llbracket \mathrm{S} \rrbracket_{E}, \quad$ where $\mathrm{R}, \mathrm{s} \in \mathcal{R}$

Conjuncts correspond with $冈$ from before. Conjunctions are often constraints themselves as if both conjuncts $R$ and $S$ are a constraint, then a conjunction is also a constraint. The typewriter font multiplication symbol ' $x$ ' is only used to represent conjunctions. In the written presentation in this dissertation, I will follow the standard order-of-operations with ' + ' and ' $x$ ', with ' $x$ ' binding more tightly than ' + '. Further note that we define all conjunctions with a 0 as 0 , even in the case of an $\infty$ multiplicity. Hence $0 \times \infty=0$.

Just like with disjunctions, the free variables are the union of the free variables in the conjunctive sub-R-exprs: $\operatorname{vars}(R * S)=\operatorname{vars}(R) \cup \operatorname{vars}(S)$.

### 5.2.2.7 Multiplicities

Multiplicities R-expr kinds are defined as the multiplicity they represent.
6. $\llbracket M \rrbracket_{E}=M, \quad$ where $M \in \mathcal{M}$

Multiplicities always return the same value $M$, and are not influenced by the environment $E$. For example, the multiplicity 1 represents the bag that contains all environments exactly once, and 0 represents an empty bag. E.g. $\llbracket 1 \rrbracket_{E}=1$ and $\llbracket 0 \rrbracket_{E}=0$.

### 5.2.2.8 Conditionals

We allow conditional or if-expressions in our relational algebra as follows:
7. $\llbracket \mathrm{if}(\mathrm{Q}, \mathrm{R}, \mathrm{S}) \rrbracket_{E}=$ if $\llbracket \mathrm{Q} \rrbracket_{E}>0$ then $\llbracket \mathrm{R} \rrbracket_{E}$ else $\llbracket \mathrm{S} \rrbracket_{E}, \quad$ where $\mathrm{Q}, \mathrm{R}, \mathrm{S} \in \mathcal{R}$

Notice that this construction can be used to implement anti-joins, set differences, and priority unions on sets. It does not support bag differences, which would involve subtracting multiplicities, but instead uses $Q$ to choose between two multiplicities (which can be 0). Furthermore, an if-expression is a constraint if $R$ and $S$ are both constraints. ${ }^{77}$

As with disjunctions and conjunctions, the free variables are the union of the free variables from the sub-R-exprs: $\operatorname{vars}(i f(Q, R, S))=\operatorname{vars}(Q) \cup \operatorname{vars}(R) \cup \operatorname{vars}(S)$.

### 5.2.2.9 Projection

Next, we define projection, which removes a named column (variable) from a bag relation, summing the multiplicities of tuples that have become equal. When translating a Dyna program into an R-expr (chapter §7), projections will be used to eliminate a rule's local variables.
8. $\llbracket \operatorname{proj}(\mathrm{X}, \mathrm{R}) \rrbracket_{E}=\sum_{x \in \mathcal{G}} \llbracket \mathrm{R} \rrbracket_{E[\mathrm{X}=x]}$, where $\mathrm{X} \in \mathcal{V}, \mathrm{R} \in \mathcal{R}$, and where $E[\mathrm{X}=x]$ means a version of $E$ that has been modified such that $E(\mathrm{x})=x,^{78} 79$ and $\sum_{x \in \mathcal{G}}$ is a summation of the multiplicities over all possible values in $\mathcal{G} .{ }^{80}$

In the case that the variable $x$ is not contained in $\operatorname{vars}(R)$, and the multiplicity of R is non-zero $\left(\llbracket \mathrm{R} \rrbracket_{E}>0\right)$, then the multiplicity of projection is defined as infinity $(\infty) .{ }^{81}$ The reason is that this is equivalent to summing a non-zero value an infinite number of times because $\mathcal{G}$ is an infinite set.

The free variables of projection are again based on the sub-R-expr's free variables, but this time, we are removing the variable $x$ from the set of free variables:

[^50]$\operatorname{vars}(\operatorname{proj}(X, R))=\operatorname{vars}(R)-\{X\}$.

### 5.2.2.10 Aggregation

Aggregation does the same as projection, but instead of adding multiplicities, it aggregates the $X$ value and computes a new value $Y$ as the result of aggregation. Thus, it removes column $X$ from the bag but introduces a new column $Y$.

Throughout this dissertation, I will use the 'sum' aggregator as a prototypical aggregator. However, there are other aggregators which have similar semantic denotations and rewrite rules.
9. $\llbracket \mathrm{A}=\operatorname{sum}(\mathrm{X}, \mathrm{R}) \rrbracket_{E}=$ if $E(\mathrm{~A})=\sum_{x \in \mathcal{G}}\left(x * \llbracket \mathrm{R} \rrbracket_{E[\mathrm{X}=x]}\right)$ then 1 else 0 ,
where $A, X \in \mathcal{V}, R \in \mathcal{R}$. The $*$ in the summand means that the summation includes $\llbracket \mathrm{R} \rrbracket_{E[\mathrm{X}=x]}$ copies of the value $x$ (possibly $\infty$ copies $^{82}$ ). If there are no summands (meaning $\llbracket \mathrm{R} \rrbracket_{E}=0$ for all $E$ ), then the result of $\sum \cdots$ is defined as null, which is a value outside of $\mathcal{G}$. Hence, it has the multiplicity of 0 . In other words, when R is 0 , we have $\llbracket A=\operatorname{sum}(X, 0) \rrbracket_{E}=0$. If there are non-numeric summands, or the sum is not well defined, ${ }^{83}$ then the result of the sum is defined to be an error value in $\mathcal{G}$, such as NaN .

Notice that the aggregator $\mathbf{R}$-expr is a constraint as its multiplicity is either 1 or 0 . Additionally, observe we do not define $\llbracket \operatorname{sum}(X, R) \rrbracket_{E}$, whose value would not be a multiplicity. Rather, the aggregator is written as $A=\operatorname{sum}(X, R)$, which denotes where the result of aggregation will be saved (variable $A$ in this case) and is considered a more readable notation for what could have been written as sum (A, $X, R$ ).

The free variables of aggregation project out the input variable $X$ and add the variable $A$ to the free variables: $\operatorname{vars}(A=\operatorname{sum}(X, R))=(\operatorname{vars}(R)-\{X\}) \cup\{A\}$

### 5.2.2.11 User-defined R-exprs

Finally, it is convenient to augment the built-in relations with user-defined relation kinds. Choose an identifier $f \in \mathcal{F}$ and choose some $\mathbf{R}$-expr $R_{f}$ with $\operatorname{vars}\left(R_{f}\right) \subseteq$

[^51]$\left\{x_{1}, \ldots, x_{n}\right\}$ (which are $n$ distinctly named variables) to serve as the definition or $\mathbf{R}$-expr expansion of $f$. Now define
10. $\llbracket \mathrm{f}\left(\mathrm{T}_{1}, \ldots, \mathrm{~T}_{n}\right) \rrbracket_{E}=\llbracket \mathrm{R}_{\mathrm{f}}\left\{\mathrm{X}_{1} \mapsto \mathrm{~T}_{1}, \ldots, \mathrm{X}_{n} \mapsto \mathrm{~T}_{n}\right\} \rrbracket_{E}$, where $\mathrm{T}_{1}, \ldots, \mathrm{~T}_{n} \in \mathcal{V}$

The $\{\mapsto\}$ notation denotes substitution for variables, where free variables of $R_{f}$ are renamed, and variables captured by other R-exprs, such as projection or aggregation, are first renamed to avoid accidental capture. ${ }^{84}$

User-defined $\mathbf{R}$-exprs make it possible to define $\mathbf{R}$-exprs which are circularly defined in terms of themselves. ${ }^{85}$ In fact, a Dyna program will normally do this. In this case, the definition of $\llbracket \cdot \rrbracket_{E}$ is not an inductive function. Rather, $\llbracket \cdot \rrbracket_{E}$ is interpreted as a variable in a system of equations, where we solve for values of the $\llbracket \cdot \rrbracket_{E}$ that satisfies these constraints. This is allowed by Dyna, as in section §2.5, I stated that the semantics of Dyna allow us to find any consistent assignment. In practice, there can be multiple possible assignments, and there is no guarantee that the system will find a particular assignment. ${ }^{86}$

### 5.3 Example R-exprs

Given that we have now defined the basic $\mathbf{R}$-expr kinds, let us look at some simple $\mathbf{R}$-exprs to see how they can represent bags and simple Dyna programs. This is not intended as a complete introduction to how $\mathbf{R}$-exprs can represent Dyna programs or bag relations, as we will see more in chapter $\S 7$.

### 5.3.1 Finite Materialized Relation

As a first example, we can express a finite bag relation shown in figure 5-2a as an $\mathbf{R}$-expr using conjunctions, disjunctions, and equality assignments as in figure 5-2b.

[^52]\[

\left\{$$
\begin{array}{l}
\langle X=1, Y=1\rangle @ 1 \\
\langle X=2, Y=6\rangle @ 1 \\
\langle X=2, Y=7\rangle @ 2 \\
\langle X=5, Y=7\rangle @ 1
\end{array}
$$\right\} \quad $$
\begin{aligned}
& (X=1) *(Y=2)+ \\
& (X=2) *(Y=6)+ \\
& (X=2) *(Y=7) * 2+ \\
& (X=5) *(Y=7)
\end{aligned}
$$
\]

(a) Bag
(b) R-expr

Figure 5-2. Bag of ground assignments to the variables $X$ and $Y$. The ( $X=1$ ) R-exprs are each individual $\mathbf{R}$-expr equality constraints (as previously defined, section §5.2.2.1). The equality constraints are combined into conjunctions using ' *' (section §5.2.2.6). Finally, the conjunctions are combined into a big disjunction using '+' (section §5.2.2.5).

Each individual tuple of the bag translates into a product (*) of several (Variable $=$ value) expressions. Encoding the multiple tuples contained in the bag is done by simply adding (+) the $\mathbf{R}$-exprs for different tuples together using a disjunction.
$\mathbf{R}$-exprs of this form can be used as the basis for user-defined $\mathbf{R}$-exprs when representing external data as an $\mathbf{R}$-expr (section §5.2.2.11). Finite materialized $\mathbf{R}$-exprs are also desirable representations to return to the user when making a query, as they can be easily understood by reading the relevant values out of the R-expr (section §2.3).

### 5.3.2 Bag with Constraints

$$
\begin{gathered}
\left\{\langle X\rangle @ c: c= \begin{cases}1 & \text { if } \exists y \in \mathbb{Z} \text { s.t. } 2 * y=\mathrm{x} \\
0 & \text { otherwise }\end{cases} \right. \\
\text { (a) Bag } \\
\operatorname{if(\operatorname {proj}(Y,\operatorname {int}(Y)*\operatorname {times}(2,Y,X)),1,0)} \\
\text { (b) } \mathbf{R}-\operatorname{expr}
\end{gathered}
$$

Figure 5-3. Bag of even integers.

The bag expression in figure 5-3a checks whether there exists an integer $y$ such that $2 * y=\mathrm{X}$. In figure 5 -3b we introduce a new variable Y using $\operatorname{proj}(\mathrm{Y}, \cdot)$. We ensure that the multiplicity of the $\mathbf{R}$-expr is either 0 or 1 by using the if-expression to match the $\exists$ in the bag definition. Otherwise, the projection could have a multiplicity greater than 1 , though this does not actually happen in this case.

### 5.3.3 Simple Dyna Rule

As a preview for chapters 6 and 7, figure 5-4 shows a single-rule Dyna program and the equivalent $\mathbf{R}$-expr representation.

$$
\begin{aligned}
& { }_{358} \mid a(X)+=(X+7) * X . \quad a(X, \text { Result }) \xrightarrow{1} \text { (Result=sum(Inp, } \\
& \text { proj(Tmp,plus(X,7,Tmp)*times(Tmp,X,Inp)))) }
\end{aligned}
$$

(a) Dyna

## (b) R -expr

Figure 5-4. Simple Dyna program written as an R-expr.

Dyna user-defined terms, which are comprised of rules such as line 358, define their own user-defined R-expr kind (section §5.2.2.11) and rewrite rule (shown in figure 5-4b).

The user-defined $\mathbf{R}$-expr has the variable $x$ for the arguments to the expression, as well as a Result variable for the returned value of the expression. A rewrite rule is defined to rewrite the user-defined $\mathbf{R}$-expr to its corresponding definition. Rewrite rules are represented as an arrow $(\rightarrow)$, with important rewrite rules having a reference number over the arrow that will appear in the text and are clickable (e.g. rewrite rule 1). ${ }^{87}$

Calls to built-ins, written as ' + ' and ' $*$ ' in the Dyna source on line 358 are converted into the R-exprs plus $(\cdot, \cdot, \cdot)$ and times $(\cdot, \cdot, \cdot)$. We reserve the ' $*$ ' and ' + ' symbols in the $\mathbf{R}$-expr representation for conjunction and disjunction, which are used much more frequently.

All local variables and intermediate variables are projected out in user-defined $\mathbf{R}$-exprs. All free variables must appear in the user-defined R-exprs. Intermediate

[^53]variables are introduced as needed when translating from Dyna into R-exprs, as exemplified by the variable Tmp in this example.

### 5.4 Conclusion of R-expr Semantics

The R-expr kinds defined in this chapter form the basis for our internal representation. In the next three chapters, I will define semantic preserving rewrite rules for R-exprs (chapter §6), as well as show in further detail how Dyna is translated into R-exprs (chapter §7) and how we can implement a simple term rewriting system using R-exprs (chapter §8).

## Chapter 6

## Rewrites Rules for R-exprs

The previous chapter gave the denotational semantics for R-exprs. I will now provide the operational semantics. The basic idea is that we can use rewrite rules to simplify, an R-expr until it is either a finite materialized relation-a list of tuples-or has some other convenient form. All of our rewrite rules are semantics-preserving by construction, so our rewrite system is sound, although it is not and cannot be complete, as chapter $\S 15$ will discuss. Our rewrites can be applied to any sub-R-expr contained within a larger R-expr. Chapter $\S 8$ will discuss the procedure we use to apply rewrites. Our rewrite system is not confluence, meaning that the order in which rewrites are applied can result in different semantically equivalent $\mathbf{R}$-exprs. However, our rewrite system is normalizing, in that it eventually reach a state where no more directional rewrites that can be applied.

### 6.1 Equality Constraints and Multiplicity

We start with rewrites for basic equality constraints.

$$
\begin{aligned}
& (\mathrm{X}=\mathrm{X}) \xrightarrow{2} 1 \quad \triangleright \text { trivially true when the variables/values are the same } \\
& (X=Y) \xrightarrow{3} 0 \quad \text { if } X, Y \in \mathcal{G} \text { and } X \neq Y \quad \triangleright \text { values not equal } \\
& (G=X) \xrightarrow{4}(X=G) \quad \text { if }(G \in \mathcal{G} \text { and } X \in \tilde{\mathcal{V}}) \text { or }(G \prec X) \quad \triangleright \text { canonical order } \\
& (X=Y) * R \xrightarrow{5}(X=Y) * R\{X \mapsto Y\} \quad \text { if } X \in \tilde{\mathcal{V}} \quad \text { - equality propagation }
\end{aligned}
$$

First, rewrite rule 2 checks for equivalence between its two arguments. When a

R-expr constraint has been successfully and completely "checked", it is rewritten as a 1, as in this case. Conversely, rewrite rule 3 handles the case where $X$ and $Y$ are not equal. It requires that both $X$ and $Y$ be known ground values. When both values are known, we can use the equality operator defined on ground values $\mathcal{G}$ to check if $X$ and $Y$ are equal. Constraints whose check fails are rewritten as 0 , which indicates that the current assignment of variables in the environment $E(\cdot)$ is inconsistent and, therefore, not contained in the bag represented by the R-expr.

For equality, we prefer to have a canonical ordering where we place any ground value as the second argument. If we have two variables, rewrite rule 4 reorders the equality constraint according to an arbitrarily chosen $\prec$ ordering on variables. ${ }^{88}$

The most common operation with equality constraints is performed by rewrite rule 5, which propagates a variable binding to other conjoined $\mathbf{R}$-exprs. The $Y$ is likely a ground value, in which case propagating the value of $Y$ in place of $X$ enables other rewrites, as we shall see.

### 6.1.1 Structured Term Equality Rewrites

Structured terms are similar to equality constraints; however, they involve multiple variables and a named term $f, g \in \mathcal{F}$.

$$
\begin{aligned}
& \left(\mathrm{X}=\mathrm{f}\left[\mathrm{Y}_{1}, \ldots, \mathrm{X}, \ldots, \mathrm{Y}_{n}\right]\right) \xrightarrow{6} 0 \quad \text { Doccurs check, not true for any ground value of } \mathrm{X} \\
& \left(X=f\left[Y_{1}, \ldots, Y_{n}\right]\right) *\left(X=f\left[Z_{1}, \ldots, Z_{n}\right]\right) \xrightarrow{7}\left(X=f\left[Y_{1}, \ldots, Y_{n}\right]\right) *\left(Y_{1}=Z_{1}\right) * \cdots *\left(Y_{n}=Z_{n}\right) \\
& \left(X=f\left[Y_{1}, \ldots, Y_{n}\right]\right) *\left(X=g\left[Z_{1}, \ldots, Z_{m}\right]\right) \xrightarrow{8} 0 \text { if }(f \neq g \text { or } n \neq m) \\
& \left(G=f\left[Y_{1}, \ldots, Y_{n}\right]\right) \xrightarrow{9}\left(Y_{1}=Z_{1}\right) * \cdots *\left(Y_{n}=Z_{n}\right) \text { if } G \in \mathcal{G} \text { and } G=f\left[Z_{1}, \ldots, Z_{n}\right] \\
& \left(G=f\left[Y_{1}, \ldots, Y_{n}\right]\right) \xrightarrow{10} 0 \text { if } G \in \mathcal{G} \text { and } G \neq f\left[Z_{1}, \ldots, Z_{n}\right]
\end{aligned}
$$

Rewrite rule 6 implements the occurs check, looking for expressions of the form $X=f[Y, Z, X]$. In Dyna, we do not allow for cyclic data structures (which is allowed in some Prolog implementations). Hence, there is no value $x \in \mathcal{G}$, so this expression is unsatisfiable, thus can be rewritten as 0 , without having to identify the value of $x$.

Rewrite rules 7 and 8 perform unification between different structured term rewrites. If two incompatible structured terms are unified together (in this case,

[^54]$\mathrm{f}, \mathrm{g} \in \mathcal{F}$ and $\mathrm{f} \neq \mathrm{g}$ ), then this can be rewritten as 0 as seen in rewrite rule $8 .{ }^{89}$ Similarly, when the functor names on two structured term R-exprs are equivalent to each other, all of the arguments of the terms are unified together, as in rewrite rule 7 . Observe that the result of rewrite rule 7 still includes at least one structured unification $\mathbf{R}$-expr. The reason is that we still need the constraint on the variable $x$. Without this constraint, the bag semantics of the $\mathbf{R}$-expr would change, and there are likely other conjunctive sub-R-exprs in the containing $\mathbf{R}$-expr that depend on the value of $x$.

Rewrite rules 9 and 10 replicate rewrite rules 7 and 8 but are for when the structured term is a ground, fully known value. For example, $f[1,2,3]$ is ground and can be entirely represented by the value variable $G$, and can be handled by rewrite rule 9 or 10 . Whereas, $f[1,2, x]$ is not ground, as it contains the variable $x$, and must be handled with rewrite rule 7 or 8 .

It can be seen that rewrite rules $2-10$ is equivalent to the well-known unification algorithm commonly seen in logic programming [102].

### 6.1.2 Multiplicity Rewrites

Given the rewrite rules that we have so far, we may end up creating an arithmetic expression like $0 * 0+1 * 0+1 * 1+1 * 1+0 * 1$, which can be reduced to the multiplicity 2. This is done using rewrites that implement basic arithmetic on multiplicities $\mathcal{M}$ :

```
M+N \xrightarrow{}{11}L\mathrm{ if }M,N\in\mathcal{M}\mathrm{ and }M+N=L
M*N \xrightarrow{}{12}L\mathrm{ if }M,N\in\mathcal{M}\mathrm{ and }M*N=L
```


### 6.2 Joining Relations

So far, we have the rewrites on the equality $\mathbf{R}$-expr kinds. To represent a bag with multiple variables, we need to combine multiple equality constraints together, as in section §5.3.1. Furthermore, we need to be able to rearrange the $\mathbf{R}$-expr to bring relevant sub-R-exprs together so that we can carry out the aforementioned rewrites.

[^55]To carry out such rewrites, we use the fact that multiplicities form a commutative semiring under + and *. Since any $\mathbf{R - e x p r}$ evaluates to a multiplicity, these rewrites can be used to rearrange unions and joins of $\mathbf{R}$-exprs $Q, R, S \in \mathcal{R}$ :
$1 * R \xrightarrow{13} R$
$0 * R \xrightarrow{14} 0$
$0+R \xrightarrow{15} R$
$\infty * R \xrightarrow{16} \infty$ if $R \in \mathcal{M}$ and $R>0$
$\infty+\mathrm{R} \xrightarrow{17} \infty$
$\mathrm{R}+\mathrm{S} \stackrel{18}{\longleftrightarrow} \mathrm{~S}+\mathrm{R} ; \quad \mathrm{R} * \mathrm{~S} \stackrel{19}{\longleftrightarrow} \mathrm{~S} * \mathrm{R} \quad$ - commutativity
$\mathrm{Q}+(\mathrm{R}+\mathrm{S}) \stackrel{20}{\longleftrightarrow}(\mathrm{Q}+\mathrm{R})+\mathrm{S} ; \quad \mathrm{Q} *(\mathrm{R} * \mathrm{~S}) \stackrel{21}{\longleftrightarrow}(\mathrm{Q} * \mathrm{R}) * \mathrm{~S} \quad \triangleright$ associativity
$\mathrm{Q} *(\mathrm{R}+\mathrm{S}) \stackrel{22}{\longleftrightarrow} \mathrm{Q} * \mathrm{R}+\mathrm{Q} * \mathrm{~S} \quad \triangleright$ distributivity
$R * M \stackrel{23}{\longleftrightarrow} R+(R * N)$ if $M, N \in \mathcal{M}$ and $(1+N=M) \quad$ Dimplicitly does $M=1+N$
$R \stackrel{24}{\longleftrightarrow} R * R$ if $R$ is a constraint $\quad \triangleright$ as defined in section $\S 5$.2.2.4
$\mathrm{R} \stackrel{25}{\longleftrightarrow} \mathrm{R} * \mathrm{~S}$ if S is a constraint and $\forall_{E} \llbracket \mathrm{R} \rrbracket_{E}=\llbracket \mathrm{R} * \mathrm{~S} \rrbracket_{E} \quad \triangleright S$ is redundant with $R$

Note, rewrite rules 18 to 25 are bidirectional rewrites, which may appear contradictory to statement at the start of this chapter that our rewrite system will be normalizing. As will be discussed in chapter $\S 8$, our rewrite system is normalizing with respect to the directional rewrite rules, and bidirectional rewrites, such as rewrite rules 18 to 25 , will be handled specially and will only be used to rearrange the $\mathbf{R}$-expr to enable directional rewrites.

### 6.2.1 Simple Example

Let us try applying the rewrites we have to the finite materialized bag $\mathbf{R}$-expr, like the $\mathbf{R}$-expr from section §5.3.1. Let us compute the intersection between two finite relations R-exprs.

$$
\binom{(X=2) *(Y=6) * 3+}{(Y=7)} *\left(\begin{array}{l}
(X=1) *(Y=2)+  \tag{6.1}\\
(X=2) *(Y=6)+ \\
(X=2) *(Y=7) * 2+ \\
(X=5) *(Y=7)
\end{array}\right)
$$

We first expand out the first disjunct using the distributive rewrite (rewrite rule 22):

$$
\rightarrow \quad(X=2) *(Y=6) * 3 *\left(\begin{array}{l}
(X=1) *(Y=2)+  \tag{6.2}\\
(X=2) *(Y=6)+ \\
(X=2) *(Y=7) * 2+ \\
(X=5) *(Y=7)
\end{array}\right)+(Y=7) *\left(\begin{array}{l}
(X=1) *(Y=2)+ \\
(X=2) *(Y=6)+ \\
(X=2) *(Y=7) * 2+ \\
(X=5) *(Y=7)
\end{array}\right)
$$

The assignments to $X$ and $Y$ can be propagated through the conjunctive $\mathbf{R}$-expr using rewrite rule 5:

$$
\rightarrow \quad(X=2) *(Y=6) * 3 *\left(\begin{array}{l}
(2=1) *(6=2)+  \tag{6.3}\\
(2=2) *(6=6)+ \\
(2=2) *(6=7) * 2+ \\
(2=5) *(6=7)
\end{array}\right)+(Y=7) *\left(\begin{array}{l}
(X=1) *(7=2)+ \\
(X=2) *(7=6)+ \\
(X=2) *(7=7) * 2+ \\
(X=5) *(7=7)
\end{array}\right)
$$

Using rewrite rules 2 and 3 equality constraints which have the same value can be rewritten as 1 and different values are rewritten as 0 :

$$
\rightarrow \quad(X=2) *(Y=6) * 3 *\left(\begin{array}{l}
0 * 0+  \tag{6.4}\\
1 * 1+ \\
1 * 0 * 2+ \\
0 * 0
\end{array}\right)+(Y=7) *\left(\begin{array}{l}
(X=1) * 0+ \\
(X=2) * 0+ \\
(X=2) * 1 * 2+ \\
(X=5) * 1
\end{array}\right)
$$

The zeros in the R-expr cause branches of the disjunction to be eliminated (rewrite rule 14), and multiplicities can be combined together using multiplication (rewrite rule 12):

$$
\begin{equation*}
\rightarrow \quad(X=2) *(Y=6) * 3 *(1)+(Y=7) *\binom{(X=2) * 2+}{(X=5) * 1} \tag{6.5}
\end{equation*}
$$

### 6.2.2 Structuring Disjuncts and Conjuncts as Tries

Notice that the factored form $(Y=7) *((X=2) * 2+(X=5) * 1)+(Y=6) *((X=2) * 3)$ in example (6.5) is a more compact $\mathbf{R}$-expr than the sum-of-products form from example (6.1) that we started with, and it is preferable in some settings. In fact, it is an example of a trie representation of a bag relation. ${ }^{90}$ Like the root node of a

[^56]trie, the expression partitions the bag of $\langle X, Y\rangle$ tuples into disjuncts according to the value of $Y$, and then partitions further by the value of $X$.

A trie-shaped $\mathbf{R}$-expr generally has a smaller branching factor than a sum-ofproducts $\mathbf{R}$-expr. As a result, it is comparatively fast to query it for all tuples that match on a given value of $Y$, or on a ( $Y, X$ ) pair, by narrowing down to the matching sub-R-exprs at each level of the trie. For example, suppose that we query the $\mathbf{R}$-expr from example (6.5) by conjoining it with the query $(Y=5)$, which is represented as an $\mathbf{R}$-expr. In this case, we can rewrite $(Y=5) *(Y=7) \rightarrow 0$ and $(Y=5) *(Y=2) \rightarrow 0$ without having to check the part of the trie for the variable $X$.

The example query ( $\mathrm{Y}=5$ ) provides an opportunity for a larger point. The trie has the form $(Y=7) *(X=2) * R$, where this $\mathbf{R}$-expr can be rewritten to 0 based on the first sub-R-expr ( $Y=7$ ), without spending any effort to rewrite $R$, which may represent a large $\mathbf{R}$-expr. This is an example of short-circuiting evaluation and is the same logic that allows a SAT solver or Prolog solver to backtrack immediately upon detecting a contradiction.

### 6.3 Built-in R-expr Rewrites

Built-in constraints are an important ingredient in representing bag relations. While they are not the only ingredient, ${ }^{91}$ they have the advantage that libraries of built-in constraints, such as plus ( $I, J, K$ ) (section $\S 5.2 .2 .3$ ), usually come with rewrite rules for reasoning about these constraints [72]. Some of the rewrite rules invoke opaque procedural code.

Recall that the arguments to a plus constraint are either variables or ground constants. Not all plus constraints can be rewritten, but a library should provide at least the cases:

[^57]```
plus \((I, J, K) \xrightarrow{26} 0\) if \((I, J, K \in \mathbb{R}\) and \(I+J \neq K\) ) or \(I \notin \mathbb{R}\) or \(J \notin \mathbb{R}\) or \(K \notin \mathbb{R}\)
plus \((I, J, K) \xrightarrow{27} 1\) if \(I, J, K \in \mathbb{R}\) and \(I+J=K \quad \triangleright\) Note that \(\mathbb{R} \subset \mathcal{G}\)
plus \((I, J, X) \xrightarrow{28}(X=I+J) \quad\) if \(I, J \in \mathbb{R}\) and \(X \in \mathcal{V}\)
\(\operatorname{plus}(I, X, K) \xrightarrow{29}(X=K-I) \quad\) if \(I, K \in \mathbb{R}\) and \(X \in \mathcal{V}\)
plus \((X, J, K) \xrightarrow{30}(X=\underbrace{K-J}_{\in \mathbb{R}}) \quad\) if \(J, K \in \mathbb{R}\) and \(X \in \mathcal{V}\)
```

As an example, the $\mathbf{R}$-expr $R=\operatorname{proj}(J, \operatorname{plus}(I, 3, J) \star \operatorname{plus}(J, 4, K))$ represents the infinite set of $(I, K)$ pairs such that $K=(I+3)+4$ arithmetically. (The intermediate temporary variable $J$ is projected out.) The rewrite rules already presented (plus a rewrite rule from section $\S 6.4$ below to eliminate proj) suffice to obtain a satisfactory answer to the query ( $I=2$ ) or $(K=9)$, by rewriting either ( $I=2$ )*R or $R *(K=9)$ to ( $\mathrm{I}=2$ ) * ( $\mathrm{K}=9$ ).${ }^{92}$

On the other hand, if we wish to reduce the $\mathbf{R}$-expr R on its own, the above rules do not apply. In the jargon, the two plus constraints within $R$ remain as delayed constraints, which cannot do any work until more of their variable arguments are replaced by constants (e.g., due to equality propagation from a query, as above).

We can do better in this case with a library of additional rewrite rules that implement standard theorems of arithmetic [72]. With these, the R from the same example reduces to plus ( $I, 7, K$ ), which is a simpler description of this infinite relation. Such rewrite rules are known as idempotent constraint handling rules. Other useful examples concerning plus include plus ( $0, \mathrm{~J}, \mathrm{~K}$ ) $\xrightarrow{31}(\mathrm{~K}=\mathrm{J})$ and plus $(I, J, J) \xrightarrow{32}(I=0)$, since, unlike the rules at the beginning of this section, they can make progress even on a single plus constraint whose arguments include more than one variable.

Similarly, some useful constraint propagators for the lessthan relation include lessthan $(\mathrm{J}, \mathrm{J}) \xrightarrow{33} 0 ;$ the transitivity rule lessthan(I, J)*lessthan(J,K) $\xrightarrow{34}$

[^58]lessthan (I, J)*lessthan (J, K)*lessthan(I, K); and lessthan(0, I)*plus(I, J, K) $\xrightarrow{35}$ lessthan( $0, I$ ) *plus(I, J, K)*lessthan(J,K). The integer domain can be split by rules such as int(I) $\xrightarrow{36} \operatorname{int}(I) *($ lessthan $(I, 1)+$ lessthan $(0, I))$ in order to allow case analysis of, for example, int(I)*myconstraint(I). Rewrite rules can also be used to propagate the domain information about variable: lessthan( $I, A$ ) *lessthan ( $I, B$ ) *plus $(A, B, C) \xrightarrow{37}$
$\operatorname{proj}(K, l e s s t h a n(I, A) * l e s s t h a n(J, B) * p l u s(A, B, C) * p l u s(I, J, K) * l e s s t h a n(K, C)) . A l l$ of these rules apply even if their arguments are variables, so they can apply early in a reduction before other rewrites have determined the values of those variables. Indeed, they can sometimes short-circuit the work of determining those values.

Like all rewrites, built-in rewrites $R \rightarrow S$ must not change the denotation of $R$ : they ensure $\llbracket \mathrm{R} \rrbracket_{E}=\llbracket \mathrm{S} \rrbracket_{E}$ for all $E$. For example, lessthan $(\mathrm{X}, \mathrm{Y}) *$ lessthan $(\mathrm{Y}, \mathrm{X}) \rightarrow^{*} \theta$ is semantics-preserving because both forms denote the empty bag relation.

### 6.4 Projection

Projection is implemented using the following rewrite rules. The first two rewrites make it possible to push the $\operatorname{proj}(X, \cdots)$ R-expr down through the sums and products of $R$, so that it applies to smaller subexpressions that mention $X$. The third rewrite allows for projections of different variables to commute with each other:
$\operatorname{proj}(X, R+S) \stackrel{38}{\longleftrightarrow} \operatorname{proj}(X, R)+\operatorname{proj}(X, S) \quad \triangleright$ distributivity over + $\operatorname{proj}(X, R * S) \stackrel{39}{\leftrightarrows} R * \operatorname{proj}(X, S) \quad$ if $X \notin \operatorname{vars}(R) \quad$ see also rewrite rule 45 below $\operatorname{proj}(X, \operatorname{proj}(Y, R)) \stackrel{40}{\longleftrightarrow} \operatorname{proj}(Y, \operatorname{proj}(X, R)) \quad \triangleright \operatorname{projections}$ commute with each other

We can then use the following rewrite rules to eliminate the projection operator from $\mathbf{R}$-exprs whose projection is easy to compute. (In other cases, it must remain delayed.)

```
\(\operatorname{proj}(X,(X=Y)) \xrightarrow{41} 1\) if \(X \neq Y\)
\(\operatorname{proj}\left(X,\left(X=f\left[Y_{1}, \ldots, Y_{n}\right]\right)\right) \xrightarrow{42} 1\) if \(X \notin\left\{Y_{1}, \ldots, Y_{n}\right\}\)
\(\operatorname{proj}(X, \operatorname{bool}(X)) \xrightarrow{43} 2 \quad \triangleright\) cardinality of a variable given bool constraint
\(\operatorname{proj}(X, \operatorname{int}(X)) \xrightarrow{44} \infty \quad \triangleright\) cardinality of a variable given int constraint
\(\operatorname{proj}(X, R) \xrightarrow{45} R * \infty \quad\) if \(X \notin \operatorname{vars}(R) \quad \triangleright\) cardinality of an unconstrained variable
```

$\operatorname{proj}(X, \operatorname{proj}(Y$, boolean_or $(X, Y, t r u e))) \xrightarrow{46} 3 \quad \triangleright$ cardinality of a variable pair given $a$ certain constraint (bOOLEAN_OR has 3 satisfying assignments that result in true)
$\operatorname{int}(X) * \operatorname{int}(Y) * \operatorname{proj}(Z, \operatorname{plus}(X, Y, Z)) \xrightarrow{47} \operatorname{int}(X) * \operatorname{int}(Y) \quad \triangleright$ cardinality of integer addition (a unique answer always exists)

How are these rewrites justified? Observe that $\operatorname{proj}(X, R)$ in an environment $E$ denotes the number of $X$ values that are consistent with $E$ 's binding of R's other free variables. Thus, we may safely rewrite it as another expression that achieves the same denotation for every $E$. For example, in the case that a variable is restricted to be a boolean value (e.g. $(X=$ true $)+(X=f a l s e)$ ), then there are two possible assignments to the variable $\times$ (rewrite rule 43).

### 6.4.1 Example Projection

As a simple example, let us project the variable $K$ out of simple $\mathbf{R}$-expr comprised of equality constraints.

$$
\begin{aligned}
& \operatorname{proj}(K,((J=1) *(K=1)+\quad(J=1) * \operatorname{proj}(K,(K=1)) \\
& (J=2) *(K=6)+\quad+(J=2) * \operatorname{proj}(K,(K=6)) \\
& (\mathrm{J}=2) *(\mathrm{~K}=7)+\rightarrow^{*}+(\mathrm{J}=2) * \operatorname{proj}(\mathrm{~K},(\mathrm{~K}=7)) \\
& \text { (J=2) * (K=7) + + (J=2) * proj(K, (K=7)) } \\
& (\mathrm{J}=5) *(\mathrm{~K}=7) \quad) \quad+(\mathrm{J}=5) * \operatorname{proj}(\mathrm{~K},(\mathrm{~K}=7))) \\
& \rightarrow^{*}(\mathrm{~J}=1)+(\mathrm{J}=2) * 3+(\mathrm{J}=5)
\end{aligned}
$$

Figure 6-1. Example Rewrites with Projection

The second step uses rewrite rules 38 and 39, while the third step involves rewrite rule 41.

### 6.5 Aggregation

We give the following rewrite rules only for the aggregator sum, as isomorphic rewrites apply to the other aggregators. They rewrite $A=\operatorname{sum}(X, R)$ as a chain of plus constraints that maintain a running total. The three rewrites handle cases where
$R$ is expressed as a disjunction of 0 , 1 , or 2 bag relations, respectively. A larger disjunction such as $(Q+R)+S$ is handled recursively.

$$
\begin{aligned}
& \begin{array}{l}
A=\operatorname{sum}(X, 0) \xrightarrow{48} 0 \\
A=\operatorname{sum}(X,(X=Y))
\end{array} \xrightarrow{\text { 49 }}(A=Y) \\
& A=\operatorname{sum}(X, R+S) \xrightarrow{50} \operatorname{proj}(B, \operatorname{proj}(C,(B=\text { sum }(X,(X=\text { agg_null })+R)) * \\
& (C=\text { sum }(X,(X=\text { agg_null })+S)) * \\
& \quad \text { plus }(B, C, A) * \text { not_equal }(A, \text { agg_null })))
\end{aligned}
$$

The first two rewrites, 48 and 49 , handle the case where there are 0 or 1 assignments to the variable $x$ in the bag relation. If there is nothing contained in the bag being aggregated, then the result of aggregation was defined "as a value not contained in $\mathcal{G} "$, which is equivalent to rewriting the expression as 0 . In the case that there is only one value to sum over, then the aggregation does not change the resulting value, hence it can be eliminated.

### 6.5.1 Rewrite Rule 50 for Handling Disjunctions

The third rewrite rule requires some explanation. Here, we are handling the case where there is a disjunction $\mathrm{R}+\mathrm{S}$ that has been brought to the top of the aggregator. In this case, we can aggregate over R and S independently of each other and then combine the result using the aggregator's associated operator. In this case, using the built-in plus( $B, C, A$ ). To correctly handle the disjunction, there are a few scenarios that we need to handle carefully, namely the cases where $R \rightarrow^{*} 0$ or $S \rightarrow^{*} 0$.

First, let us consider the easy case where both $R$ and $S$ are rewritten to have some contributed value. For demonstration purposes, suppose that we have rewritten $R$ into $R \rightarrow(X=5)$ and rewritten $S$ into $S \rightarrow(X=7)$. In this case, the entire $R$-expr would be $A=\operatorname{sum}(X, \quad(X=5)+(X=7))$. We need to rewrite this $\mathbf{R}$-expr so that the values 5 and 7 are arguments to plus so we can compute their sum. Using rewrite rule 50 , the resulting $\mathbf{R}$-expr includes the sub-R-exprs $B=$ sum ( $X,(X=$ agg_null $)+(X=5))$ $C=\operatorname{sum}(X,(X=$ agg_null $)+(X=7)) .{ }^{93}$ In which case we can then use rewrite rule 51 to

[^59]work around the ( $\mathrm{X}=$ agg_null) disjunct that has been added to this expression.
$$
A=\operatorname{sum}(X, \quad(X=\text { agg_null })+(X=Y)) \xrightarrow{51}(A=Y)
$$

This will assign the value of $(B=5)$ and $(C=7)$. Using equality propagation, we get the R-expr plus ( $5,7, A$ ), which can be rewritten using rewrite rule 28 to get the final result ( $A=12$ ).

Now, let us consider a harder case where one of $R$ or $S$ is rewritten as 0 . Without loss of generality, we will assume that R can be rewritten as 0 . And for demonstration purposes, we will say that $S$ can be rewritten as $S \rightarrow(X=7)$. Hence, we could have the rewrite sequence $(A=\operatorname{sum}(X, R+S)) \rightarrow(A=\operatorname{sum}(X, 0+S)) \rightarrow(A=\operatorname{sum}(X, S)) \rightarrow(A=\operatorname{sum}(X$, $(X=7))) \rightarrow(A=7)$, where the aggregation is only performed only on $S$. However, using this sequence of rewrite requires that we defer rewrite rule 50 until we are sure that all remaining disjuncts cannot be rewritten as 0 . This is contradictory to our design, where we allow rewrites to apply whenever they can be matched. As such, rewrite rule 50 can be applied when we match a disjunction inside of the aggregator. To continue with the scenario, for the sake of demonstration, let us consider what happens if we leave off the ( $X=$ agg_null) disjunct on rewrite rule 50 . In which case we will have ( $B=\operatorname{sum}(X, R)) *(C=s u m(X, S))$ as a sub-R-expr on the right-hand side of rewrite rule 50 . As already stated, $R$ can be rewritten as 0 . Hence, the resulting rewrite sequence would therefore become $(B=\operatorname{sum}(X, R)) *(C=\operatorname{sum}(X, S)) \rightarrow(B=\operatorname{sum}(X, 0)) *(C=\operatorname{sum}(X, S)) \xrightarrow{48} 0 *(C=\operatorname{sum}(X, S)) \xrightarrow{14} 0$. Clearly, this is incorrect. It does not match $(A=7)$, which we know is the correct answer as $S \rightarrow(X=7)$. The issue here is that rewrite rule 48 takes $(B=\operatorname{sum}(X, 0))$ and rewrites it as 0 . As such, we need to suppress the behavior of rewrite rule 48 while still allowing rewrite rule 49 to rewrite the aggregator so that we can get the contributed value. ${ }^{94}$

[^60]To suppress rewrite rule 48 we are going to introduce the disjunct ( $X=$ agg_null), so the $\mathbf{R}$-expr now be $(B=\operatorname{sum}(X,(X=$ agg_null $)+R)) *(C=\operatorname{sum}(X,(X=$ agg_null $)+S))$. Now when $R$ is rewritten as 0 , this causes the first conjunct to become ( $B=\operatorname{sum}(X$, $\left.\left.\left(X=a g g \_n u l l\right)+0\right)\right) \xrightarrow{15,49}(B=$ agg_null). Here, the interpretation of the value agg_null is that the aggregated $\mathbf{R}$-expr $R$ returned nothing. Given that this no longer rewrites as 0 , this means the $\mathbf{R}$-expr conjunctive ( $B=$ agg_null) $) ~(C=\operatorname{sum}(X,(X=$ agg_null $)+S)$ ) is not rewritten as 0 but instead rewritten as ( $B=$ agg_null $) *(C=7)$. Now, we have to work around the value agg_null again. This is done by extending the definition of plus so that it ignores agg_null and treats it like a numerical 0 . This is done using rewrite rule 52.

$$
\begin{aligned}
& \text { plus(agg_null, } A, B) \xrightarrow{52}(A=B) \\
& \text { plus(A, agg_null, B) } \xrightarrow{52}(A=B)
\end{aligned}
$$

This means that we will get plus(agg_null, $7, A) \rightarrow(A=7)$, which is the correct result for this example.

Finally, there is one last case that we need to handle. This is the case where both $R$ and $S$ are rewritten as 0 . In this case, we have that $(A=\operatorname{sum}(X, R+S)) \rightarrow(A=\operatorname{sum}(X, 0))$ $\xrightarrow{49} 0$. With $R$ and $S$ being rewritten as 0 after rewrite rule 50 was applied, we will get the $\mathbf{R}$-expr ( $\left.B=a g g \_n u l l\right) *\left(C=a g g \_n u l l\right) * p l u s(B, C, A)$. The value agg_null will be assigned to the variable $A$ because of rewrite rule 52 passing the value through the plus. Hence, we will get (A=agg_null), which is not the 0 we want. We can fix this by adding in the constraint not_equal(A, agg_null), which checks that the final value from the aggregator is not agg_null. In the case where the value is agg_null, this $\mathbf{R}$-expr is rewritten as 0 , which causes the conjunction to become 0 which is the behavior we want. ${ }^{95}$

We can further work around the value agg_null by also introducing rewrite rule 53 , which mimics rewrite rule 50 :

$$
\begin{aligned}
A=\operatorname{sum}(X, \quad(X=\text { agg_null })+R+S) \xrightarrow{53} \operatorname{proj}(B, \operatorname{proj}(C, & (B=\operatorname{sum}(X,(X=\text { agg_null })+R)) * \\
& (C=\operatorname{sum}(X,(X=\text { agg_null })+S)) * \\
& \text { plus }(B, C, A))) \\
& \text { if } R \neq(X=\text { agg_null }) \text { and } S \neq(X=\text { agg_null })
\end{aligned}
$$

[^61]
### 6.5.2 Other Aggregation Rewrites

In addition to rewrite rules 48 to 50, we have rewrite rules to handle other common R-exprs seen inside of an aggregator. For example, conjunctions are handled with the following rewrite rules:

$$
\begin{aligned}
& A=\operatorname{sum}(X, R * S) \stackrel{54}{\hookrightarrow} R *(A=\operatorname{sum}(X, S)) \quad \text { if } R \text { is a constraint, and } X \notin \operatorname{vars}(R) \\
& A=\operatorname{sum}(X, M *(X=T)) \xrightarrow{55} \quad(A=(T \times \llbracket M \rrbracket)) \text { if } M \in \mathcal{M} \text { and } M>0 \\
& A=\operatorname{sum}(X,(X=\text { agg_null })+M *(X=T)) \xrightarrow{56} \quad(A=(T \cdot \llbracket M \rrbracket)) \text { if } M \in \mathcal{M} \text { and } M>0
\end{aligned}
$$

First, rewrite rule 54 allows for aggregators to be permeable to constraints that do not mention the variable $X$. For example, in the $\mathbf{R}$-expr $A=\operatorname{sum}(X$, lessthan $(Z, 5) *(X=7))$, the constraint lessthan $(Z, 5)$ does not impact $X$ directly, besides controlling if the entire expression has a non-zero multiplicity. Hence, we can move it out of the aggregator without influencing the overall expression. Additionally, if we have an $\mathbf{R}$-expr like $(Y=3) *(A=\operatorname{sum}(X$, plus $(4, Y, X)))$, then the constraint $(Y=3)$ can be brought into the aggregator so that we can evaluate the built-in plus.

Rewrite rules 55 and 56 handle the case that there are other non-constraint $\mathbf{R}$-exprs contained inside of the aggregator that does not influence the aggregated value $X$. For example, if we have an $\mathbf{R}$-expr like $A=\operatorname{sum}(X$, (lessthan $(Z, 5)+$ lessthan $(1, Z)) *(X=3)$ ), then once $Z$ is rewritten as a ground numerical value, then each of the lessthan constraints can be rewritten as either 0 or 1 . Summing their multiplicities together, this will result in either 1 or 2 . In the case of 1 , we can simply remove the 1 from the expression using rewrite rule 13 . However, when the multiplicity is 2 , we need to count two copies of 3 into the aggregator. ${ }^{96}$ Additionally, note that rewrite rules 55 and 56 are generalizations of previously presented rewrite rules 49 and 51.

The way non-constraints R-exprs are handled differs between different aggregators. For example, rewrite rules 57 and 58 are the same as rewrite rules 55 and 56, but they are defined for the max aggregator instead of sum.

$$
A=\max (X, M *(X=T)) \xrightarrow{57}(A=T) \text { if } M \in \mathcal{M} \text { and } M>0
$$

[^62]$$
A=\max (X,(X=\text { agg_null })+M *(X=T)) \xrightarrow{58} \quad(A=T) \text { if } M \in \mathcal{M} \text { and } M>0
$$

The max aggregator does not care about how many copies a particular value it has, as long as there is at least one copy. In this case, the aggregator ignores the multiplicity M as long as it is greater than 0 .

We also define rewrites for handling general constraints involving aggregators.

$$
A=\operatorname{sum}(X, R) \xrightarrow{59} R\{X \mapsto A\} \text { if } R \text { is a constraint }
$$

With rewrite rule 59, if we are aggregating over a constraint, then we can remove the aggregator (subject to variable renaming). This can be useful when there are nested aggregators, as aggregators themselves are constraints. For example, the R-expr $(A=\operatorname{sum}(X,(X=\min (Y, R))))$ can be rewritten as $(A=\min (Y, R))$ as the outer sum aggregator does nothing.

### 6.5.3 Rewriting Aggregation With Partial Information

Some aggregators have special behavior, which is expressed using rewrites. One such case in logic programming is an exists aggregator (:-), which checks that there is some true input. As such, once we have found a positive result, we can immediately stop rewriting other disjunctive $\mathbf{R}$-exprs contained in the aggregator.

$$
A=\text { exists }(X,(X=\text { true })+R) \xrightarrow{60} A=\text { true }
$$

Another case often seen in AI tree search algorithms is alpha-beta pruning with the $\mathrm{min} / \mathrm{max}$ value of an aggregator. Alpha-beta pruning prunes useless computation once an upper/lower bound is known. This can be expressed using a rewrite rule, which infers lessthan constraints once some value is known.
$A=\min (X,(X=T)+R) \xrightarrow{61}$ lessthaneq $(A, T) *(A=\min (X,(X=T)+$ lessthan $(X, T) * R)$ ) if $T \in \mathcal{G}$
$A=\max (X,(X=T)+R) \xrightarrow{62}$ lessthaneq $(T, A) *(A=\max (X,(X=T)+$ lessthan $(T, X) * R))$ if $T \in \mathcal{G}$
The lessthaneq on the outside of the aggregator can be used with rewrite rules 34 and 35 to infer if the result of this aggregation will be useful. Similarly, the lessthan added to the other disjunctive branch $R$ informs that $\mathbf{R}$-expr what the value must be for it to influence the result of aggregation. This allows for the elimination of useless disjunctive branches of the min/max aggregator.

### 6.6 Conditional if-Expression Rewrites

An if-expression switches between two different $\mathbf{R}$-exprs. This is used to maintain $\mathbf{R}$-exprs, which are only used in a few scenarios to override other $\mathbf{R}$-exprs. To rewrite an if-expression, the rewrite system must determine if the conditional sub-R-expr is "true" (has multiplicity greater than zero). Here, the goal is to eventually match the conditional expression using rewrite rules 63 and 64 such that the expression can be rewritten as the true or false branch.
if $(M+Q, R, S) \xrightarrow{63} R \quad$ if $M \in \mathcal{M}$ and $M>0 \quad \triangleright$ Return True $\mathbf{R}$-expr branch
if $(0, R, S) \xrightarrow{64} \mathrm{~S} \quad \triangleright$ Return False $\mathbf{R}$-expr branch
Following from the definition of if-expression in section $\S 5.2 .2 .8$, the multiplicity of the if-expression R-expr is only influenced by the then $(\mathrm{R})$ and else ( S ) sub-$\mathbf{R}$-exprs. The multiplicity of the condition does not influence the multiplicity besides switching between the two R-exprs. However, to be able to rewrite the condition into a form that can be matched by rewrite rules 63 and 64, the conditional sub-R-expr needs to be able to "read" from other conjunctive R-exprs so it can be rewritten. This is accomplished using rewrite rule 65 . Here, any conjunctive R-expr can be copied into the condition sub-R-expr, allowing other rewrites to simplify the conditional R-expr. Note, because we only care about the multiplicity of an expression being zero vs. non-zero, duplicating a $\mathbf{R}$-expr may change the magnitude (e.g., 2 vs 4) but does not change the result of comparing zero vs. a nonzero multiplicity.

$$
T * i f(Q, R, S) \stackrel{65}{\leftrightarrows} T * i f(T * Q, R, S)
$$

If-expressions provide many opportunities for potential rewrites that may be useful in various contexts. Observe that if-expressions allow us to switch between two different $\mathbf{R}$-exprs. As such, a rewrite can be used to conditionally pre-rewrite some portion of the $\mathbf{R}$-expr. This is a central insight to enabling memoization, as I will discuss in chapter $\S 10$. Doing this as a rewrite, observe that we can introduce an if-expression anywhere in an R-expr as long as both the true and false branches are identical, as in rewrite rule 66.
if $(\mathrm{Q}, \mathrm{R}, \mathrm{R}) \stackrel{66}{\leftrightarrows} \mathrm{R} \quad \triangleright$ Regardless of $Q$ the same $\mathbf{R}$-expr is returned
Furthermore, we can use the conditional branch of an if-expression to rewrite the true branch of the if-expression. This is allowed because the true branch is only
returned when $Q$ is rewritten as nonzero. Hence, the true branch only needs to be conditionally semantic preserving under Q :
if $(Q, R, S) \xrightarrow{67}$ if $(Q, Q * R, S)$ if $Q$ is a constraint $\triangleright$ Allow possible simplifications of true-branch sub-R-expr

Many if-expressions can be introduced by the rewriting engine when handling memoization. It is, therefore, beneficial to have a number of rewrites that allow us to rearrange the if-expressions to keep the R-expr tidy and efficient:

T*if $(Q, R, S) \stackrel{68}{\hookrightarrow}$ if $(Q, T * R, T * S) \quad \triangleright$ The if-expression is permeable if $(\mathrm{Q}, \mathrm{R}$, if $(\mathrm{Q} 2, \mathrm{R} 2, \mathrm{~S})) \xrightarrow{69}$ if $(\mathrm{Q}+\mathrm{Q} 2, \mathrm{R}+\mathrm{R} 2, \mathrm{~S})$ if $\forall_{E} \llbracket \mathrm{Q} * \mathrm{Q} 2 \rrbracket_{E}=0$ and $\llbracket \mathrm{R} * \mathrm{R} 2 \rrbracket_{E}=0^{97}$
$\triangleright$ Combine two non-overlapping nested if-expressions
if $(Q, R$, if $(Q 2, R 2, S)) \stackrel{70}{\longleftrightarrow}$ if(Q+Q2, if(Q, R, R2), S) $\triangleright$ Nested if-expressions can be rearranged
if $(\mathrm{Q}, \mathrm{R}, \mathrm{S}) \xrightarrow{71} \mathrm{~S}$ if Q is a constraint and $\llbracket \mathrm{Q} * \mathrm{R} \rrbracket_{E}=\llbracket \mathrm{Q} * \mathrm{~S} \rrbracket_{E} \quad \triangleright$ if-expressions can be eliminated
if $(Q 1+Q 2, R, S) \xrightarrow{72}$ if $(Q 1, R, i f(Q 2,0,1) * S)+i f(Q 1,0,1) * i f(Q 2, R, S)$
$\triangleright$ if-expressions can be split
$\mathrm{R} * \mathrm{if}(\mathrm{R}, 1,0) \stackrel{73}{\leftrightarrows} \mathrm{R} \quad \triangleright$ Redundant if-expressions constraints can be removed/introduced

### 6.7 User-Defined R-exprs Rewrites

User-defined rewrites are used to implement user-defined relations as R-exprs, as we will see in chapter $\S 7$. These rewrites match with a named functor $f \in \mathcal{F}$ that contain $n$ value types $Y_{1}, \ldots Y_{n} \in \mathcal{V}$. The value typed $Y_{i}$ correspond to the values/variables used where the user-defined R-expr kind $f(\cdots)$ appears within a large R-expr. Each functor $f$ has at most one rewrite associated with it, and it always rewrites as another $\mathbf{R}$-expr $R_{f}$.

$$
\mathrm{f}\left(\mathrm{Y}_{1}, \ldots, \mathrm{Y}_{n}\right) \xrightarrow{74} \mathrm{R}_{\mathrm{f}}\left\{\mathrm{X}_{1} \mapsto \mathrm{Y}_{1}, \cdots, \mathrm{X}_{n} \mapsto \mathrm{Y}_{n}\right\}
$$

The variables $X_{1}, \cdots, X_{n}$ are identifiers for the free variables contained in the userdefined $R_{f}$. The free variables of $R_{f}$ must be a subset of $X_{1}, \cdots, X_{n}$, otherwise the

[^63]R-expr will become ill-formed $\left(\operatorname{vars}(R) \subseteq\left\{X_{1}, \ldots, X_{n}\right\}\right)$. The names of the variables in $R_{f}$ are renamed to match the names $Y_{i}$ where $f$ originally appeared.

### 6.8 Incompleteness of Included Rewrites

It is impossible to build a sound and complete set of rewrites for all R-exprs given that $\mathbf{R}$-exprs are sufficiently powerful to express mathematical proofs. This is in contrast to Datalog (section §3.1.2) and SLD resolution [101] that have a more limited representation for terms and can therefore have complete implementations. In other words, a rewrite system that can rewrite all R-exprs would be capable of solving all mathematical proofs-which is impossible under Gödel's incompleteness theorem. This will be shown in section §15.1.2. Rather than attempting to make R-exprs and their rewrite complete, we hope to provide a useful but incomplete set of rewrites, much like compute algebra systems such as Mathematica [150] or SymPy [103].

## Chapter 7

## Conversion of Logic Programs to Relational Expressions

A Dyna program is mechanically converted into $\mathbf{R}$-exprs by our front-end parser. I will illustrate the translation to R-exprs with a few examples. In general, this chapter should not be surprising for those who are familiar with the design of front-ends for compilers or interpreters.

### 7.1 Dyna Programs Represent a Key-Value Map

A Dyna program can be thought of as defining a map between keys and valves. Keys are the names of terms in the program, and the value is the returned value computed from a user-defined term's definition in the Dyna program. For example, consider the single rule defined on line 359:

$$
\begin{aligned}
& \operatorname{a(Arg1,Result)} \rightarrow(\text { Result=sum(Inp, } \\
& \operatorname{proj}(X,(X=\operatorname{Arg} 1) * \operatorname{times}(X, X, \text { Inp }))))
\end{aligned}
$$

(a) Dyna
(b) The Program Represented through the is relation as an $\mathbf{R}$-expr.

Figure 7-1. Single rule Dyna program.

The user-defined term $a(X)$ is defined by the rule ' $a(X)+=X * X$.' on line 359 , which defines a relation between keys of the form $a(X)$, such as a(5), and the returned values, such as 25 in the case of a(5). We represent this relation as an R-expr we denote as a(X,Result). ${ }^{98}$ All user-defined terms in Dyna return some value. When representing user-defined logic programming expressions, a "dummy" return value of true is used.

All R-exprs created from Dyna terms have an aggregator at the top of the R-expr. The aggregator enforces the functional dependency between the arguments of the user-defined term and its return value. Functional dependencies are not otherwise enforced by $\mathbf{R}$-exprs.

### 7.1.1 Grouping User-Defined Rules by Name

User-defined terms are grouped according to the outer functor name and arity (number of arguments). This grouping is usually referred to using a slash. For example, we write ' $a / 1$ ' to reference the term defined in figure $7-1$. This grouping is natural and aligns with the user's expectations in how terms are identified by meta-Dyna rules like \$memo that controls memoization (section §2.7).

All contributed rules are grouped under an aggregator R-expr. For example, b(X) in figure 7-2 shows two overlapping rules defined on lines 360 and 361:

[^64]```
b(Arg1,Result) -> (Result=sum(Inp,
    proj(X,(Arg1=X)*times(X,X,Inp))+
| b(X) += X*X
```

(a) Dyna
(b) Program with two user-defined rules combined using a disjunction +.

Figure 7-2. Two user-defined rules with overlap.

Arguments to user-defined terms are given placeholder variable names, such as Arg1 used in these examples. Local variables are introduced using projection and are unified with the placeholder argument variables. This allows us to support expression appearing on the left-hand side of an aggregator without special handling:
sumlist([]) = 0.
sumlist([H|T]) =
H+sumlist(T).
(a) Dyna

```
```

sumlist(Arg1,Result) -> (Result=only(Inp,

```
```

sumlist(Arg1,Result) -> (Result=only(Inp,
(Arg1=list_end[])*(Inp=0)+
(Arg1=list_end[])*(Inp=0)+
proj(H,proj(T,proj(Tmp,
proj(H,proj(T,proj(Tmp,
(Arg1=list_cons[H,T])*
(Arg1=list_cons[H,T])*
sumlist(T,Tmp)*
sumlist(T,Tmp)*
plus(H,Tmp,Inp))))))

```
```

        plus(H,Tmp,Inp))))))
    ```
```

(b) R-expr

Figure 7-3. The sumlist rule (lines 362 to 364 is converted into a single R-expr.

This might appear inefficient at first glance, but we can use the rewrite rules that we have already defined to rewrite these $\mathbf{R}$-exprs to eliminate unneeded projections (e.g. rewrite rules 5, 39 and 41).

### 7.1.2 Different Aggregators

Dyna allows for different aggregators to be defined for the same functor name-arity grouping. To support this as an R-expr, we first group by the aggregator's name and then nest different aggregators under the 'only' aggregator. The 'only' aggregator corresponds to the "equal-sign aggregator" (=) which ensures that there is only one result; otherwise it will return an error (e.g. $(A=o n l y(X,(X=1)+(X=2))) \rightarrow(A=e r r o r))$.

```
c(X) += X*X c(Arg1,Result) -> (Result=only(Inp1,
    for }X>0.\quad(Inp1=sum(Inp2
c(X) min= exp(X) proj(X,(X=Arg1)*lessthan(0,X)*times(X,X,Inp2))))+
    for }x<0
    (Inp1=min(Inp2,
c(X) min= sin(X) proj(X, (X=Arg1)*lessthan(X,0)*exp(X,Inp2))+
    for }X<0.\quad\operatorname{proj}(X,(X=Arg1)*lessthan(X,0)*\operatorname{sin}(X,\operatorname{Inp2))))))
```

(a) Dyna
(b) only is used to ensure there is only one contribution.

Figure 7-4. Multiple aggregators can co-exist as long as they do not overlap.

### 7.1.3 Additional Metadata for Aggregators

Some aggregators require additional steps when translating from Dyna into R-exprs. For example, the := aggregator allows us to override the value by defining additional rules. To make last_override a commutative operator, we annotate each aggregated value with its line number and then select the value contributed from the last defined line number.

```
    d(Arg1,Result) -> (Result=last_overrides(Inp,
    proj(X,proj(Tmp,
        (X=Arg1)
        plus(X,1,Tmp)*
371 |d(X) := X + 1. (Inp=value_from_line[371,Tmp])))+
372 d(X) := X * 2 proj(X,proj(Tmp,
373 for }X<=0
        (X=Arg1)
        times(X, 2,Tmp)*
        (a) Dyna (Inp=value_from_line[372,Tmp])*
        lessthaneq(X,0)))))
```


## (b) R-expr

Figure 7-5. The translation of := aggregator annotates every contribution with its line number so it can determine which value should be returned.

### 7.1.4 Built-ins

Built-ins such as plus and times are mapped from their source code infix representation of ' + ' and ' $*$ '. Other infix operators, such as subtraction '-' do not have their own R-expr representation but instead are written in terms of the plus relation by reordering its arguments.

$$
\begin{array}{ll}
" \text { subtract" }(A, B, C) \rightarrow \text { plus }(A, C, B) & \triangleright \text { subtract does not exist, it is represented with plus } \\
" \text { "divide" }(A, B, C) \rightarrow \text { times }(A, C, B) & \triangleright \text { divide does not exist, it is represented with times } \\
\text { "greaterthan" }(A, B) \rightarrow \text { lessthan }(B, A) \triangleright \text { greaterthan does not exist, represented as lessthan }
\end{array}
$$

Figure 7-6. There do not exist R-expr kinds for subtract, divide, or greaterthan, hence the "scare quotes" around their name. Instead, they are represented using equivalent built-in R-expr kinds by rearranging their arguments.

All built-in callable that are from the source code (such as $C=A+B$ in the case of plus(A, B, C)) return some value, just like the user-defined terms. This includes binary operators expressed as ternary relations like logical and, logical or, less than, and greater than, which all have three value slots. This is done to support logical expressions using and '\&\&', or ' $|\mid$ ', and not '!', (e.g. the Dyna expression '(A<B) \|\| ! $C>D$ )'). That said, in this dissertation, I will often write relations like lessthan using only two of the three slots, as the third argument is almost always the constant value true, and this makes the presentation more concise.
lessthan $(A, B$, true $) \equiv$ lessthan $(A, B) \quad \triangleright$ Third argument omitted in presentation

Built-ins are also used to implement language features such as indirect function calls and dictionaries.

```
    e(Arg1,Arg2,Result) -> (Result=only(Inp,
    proj(Func,proj(A,proj(B,proj(Tmp1,
e(Func,{A, B}) =
Func(A, B).
    (Func=Arg1)*
    map_access(Arg2, "A", A ,Tmp1)*
    map_access(Tmp1,"B",B,empty_map)*
    indirect_call(Func,A,B,Inp)))))))
(a) Dyna, (§2.8)
```

(b) R-expr

Figure 7-7. Dyna syntactic features with their semantics backed by built-ins

### 7.1.5 Dynabases

Dynabases and how they are converted into R-exprs are deferred until chapter $\S 13$.

## Chapter 8

## A Basic Implementation of R-expr Rewriting

In this chapter, I will discuss a minimal and simple implementation for rewriting $\mathbf{R}$-exprs. The system described in this chapter can be implemented in approximately 1000 lines of Python, and the pseudocode for this algorithm can be found at the end of this chapter in section §8.A.

In chapter §11, I will discuss a "more realistic", "fully featured" implementation of term rewriting built on R-exprs. I note that the implementation described in the present chapter $\S 8$ is sufficient to execute many complicated programs represented as R-exprs but not all. Additionally, programs are likely to run very slowly and have a suboptimal asymptotic runtime.

The implementations of Dyna using R-exprs in this chapter and in chapter $\S 11$ are designed to be as close as possible to the "mathematical" design of the R-expr semantics and rewrite rules presented in chapters 5 and 6. Internally, Dyna is as homogeneous as possible, representing almost all state using R-exprs. This includes the program itself, as in chapter $\S 7$, queries, query results (as discussed in chapter §2), and memoized and compiled representations (as we will see in chapters 10 and 12).

As discussed in section $\S 2.3$, a user's interaction with the Dyna system consists of interleaved queries and updates, much like a database. Queries observe the state of the Dyna program, and updates modify the state of the program. When the

Dyna system is queried, it translates the query into an $\mathbf{R}$-expr (as in chapter $\S 7$ ), and rewrites the query's R-expr into a semantically equivalent R-expr. Ideally, the resulting R-expr will be "simpler" and provide a useful answer. I will give an intuitive definition of "simpler" shortly and a formal definition in chapter $\S 15$. In principle, we are allowed to return any semantically equivalent $\mathbf{R}$-expr, including the initial query-though echoing back the initial query is undesirable as it is not useful. In this way, Dyna resembles a computer algebra system such as Mathematica [150] or SymPy [103], which has an incomplete but useful collection of identities and proof strategies.

I will start with a high-level description of R-exprs, the data structure, and then discuss our core rewriting procedure, a pair of functions called SIMPLIFY and SimplifyNormalize.

### 8.1 R-exprs, The Data Structure

An $\mathbf{R}$-expr is an immutable recursive data structure. The $\mathbf{R}$-expr is rewritten via recursive functions that return a new $\mathbf{R}$-expr or the exact same $\mathbf{R}$-expr unmodified. The two most commonly used recursive functions used on an $\mathbf{R}$-expr are variable renaming and rewriting for execution via Simplify described in section §8.2. This design should be familiar to those who have used purely functional data structures and programming languages.

We will usually think of an $\mathbf{R}$-expr as a tree data structure, though, in practice, it can be a DAG (directed acyclic graph) with the same sub-R-expr pointed to multiple times. This does not matter as the $\mathbf{R}$-expr is an immutable data structure. Additionally, sharing internal sub-R-exprs can speed up equality checks with pointer equality.

An $\mathbf{R}$-expr is always bounded in size. This property ensures that the $\mathbf{R}$-expr can be represented in memory and that we can traverse the $\mathbf{R}$-expr using recursive depth-first procedures. User-defined $\mathbf{R}$-exprs can represent recursive functions (terms that are defined recursively, section §5.2.2.11). At any given point in time, these user-defined $\mathbf{R}$-exprs are only expanded up to some bounded depth of the
recursion. ${ }^{99}$
The class that implements an $\mathbf{R}$-expr kind contains named fields that can contain an $\mathbf{R}$-expr, value type, or an array of $\mathbf{R}$-exprs or value types. A value type can be either a variable identifier ${ }^{100}$ or a constant value. This allows us to write both plus( $\mathrm{X}, \mathrm{Y}, \mathrm{Z}$ ) and plus( $1,2,3$ ) using the same plus(.,.,.) $\mathbf{R}$-expr kind. Some classes that implement $\mathbf{R}$-exprs contain metadata fields. This allows a single class to implement different $\mathbf{R}$-exprs from the Dyna language. For example, equality constraints with structural terms (sections § 2.1.1 and 5.2.2.2 e.g. $\mathrm{X}=$ =name[Var1, Var2]) are all handled by the same $\mathbf{R}$-expr kind, where the $\mathbf{R}$-expr kind contains array that track the variables (e.g. [Var1,Var2]) as well as a string field that tracks the functor's name.

Lifetime management of an $\mathbf{R}$-expr is handled using the usual garbage collection mechanism in the host programming language.

R-exprs support syntactic equality checking, which requires that all variable names are the same and all sub-R-exprs are in the same order. In many cases, this check is bypassed due to shared internal sub-R-exprs and by checking pointer equality. All $\mathbf{R}$-exprs are hashable and cache their hash code internally. The hash code speeds up equality checking when $\mathbf{R}$-exprs are not equal. It also allows $\mathbf{R}$-exprs to be used as the key (rather than just the value) in a hash table.

### 8.2 Evaluation by Simplifying an R-expr

Dyna programs and user queries are represented as an $\mathbf{R}$-expr. Evaluation of the program is performed through the application of the rewrite rules from chapter §6. We call this process simplification, and have accordingly named the function that applies these rewrites Simplify.

The reason we have called this process simplification is that our rewrite rules

[^65]attempt to rewrite an R-expr into another $\mathbf{R}$-expr that is "simpler" and "easier to understand". For example, the $\mathbf{R}$-expr plus $(1,2, X)$ is rewritten into ( $X=3$ ) using rewrite rule 28. The $\mathbf{R}-\operatorname{expr}(X=3)$ is simpler as no computation is required to identify the value of $x$. Whereas plus $(1,2, x)$ requires adding 1 and 2 to get the value 3 . In section §15.3, I will formalize the notion of simpler, but for now, an intuitive definition of "easier to understand" and "requires fewer steps of computation to get a final answer" will suffice.

### 8.2.1 Properties of Simplify

The function Simplify: $\mathcal{R} \times \mathcal{C} \rightarrow \mathcal{R}$ performs a pass of rewriting and takes an $\mathbf{R}$-expr $R$ and a context $\mathcal{C}$ (to be defined shortly) and rewrites the $\mathbf{R}$-expr as another semantically equivalent $\mathbf{R}$-expr using the rewrite rules from chapter $\S 6 .{ }^{101}$ SIMPLIFY is allowed to selectively apply rewrites as long as it guarantees that no rewrite is completely starved. ${ }^{102}$ SIMPLIFY is not required to apply all possible rewrites every time it is invoked. In fact, to run faster, it will usually only apply a few rewrites each time. Simplify will return an identical R-expr if-and-only-if there are no directional ${ }^{103}$ rewrites that can be applied. This property allows SimplifyNormalize to identify that it has finished rewriting ${ }^{104}$ the R-expr (algorithm 1).

Simplify recursively invokes itself on the $\mathbf{R}$-expr tree structure depth-first. For example, when Simplify encounters an $\mathbf{R}$-expr like $R * S$, it rewrites the $\mathbf{R}$-expr as "Simplify $(R, \mathcal{C})$ "*"Simplify $(S, \mathcal{C})$ "-recursively invoking Simplify on $R$ and $S$ and combining the results using a conjunction ' $*$ '. We require that Simplify runs in a bounded amount of time. This means that Simplify can scan the entire R-expr tree data structure and apply some (but not all) rewrites. ${ }^{105}$ This property prevents Simplify from getting stuck in an infinite loop when rewriting an R-expr. This

[^66]```
function SimplifyNormalize(R)
    \mathcal { C } \leftarrow \text { AlLConjunctiveRexprs(R) D The context, section \&8.2.2.1}
    repeat
        R prev }\leftarrowR\quad\triangleright The \mathbf{R}\mathrm{ -expr R is a pointer to an immutable, recursive structure.
        R\leftarrowSIMPLIFY}(R,\mathcal{C})\quad\triangleright\mathrm{ Attempt to make progress by rewriting R
    until R prev == R DIf no rewrites are performed, then return R
    return R
```

Algorithm 1. SimplifyNormalize invokes Simplify on the R-expr until no more rewrites can be performed by Simplify. SimplifyNormalize is our Turing-complete rewrite processes, whereas Simplify only performs a bounded number of rewriting steps (section §15.2). SimplifyNormalize is only invoked a the root of an R-expr. This is done because SimplifyNormalize does not guarantee fairness to all potential rewrites, whereas Simplify ensures that it returns in a bounded amount of time.
property also ensures that we do not starve a rewrite. For example, suppose that we have the $\mathbf{R}$-expr $Q * R * S$, where there exists a rewrite sequence such that $R \rightarrow{ }^{*} 0$. We do not want to starve the rewrites on $R$ by spending too much time rewriting $Q$.

To ensure that an R-expr is completely rewritten into the "simplest form possible", we invoke the function SIMPLIFY many times until it reaches a normal form, meaning there are no more applicable rewrites, which is indicated by Simplify returning an identical R-expr to the one passed as an argument. Therefore, the normal form is a fixed-point of Simplify, and we call the function that finds this fixed-point SimplifyNormalize (algorithm 1).

SimplifyNormalize is only invoked at the root of an R-expr. The reason for this is that SimplifyNormalize is not guaranteed to return, and on programs that contain infinite loops will never return. ${ }^{106}$ Conversely, Simplify is a single pass ${ }^{107}$ of rewriting, and therefore will not get stuck performing rewrites endlessly and is therefore guaranteed to return.

[^67]
### 8.2.2 Finding Applicable Rewrites

Rewrites are grouped by the R-expr kind (implementation class) that they match. $\operatorname{Simplify}(R, \mathcal{C})$ will use this grouping and checks all rewrites which might match. Simplify is allowed to perform rewrites on any sub-R-expr contained in a larger $\mathbf{R}$-expr. For example, if we have the $\mathbf{R}$-expr $\operatorname{proj}(X, \operatorname{times}(Y, X, Z) * p l u s(1,2, X))$, then Simplify can rewrite plus $(1,2, X) \rightarrow(X=3)$ using rewrite rule 28 . The way this is implemented is that recursive $\mathbf{R}$-expr kinds, such as such as conjunction ( $\cdot * \cdot$ ), disjunction $(\cdot+\cdot)$, projection $(\operatorname{proj}(X, \cdot))$, or aggregation $(X=\operatorname{sum}(A, \cdot))$, have special rules that recursively invoke Simplify on their sub-R-exprs. This design is known as a visitor pattern and ensures that Simplify will visit the sub-R-exprs and ensures that Simplify attempts to rewrite all visited sub-R-exprs.

Some rewrite rules only require local syntactic information to match. For example, the R-expr plus ( $1,2, X$ ) can be matched with rewrite rule 28 and rewritten as $(X=3)$. The numerical values of the first two arguments are constant value types and are embedded in the R-expr. However, this does not work for all rewrites. Many rewrite rules require additional context to match. For example, rewrite rules 5, 34, 35 and 37 combine two conjunctive $\mathbf{R}$-exprs to infer a third. As presented in chapter $\S 6$, these rewrite rules require that the two $\mathbf{R}$-exprs combined together must be next to each other to be triggered. Now, the R-exprs can be rearranged using the bidirectional rewrites from section $\S 6.2$, which handle commutativity, associativity, and distributivity of conjunction and disjunction. ${ }^{108}$ However, this also has problems. To efficiently rearrange the $\mathbf{R}$-expr would require an "oracle" to identify how best to rearrange the R-expr. Additionally, these rewrites are bidirectional, meaning that we could get stuck in an unproductive cycle such as in figure 8-1 (e.g. $R * S \rightarrow S * R \rightarrow R * S$ ). To fix this, we are going to use the context.

### 8.2.2.1 The Context (Filled With Conjunctive R-exprs)

To avoid the issue of rearranging the $\mathbf{R}$-expr, we observe that many rewrite rules that apply to conjunctions only change a part of the $\mathbf{R}$-expr, rather than the entire conjunction. In other words, these rewrites are of the form $R * S 1 \rightarrow R * S 2$ where the $R$ remains unchanged and only serves as a "license" for the rewrite on S1. As such,

[^68]
## Rewriting Without a Context and Explicit Bidirectional Rules

| Step | Rewrite rule used | R-expr |
| :---: | :---: | :---: |
| 1) | Initial R-expr | $(\mathrm{X}=1) \times \operatorname{proj}(\mathrm{Y},(\mathrm{Z}=2) * \operatorname{plus}(\mathrm{X}, 3, Y))$ |
| 2) | ( $X=1$ ) is moved into proj via rewrite rule 39 | $\operatorname{proj}(Y,(X=1) *(Z=2) * \operatorname{plus}(X, 3, Y))$ |
| 3) | ( $X=1$ ) and ( $Z=2$ ) are flipped via commutativity (rewrite rule 19) | $\operatorname{proj}(Y,(Z=2) *(X=1) * \operatorname{plus}(X, 3, Y))$ |
| 4) | ( $X=1$ ) propagates into plus $(X, 3, Y$ ) via rewrite rule 5 | $\operatorname{proj}(Y,(Z=2) *(X=1) * \operatorname{plus}(1,3, Y))$ |
| 5) | plus $(1,3, Y)$ is computed using the built-in rewrite rule 28 | $\operatorname{proj}(Y,(Z=2) *(X=1) *(Y=4))$ |
| 6) | ( $Z=2$ )* $(X=1)$ are moved out of proj via rewrite rule 39 | $(Z=2) *(X=1) * \operatorname{proj}(Y,(Y=4))$ |
| 7) | $\operatorname{proj}(Y,(Y=4))$ is simplified as $Y$ is known, via rewrite rule 41 | $(Z=2) *(X=1) * 1$ |
| 8) | Useless application of commutativity (rewrite rule 19) | $(X=1) *(Z=2) * 1$ |

Figure 8-1. If we directly apply rewrite rules without knowledge of other conjunctive R-exprs, then we need to use the bidirectional rewrite rules on steps 1,2 and 5 to rearrange the R-expr first. Without an "oracle" to tell us when to apply a bidirectional rule, this can result in unproductive rearranging of the $\mathbf{R}$-expr or cycles such as on steps 7 and 8.
tracking the presence of $R$ is sufficient to permit rewriting $S 1 \rightarrow S 2 .{ }^{109}$ We perform this tracking of conjunctive $\mathbf{R}$-exprs using the context $\mathcal{C}$.

The Context $\mathcal{C}$ is a set ${ }^{110}$ that tracks all $\mathbf{R}$-exprs that are conjunctive with the R-expr $R$ that passed as an argument to Simplify. When Simplify recurses through a disjunct, it adds all sub-R-exprs that become conjunctive, as seen in figure 8-2. This allows Simplify to find the conjunctive sub-R-exprs without scanning the entire R-expr. The context, unlike R-exprs, can be mutated in place, allowing for efficient tracking of conjuncts. ${ }^{111}$ When SIMPLIFY recurses through the R-expr, shallow copies ${ }^{112}$ of the context are made to prevent sub-R-exprs disjuncts from being added into the parent's context. ${ }^{113}$

The context allows us to avoid using bidirectional rewrites. This means that we do not need an "oracle" to determine how to arrange an R-expr and that we can avoid unproductive rewrites. For example, if we use SIMPLIFY to rewrite the example from figure 8-1, then we do not have to explicitly pull down the conjunct ( $X=1$ ) into the projection, and propagation of $X$ 's assignment can be handled via the context, as shown in figure 8-3.

The context is used frequently when checking for matches for $\mathbf{R}$-exprs. As such, it is important that looking up an $\mathbf{R}$-expr in the context is as efficient as possible. To enable this, the context includes indexes that allow for efficient retrieval of relevant $\mathbf{R}$-exprs. For example, a particularly frequent case that we have to handle is $\mathbf{R}$-exprs that assign a value to a variable (e.g. ( $X=7$ ), section §5.2.2.1). These $\mathbf{R}$-exprs can be efficiently represented using an associative map (hash-map) from the variable's

[^69]Context $\mathcal{C}$ Tracking Conjunctive R-exprs During Recursive Simplify Calls


Figure 8-2. The context $\mathcal{C}$ tracks the sub-R-exprs that are conjunctive (shown in color and with an underline) when Simplify is invoked on a sub-R-expr shown with a box. The $\mathbf{R}$-exprs $R, S 1, S 2, Q 1, Q 2$ represent $\mathbf{R}$-exprs which are leaves of the $\mathbf{R}$-expr expression, such as built-in constraints or user-defined $\mathbf{R}$-exprs which have not been expanded yet.
name (e.g. X ) to the assigned value (e.g. the number 7).

### 8.2.3 Canonical Ordering of an R-expr

Although bidirectional rewrites such as rewrite rules 18 to 24 are not used for finding conjunctive $\mathbf{R}$-exprs, there are still some $\mathbf{R}$-expr orderings/arrangements that are preferable. For example, placing $\mathbf{R}$-exprs that are more likely to be rewritten as 0 earlier can make Simplify more efficient, as it can stop rewriting earlier, such line 29 of the pseudocode in section §8.A that stops the evaluation of a conjunction once one of the conjuncts is rewritten as $0 .{ }^{114}$

As such, we have a canonical form for the $\mathbf{R}$-expr and do use bidirectional rewrites to attempt to rearrange the $\mathbf{R}$-expr into this form. We perform these rearrangement rewrites on an "is convenient basis", and they are not considered necessary. This means that we perform these rearrangements as long as it is not too expensive to match the rewrite and rearrange the R-expr. ${ }^{115}$

[^70]
## Rewriting With a Context and No Bidirectional Rules

| Step | Rules used / Explanation | R-expr | Context ( $\mathcal{C}$ ) |
| :---: | :---: | :---: | :---: |
| 1) | Initial R-expr and empty Context | $(\mathrm{X}=1) * \operatorname{proj}(\mathrm{Y},(\mathrm{Z}=2) * \operatorname{plus}(\mathrm{X}, 3, Y))$ | \{\} |
| 2) | Conjunctive constraints are added to $\mathcal{C}$ | $(X=1) * \operatorname{proj}(\mathrm{Y},(\mathrm{Z}=2) * \operatorname{plus}(\mathrm{X}, 3, Y))$ | $\begin{aligned} & \{(X=1),(Z=2), \\ & \operatorname{plus}(X, 3, Y)\} \end{aligned}$ |
| 3) | X is propagated using the context | $(\mathrm{X}=1) * \operatorname{proj}(\mathrm{Y},(\mathrm{Z}=2) * \operatorname{plus}(1,3, Y))$ | $\begin{aligned} & \{(X=1),(Z=2), \\ & \operatorname{plus}(X, 3, Y)\} \end{aligned}$ |
| 4) | plus is rewritten using rewrite rule 28 | $(X=1) * \operatorname{proj}(Y,(Z=2) *(Y=4))$ | $\begin{aligned} & \{(X=1),(Z=2), \\ & \operatorname{plus}(X, 3, Y)\} \end{aligned}$ |
| 5) | ( $Z=2$ ) is lifted out of proj via rewrite rule 39 | $(X=1) *(Z=2) * \operatorname{proj}(Y,(Y=4))$ | $\begin{aligned} & \{(X=1),(Z=2), \\ & \text { plus }(X, 3, Y), \\ & (Y=4)\} \end{aligned}$ |
| 6) | $\operatorname{proj}(Y,(Y=4))$ is simplified as $Y$ is known, via rewrite rule 41, constraints that reference $Y$ are removed from $\mathcal{C}$ | $(X=1) *(Z=2) * 1$ | $\{(\mathrm{X}=1),(\mathrm{Z}=2)\}$ |

Figure 8-3. With a context, all conjunctive sub-R-exprs are tracked in $\mathcal{C}$. This eliminates the need to rearrange the $\mathbf{R}$-expr to apply rewrites such as rewrite rule 5 on step 3 and can be used to implement lifting out of the projection on step 5.

The general principle is that we want the $\mathbf{R}$-expr as factored as possible. Sub-R-exprs are pulled as high up as possible in the R-expr and out of projection, aggregation, and disjunctions. For example, rewrite rule 22 is used to rewrite $R * S+R * Q \rightarrow R *(S+Q)$, and rewrite rules 39 and 54 are used to pull out of projection and aggregation: to repeat for convenience, $\operatorname{proj}(X, R * S) \xrightarrow{39} R * \operatorname{proj}(X, S)$ and ( $A=\operatorname{sum}(X, R * S)) \xrightarrow{54} R *(A=\operatorname{sum}(X, S))$ where $X \notin \operatorname{vars}(R)$, and $R$ must be a constraint when using rewrite rule 54.

### 8.2.3.1 Why factored R-exprs are Preferable

In logic programming without aggregation, creating factored $\mathbf{R}$-exprs would only be a nice feature. However, aggregation can make it essential that we factor $\mathbf{R}$-exprs. The reason is that pulling a constraint from an aggregator requires that it is conjunctive with the entire body of the aggregator-meaning it has been factored from any nested disjunctions.

To see this, let us start with a Prolog-style program and translate it into R-exprs with and without Dyna's aggregators.

```
a(X) :- q(X) for X > 5.
a(X) :- s(X) for X > 5.
b(Y) :- r(Y) for Y < 5.
b(Y) :- t(Y) for Y < 5.
should_be_nothing :- a(Z), b(Z).
```

Figure 8-4. Dyna program which requires solving the intersection of $Z>5$ and $\mathrm{Z}<5$ to identify should_be_nothing is unsatisfiable (rewrites as 0 ).

The rule should_be_nothing represents the intersection of a set of values greater than 5 and another set of values less than 5 , which is empty, as such this rule results in nothing: lessthan $(Z, 5) *$ lessthan $(5, Z) \xrightarrow{34,33} 0$.

First, let us consider what should_be_nothing looks like as an R-expr when there are no aggregators:

```
((lessthan (5, Z)*q(Z,ARes))+ Drule a(X)
    (lessthan(5,Z)*s(Z,ARes)))* Drule a(X)
((lessthan(Z,5)*r(Z,BRes))+ Drule b(Y)
    (lessthan(Z,5)*t(Z,BRes))) Drule b(Y)
```

Figure 8-5. Translation of figure 8-4 without aggregators.

We can prove that the $\mathbf{R}$-expr in figure $8-5$ is empty by combining lessthan $(5, Z)$ and lessthan $(Z, 5)$ into the same conjunction. Without aggregation, this can be accomplished using the distributive rewrites to expand this into four different conjunctive cases:

```
(lessthan(5,Z)*q(Z,ARes)*lessthan(Z,5)*r(Z,BRes))+
(lessthan (5,Z)*q(Z,ARes)*lessthan(Z,5)*t(Z,BRes))+
(lessthan(5,Z)*s(Z,ARes)*lessthan(Z,5)*r(Z,BRes))+ ->0
(lessthan(5,Z)*s(Z,ARes)*lessthan(Z,5)*t(Z,BRes))
```

Figure 8-6. Figure 8-5 expanded into four conjunctive cases, and each individually rewritten as 0 .

Each conjunction can be individually rewritten as 0 , which allows the entire R-expr to be rewritten as 0 , proving that should_be_nothing is empty. ${ }^{116}$

Now, let us again consider the program from figure 8-4, but this time translate it into an R-expr with aggregators, as in figure 8-7:

```
(ShouldBeNothing=exists(true,
    proj(Z,
        (true=exists(true, Drule a(X)
            (proj(ARes, (lessthan(5,Z)*q(Z,ARes)))+
            proj(ARes, (lessthan(5,Z)*s(Z,ARes))))))*
        (true=exists(true, \trianglerightrule b(Y)
            (proj(BRes, (lessthan(Z,5)*r(Z,BRes)))+
            proj(BRes, (lessthan(Z,5)*t(Z,BRes)))))))))
```

Figure 8-7. Figure 8-4 translated into R-exprs with aggregators.

In figure 8-7, the aggregator exists prevents us from using the distributive property to expand this R-expr into the four conjunctive cases. Furthermore, the four lessthan constraints are nested under a disjunction, as the lessthan constraint was contributed by each rule from the original Dyna program contributing its own lessthan constraint rather than having a "global" lessthan constraint on $a(X)$ or $b(Y)$. As such, the only way to solve this program is to factor the lessthan constraint out of the $\mathbf{R}$-expr and pull it out of the projection, the disjunction, and then the aggregator. The resulting $\mathbf{R}$-expr is shown in figure 8-8.

[^71]```
(ShouldBeNothing=exists(true,
    proj(Z,
        lessthan(5,Z)*
        (true=exists(true, \trianglerightrule a(X)
            (proj(ARes, q(Z,ARes))+
            proj(ARes, s(Z,ARes)))))*
        lessthan(Z,5)*
        (true=exists(true, Drule b(Y)
            (proj(BRes, r(Z,BRes))+
            proj(BRes, t(Z,BRes))))))))
```

Figure 8-8. Figure 8-7 after factoring lessthan constraint out of the projections, disjunctions and aggregators.

Once we have constructed the $\mathbf{R}$-expr in figure 8-8, we have that both lessthan ( $5, \mathbf{Z}$ ) and lessthan $(Z, 5)$ are in the same conjunct. This means that we can use rewrite rules 33 and 34 to rewrite this conjunction as 0 . This proves that should_be_nothing is false.

## 8.A Appendix: Basic Simplify Pseudocode

Here is the high-level pseudocode of Simplify. The Simplify function recursively calls itself on the $\mathbf{R}$-expr. Simplify starts by matching against the kind of the $\mathbf{R}$-expr. Then, a more detailed matching is performed using the context. This includes checking if a variable is assigned some value (using isGround) and getting the value (using getValue). The context is mutated in place to track conjuncts that are added. When Simplify recurses through recursive R-exprs like projection, disjunction, or aggregation, it makes a "copy" ${ }^{112}$ of the context to scope conjunctions that are added to the context. All new conjuncts are immediately added to the context using allConjunctiveRexprs to find all conjunctive R-exprs before applying the rewrites.

```
function Simplify(\mathcal{C},R)
    if R matches ( }X=Y\mathrm{ ) : }\quad\triangleright\mathrm{ Equality constraints, section §5.2.2.1
        if isGround (\mathcal{C},X) and isGround (\mathcal{C},Y):
```

if $\operatorname{GetValue}(\mathcal{C}, X)==\operatorname{GetValue}(\mathcal{C}, Y):$
return $1 \quad \triangleright$ Return the multiplicity $\mathbf{R}$-expr 1 to indicate success
else
return $0 \quad \triangleright$ Return the multiplicity $\mathbf{R}$-expr 0 to indicate failure
else if isGround $(\mathcal{C}, Y)$ :
$\mathcal{C}[\mathrm{X}] \leftarrow \operatorname{GetValue}(\mathcal{C}, Y) \quad \triangleright$ Record value of X into the context return $1 \triangleright$ Remove equality constraint from $\mathbf{R}$-expr
else if other similar rewrites omitted for brevity : other similar rewrites omitted for brevity

```
else if R matches (X=f[Y},\cdots\mp@subsup{Y}{n}{}]):\quad\triangleright : Structured terms, section $5..2.2.2
```

    if isGround \((\mathcal{C}, \mathrm{x})\) :
        \(\mathrm{x} \leftarrow \operatorname{GetValue}(\mathcal{C}, \mathrm{x})\)
        if \(\mathbf{x}\) matches \(f\left[z_{1}, \cdots, z_{n}\right]\) :
            return \(\left(Y_{1}=Z_{1}\right) * \cdots *\left(Y_{n}=Z_{n}\right) \quad \triangleright\) Unpack the value, rewrite rule 9
        else
            return \(0 \quad \triangleright\) Incompatible value, rewrite rule 10
    else if \(\operatorname{isGround}\left(\mathcal{C}, Y_{1}\right)\) and \(\cdots\) and \(\operatorname{isGround}\left(\mathcal{C}, Y_{n}\right)\) :
        \(z_{1} \leftarrow \operatorname{GetValue}\left(\mathcal{C}, \mathrm{Y}_{1}\right), \ldots, z_{n} \leftarrow \operatorname{GetValue}\left(\mathcal{C}, Y_{n}\right)\),
        return ( \(\mathrm{X}=\mathrm{f}\left[z_{1}, \ldots, z_{n}\right]\) )
    else if \(\left(X=f\left[z_{1}, \cdots, z_{n}\right]\right) \in \mathcal{C}\) :
        return \(\left(X_{1}=Z_{1}\right) * \cdots *\left(X_{n}=Z_{n}\right) \quad \triangleright\) Unify with another \(\mathbf{R}\)-expr that is the same
    structural term, rewrite rule 7
else if $\left(x=g\left[z_{1}, \cdots, z_{m}\right]\right) \in \mathcal{C}$ :
return $0 \quad \triangleright$ Incompatible with different term name, rewrite rule 8
else if $R$ matches $Q * S$ :
$Q^{\prime} \leftarrow \operatorname{Simplify}(\mathcal{C}, Q)$
if $Q$ ' matches 0 : $\triangleright$ Avoid unnecessary rewriting when we get a 0 , rewrite rule 14
return 0
$S^{\prime} \leftarrow \operatorname{Simplify}(\mathcal{C}, \mathrm{s})$
if $Q^{\prime} \in \mathcal{M}$ and $S^{\prime} \in \mathcal{M}$ :
return $Q^{\prime} * S^{\prime} \quad \triangleright$ Multiply the multiplicities together, rewrite rule 12
else if $\mathrm{S}^{\prime}$ matches 0 :
return 0
else if $Q^{\prime}$ matches 1 : $\quad \triangleright$ Handle identity, rewrite rule 13
return $\mathrm{S}^{\prime}$
else if $\mathrm{S}^{\prime}$ matches 1 :
return $Q^{\prime}$

```
        else
            return Q'*S'
    else if R matches proj(X,Q):
        \mathcal{C}}\leftarrow\operatorname{Copy}(\mathcal{C})\quad\trianglerightA\mathrm{ "copy" is made to avoid interfering with the parent context
        \hat { \mathcal { C } } \leftarrow \hat { \mathcal { C } } \cup \text { AlLCONJUNCTIVEREXPRS(Q) DR-exprs that are now conjunctive are}
added to the context
        Q'}
        if IsGround}(\hat{\mathcal{C}},X): \triangleright If ground in the context used inside the projectio
```



```
        return Q'{X\mapstox} }\triangleright\mathrm{ Rename }\textrm{X}\mathrm{ to the constant x, eliminate the projection
        if Q' matches Q1*Q2 where X & vars(Q1) and isConstraint(Q1):
        return Q1*proj(X,Q2) }\triangleright\mathrm{ Pull out constraints that do not depend on X
        return proj(X,Q') \triangleright Can not eliminate the projection
    else if R matches Q+S :
        \mp@subsup{\mathcal{C}}{1}{}\leftarrow\operatorname{COPY}(\mathcal{C})}\quad\triangleright\mathrm{ Rewrite first branch of disjunct
        \mp@subsup{\mathcal{C}}{1}{}}\leftarrow\mp@subsup{\hat{\mathcal{C}}}{1}{}\cup\mathrm{ ALLCONJUNCtIVEREXPRS(Q)
        Q'}\leftarrow\operatorname{Simplify}(\mp@subsup{\hat{\mathcal{C}}}{1}{},Q
        \mp@subsup{\hat{\mathcal{C}}}{2}{*}}\leftarrow\operatorname{COPY}(\mathcal{C})\quad\triangleright Rewrite second branch of disjunc
        \mp@subsup{\mathcal{C}}{2}{}}\leftarrow\mp@subsup{\hat{\mathcal{C}}}{2}{}\cup\mathrm{ AlLCONJUNCtIVEREXPRS(S)
        Q'}\leftarrow\operatorname{SimpLiFY}(\mp@subsup{\hat{\mathcal{C}}}{2}{},\textrm{S}
        if Q' matches 0:
            C}\leftarrow\mathcal{C}\cup\mp@subsup{\hat{\mathcal{C}}}{2}{}\quad\triangleright\mathrm{ One disjunct is 0, so return the other
            return S'
        else if S' matches 0:
            \mathcal{C}}\leftarrow\mathcal{C}\cup\mp@subsup{\hat{\mathcal{C}}}{1}{
            return Q'
        else
            \mathcal { C } \leftarrow \mathcal { C } \cup ( \hat { \mathcal { C } } _ { 1 } \cap \hat { \mathcal { C } } _ { 2 } ) \quad \triangleright ~ C o m m o n ~ c o n s t r a i n t s ~ a r e ~ p u l l e d ~ o u t ~
        \triangleright \text { Constraints not kept in C are encoded back into the R-expr}
        \ (Equality constraints are removed from the \mathbf{R}\mathrm{ -expr by line 9 and saved back)}
            return (( (\hat{\mathcal{C}}
        else if R matches (A=sum(X,Q)) :
        \hat { \mathcal { C } } \leftarrow \operatorname { c o p y } ( \mathcal { C } ) \quad \triangleright A \text { "copy" is made to avoid interfering with the parent context}
        \hat { \mathcal { C } } \leftarrow \hat { \mathcal { C } } \cup \text { ALLCONJUNCTIVEREXPRS(Q) }
    added to the context
        Q'}
```

Rule 28

```
    if Q' matches (X=x):
        return (A=x) \triangleright Aggregation "complete" return, rewrite rule 49
    else if Q' matches Q1+Q2 and Q1 & Q2 not match ( }X=\mathrm{ agg_null) :
        return proj(B,proj(C,(B=sum(X,(X=agg_null)+Q1))* }\triangleright\mathrm{ Rewrite rule 50
        (C=sum(X,(X=agg_null)+Q2))*plus(B, C,A)*not_equal(A, agg_null)))
    else if Q' matches Q1*Q2 where X\not\in\operatorname{vars(Q1) and isConstraint(Q1):}
        return Q1*(A=sum(X,Q2)) \triangleright Pull out constraints that do not influence }
        return (A=sum( }X,\mp@subsup{Q}{}{\prime}))\quad\triangleright Return the R-expr with newly rewritten Q'
    else if R matches if (Q,R,S):
    \mathcal{C}}\leftarrow\operatorname{copy}(\mathcal{C})\quad\trianglerightA\mathrm{ "copy" is made to avoid interfering with the parent context
    Q'}\leftarrow\operatorname{Simplify(Q,\hat{C})
    if Q' matches 1+T (for some T): \triangleright rewrite rule 63
        return R }\triangleright\mathrm{ Return true branch of if-expression
    else if Q' matches 0: \triangleright rewrite rule 64
        return S }\triangleright\mathrm{ Return false branch of if-expression
    else
        return if (Q',R,S) \triangleright Unable to evaluate conditional, return delayed if-expr
    else if R matches f(Y, Y , , , , Y ( ) : 
    R}f\leftarrowL\mp@code{LookupUserDefinition(f, n)
    return R }\mp@subsup{f}{f}{}{\mp@subsup{X}{1}{}\mapsto\mp@subsup{Y}{1}{},\mp@subsup{X}{2}{}\mapsto\mp@subsup{Y}{2}{},\cdots,\mp@subsup{X}{n}{}\mapsto\mp@subsup{Y}{n}{}}\triangleright\mathrm{ Rename variables to names used
    by the }\mathbf{R}\mathrm{ -expr currently being rewritten, }\textrm{R
    else if R matches plus(I,J,K):
    if isGround(\mathcal{C},I) and isGround}(\mathcal{C},\textrm{J})\mathrm{ and isGround}(\mathcal{C},K):\quad\triangleright Rule 27
    and 26
        if GetValue(\mathcal{C},I)+GetValue(\mathcal{C},\textrm{J})== GetValue(\mathcal{C},K):
                return 1 \triangleright Assignment to I, J and K is consistent with plus
        else
            return 0 D Assignment to I, J and K is inconsistent with plus
        else if isGround}(\mathcal{C},I)\mathrm{ and isGround}(\mathcal{C},J)\mathrm{ and not isGround (C, K) : }
    Rule 28
        k\leftarrow\operatorname{GetValue(\mathcal{C}, I) + GetVAlue(\mathcal{C},\textrm{J})}\quad\triangleright\mathrm{ Compute using addition}
        return (K=k)
    else if isGround(\mathcal{C},I) and not isGround}(\mathcal{C},J)\mathrm{ and isGround (C, K) : }
Rule 29
        j}\leftarrow\operatorname{GetValue(\mathcal{C}, I) - GetValue(\mathcal{C},\textrm{K})\quad\triangleright Compute using subtraction
        return (J= j)
            else if not isGround}(\mathcal{C},I)\mathrm{ and isGround (C, J) and isGround (C, K) : }
```


## Rule 30

107:
108:
109:
110:
111:
112:
113:
114:
115:
116:
117:
118:
119:
120:
121:
122:
123:
124:

```
            i\leftarrowGetValue(\mathcal{C},J) - GetValue(\mathcal{C},\textrm{K})\quad\triangleright Compute using subtraction
            return (I=i)
        return R }\triangleright\mathrm{ Return the R-expr unmodified
    else if R matches lessthan(I,J) :
        if isGround}(\mathcal{C},I)\mathrm{ and isGround}(\mathcal{C},\textrm{J})
            if GetValue(\mathcal{C}, I) < GetValue(\mathcal{C},\textrm{J}):
                return 1 \triangleright Assignment to I and J is consistent with lessthan
        else
            return 0 \triangleright Assignment to I and J is inconsistent with lessthan
    else if lessthan (J,K) \in\mathcal{C}\mathrm{ and lessthan (I,K) &C : }\quad\triangleright\mathrm{ Rewrite rule 34}
        \mathcal { C } \leftarrow \mathcal { C } \cup \{ l e s s t h a n ( I , K ) \} \triangleright ~ T r a c k ~ v i a ~ c o n t e x t ~ t h a t ~ a ~ n e w ~ l e s s t h a n ~ i s ~ i n f e r r e d ~
        return R*lessthan(I,K)
    else if lessthan (K,I) \in\mathcal{C}\mathrm{ and lessthan (K,J) &C C: }\, \triangleright Rewrite rule 34
        \mathcal { C } \leftarrow \mathcal { C } \cup \{ l e s s t h a n ( K , J ) \} \triangleright ~ T r a c k ~ v i a ~ c o n t e x t ~ t h a t ~ a ~ n e w ~ l e s s t h a n ~ i s ~ i n f e r r e d ~
        return R*lessthan(K,J)
    return R }\triangleright\mathrm{ Return the R-expr unmodified
    else if R matches .. . :
        Omitted for brevity: Many other rewrites and matching rules.
```

Algorithm 2. An example of how a few rules of Simplify are implemented.

```
function isGround \((\mathcal{C}, \mathrm{v})\)
    if \(\vee \in \mathcal{G}: \quad \triangleright\) A ground constant as a value type is always ground
        return true
    else if \(\mathcal{C}[\mathrm{V}]: \quad \triangleright\) Check if there exists a binding to \(\vee\) in the context \(\mathcal{C}\)
        return true
    else
        return false
function \(\operatorname{GetValue}(\mathcal{C}, \mathrm{V})\)
    if \(\vee \in \mathcal{G}\) :
        return \(\vee\)
    else if \(\mathcal{C}[\mathrm{v}]\) :
        return \(\mathcal{C}[\mathrm{V}]\)
    else
        throw Error \(\triangleright\) Should not get value if one does not exist
function AllConjunctiveRexprs(R)
    \(\triangleright\) Return all of the conjunctive \(\mathbf{R}\)-expr
    if \(R\) matches \(Q * S\) :
        return allConjunctiveRexprs \((Q) \cup\) allConjunctiveRexprs(S)
    else if \(R\) matches \(\operatorname{proj}(X, Q)\) :
        \(\mathrm{c} \leftarrow\) AllConjunctiveRexprs(Q)
        \(\mathrm{V} \leftarrow\) makeDummyVariable ()
        return \(\{\mathrm{S}\{\mathrm{X} \mapsto \mathrm{V}\}: \mathrm{S} \in c\} \quad \triangleright\) Rename variable X to a dummy variable name
    else if \(R\) matches \(A=\operatorname{sum}(X, Q)\) :
        \(\mathrm{c} \leftarrow\) AllConjunctiveRexprs(Q)
        \(\vee \leftarrow\) makeDummyVariable ()
        return \(\{\mathrm{S}\{\mathrm{X} \mapsto \mathrm{V}\}: \mathrm{S} \in c\} \cup\{\mathrm{A}=\operatorname{sum}(\mathrm{X}, \mathrm{Q})\} \quad \triangleright\) Rename variable X
    else if \(R\) matches \(Q+S\) :
        return \(\{\mathrm{Q}+\mathrm{S}\} \quad \triangleright\) Return the disjunction itself, not its children
    else
            return \(\{R\}\)
function \(\operatorname{copy}(\mathcal{C})\)
    \(\triangleright\) Make a copy of the context \(\mathcal{C}\). The copied context shares any immutable internal state
        to avoid making a deep copy.
```

Algorithm 3. Helper functions used by Simplify for interacting with the context.

## Chapter 9

## Rearranging R-exprs to Enable Further Rewriting

In chapter $\S 8$, I introduced the Simplify procedure, which is central in rewriting R-exprs. While Simplify is capable of handling a large number of programs, it is not complete. This is both in a theoretical sense, in that there will always exist programs which Simplify cannot be reduced into a "simple" representation, and in a practical sense, in that there are programs that we know how to solve but Simplify is unable to solve.

In this chapter, I will introduce a formalism for additional rewrites, which can be used to rearrange the R-expr in useful ways. These rewrites, like our bidirectional rewrites from chapter $\S 6$, are not explicitly used. Rather, these rewrites will license us to perform more complicated sequences of rewrites in chapter $\S 11$, just as the rewrites in chapter $\S 6$ licensed the Simplify procedure in chapter $\S 8$.

### 9.1 The Problem with Simplify from Chapter §8

Let us first work through an example where Simplify fails to produce a "useful" and "simple" result.

Our Simplify function creates factored R-exprs so that it can pull constraints out of an aggregator (section §8.2.3). As already discussed in section §8.2.3.1, this is something we have to do to work around aggregators. However, there are many
problems that require us to use distributive rewrites to expand an R-expr into smaller bags ${ }^{117}$ to be able to solve the program. Hence, we need some way to divide the $\mathbf{R}$-expr into smaller, solvable units while still maintaining the factored $\mathbf{R}$-expr so that we can work around aggregators.

To see this, let us work through a program that requires subdividing the problem into smaller subproblems. One such program is boolean SATisfiability, commonly known as the SAT problem [41, 96]. We can write a SAT formula as a Dyna program, as I have done in figure 9-1. A SAT formula is comprised of boolean variables and clauses. Variables can take on the value true or false. Each clause is comprised of one or more literals, which are either variables or the negation of variables. Each clause must have at least one literal that is true. Solving a SAT formula requires finding an assignment to all of the variables such that all clauses are satisfied or proving that there is no assignment that can satisfy all clauses.

```
boolean(1). % true value
boolean(0). % false value
negate(1) = 0. % negate the argument by mapping true to false
negate(0) = 1. % and vice versa
one_true(1, _, _). % check that at least one argument is true
one_true(_, 1, _).
one_true(_, _, 1).
one_true(1, _). % check one argument is true for two variables
one_true(_, 1).
one_true(1). % check that the one argument is true
    % represent a SAT problem as a conjunct of disjuncts (using one_true)
is_sat :- boolean(A), boolean(B), boolean(C), % variables are boolean
    one_true(A, B, negation(C)), % SAT clause
    one_true(negation(A), negation(B)), % SAT clause
    one_true(C), % SAT clause
    one_true(A, negation(B)). % SAT clause
```

Figure 9-1. The Boolean SAT formula $(A \vee B \vee \neg C) \wedge(\neg A \vee \neg B) \wedge C \wedge(A \vee \neg B)$ expressed as a Dyna program

[^72]The SAT problem is known to be NP-complete and, therefore, difficult to solve. SAT can be solved using the DPLL algorithm [45]. ${ }^{118}$ The DPLL algorithm works by searching for possible assignments to variables in the problem. The DPLL does this by alternating between branching and propagating. When branching, the DPLL algorithm will arbitrarily pick an unassigned variable ${ }^{119}$ and loop over the variable's domain (in this case, the domain is $\{1,0\}$ ), trying each assignment in turn. When a variable is assigned, the DPLL algorithm propagates the variable's assignment to all clauses in the program. If an assignment to a variable is found to be inconsistent, the DPLL algorithm backtracks to the last branching location. Propagation can also cause other variables to be assigned, which also need to be propagated. If propagation completes without any inconsistencies, then another variable will be selected for branching. The DPLL algorithm will continue this process of alternating between branching and assignment.

Given the above explanation of the DPLL algorithm, let us take the SAT formula from figure 9-1 and see what happens when we run it under the Simplify function defined in chapter §8. The first step is to translate figure $9-1$ into an $\mathbf{R}$-expr as shown in figure 9-2 using the method described in chapter $\S 7$.

[^73]```
IsSat=exists(true,
    proj(A, proj(B, proj(C,
        (true=exists(true, \((A=1)+(A=0))) * \quad \triangleright\) boolean constraint on \(A\), line 393
        (true \(=\) exists \((\) true,\((B=1)+(B=0))) *\) boolean constraint on \(B\), line 393
        (true=exists(true, \((C=1)+(C=0))) * \quad \triangleright\) boolean constraint on \(C\), line 393
        (true=exists(true, \((\mathrm{A}=1)+\quad \triangleright\) Clause from line 394
            (B=1)+
            proj(NegC, \(\quad \triangleright\) Negation expanded from line 394
                            (NegC=only (Inp, (Inp=0)* (C=1) +
                                    \((\operatorname{Inp}=1) *(C=0))) *\)
                    (NegC=1))))*
    (true=exists(true, proj(NegA, \(\triangleright\) Clause from line 395
            (NegA=only (Inp, \((\operatorname{Inp}=0) *(A=1)+\)
                                    \((\operatorname{Inp}=1) *(A=0))) *\)
            (NegA=1)) +
            proj(NegB,
            (NegB=only (Inp, (Inp=0)*(B=1)+
                                    \((\operatorname{Inp}=1) *(B=0))) *\)
                            (NegB=1))))*
    (true=exists(true, (C=1)))* \(\triangleright\) Clause from line 396
    (true=exists(true, (A=1)+ \(\triangleright\) Clause from line 397
proj(NegB,
                            (NegB=only (Inp, (Inp=0)*(B=1)+
                        \((\operatorname{Inp}=1) *(B=0))) *\)
                    (NegB=1))))))))
```

Figure 9-2. Boolean SAT formula (figure 9-1) translated into an $\mathbf{R}$-expr. Userdefined $\mathbf{R}$-exprs have been expanded (section $\S 6.7$ ). Some variables (such as NegA) can be rewritten away due to their assignment (e.g. with (NegA=1)), but were left in for clarity about the return value of the negation function (lines 383 to 384).

Once we have the $\mathbf{R}$-expr in figure 9-2, the Simplify function will look for opportunities to apply its rewrite rules. The relevant rewrite rule in this case is lifting ( $C=1$ ) out of the aggregator (true=exists(true, ( $C=1$ ))) from line 396 using rewrite rule 54. This rewrite leaves us with (C=1)*(true=exists(true, 1)). The (true=exists(true, 1)) can be rewritten as 1 , therefore removed from the $\mathbf{R}$-expr, and $(C=1)$ is now at the top level of the $\mathbf{R}$-expr and is therefore used to propagate the assignment of true (represented as 1 ) to the variable $C$.

```
IsSat=exists(true,
    proj(A, proj(B, proj(C,
        (C=1)* \(\triangleright\) Assignment from propagation
        (true=exists(true, \((A=1)+(A=0))) * \quad \triangleright\) boolean constraint on A, line 393
        (true \(=\) exists \((\) true,\((B=1)+(B=0))) * \quad \triangleright\) boolean constraint on \(B\), line 393
        (true=exists(true, \((A=1)+\quad \triangleright\) Clause from line 394
        ( \(B=1\) )) )*
        (true=exists(true, (1=only(Inp, (Inp=0)*(A=1)+ \(\triangleright\) Clause from line 395
                        (Inp=1)*(A=0)))
                            (1=only (Inp, \((\operatorname{Inp}=0) *(B=1)+\)
                                    \((\operatorname{Inp}=1) *(B=0))))) *\)
    (true=exists(true, \((A=1)+\quad \triangleright\) Clause from line 397
    (1=only (Inp, \((\operatorname{Inp}=0) *(B=1)+\)
                                    \((\operatorname{Inp}=1) *(B=0)))))\) )) )
```

Figure 9-3. SAT formula from figure 9-2 after propagation and some rewrites. The Simplify function from chapter $\S 8$ does not perform further rewrites on this R-expr.

The propagation of $(C=1)$ is done by adding it to the context $\mathcal{C}$ and will cause all R-exprs of the form ( $\mathrm{C}=0$ ) to be rewritten as 0 . Hence, the negation( $C$ ) ) from line 394 , which is equivalent to ( $\mathrm{C}=0$ ) is rewritten as 0 and removed from this clause entirely.

Once propagation is completed, we are left with the R-expr in figure 9-3 and there are no more rewrites that Simplify will apply. ${ }^{120}$ Continuing from this point requires that one of the remaining unassigned variables, $A$ or $B$, is assigned some value. This requires branching, which Simplify from chapter $\S 8$ does not support. In this way, we can see that Simplify is capable of propagation but does not support branching.

### 9.2 Using Nested Constraints

When we have an R-expr similar to the one in figure 9-3, we want to expand the $\mathbf{R}$-expr by branching to subdivide the $\mathbf{R}$-expr into smaller $\mathbf{R}$-exprs which are solved

[^74]individually. In other words, given an $\mathbf{R}$-expr of the form $R *(Q 1+Q 2)$, it should rewrite into $R * Q 1+R * Q 2 .{ }^{121}$ However, before we can get to an $\mathbf{R}$-expr of the form $R *(Q 1+Q 2)$, there is another problem that we must solve. Observe that in figure 9-3, there is no top-level disjunction of the form $\mathrm{Q} 1+\mathrm{Q} 2$. Hence, there is no top-level disjunction that can be expanded. Therefore, we need to make a top-level disjunction.

To make a top-level disjunction, observe that in figure 9-3, there are disjuncts nested under the aggregator that could be used if we could bring them to the top of the R-expr. To accomplish this, I will introduce the idea of an optional constraint. An optional constraint is a constraint that we are allowed to be ignored or relaxed. In essence, an optional constraint is something that must be true but can be ignored when it is inconvenient (expensive to compute) or can be modified to make it easier to work with.

An optional constraint is denoted with opt(•) and is defined using the following semantic definition:

## 11. $\llbracket \operatorname{opt}(\mathrm{R}) \rrbracket_{E}=$ if $\llbracket \mathrm{R} \rrbracket_{E}=0$ then (nondeterminstic choice of 1 or 0 ) else 1

From this semantic definition, we have that in the case of opt(0), we "allow" this $\mathbf{R}$-expr to nondeterministically represent the multiplicity of 1 or 0 . The reason we allow this nondeterminism with optional constraints is that optional constraints are never used on their own. Instead, they are used in the context of a larger $\mathbf{R}$-expr, where regardless of the semantic interpretation used ( 1 or 0 ), the larger R-expr's semantic interpretation will not change. In other words, if we find opt(0) in an R-expr, then we know that the larger context will look something like opt ( 0 ) $* \cdots * \cdots * 0$. Hence, we can rewrite opt ( 0 ) as 0 , stopping the evaluation early, knowing that some other conjunctive sub-R-expr in the R-expr would have eventually been rewritten as 0 .

Additionally, note that the $\mathbf{R}$-expr R wrapped contained in the optional constraint does not need to be a constraint, as being a constraint, with a multiplicity of at most 1 , is enforced by the optional constraint $\mathbf{R}$-expr.

The nondeterministic behavior of the optional constraint can be defined with the following rewrite rules:

[^75]$R \xrightarrow{75} R *$ opt $(R)$
$\operatorname{opt}(R) \xrightarrow{76}$ if(R,1,0) $\triangleright$ make optional constraint, non-optional
$\operatorname{opt}(0) \xrightarrow{77} 0 \quad \triangleright$ Short cut for rewrite rule 76 when zero
$\operatorname{opt}(R) \xrightarrow{78} 1$
$T * \operatorname{opt}(R) \xrightarrow{79} T * \operatorname{opt}(T * R)$
$\operatorname{opt}(\mathrm{R}) \xrightarrow{80} \operatorname{opt}(\mathrm{R}+\mathrm{S}) \quad \triangleright$ Weaken the constraint with any disjunct S
$\operatorname{opt}(R) \xrightarrow{81} \operatorname{opt}(\operatorname{proj}(X, R)) \quad \triangleright$ Weaken the constraint by ignoring any variable $X$
$\operatorname{opt}(R * S) \xrightarrow{82} \operatorname{opt}(R)$
$\triangle$ Introduce optional constraint
$\triangleright$ Ignore optional constraint $\triangleright$ Enable rewrites by reading from context $\triangleright$ Weaken the constraint by removing a conjunct

Observe that rewrite rules 76 to 78 are for solving the optional constraint. If we have opt(0), we are allowed to apply any of these rewrites nondeterminically. Rewrite rule 78 allows us to rewrite the optional constraint as 1 before we know the multiplicity of $R$. This allows us to ignore the optional constraint entirely; hence, it does not affect containing R-expr's multiplicity. Rewrite rule 76 turns the optional constraint back into a non-optional constraint using the if-expression. The if-expression can be removed in the case that $R$ is a constraint using rewrite rule 83.

$$
\text { if }(R, 1,0) \xrightarrow{83} R \quad \text { if } R \text { is a constraint }
$$

Optional constraints can be introduced from any $\mathbf{R}$-expr using rewrite rule 75. As we can see in rewrite rule 75, the optional constraint is conjunctive with the $\mathbf{R}$-expr $R$ which will enforce the multiplicity of the $\mathbf{R}$-expr, meaning the optional constraint is redundant. As a brief reminder of what I said at the beginning of the chapter, the optional constraint does not "actually exist". Rather, optional constraints and rewrites on optional constraints serve as a license for complicated sequences of rewrites that we will see in chapter $\S 11$. As such, we do not have unproductive rewriting cycles like $R \xrightarrow{75} R *$ opt $(R) \xrightarrow{78} R * 1 \xrightarrow{13} R$ with optional constraints.

Optional constraints can also "read" from the surrounding context using rewrite rule 79 to copy in the conjunctive $\mathbf{R}$-exprT, allowing rewrites to be performed internally-just like the conditional in an if-expression with rewrite rule 65.

Finally, optional constraints can be weakened using rewrite rules 80 to 82. This means that the constraint will be true more often. For example, rewrite rule 80 adds in the additional disjunct $S$, which may be true (nonzero) in cases where $R$ is false. Similarly, rewrite rule 81 projects a variable, allowing it to ignore the value that is
assigned to some variable. To see an example of weakening an $\mathbf{R}$-expr constraint, consider the R-expr int $(X) *$ times $(3, X, Y)$, which defines a constraint on both $X$ and $Y$, as well as a bag relation on two variables. If we project out the variable $X$ from this $\mathbf{R}$-expr, we will have $\operatorname{proj}(X, \operatorname{int}(X) * t i m e s(3, X, Y)$ ). This $\mathbf{R}$-expr is only a constraint on the variable $Y$. It defines that $Y$ must be some multiple of the number 3, but otherwise has no influence on the variable $X .{ }^{122}$

Given this, we can think of an optional constraint as defining a valid upper bound on the support ${ }^{123}$ of an $\mathbf{R}$-expr. As such, the optional constraint never eliminates something that is true. They allow us to use useful constraints that are nested deep inside of an $\mathbf{R}$-expr.

### 9.2.1 Example Using Optional Constraints

To see how optional constraints can be used to split the domain of a problem, suppose that we have the following R-expr

$$
\begin{array}{r}
(A=\operatorname{sum}(Y,(X=1) *(Y=1)+ \\
(X=1) *(Y=3)+ \\
(X=2) *(Y=5)))
\end{array}
$$

Figure 9-4. R-expr with aggregation over a disjunction

We can create the optional constraint from any $\mathbf{R}$-expr in this expression. For this example, we are going to go ahead and create optional constraints from ( $X=1$ ) and ( $X=2$ ) using rewrite rule 75 as this will allow us to illustrate how optional constraints are usually used.

[^76]```
(A=sum(Y, (X=1)*(Y=1)*opt ((X=1))+
    (X=1)*(Y=3)*opt ((X=1))+
    (X=2)*(Y=5)*opt((X=2))))
```

Figure 9-5. Optional constraint introduced for $X$.

In figure 9-5, the optional constraints that were introduced are not too useful at first. We will use rewrite rule 80 to weaken these optional constraints with a disjunction.

```
(A=sum(Y, (X=1)*(Y=1)*opt ((X=1)+(X=2))+
    (X=1)*(Y=3)*opt((X=1)+(X=2))+
    (X=2)*(Y=5)*opt((X=2)+(X=1))))
```

Figure 9-6. Optional constraint weakened with a disjunct.

The weakening with the new disjunctive $\mathbf{R}$-expr can be done using any $\mathbf{R}$-expr. However, we have intelligently selected the $\mathbf{R}$-expr that we add in this case. Observe that now we have the constraint opt $((X=1)+(X=2))$ under all branches of the disjunction. Hence, we can now factor the optional constraint out of the disjunct using rewrite rule $22 .{ }^{124}$

$$
\begin{aligned}
(A=\operatorname{sum}(Y, \operatorname{opt}((X=1)+(X=2)) * & ((X=1) *(Y=1)+ \\
& (X=1) *(Y=3)+ \\
& (X=2) *(Y=5))))
\end{aligned}
$$

Figure 9-7. Optional constraints from figure 9-6 factored out of the disjunct.

Because the optional constraint in figure 9-7 does not interact with $Y$, we can further lift it out of the aggregator to the top of the R-expr. We can further use

[^77]rewrite rule 76 and rewrite rule $83^{125}$ to convert the optional constraint back into a normal constraint.
\[

$$
\begin{aligned}
& ((X=1)+(X=2)) *(A=\operatorname{sum}(Y,((X=1) *(Y=1)+ \\
& (X=1) *(Y=3)+ \\
& (X=2) *(Y=5))))
\end{aligned}
$$
\]

Figure 9-8. The optional constraint has been lifted to the top of the $\mathbf{R}$-expr and turned into a normal constraint.

From this point, we use the distributive rewrites on figure 9-8 to expand the R-expr and handle each of these cases separately. For this example, we expand this out using the distributive rewrite and have the case where $(X=1)$ and $(X=2)$.

$$
\begin{aligned}
(X=1) *(A=\operatorname{sum}(Y, & (X=1) *(Y=1)+ \\
& (X=1) *(Y=3)+ \\
& (X=2) *(Y=5))))+\quad \rightarrow^{*} \begin{array}{l}
(X=1) *(Y=4)+ \\
(X=2) *(Y=5)
\end{array} \\
(X=2) *(A=\operatorname{sum}(Y, & (X=1) *(Y=1)+ \\
(X=1) *(Y=3)+ & \\
(X=2) *(Y=5)))) &
\end{aligned}
$$

Figure 9-9. Aggregator evaluated and rewritten.

Admittedly, the use of the distributive rewrite in figure 9-9 goes against our philosophy of using the distributive rewrite (rewrite rule 22) to split the R-expr into $(X=1)$ and ( $X=2$ ) cases. As such, we can see that the "only factor" philosophy is not sufficient in all cases. In section $\S 11.7$, I will continue to build on optional constraints so that an $\mathbf{R}$-expr can be split into a disjunction of finitely many smaller R-exprs where each will be individually handled.
Solving the SAT formula To solve the SAT formula at the beginning of this chapter (figure 9-2), the system will use the same lift nested boolean assignments trick. The SAT formla contains $\mathbf{R}$-exprs of the form (true=exists(true, ( $A=1$ ) $+(A=0)$ )).

[^78]Using optional constraints, the constraint $(A=1)+(A=0)$ can be lifted to the top of the R-expr, eventually resulting in (true=exists(true, $(A=1)+(A=0))) *((A=1)+(A=0))$. The $(A=1)+(A=0)$ can be used to split the $\mathbf{R}$-expr, and perform case analysis (propagation of assignments and further branching).

### 9.2.2 Can Optional Constraints Solve all Disjunctive R-exprs?

No. Optional constraints must represent a valid upper bound on the support of an R-expr. They accomplish this by lifting up existing constraints through disjunctions and aggregations. Unfortunately, sometimes even the tightest upper bound is useless. For example, consider the program in figure 9-10.


Figure 9-10. The lack of a constraint on line 398 prevents us from creating a useful optional constraint for the variable X .

Because of line 398, the upper bound on the support for values of $X$ is any value. The reason is that without a constraint on X , line 398 always matches. Therefore, any optional constraint derived from a/1 will be always be true, represented as opt(1). ${ }^{126}$ If we have an optional constraint that is always true combined with any other optional constraint through a disjunction, then the resulting optional constraint is also going to be always true. In this case, the rewrite sequence is opt $(1+\cdots) \xrightarrow{76} \mathrm{if}(1+\cdots, 1,0) \xrightarrow{63} 1$.

As such, the best we can do in the case of a query $a(X)$ ? against the program in figure 9-10 would be to return the $\mathbf{R}$-expr in figure 9-10 (b).

[^79]
## Chapter 10

## Memoization, Reactivity, Cycles, and Updates

This chapter covers how we implement memoization, updates, and cyclic programs with R-exprs. For this chapter, I only assume the implementation described in chapter $\S 8$. Therefore, I will defer some of the discussion around being efficient till section §11.6.2, and instead focus on the fundamental ideals of memoization as it pertains to R-exprs and rewriting.

I should note ahead of time that memoization, updates, and cyclic programs were previously studied by Nathaniel Filardo in work on the Dyna project. As such, I recommend that anyone interested in the memoization topic in general should also read Filardo's dissertation [66]. In [66], Filardo developed theoretical algorithms that allowed the values associated with ground terms to be updated in different ways using message passing given a known computation graph. Filardo's work focused on making recomputation efficient and allowed for updates to be processed at different points in the algorithms running. ${ }^{127}$ In this chapter, I will instead focus on the details necessary to introduce memoization into an R-expr-based rewrite system. The system presented here uses a more complex representation for memos in that we allow R-exprs to be memoized vs. Filardo's algorithm, which only supports ground values. Because we are memoizing R-exprs, we support memos of non-ground terms and partially evaluated computation (represented

[^80]as an R-expr)-a capability that we have been unable to find elsewhere in the memoization literature. Given that the focus here is on the implementation of memos and R-exprs, the kinds of update messages that our system currently supports are much more limited than what Filardo's algorithm supported. We will only use invalidation messages that cause our system to run a recomputation. Hence, the approach for handling updates may be conceptually closer to a simple implementation of reactive programming [14]. Future work may wish to investigate how the different message types supported by Filardo can be integrated into the R-expr-based memoization presented here.
Chapter Outline The presentation in this chapter is ordered as follows. I will start by reviewing the necessary background on memoization and its procedural implementation. Then, I will transition to discussing the same abstract ideas of memoization as they apply to R-exprs by focusing on "simple" Dyna programs without external or internal updates. ${ }^{128}$ This means that the result of any computation will not change, allowing us to temporarily ignore the issue of sending invalidation messages (cache invalidation). Once we have developed the necessary background on unchanging Dyna programs with memos, I will introduce the additional mechanisms that are needed to handle external ${ }^{129}$ and internal updates (section §10.4). Finally, once we have developed the necessary background on the ideas underlying memoization and updates, I will present further clarifications about how this is actually implemented with $\mathbf{R}$-exprs (section $\S 10.6$ ), how updates are handled (section $\S 10.7$ ) and how $\$ \operatorname{memo}(\cdot)^{130}$ is implemented and used to control memoization (section §10.8).

In this chapter, you will notice that the memoization approach presented here is capable of memoizing any R-expr expression. However, the \$memo(•) memoization control mechanism defined in section $\S 2.7$ is only capable of defining memoization policies for user-defined terms. The discrepancy between \$memo(•) and this chapter is not a mistake. The reason is that controlling the "full power" of memoization

[^81]mechanisms presented in this chapter is difficult and requires direct manipulations of $\mathbf{R}$-exprs. Instead, our implementation of Dyna provides automated manipulations of $\mathbf{R}$-exprs that are limited to adding memoization to user-defined terms and are controlled using \$memo(•) (these "automated manipulations" will be discussed in section §10.8). Future work may consider extending \$memo(•) to expose more control or may consider researching automatic control of the memoization mechanisms presented here (section §16.4).

### 10.1 What is Memoization?

Let us review what memoization is before we jump into how it will work with $\mathbf{R}$-exprs. Memoization is a common concept that is taught to virtually all programmers. Hence, most readers will probably already have some preconceived notion of what memoization is. In this section, we are going to distill memoization down to a few simple conceptual elements. Then, we will look for those same elements in our R-expr rewrite system.

Memoization is a technique to avoid redundant computation by storing in memory a memo, which is the result of a computation, and retrieving it later to avoid redoing the same computation. The technique of storing memos was in fact called "Machine Learning" in the original memoization paper by Michie in 1968 [104]. To be able to store and later retrieve a memo, there needs to be a "signature" of the computation. ${ }^{131}$ Ideally, the signature should be the same every time that the same computation is performed-however, this is not a requirement, in which case a memoized algorithm will redo work and be less efficient but is still "correct". Finally, we need to modify the program to "intercept" relevant calls to the computation with a check to see if an equivalent computation was previously done and intercept the returned result of a computation so that it can be retrieved later.

[^82]
### 10.1.1 Example of Memoization in a Procedural Programming Language

I have chosen to start this discussion with an example in Python, a procedural programming language. The reason for this is that it will allow us to discuss the procedural steps required for memoization that we need to replicate with $\mathbf{R}$-exprs.

To illustrate this, we start with the Fibonacci program in figure 10-1. The Fibonacci program was chosen as it is an iconic example used when discussing memoization. The Fibonacci sequence is defined as the sequence (computed by the function in figure 10-1) where each number in the sequence is the sum of the previous two numbers: $0,1,1,2,3,5,8,13,21,24,55, \ldots$.. When the previous two numbers in the sequence are available, the next number in the sequence can be efficiently computed in $O(1)$ time. However, if the Fibonacci program is written such that every value is recomputed entirely from scratch each time it is used, then the program will take $\Omega\left(2^{n / 2}\right)$ time ${ }^{132}$ to compute the $n$-th element in the sequence.

```
def fib(N):
    if N == 0:
        return 0
    elif N == 1:
        return 1
    else:
        return fib(N-1) + fib(N-2)
```

Figure 10-1. Fibonacci program written in Python so that we can discuss the procedural processes of how memoization is implemented. On line 407, the Fibonacci function calls itself twice. Hence, if we run Fibonacci as written here, it will take $\Omega\left(2^{n / 2}\right)$ time to compute the result.

To add memoization to the Python fib function (line 401), we first need to identify the "signature", which sufficiently identifies the computation to be performed. In this case, we can use the argument for the fib function N. Note: The arguments to

[^83]the function are sufficient in the case that the function is a pure function ${ }^{133}$, as here, meaning that the result of the function is entirely determined by its arguments. In general, Python does not require that functions are pure, so the arguments might not be a sufficient signature, and it is a Python programmer's responsibility to ensure that a memoized function is pure. ${ }^{134}$ Dyna automatically tracks the necessary dependencies and changes for a memo without having to concern the Dyna developer further.

Now that we have identified the signature of the Fibonacci function as its argument, we need to create some "global" storage associated with the function where the results of the computation will be stored.

Finally, we need to check upon being called if the requested computation was previously performed and retrieve it from storage. Or, if the computation was not previously performed, do the computation and save it to storage for later use.

Adding this to the Python program from figure 10-1, we get the program shown in figure 10-2.

[^84]```
def fib_original(N):
    if N == 0:
        return 0
    elif N == 1:
        return 1
    else:
        return fib(N-1) + fib(N-2)
memoized_values = {}
def fib(N_signature): # override fib, so it is now memoized
    if N_signature in memoized_values:
        return memoized_values[N_signature]
    r = fib_original(N_signature)
    memoized_values[N_signature] = r
    return r
```

Figure 10-2. The Fibonacci program in Python with memoization manually added by the programmer. On line 416, a global hash map is created that is used to hold the memoized values. The hash map is checked at the top of the memoized fib function on line 418 before the original fib_original is called (performing computation). In the original fib_original function, the memoized version of the function is called on line 414 as we have overridden the fib function. After the original fib_original function performs the computation, the result is saved by the memoizing version of fib on line $421 .{ }^{135}$

### 10.1.2 The Facets of Memoization

Taking a step back from the Fibonacci example, we can see that there are four things that we need to identify when designing memoization for $\mathbf{R}$-exprs:

1. A location (in memory) for a data structure to store a memo (short for memoized result of computation). This can either be a data structure that exists globally for the entire running of the program (as in figure 10-2), or a data

[^85]structure that only exists within the context of a single computation (e.g. within the context of a single query handled by SimplifyNormalize).
2. A signature to identify the computation requested. The signature must be sufficient to identify the computation, otherwise the memo will be incorrect. ${ }^{134}$ The signature should occur many times, otherwise there will be no advantage when memoizing the computation under an ineffective signature. Usually, the signature is the arguments to a function; however, a backed off version of the arguments could also be used. For example, if we have the arguments $x=-3$, then we could choose to memoize for the negative interval $x<0$, as a backed-off version (section §10.2.3).
3. A way to intercept the calls to function and identify the signature of what computation is requested. This is done to check the data structure that contains the stored results before attempting to perform further computation.
4. A way to store the result of a computation so that it can be reused later. Generally, in a language like Python, the result is the value returned from a function (as on line 421); however, as we will see, this can be generalized with R-exprs.
5. A way to control what is memoized. In the case of Python, this is how the programmer has modified their program to enable memoization (as in figure 102). In Dyna, the "program" itself does not require any changes. Instead, we will control memoization using \$memo.

By the end of this chapter, we will be able to identify these same elements (and more) in R-exprs-based memoization.

### 10.2 First Steps Towards R-expr Memoization

Given the requirements of memoization in section $\S 10.1 .2$, let us identify where these requirements exist in the context of $\mathbf{R}$-exprs. To keep things "simple" in this section, I will only focus on creating simple memos, and I will defer modifications, cyclic programs, internal \& external updates, and change propagation until section §10.4.

$$
\begin{aligned}
& \mathrm{m}(\mathrm{X}, \text { Res }) \rightarrow \quad \triangleright \text { user-def' with memo } \\
& \text { if }((X=1)+(X=2), \quad \triangleright \text { Disjunction of signatures } \\
& \left((X=1) *(\text { Res }=5)+\quad \triangleright \text { The memos }{ }^{136}\right. \\
& { }_{423} \mid m(X)=2 * X+3 . \\
& \text { (a) Simple Dyna rule. } \\
& \text { ( } X=2 \text { ) * (Res=7) ), } \\
& \text { (Res=only(Inp, } \triangleright \text { Original user-defined } \mathbf{R} \text {-expr } \\
& \text { proj(Tmp, } \\
& \text { times(2,X,Tmp)*plus(Tmp,3,Inp))))) }
\end{aligned}
$$

Figure 10-3. Simple R-expr with the memoized values for 1 and 2 stored inside of the $\mathbf{R}$-expr. X is the argument and Res is the returned result of the user-defined R-expr.

Now, the goal of memoization is to avoid redoing computation. Computation, in the $\mathbf{R}$-expr setting, are rewrites performed by $\operatorname{Simplify}(R, \mathcal{C})$ against an $\mathbf{R}$-expr, to get another semantically equivalent $\mathbf{R}$-expr.

As such, a first attempt would be to apply memoization to the $\operatorname{Simplify}(R, \mathcal{C})$ function. Unfortunately, memoizing Simplify does not work that well. The reason is that the arguments for Simplify are both a large $\mathbf{R}$-expr R and the entire context $\mathcal{C}$, which contains many irrelevant details. Therefore, a signature match against $R$ and $\mathcal{C}$ would be inefficient and ineffective. Furthermore (and more importantly), memoizing Simplify does not work with handling internal or external updates, as will be discussed in section §10.5.3.1.

Rather, we are going to create memos for sub-R-exprs using an approach based on the previously introduced if-expression kind $\mathbf{R}$-expr (section §5.2.2.8). Recall that given any $\mathbf{R}$-expr $R$, for which we want to create memos for, the system can use rewrite rule 66 to introduce an if-expression as $R \xrightarrow{66}$ if $(S, R, R$ ) (as the same $\mathbf{R}$-expr $R$ is returned regardless of the $S$ that the system picks). As we will see, $S$ is the signature of the memo. I will get into details of how $S$ is selected later, but for now, we will just assume that $S$ is selected to be "useful for memoization".

Once we have this R-expr, the system can use rewrite rule 67 to rewrite the

[^86]true branch as if $(S, R, R) \xrightarrow{67}$ if $(S, S * R, R)$ provided that $S$ is a constraint. We can perform $S * R \rightarrow{ }^{*}$ S*RMemo using Simplify. ${ }^{137}$

The $\mathbf{R}$-expr now has a structure like if ( $S, S * M e m o, R$ ). A more complete example of if(S,S*Memo,R) R-expr is shown in figure 10-3. Let us look at this R-expr and check which facets of memoization are satisfied.

First, memoization requires that there is some location in memory where the result of a computation is stored. Here, RMemo is the result of Simplify applied to $R$ in the context of $S$. Additionally, RMemo is stored inside of the $\mathbf{R}$-expr if ( $S, S * R M e m o, R$ ).

Second, there needs to be a signature of the computation. The signature needs to be checked and return the memo in the case it matches, or bypass the memo and perform computation in the case the memo does not exist. Here, S is the signature of the computation. To see that $S$ is indeed the memo's signature, let us consider the third requirement of memoization, that we can intercept "calls" to compute and instead return the memoized result. The way in which we perform computation with $\mathbf{R}$-exprs is by rewriting an $\mathbf{R}$-expr with Simplify in the context of other conjunctive $\mathbf{R}$-exprs tracked via the context $\mathcal{C}$. For example, suppose that we have the $\mathbf{R}$-expr $R$, which is being rewritten in the context of $Q$. Hence, the system is rewriting the $\mathbf{R}$-expr $Q * R .{ }^{138}$ Now, let us consider a version $R$ with a memo. The $\mathbf{R}$-expr $R$ with a memo has the form if ( $S, S * R M e m o, R$ ) (where the if-expression was introduced using rewrite rules 66 and 67 ). Now, when rewriting if ( $S, S * R M e m o, R$ ) in the context of $Q$, the system is rewriting the $\mathbf{R}$-expr $Q * i f(S, S * R M e m o, R)$. Using rewrite rule 65, the if-expression's conditional is allowed to "read" the context $Q$, resulting in the $\mathbf{R}$-expr $Q * i f(Q * S, S * R M e m o, R)$. If the signature $S$ matches the context, then there will exist a sequence of rewrites such that $Q * S \rightarrow^{*} Q *(1+T)$ for some $T .{ }^{139}$ This means that the true-branch of $\mathrm{Q} *$ if $(\mathrm{Q} * \mathrm{~S}, \mathrm{~S} * \mathrm{RM}$ momo, R ) will be returned ( $\mathrm{Q} * \mathrm{if}(\mathrm{Q} *(1+\mathrm{T}), \mathrm{S} * \mathrm{RMemo}, \mathrm{R})$ $\xrightarrow{65,63} \mathrm{Q} * S *$ RMemo). Hence, we have intercepted the "call" to $R$ and substituted in the memo. Retrieving the memo was previously called Lookup in Filardo's work [66, 67].

Similarly, if $Q$ is not memoized, then $Q * S \rightarrow^{*} 0$, causing the false-branch of the

[^87]if-expression to be returned: $\mathrm{Q} *$ if $(\mathrm{Q} * 0, \mathrm{~S} * \mathrm{RM}$ emo, R$) \xrightarrow{64} \mathrm{Q} * \mathrm{R}$. This corresponds with falling back to the original computation (known as COMPUTE in Filardo's work [66, 67]).

This way in which this R-expr-based if-expression works can be seen as conceptually the same as the if-expression in the Python program (line 418). ${ }^{140}$ When the signature in the Python program matches, the true branch, which reads from the hash-table, is used. If the signature fails to match, then the false branch is used, which falls back to the original definition. However, by representing the signature $S$ and the query $Q$ as $\mathbf{R}$-exprs, we can memoize non-ground relations and condition memos on general expressions rather than ground terms used as keys for a memo table.

### 10.2.1 Advantages of Homogeneity

The advantage of a homogeneous system which represents all state as $\mathbf{R}$-exprs shows through in our representation of memos.

For example, in the procedural Fibonacci example (section §10.1.1), the signature of a memo was the argument to the function-which is an integer value like the number 7. Using a value like this is quite typical of memoization. However, with our R-expr if-expression representation, we can further generalize memos to use any executable $\mathbf{R}$-expr as the signature. For example, we can choose to memoize all negative values by using lessthan $(X, 0)$ as a signature. Admittedly, a programmer could manually introduce the relevant test and if-expression in another programming language. However, the R-expr backed formalism allows us to do this automatically.

A second advantage of homogeneity is that the result returned from the memo is an $\mathbf{R}$-expr. This means that we are not limited to memoizing values, but instead, we can memoize partially completed computation. For example, suppose that we have the sum3 rule defined in figure 10-4 that sums up its three arguments.
${ }^{140}$ Additionally, note that accessing the memo contained in the if-expression required no modifications to Simplify. In this way, Simplify is comparable to a bytecode interpreter, and the addition of the if-expression to the $\mathbf{R}$-expr can be considered a modification to the program, just like how we modified the Python program (not the Python interpreter).
${ }_{424} \mid \operatorname{sum} 3(A, B, C)=A+B+C . \quad \operatorname{proj}(T m p, \operatorname{plus}(A, B, T m p) * p l u s(T m p, C, R e s u l t))$
(a) Sum3 in Dyna
(b) Body of Sum3 as an R-expr

Figure 10-4. Rule defining sum of its 3 arguments.

In memoizing sum 3 with only 2 of its three arguments (say A and B), we will still have one plus $\mathbf{R}$-expr remaining. The memo can memoize the intermediate sum of $A+B$ but must still represent the remaining plus as shown in figure 10-5.

```
if((A=1)*(B=2), (A=1)*(B=2)*plus(3, C, Result),
    proj(Tmp, plus(A,B,Tmp)*plus(Tmp,C,Result)))
```

Figure 10-5. sum 3 with a memo created where only two of its arguments ( $A, B$ ) are known. Computation to compute $1+2=3$ was performed, leaving an $\mathbf{R}$-expr which represents the computation $3+C$ that still needs to be completed. Admittedly, in this example, we are not saving that much "work" by memoizing $1+2$, however, in general, we would memoize a computation that takes more time to complete.

### 10.2.2 Persisting Memos to Make them Globally Usable

In our design, so far, a partial memo is represented as another R-expr, namely $R \xrightarrow{66,67}$ if $(S, S * R, R)$. This is a very general approach to creating and storing a memo, as there always exists some $\mathbf{R}$-expr, and therefore the system can create a memo whenever it wants. However, "whenever it wants" is a double-edged sword, in that some points where a memo could be created may be useless. For example, suppose that a memo is created inside of an R-expr that was created to evaluate a user's query (section §2.3). In this case, the memo is usable only within the context of that query. The R-expr with the memo does not persist within the Dyna system after the query has been completed.

Ideally, a memo should be usable throughout the entire running of the programmeaning that it should persist between queries initiated by the user. This can be accomplished by carefully choosing when and where a memo is constructed. Recall from chapter $\S 7$ that the user's program is defined in terms of user-defined named R-exprs, recreated in figure 10-6 for convenience.

$$
\begin{array}{ll}
{ }_{425} \mid f(X)+=X * X . & f(X, \operatorname{Res}) \xrightarrow{74}(\text { Res }=\operatorname{sum}(\text { Inp }, \\
\operatorname{times}(X, X, \text { Inp })))
\end{array}
$$

(a) Dyna
(b) R-expr

Figure 10-6. Recall how Dyna is translated to R-exprs, chapter $\S 7$.

Every time that a user-defined $\mathbf{R}$-expr is referenced, it indirects through rewrite rule 74 . This means that if we modify rewrite rule 74 , then the resulting $\mathbf{R}$-expr will be accessible globally. The resulting $\mathbf{R}$-exprs will look something like figure 10-7.

```
f_original(X,Res) -> (Res=sum(Inp, }\quad\mathrm{ Original R-expr (like fib_original
    times(X,X,Inp))) in figure 10-2)
f(X, Res) }\xrightarrow{}{74}\mathrm{ if((X=3), }\triangleright\mathrm{ Disjunction of all signatures
    (X=3)*(Res=9), \trianglerightMemo
    f_original(X,Res)) \trianglerightCall to original definition
```

Figure 10-7. The user-define f R-expr with a memo for the value of 3.

### 10.2.3 Things to Consider When Choosing a Signature for Memoization

So far, we have that the memo is contained inside of the $\mathbf{R}$-expr that user-defined R-exprs are rewritten as. However, we need to discuss how the system determines what to memoize. More concretely, we need to pick the signature $S$ used by rewrite rule 66 when introducing an if-expression ( $R \xrightarrow{66,67} \operatorname{if}(S, S * R, R)$ ). The common approach we saw with the Python example in section $\S 10.1 .1$ is to wait until the function is called with an unmemoized query and then use the value that was used to check the memo table as the signature (as in figure 10-2 lines 418, 420 and 421). However, the Python program hard coded that the signature was the argument of the Fibonacci function. Whereas with R-exprs, we have more flexibility in choosing the signature as we are allowed to choose any R-expr, and we do not have to commit to a particular $S$ until just before the memo is created by rewrite rule 66. Further complicating this, the choice of $S$ will have a significant impact on
the usefulness and efficiency of our memo.
As an example of choosing a signature, suppose that we are performing a query $Q$ against the $\mathbf{R}$-expr $R$. If the system chooses $S$ as having no relation to $Q$, then it is likely that $S$ and $Q$ will be incompatible (e.g. $Q * S \rightarrow^{*} 0$ ). This means that spending time constructing a memo using a bad signature $S$ would actually delay the computation with useless rewrites: $Q * R \rightarrow Q * i f(S, S * R, R) \rightarrow Q * i f(S, S * R M e m o, R) \rightarrow Q * R$.

This gives us our first hint of how to pick a signature. A useful signature must have a non-empty intersection with the query $Q$.

A second thing to consider when choosing a signature is the amount of reuse we will get from a memo. A signature that is more general can be used in more cases. For example, suppose that we are making the query lessthan $(5, W) *(X=1) *(Y=2) *(Z=3)$ against the $\mathbf{R}$-expr $R$. We can choose to use the signature $(X=1) *(Z=3)$, ignoring lessthan $(5, W) *(Y=2)$. This will mean that we get an $\mathbf{R}$-expr with a memo like if $((X=1) *(Z=3),(X=1) *(Z=3) * R M e m o, R)$. Now, if we get a second query $(X=1) *(Z=3)$ * $(Y=7)$ against this $\mathbf{R}$-expr, we can reuse $R M e m o$ as the signature $(X=1) *(Z=3)$ matches this query.

A third consideration when choosing the signature $S$ is the data structure that will be used to store the memo. For example, ground assignments to variables (e.g. $(X=1) * R 1+(X=7) * R 2$ ), can be represented efficiently using a hash table ( $\mathrm{sec}-$ tion §11.6.1.2). Whereas a signature like $\operatorname{int}(X) *$ lessthan $(0, X) *$ lessthan $(X, 10)$ $+\operatorname{int}(X) *$ between $(X, 15,18)+\cdots+$ int $(X) *$ lessthan $(321, X) *$ lessthan $(X, 325)$ requires that the system scan through all of the lessthan constraints to check if any of them match the current query.

Finally, I note that when selecting the signature S , the system has access to the context $\mathcal{C}$ for the current query. This means that the system can efficiently retrieve conjunctive R-exprs when accessing the memo. Hence, the signature $S$ will be chosen as a relevant subset of $\mathcal{C}$ that satisfies the considerations listed above.

### 10.3 Example: Memoization of an R-expr

Let us use the representation in memoization discussed so far to work through an example on the Fibonacci program, this time represented as a Dyna program and an $\mathbf{R}$-expr in figure 10-8. Note that the $\mathbf{R}$-expr here is semantically equivalent to
the Dyna program, though not identical to the $\mathbf{R}$-expr that would result from the mechanical translation of Dyna to R-exprs (as defined in chapter §7).

```
fib(0) += 0. fib(N, Res) -> (Res=sum(Inp,
fib(1) += 1. (N=0)*(Inp=0)+
fib(N) += fib(N-1) (N=1)*(Inp=1)+
    for N > 1. proj(Tmp, lessthan(1,N)*plus(Tmp,1,N)*
fib(N) += fib(N-2) fib(Tmp,Inp))+
    for N > 1. proj(Tmp, lessthan(1,N)*plus(Tmp,2,N)*
(a) Dyna
```

```
    fib(Tmp,Inp))))
```

```
    fib(Tmp,Inp))))
```

(b) R-expr

Figure 10-8. Fibonacci program shown as a Dyna program and R-expr.

First, suppose that a query is initiated against the fib $\mathbf{R}$-expr. For example, this might be fib(5,Res) where the system is computing the Fibonacci value for the number 5. The Fibonacci $\mathbf{R}$-expr will be simplified by expanding the user-defined sub-R-expr until it reaches the base cases of $(N=0)$ and $(N=1)$. At this point, the system can identify ( $\mathrm{N}=0$ ) and ( $\mathrm{N}=1$ ) as two queries made against the Fibonacci relation. Both are combined into a signature to indicate that the system will memoize the values for both $(\mathrm{N}=0)$ and $(\mathrm{N}=1)$. The resulting $\mathbf{R}$-expr is shown in figure 10-9:

```
fib(N, Res) -> if((N=0)+(N=1),
    ((N=0)*(Res=0)+ \trianglerightThe memo
    (N=1)*(Res=1) ),
    (Res=sum(Inp, \trianglerightThe original R-expr
        (N=0)*(Inp=0)+
        (N=1)*(Inp=1)+
        proj(Tmp, lessthan(1,N)*plus(Tmp,1,N)*
        fib(Tmp,Inp))+
        proj(Tmp, lessthan(1,N)*plus(Tmp,2,N)*
                fib(Tmp,Inp)))))
```

Figure 10-9. Fib R-expr with memo for 0 and 1

Next, the value for $\mathrm{fib}(2, R e s)$ can be computed and stored. The current disjunc-
tion of the signatures $(N=0)+(N=1)$ indicates that the $(N=2)$ value is not currently memoized, ${ }^{141}$ and therefore will fall back to the original $\mathbf{R}$-expr. A new memoizing if-expression is nested under the false-branch of the original memo. This is consistent with the idea that we are creating memos on any $\mathbf{R}$-expr, rather than modifying the existing memo. The resulting R-expr is shown in figure 10-10.

```
fib(N, Res) -> if((N=0)+(N=1),
    ((N=0)*(Res=0)+ \trianglerightThe previous memos
        (N=1)*(Res=1) ),
    if((N=2), }\quad\triangleright\mathrm{ The new memo nested under the false-branch of the first memo
        (N=2)*(Res=1),
        (Res=sum(Inp, }\quad\mathrm{ The original R-expr deeply nested under the false-branches of all
        memos
            (N=0)*(Inp=0)+
            (N=1)*(Inp=1)+
            proj(Tmp, lessthan(1,N)*plus(Tmp,1,N)*
                    fib(Tmp,Inp))+
            proj(Tmp, lessthan(1,N)*plus(Tmp,2,N)*
                        fib(Tmp,Inp))))))
```

Figure 10-10. Fib R-expr with a new memo for ( $\mathrm{N}=2$ ), which was introduced using rewrite rules 66 and 67. The newly added memo is under a second if-expression, and does not modify the previously created memo for $(\mathrm{N}=0)+(\mathrm{N}=1)$.

Observe that falling through several if-expressions will be inefficient when there are a large number of memoized $\mathbf{R}$-exprs and $\mathbf{R}$-expr signatures. Therefore, the system can use rewrite rule 69 to merge the if-expressions together, which avoids creating a nested chain of if-expressions:

[^88]```
fib(N, Res) -> if((N=0)+(N=1)+(N=2),
    ((N=0)*(Res=0)+ \trianglerightThe memos (now combined)
        (N=1)*(Res=1)+
        (N=2)*(Res=2) ),
    (Res=sum(Inp, \trianglerightThe original R-expr
        (N=0)*(Inp=0)+
        (N=1)*(Inp=1)+
        proj(Tmp, lessthan(1,N)*plus(Tmp,1,N)*
        fib(Tmp,Inp))+
        proj(Tmp, lessthan(1,N)*plus(Tmp,2,N)*
        fib(Tmp,Inp)))))
```

Figure 10-11. Fibonacci with if-expressions merged using rewrite rule 69

This process of creating nested if-expression for new memos and merging the if-expressions will continue all of the way up to fib(5,Res), which was initially requested by the user's query. Once the query fib( 5, Res) has been answered, the user-defined Fibonacci R-expr fib(N,Res) will only contain the memos for $0,1,2,3,4,5$, and the system will stop modifying the definition $f i b(N, R e s)$.

Throughout this entire process, the rewrite that defines the user-defined fib(N,Res) was having its right-hand side modified. However, the semantic interpretation of fib(N,Res) never changed. This is because all of the rewrites performed against fib's definition are semantics preserving. The only difference is that we are now clever about where and when we apply the rewrite rules from chapter $\S 6$.

### 10.3.1 Conclusion of Basic Memos

So far, we have developed a sufficient understanding to implement memoization in the case of unchanging programs. The if-expression can represent the memos inside of $\mathbf{R}$-exprs. If we think of the $\mathbf{R}$-expr itself as representing the program, then modifications of the $\mathbf{R}$-expr with the if-expression are conceptually equivalent to the modifications done to the Python program when adding memoization to the Fibonacci function (figure 10-2). ${ }^{140}$ With R-exprs, the true-branch of the if-expression represents the memo itself, the signature is used to indicate what is memoized, and the false-branch is used to represent falling back perform a computation using the original definition of a user-defined $\mathbf{R}$-expr.

```
\({ }_{432} \mathrm{e}+=1 . \quad 435 \mid \mathrm{a}+=1\).
433 e += e/2.
434 print e. \% prints 2
(a) A cyclic Dyna program which converge to \(\mathrm{e}=2\).
(b) Dyna program with an update (line 437) which modifies the program.
```

Figure 10-12. Example Dyna Programs which demonstrate rules (e and a) can change. Therefore, any downstream memoized values must be updated accordingly.

### 10.4 Handling Change

So far, we have focused only on programs where the result does not change, ignoring issues such as updates (cache invalidation). However, Dyna allows the rules to change. This can be the result of externally driven updates (such as new rules added at the REPL), or due to cycles in the program iterating until a fixed point (section §2.5, e.g. figure 10-12a). To properly handle this, we must track any memo that depends on a value that has changed and needs to be updated or invalidated. We accomplish this by adapting a design similar to that of other reactive programming languages and libraries [14].

### 10.4.1 Assuming Reads Never Change

First, we have that all values and expressions in the language which can change are protected by an assumption ${ }^{142}$ that the values have not changed. Some examples of changeable values include all the R-expr definitions of user-defined terms and the mutable memo tables (which I will define shortly).

An assumption (figure 10-13) is an object within the Dyna implementation that tracks the immediate downstream dependencies of a particular value. When a value is changed, its associated assumption is invalidated and will be replaced with a new assumption for the new value. An invalidated assumption sends notification messages to all downstream dependencies that it has been invalidated. An invalid assumption can never become valid again. This design was made with the intention

[^89]```
final class Assumption {
    private boolean isValid = true; // Starts valid
    private Set<MessageListener> subscribers;
    public void subscribe(MessageListener); // Dependents are subscribed
    public void sendMessageToAll(Message); // Send to all subscribers
    public boolean getIsValid() { return isValid; }
    public void invalidate() { isValid = false; // Can only invalidate
                    sendMessageToAll(..); }
}
```

Figure 10-13. Class design for an Assumption
that future work on parallel/concurrent processing can take advantage of this design to avoid race conditions.

When computation is performed in the Dyna implementation, assumptions that protect read values will be automatically subscribed to. ${ }^{143}$ Downstream dependencies are allowed to run "any code they want" when they receive a notification that an upstream assumption has become invalidated. This allows us to integrate assumptions and invalidation messages into many different aspects of Dyna implementation outside of memoization. Currently, assumptions are only used to trigger recomputation of memos, however future work may which to depend on this mechanism. ${ }^{144}$

Note: Memoized values can be subscribed to their own assumption. This happens in the case of a cyclic program as in figure 10-12a. This is not a problem as long as the memoized $\mathbf{R}$-expr eventually converges. ${ }^{145}$

[^90]
### 10.4.1.1 Updating Memoized R-exprs

The most common message that a subscriber will receive is a notification that an upstream R-expr has changed and its corresponding assumption has been invalidated. ${ }^{146}$ In this case, this means that there is an $\mathbf{R}$-expr like if (S, RMemoExisting, $R$ ) and the system needs to recompute RMemoExisting. This is done by starting again from the original $\mathbf{R}$-expr definition $R$, and combining it with the memo's existing signature $S$ to get $R * S$. The system rewrites $R * S$ using SimplifyNormalize, as we did before, and will get $S * R \rightarrow^{*}$ RMemoNew. In the process of creating RMemoNew, all reads performed, and assumptions depended on will be tracked. Finally, RMemoNew and RMemoExisting are compared with each other for "semantic equivalence". ${ }^{147}$ If the system can prove that RMemoNew and RMemoExisting are semantically equivalent, then we do not have to invalidate the assumption associated with the memo RMemoExisting and can stop propagating update messages forward. In the case that the system cannot prove that RMemoNew and RMemoExisting are semantically equivalent, it must replace RMemoExisting with RMemoNew, and then invalidate the assumption associated with RMemoExisting.

This process of receiving messages about upstream assumptions being invalidation and recomputing RMemo will continue until the system converges ${ }^{148}$ and all messages have been processed.

### 10.4.2 Example: Updating a Dyna Program

To see how assumptions work in the context of a program receiving updates, let us work through the example given in figure 10-12b.

[^91]First, we have the initial R-expr for ' $a$ ', and the system will have created a memo using the if-expression around the original R-expr. Furthermore, I have shown the valid assumptions as black boxes and the subscription from the assumptions as a blue dashed arrow in figure 10-14.


Figure 10-14. A memo created for the Dyna rule a. Assumptions are shown as boxes on the right-hand side, and the subscription to the "definition of ' $a$ '" is shown as a blue dashed line. In this example, the signature is always the $\mathbf{R}$-expr 1, meaning that the entire relation is memoized.

When the definition of 'a' is updated, the original $\mathbf{R}$-expr is modified, and the assumption that is associated with the definition of 'a' is invalidated. This is shown in figure 10-15.


Figure 10-15. The original R-expr (Dyna program) has been modified. The assumption that is associated with the user's definition has been invalidated, and the subscription to the memo needs to be processed.

This will trigger the memo to be recomputed and brought up to date with the new
value as defined by the R-expr constructed from the user's program as shown in figure 10-16.


Figure 10-16. The memoized value has been updated to $\mathrm{Val}=3$ by recomputing using the original R-expr. The Memo now depends on the new valid assumption "' a' User-Def \#2". The old assumption for the memo has been invalidated and replaced with a new assumption. Anything that was subscribed to the old assumption will have received a notification of its invalidation.

### 10.5 Handling Cyclic Programs

So far, in section $\S 10.4$, I have glossed over the details of how the computation for creating a memo is performed. In general, SimplifyNormalize (algorithm 1) is invoked on the original user-defined $\mathbf{R}$-expr in conjunction with a query, and the resulting $\mathbf{R}$-expr is saved as the memo. With a non-cyclic program, we are guaranteed that the fully expanded $\mathbf{R}$-expr is bounded in size. Hence, SimplifyNormalize is able to fully rewrite the $\mathbf{R}$-expr.

Unfortunately, the same guarantee does not exist with cyclic programs. A cyclic program recurses back onto itself. For example, the definition of 'e' in figure 10-17 is cyclic as the term 'e()' depends on the term 'e()'. Conversely, the Fibonacci program from before is recursive, not cyclic, as terms in the computation graph do not depend on themselves, rather terms with the name fib(•) depend on other terms which also have the name fib(•).

When rewriting a cyclic R-expr, SimplifyNormalize will never terminate. The reason is that the $\mathbf{R}$-expr is expanded at every invocation of Simplify to increasing depths of the recursion.

$$
\begin{array}{l|l}
{ }^{453} & \mathrm{e}+=1 . \\
{ }^{454} & \mathrm{e}(\text { Val }) \rightarrow(\text { e } \mathrm{e} / 2 . \\
(\text { Inp }=1)+
\end{array} \quad \begin{array}{ll}
\text { (a) Dyna } & \operatorname{proj}(\text { Tmp }, \mathrm{e}(\text { Tmp }) \star \text { times(Tmp, Inp, 2)))) }
\end{array}
$$

(b) R-expr without Memos

Figure 10-17. The equation $e=1+\frac{e}{2}$ is represented as an $\mathbf{R}$-expr on lines 453 and 454. The rule 'e' depends on what is in both line 454 , and in the $\mathbf{R}$-expr, with $e(V a l)$ appearing on both the left and right-hand side of the rewrite rule.

### 10.5.1 Making Guesses

To avoid this issue, we must avoid expanding the same recursive call when an $\mathbf{R}$-expr is rewritten as depending on itself as a sub-R-expr, as in figure $10-17$ with e (Val). This is accomplished by placing a "marker" along the recursive calls to avoid performing the same query against a memo twice. This is akin to setting and checking a visited bit in a depth-first search algorithm.

Like in the depth-first search, when the query is re-encountered, the system avoids redoing the exact same computation by guessing the $\mathbf{R}$-expr result to a query. A guess is a memo that is created before any computation is performed. A guess can be any $\mathbf{R}$-expr, though in practice, we always guess an empty bag, hence the R-expr 0 . A guess initially depends on an invalid assumption; therefore, it must be recomputed/checked immediately after it is created. Checking a guess is crucial as it enforces the self-consistency of a value on a cycle. Furthermore, the process of guessing and then forward-chaining invalidation messages until convergence is equivalent to Datalog-style forward chaining, though with $\mathbf{R}$-exprs rather than ground values (section §3.1.2).

I should note that guessing in the case of cycles is not novel to this dissertation and has appeared in prior work that influenced our design. First, XSB-Prolog [135, 145, 151] added memoization to standard Prolog style back-chaining by detection with marking. When a cycle is found, a guess is made. Filardo and Eisner [67] focused their work on finite circuits, which can contain cycles and allow guessing and forward-chaining messages to update memoized values (also see Filardo's dissertation [66]).

### 10.5.1.1 Example: Using Guessing with a Cyclic Program

To see how guessing works, let us work through an example of 'e' defined in figure 10-17. In figure 10-18, I have written the R-expr with an initial guess. The initial guess for ' $e$ ' is 0 , and the memo is subscribed to the invalid assumption. We also already have the guess protected by its own assumption "e Memo".


Figure 10-18. Initial memo being created as guess of 0 . The guess is subscribed to an invalid assumption, so the memo must be validated by recomputing the memo.

In the process of recomputing the memo for 'e' now, the recursive call to e(Tmp) is rewritten as 0 (as that is the memo). This means that only the contribution from the (Inp=1) (line 453) branch will be counted. Hence, the new memoized value is (Val=1), as shown in figure 10-19.

(a) R -expr
(b) Assumptions
(figure 10-13)
Figure 10-19. The memo for the cyclic 'e' after one iteration of recomputation. The memo ( $\mathrm{Val}=1$ ) depends on the old memoized R-expr of 0 . However, that R-expr was replaced, and its associated memo ("'e' Memo") has been invalidated. Hence, the system is required to again recompute the memo for ' $e$ '.

Unfortunately, the new memo in figure 10-19 depended on the old memo of 0 ,
which has since been invalidated. Hence, the system must again recompute the memo, as in figure 10-20.


Figure 10-20. The memo for the cyclic 'e' after two iterations of recomputation. The memo is now ( $\mathrm{Val}=1.5$ ), as the previous memoized value was 1 , and by the definition of line 454 we have defined this as $1+e / 2$. Just like before, we have that the memo of (Val=1.5) depends on the assumption of the previous read ("'e' Memo \#2") which has been invalidated as the memoized $\mathbf{R}$-expr was replaced. This process of recomputing the memo will continue until the memo is consistent with itself as in figure 10-21.


Figure 10-21. The memo has converged ${ }^{149}$ with (Val=2.0). When the system rechecks this memo, it constructs the same $\mathbf{R}$-expr, which can be checked for semantic equality. Hence, the memo is not changed, and the assumption "'e' Memo \# $n$ " is subscribed to by the same $\mathbf{R}$-expr it protects. The assumption remains valid, so no further propagation is required.

### 10.5.2 Choice of default Guesses

As stated earlier, we are allowed to use any $\mathbf{R}$-expr as our guess. The choice of initial guess can have a significant impact on the value to which the program converges and the speed at which the program converges. For example, we could have chosen to "guess" the R-expr (Val=2.0) ${ }^{150}$ for the 'e' definition in section §10.5.1.1. In which case the system would have been able to immediately validate that the guess (Val=2.0) is correct and avoid the iterative process of converging to the value of (Val=2.0).

That said, we have made the decision to make the initial guess always the empty $\mathbf{R}$-expr 0 . This choice was made in hopes of minimizing the amount of surprise that results from guessed values. As such, a user-defined term like ' $g$ ' in figure 10-24 would have "no value".

### 10.5.2.1 User Override for Initial Guesses

The choice of the initial guess being empty can be "overridden" by the user of Dyna if they are willing to introduce a small modification into their program. Using the := aggregator, we can set an initial value for when nothing exists. An example of this is shown in figure 10-22.

[^92]```
455 g_with_guess = 2*g - g**3. % g can either be 1 or -1
456 g := .5. % User defined guess
457 g := g_with_guess.
```

Figure 10-22. Users can override the default guess using the := aggregator. Here on line 456 , the default guess is set as .5. When g_with_guess is an empty R-exprwhich is the hard-coded default for a guess-the rule on line 457 will not contribute any value. Therefore, ' $g$ ' will take the value from line 456 . Once g_with_guess has some non-null value, then ' $g$ ' will take the value from line 457. This initial guess causes ' $g$ ' to converge to value 1 instead of -1 .

### 10.5.3 Guesses are Un-bypass-able

Guesses have additional requirements on them compared to "normal" memos from section §10.1.2. Most importantly, guesses are not bypassable. The reason is a guess can change the observed semantics of a Dyna program. The semantics of Dyna only require that the system find an assignment to all terms in the program that is consistent (section §2.5). In practice, there are many different consistent assignments to the terms in the program. For example, consider the program in figure 10-23, which has a cycle involving 'a' and 'b'.

```
458 a := 1.
4 5 9 ~ a ~ : = ~ b . ~
460 b := 2.
461 b := a.
4 6 2 ~ p r i n t ~ b . ~ \% ~ p r i n t ~ e i t h e r ~ 1 ~ o r ~ 2 ~ d e p e n d i n g ~ o n ~ g u e s s
```

Figure 10-23. Dyna program which prints 1 or 2 non-deterministically depending on where the back-chaining cycle is broken and how a guess is made. If the value of ' $a$ ' is guessed as null, then the value of ' $b$ ' will be set to 2 on line 460 . This causes 'a' to take on the value 2 as well, by the definition of the := aggregator, the last line of the program which is non-null for a rule defines its value, and in this case it will be line 459 is now non-null as ' $b$ ' is defined as 2 . Conversely, if ' $b$ ' is guessed as null, then ' $a$ ' will be set to 1 due to line 458 , therefore ' $b$ ' will also take on the value of 1 by the definition of the := aggregator and line 461.

The value which is returned by 'a' and 'b' can either be 1 or 2 depending on the order in which guesses are made. The Dyna specification says that either value is acceptable, as Dyna only requires that the assignment to all expressions is consistent.

A more extreme version of figure 10-23 when considering Dyna's requirement to make an assignment consistent is shown in figure 10-24.

$$
\left.{ }_{463} \mid \mathrm{g}=\mathrm{g} . \quad \mathrm{g}(\mathrm{Val}) \rightarrow(\text { Val=only(Inp, } \mathrm{g}(\mathrm{Inp}))\right)
$$

(a) Dyna
(b) R-expr

Figure 10-24. The Dyna rule ' $g$ ' can take on any value and still be consistent with its definition on line 463. Hence, whatever value is guessed for ' $g$ ' will persist.

Given the definition of rule ' $g$ ', it is "allowed" to take on any value. For example, if ' $g=77$ ', then we would have $77=77$ which is consistent. Similarly, ' $g$ ' can take on the value $\mathrm{g}=$ "hello", as "hello" = "hello", and is also consistent with g's definition.

### 10.5.3.1 Why Guessing Requires if-expressions

In section §10.2, I claimed that memoization involving cycles does not work if we are memoizing Simplify directly. The reason is that we need to ensure that we intercept user-defined $\mathbf{R}$-exprs with guesses. If we instead memoized Simplify, then we might encounter something that needs to be overridden by a guess but cannot be easily recognized due to the difficulty of comparing $\mathbf{R}$-exprs for semantic equivalence. Essentially, this means that we would either need to have a check for semantic equivalence between R-exprs (which is impossible in general because checking semantic equivalence is Turing-complete), or accept that guesses sometimes end up getting bypassed. Neither of these is acceptable, as the semantics of Dyna requires consistent set of assignments to the values associated with terms (section §2.5).

```
class MemoizationContainer implements MessageListener {
    final Rexpr original_rexpr; // The original R-expr (false if branch)
    Rexpr signature_have_memoized; // what is currently memoized
    final Rexpr what_want_memoized; // $memo representation
    Assumption assumption; // replaced when memo changes (§10.4.1)
    Rexpr the_memo; // The memo itself (true if branch)
}
```

Figure 10-25. The memoization container

### 10.6 Cleaning up the R-expr Presentation of Memoization-Memos Held Outside of the R-expr

So far, I have waved my hands when it comes to how memos are managed in an $\mathbf{R}$-expr. I have demonstrated that an if-expression can be used to override the original R-expr (as needed) for memoization, and that the system can store the memoized $\mathbf{R}$-expr inside the true-branch of the if-expression. However, we still lack details on how the $\mathbf{R}$-expr is modified and how the relevant assumptions are tracked with each R-expr.

In this section, I will lessen the abstraction and talk more specifically about how these different pieces of memoization fit together.

First, we will move the memo out of the $\mathbf{R}$-expr. To do this, we are going to replace the if-expression $\mathbf{R}$-expr and use a memoRead ( $\cdots$ ) R-expr to represent the memo. The memoRead $\mathbf{R}$-expr maintains a pointer/reference to the relevant memoization container and has a list of one or more local variable names: memoRead ( memo_container_pointer, $Y_{1}, Y_{2}, \ldots, Y_{n}$ ). The memoization container (figure 10-25) is a mutable data structure that contains the memo. The memoization container can be referenced by multiple memoRead $\mathbf{R}$-exprs at the same time. This allows all memoRead $\mathbf{R}$-exprs who access the same memo to have a consistent view of the currently memoized $\mathbf{R}$-expr. The fields on the memoization container are as follows:

1. The original R-expr.-This is used when something is configured to bypass the memo or is not current computed
2. The signature represented as an R-expr.-This corresponds with the conditional of the if-expression and records what is currently stored in the memo table.
3. An R-expr that tracks what we want to be memoized.-This is the \$memo(•) memoization control mechanism, which I will discuss in section §10.8.
4. An assumption that is associated with the current memo. Anytime the memo is changed, the assumption is invalidated, and messages are sent to all downstream dependents, as previously discussed in section §10.4.1.
5. The memo itself, represented as an R-expr.

The variables $Y_{1}, Y_{2}, \ldots, Y_{n}$ on the memoRead(memo_container_pointer, $Y_{1}, Y_{2}$, $\ldots, Y_{n}$ ) refer to local variables, which allows variable renaming in the $\mathbf{R}$-expr which uses memoRead as a sub-R-expr. Inside the memoization container, variable names are normalized to known variable names like $X_{1}, X_{2}, \ldots, x_{n}$. When memoRead is rewritten as an $\mathbf{R}$-expr returned from the memoization container, the returned $\mathbf{R}$-expr's variables are renamed to $Y_{1}, Y_{2}, \ldots, Y_{n}$ to match the local context.

### 10.6.1 Memoization Container

Comparing the memoization container with the if-expression, we can see that we maintain all the same capabilities as with the if-expression. The signature in the conditional test of the if-expression is held in the signature_have_memoized variable, the true-branch corresponds to the the_memo $\mathbf{R}$-expr, and the false-branch corresponds to the original_rexpr. The new addition in the memoization container is the what_want_memoized $\mathbf{R}$-expr variable.

The what_want_memoized $\mathbf{R}$-expr is the memoization policy, which is defined using \$memo. The reason we separate the policy from the signature of what is memoized is that sometimes the policy is not helpful when it comes to the memoized R-expr.

For example, consider the case where the signature is directly derived from the memoization policy. Suppose that our memoization policy says that the system should memoize "everything". In this case, the policy can be represented as the $\mathbf{R}$-expr 1. However, this is not a useful $\mathbf{R}$-expr when it comes to our rewrite rules. Recall that we can introduce an if-expression using any R-expr (which represents
our memoization policy, rewrite rules 66 and 67 ). As such, if we use a memoization policy of 1 , we end up with $R \rightarrow i f(1, R, R) \rightarrow i f(1,1 * R, R)$ where $1 * R$ does not yield any useful rewrites.

Alternately, consider the case where the memoization signature is different from the memoization policy. Again, suppose that the memoization policy says memoize everything, which is represented as the $\mathbf{R}$-expr 1 . However, we are now going to say that the signature represents queries that we have encountered before and have managed to do useful rewriting on the $\mathbf{R}$-expr. In other words, we can now have a memoization signature like $(X=2) *(Y=3)+(X=7) *(Y=11)$. This means that the memo corresponds with an $\mathbf{R}$-expr like $R \rightarrow \operatorname{if}((X=2) *(Y=3)+(X=7) *(Y=11)$, $(X=2) *(Y=3) * R+(X=7) *(Y=11) * R, R)$. The $R-\operatorname{expr}(X=2) *(Y=3) * R$ and $(X=7) *(Y=11) * R$ are likely to have useful rewrites as we can condition those rewrites on the values of $X$ and $Y$.

### 10.6.2 Handling Memos we "Want" but do not "Have"

It is often the case that what a user wants memoized, as defined policy, and what the system has, memoized as tracked by the signature, are often not aligned. As such, when an inconsistency is detected, the memo's signature and memoized $\mathbf{R}$-expr are updated. This is done by waiting for queries to be made against the memoRead R-expr, which are then checked against the policy. When the policy indicates that a query should be memoized, that query will have a memo created, just like in sections § 10.2 and 10.5 . Conceptually, the memoization policy held in what_want_memoized can be handled by an if-expression as shown in figure 10-26.

```
memoRead(container, \(\cdots\) ) \(\approx\) if(container.what_want_memoized( \(\cdot \cdot\) ),
    readOrMakeMemo(container, \(\cdots\) ),
    container.original_rexpr( \((\cdots)\) )
```

Figure 10-26. Approximate interpretation for the memoRead( $\cdots$ ) R-expr. When the memoization policy what_want_memoized is rewritten as 0 or $1+R$ for some $R$, this allows the system to determine if something should be memoized-in the same way that an if-expression conditions on 0 or $1+R$. In the case that the policy states that something should be memoized, the "readOrMakeMemo" will check the current memoization signature to see if the requested query is contained in the memo table. In the case that the requested query is not contained, the memo and signature will be updated as discussed in sections § 10.2 and 10.5.

Just like with an if-expression, when the memoization policy cannot be rewritten as 0 or $1+R$ for some $R$, the system will have to defer accessing the memo. Hence, a bad memoization policy can cause the system to become inoperable.

### 10.7 Update Loop

It is often the case that there are many different updates that are pending at any given point in time. Dyna permits us to process updates to memos in any order we choose. Dyna only requires that all memos are consistent with each other when all updates have finished processing. Therefore, to handle this, we buffer all updates using a (priority) queue.

The update queue is a global object. The system runs a loop processing updates until the queue is completely empty. When processing an update, the system will check if the current memo is consistent by recomputing it from the original $\mathbf{R}$-expr using SimplifyNormalize (algorithm 1) and check if the resulting $\mathbf{R}$-expr is semantically equivalent to the currently memoized $\mathbf{R}$-expr. ${ }^{151}$ In the case that the memo must be updated, the memo is updated first, and then the assumption associated with the memo is invalidated, causing downstream dependencies to receive notifications.

[^93]When a notification is received, a pending update is added to the queue, ${ }^{152}$ and the update loop will continue its processing.

This is shown in algorithm 4.

```
function ProcessUpdates()
    global UpdateQueue
    while not Empty(UpdateQueue) :
        work \leftarrow & Pop(UpdATEQUEUE) }\triangleright\mathrm{ Pop from the queue according
                    to the priority function
        \triangleright "work.memo" is a pointer to mutable container which holds the memo itself. It con-
            tains the fields that track the original \mathbf{R}-\operatorname{expr, the current memoized \mathbf{R}}\mathbf{\mathrm{ expr, and the}}\mathbf{}\mathrm{ (t)}
            assumption which protects read operations performed against the memo.
        R original }\leftarrow\mathrm{ work.memo.original_rexpr
        global assumptionsTracker }\leftarrow\emptyset\quad\triangleright A set of all assumptions depended on during
                            SIMPLIFYNormalizE, automatically tracked
        R new}\leftarrow\mathrm{ SimplifyNormalize( ( original)
        if SemanticallyEquivalent(R new,work.memo.the_memo):
                No Op \triangleright No update happens if it is semantically equivalent
        else
            oldAssumption \leftarrow work.memo.assumption }\triangleright\mathrm{ Save a pointer to existing assumption
            work.memo.the_memo }\leftarrow\mp@subsup{\textrm{R}}{\mathrm{ new }}{
            work.memo.assumption }\leftarrow\mathrm{ new Assumption()
            oldAssumption.Invalidate() \triangleright Invalidate the old assumption
        for assumption \in ASSUMPTIONSTrACKER :
            assumption.SUBSCRIBE(WORK.MEMO)
```

Algorithm 4. A rough outline of the Process Updates Function. The function uses a global update queue (which can be a priority queue) to choose the next unit of work. It recomputes the $\mathbf{R}$-expr R from scratch based on the original definition. If the newly computed $\mathbf{R}$-expr is semantically equivalent, then the memo is not modified.

### 10.8 Controlling Memoization

When it comes to memoization, there are many different memoization policies, which can have a significant impact on the runtime and memory efficiency of the

[^94]overall system. For example, the order in which updates are processed can be the difference between a program running in linear time or exponential time (section §2.7.1). Furthermore, there exist programs where a bad memoization policy can cause the program to become not terminating. ${ }^{153}$

Ideally, the Dyna system would automatically figure out how to best apply memoization and prioritize updates. Unfortunately, at this time, automatic configuration of memoization is still an open research question. Instead, what we have currently is a mechanism to allow users to define their own memoization policies and update orders using \$memo and \$priority, which are described here again but were previously introduced in section §2.7.

### 10.8.1 \$memo( $)=$ "none"|"null"|"unk". ${ }^{154}$

The most significant control mechanism for memoization is \$memo, which essentially controls if memoization is used or not, and should therefore fallback to the original un-memoized R-expr. As previously discussed in section §2.7, \$memo(•) takes one argument that matches the structured-term with the same name as the user-defined term and "returns" the string "none", "null", or "unk" to indicate "how" and what should be memoized. ${ }^{154}$

Internally, in the implementation of memoization, we are instead going to think about \$memo slightly differently from the surface-level syntax. Instead, we will say that the memoization policy defined by \$memo and represented as an $\mathbf{R}$-expr is in one of three "states". These states are as follows:

1. The query is memoized (or it should be memoized),
2. The query should not be memoized, and it should fallback to the original R-expr, and
3. There is not enough information yet to determine whether the memoization policy is in state 1 or state 2.
[^95]The first state where a memo should exist corresponds with "null" being returned by the \$memo function. The second state with no memo corresponds with "none" being returned and is also the default state if nothing else is defined for \$memo. The third state of being "unsure" corresponds with having \$memo represented as an R-expr that cannot be fully rewritten yet. Notably, the keyword "unk" does not correspond to any single state (section §10.8.3). Instead, "unk" represents a common case that switches between the third state of "unsure" and the first state where a memo exists once enough variables are grounded and will be discussed further in section §10.8.3.

### 10.8.2 \$ground and \$free Annotation

When it comes to memoization policy, the system needs to determine when it must check the memo, and it is allowed to bypass the memo and fallback to the original $\mathbf{R}$-expr. Sometimes, though, we do not care about the value that is being stored but rather that there is enough information available in the query such that any memo that we create will be useful. To support different memoization policies, we have $\$ \operatorname{ground}(\cdot)$ and $\$$ free $(\cdot)$ which work as annotations on the state of the variable.

To see how \$ground and \$free work, let us consider the following memoization policy for the rule foo $(X, Y, Z)$ :
${ }_{471} \mid \$ m e m o(f o o[X, Y, Z])=" n u l l "$.
Given the policy on line 471, there is no requirement on the value for the arguments to foo, hence the memo for the user-defined foo $(\cdot, \cdot, \cdot)$ is allowed to be equivalent to the original R-expr. ${ }^{155}$

Conversely, suppose that we have the memoization policy: ${ }^{156}$
472 | \$memo(foo[X:\$ground, $Y: \$$ ground, Z]) $=" n u l l "$.
Here, we require that $X$ and $Y$ are known ground values before a memo is created. This means that our memo signatures will look something like ' $(X=1) *(Y=7)+$ $(X=5) *(Y=3)+\cdots '$, and that the memoized $\mathbf{R}$-expr can depend on the values of $X$ and

[^96]Y. Hopefully, this will result in useful rewrites being completed before the resulting $\mathbf{R}$-expr is stored in the memo.

The way that \$ground is implemented is that it requires that its argument take some value before it returns true. If the value is not known, then \$ground becomes a delayed constraint and that prevents the $\mathbf{R}$-expr from being rewritten. This can be implemented by defining a single rewrite rule that checks that a variable is ground, as in figure 10-27.

$$
\text { 12. } \llbracket \operatorname{ground}(X) \rrbracket_{E}=1 \quad \operatorname{ground}(X) \xrightarrow{85} 1 \quad \text { if } X \in \mathcal{G}
$$

Figure 10-27. Semantic definition and rewrite rule for $\$$ ground

Because \$ground prevents the memoization policy R-expr from being rewritten, it delays accessing the memo table as the memoization policy must be rewritten as a multiplicity before the system is allowed to check the memo or bypass the memo (section §10.6.2).

To complement \$ground, \$free is designed to annotate variables that may not yet know their value (which is typically called free in logic programming). Unlike \$ground, \$free is a no-op and only serves to indicate that \$ground was not forgotten.

The reason that \$free is a no-op is that we want to avoid inconsistent behavior when it comes to reading a memo. For example, suppose that we had a memo for the query foo $(1, x)$ where $x$ is a free variable. This memo can answer the query foo $(1,3)$, as it is possible to unify these two expressions together with $X=3$. However, if we explicitly matched against the variable $x$ being free, then this would be an inconsistent memoization policy and could potentially bypass guesses, which is not allowed (section §10.5.3).

### 10.8.3 "unk" Memos

The keywords "null" and "unk" were chosen to maintain continuity with prior publications on memoization and guessing in the Dyna programming language [66, 67]. That said, only the externally observable behavior of "null" and "unk" is consistent with the prior work. The conceptualization of "null" is quite similar to the prior work, but "unk" is very different.

In the prior work [66, 67], "null" and "unk" were used to represent the "default" value of a memo. With "unk" standing for "unknown value", and "null" representing that the memo has a "null valued" guess that must be validated. When an "unk" value was encountered during back-chaining, the system would immediately perform the computation using the original program and save the result. Conversely, a null guess would be eagerly computed ahead of time-as in the style of forward-chaining Datalog. Hence, the distinction between "null" and "unk" was deciding when the computation is performed.

In the prior work, without $\mathbf{R}$-exprs, it was in fact very important to make this distinction about when computation happens. Delightfully, with R-exprs we can be a bit more fast and loose about when computation happens in terms of our memoization formalism. The reason is that $\mathbf{R}$-exprs can represent computation rather than a value in the language (e.g. a memo that can contain plus ( $1,2, \mathrm{X}$ ) vs. the value 3). Therefore, we can memoize $\mathbf{R}$-exprs that represent partially evaluated computation, which was not allowed under the prior formalism.

As such, we do not have to worry about ensuring that our memoized $\mathbf{R}$-exprs are fully rewritten into a single value assignment. This leaves one question: what did an "unk" memo represent (according to the prior work), and how are we going to emulate "unk" with R-exprs. In the prior work, the memo table consisted of fully grounded terms as keys (e.g. foo ( 1,2, "hello") ), and associated a value with each key (e.g. 3, 7, "hello", bar[1, 2], etc.). Therefore, when using a memoization policy of "unk", the prior work did not allow queries of the form foo $(1, Y, Z)$ against the memo table, as the previous approach cannot make any guarantees about terms not found in the memo table, as their value is unknown. Hence, the prior work only allowed queries for fully grounded terms like foo( 1,2 , "hello") when using an "unk" memoization policy. As such, the "unk" memoization policy can be seen as equivalent to requiring that all variables grounded: \$memo(foo $[X, Y, Z]$ ) = "unk". $\equiv$ \$memo(foo[X:\$ground, Y:\$ground,Z:\$ground]) = "null".

### 10.8.4 Implementation of \$memo

With "unk" now represented in terms of "null", we now only need to handle the three cases when implementing \$memo:

1. When "null" is returned, we need to read from the memo table,
2. When "none" is returned, we need to skip the memo table and
3. When we do not yet know if it is "null" or "none" and need to defer the read operation.

Conceptually, these three cases can be handled with an if-expression like R-expr. If the conditional branch of the if-expression can be completely rewritten as either $1+R$ for some $R$ or as 0 , then we do not know which branch of the if-expression to return. This is identical to what I presented in figure 10-26, in how we use the memoization policy. As such, we can define the variable what_want_memoized as the $\mathbf{R}$-expr:

```
proj(Name,
    (Name=foo[X 
    $memo(Name, "null"))
```

Figure 10-28. The "what want memoized" R-expr (line 467 of figure 10-25), which adapts from the \$memo memoization policy represented as an $\mathbf{R}$-expr that returns "null" or "none" or not rewrite and becomes a delayed constraint, which happens if there is enough information in the context $\mathcal{C}$ rewrite the user-defined \$memo R-expr.

The Name variable holds the structured term for the user-defined term for which a memo represents. The $\$$ memo(Name, "null") is a call to the \$memo $\mathbf{R}$-expr definition. When \$memo returns "null", this $\mathbf{R}$-expr will be rewritten as 1 , causing the ifexpression in figure 10-26 to access the memo table. When \$memo returns "none", then this will be rewritten as 0 , causing the if-expression in figure 10-26 to fall through to the original R-expr. In the case that \$memo cannot be fully rewritten (possibly caused by a \$ground annotation), the if-expression in figure 10-26 will defer reading or bypassing the memo.

### 10.8.5 Ordering Updates with \$priority

The order in which updates are processed can have a significant impact on the runtime of the program. \$priority (•) supports different ordering of updates by defining a floating point number which is used to define the priority of an update.

Whenever an assumption is invalidated and an update is pushed to the update queue, the \$priority function is consulted, using the user-defined rule name and any values that are known to compute a floating point value. This value is used to sort the priority queue, which contains all pending updates. The priority is only computed when updates are pushed to the queue to avoid the complexity of reordering the queue due to values that \$priority depended on changing.

## Chapter 11

## A Realistic Implementation of R-exprs

So far we have been discussing an abstract idea of R-exprs (chapters 5 and 6), rewriting R-exprs using Simplify (chapter §8) and extensions to this design in the form of rearranging R-exprs (chapter §9) and adding memoization (chapter §10). In this chapter, I am going to transition from abstract ideas and focus on a real implementation of R-exprs and $\mathbf{R}$-expr rewriting. Note that some design choices discussed in this chapter were made with the intention of JIT compiling $\mathbf{R}$-exprs, which will be discussed in chapter $\S 12$.

### 11.1 Design Goals for our Implementation

Before diving into the details about the implementation itself, let us discuss what the goals are for this implementation of Dyna and $\mathbf{R}$-expr rewriting.

1. It should be usable and capable of supporting the kinds of Dyna programs that we expect users to write.
2. It should be "efficient" so that it can be used on "real" problems. - Recall that our target audience for Dyna is Machine Learning (ML) and Artificial Intelligence (AI) researchers. This means that the kinds of problems we expect often have the form of performing very similar computations on slightly different
data many times in a loop. An example of this kind of computation would be matrix multiplication.
3. It should make the implementation of $\mathbf{R}$-exprs and $\mathbf{R}$-expr rewrites easy. - A complete Dyna implementation requires dozens of different $\mathbf{R}$-expr kinds and hundreds of rewrite rules. Additionally, there is a lot of potential for future work to introduce new R-expr kinds and rewrite rules.
4. It should support all of Dyna. - There have been several implementations of Dyna over the years. ${ }^{157}$ However, previous implementations have only implemented a small subset of Dyna. Building a complete implementation of Dyna demonstrates that the term rewriting approach in this dissertation is the "right formalism" and sufficient to implement a complicated logic programming language like Dyna.
5. It should support (JIT) compilation of Dyna programs. - This is a secondary goal of wanting the implementation to be "efficient" and "fast", but this is a significant design constraint on the implementation, so I mention this at the start as a design goal.

### 11.1.1 What does "Efficient Implementation" Mean?

We would like Dyna to execute programs at a level that is comparable to the "first implementation (of an algorithm) that a user would write in a procedural language." In other words, we understand that a skilled programmer can make an algorithm faster through the use of clever tricks and knowledge about computer architecture. Our goal is not to compete with the skilled programmer. Instead, our goal is that an average skilled programmer will be able to develop programs in less time with Dyna when compared to a procedural language (due to Dyna's simplicity) and have the program run at a level that is comparable to what an average skilled programmer

[^97]would write in a procedural programming language. ${ }^{158}$
To illustrate this point, let us consider a program which computes a matrix-vector product:
$$
{ }_{473} \mid a(X)+=b(X, Y) * c(Y) .
$$

If we ran line 473 using the execution strategies that we have discussed so far, we would represent $b(X, Y)$ and $c(Y)$ as the disjunctions in figure 11-1.

```
b(X,Y, Val) -> c(Y, Val) }
    (X=1)*(Y=1)*(Val=3)+ (Y=1)*(Val=11)+
    (X=2)*(Y=1)*(Val=7)+ (Y=9)*(Val=23)+
    (X=1)*(Y=2)*(Val=9)+\quad(Y=5)*(Val=0)+
    \vdots
```

(a) Example $\mathbf{R}$-expr for $b(X, Y)$
(b) Example R-expr for $\mathrm{c}(\mathrm{Y})$

Figure 11-1. Example R-exprs used by line 473.

Accessing the element from the disjunctions in figure 11-1 requires many steps of rewriting. We would have to use the distributive rewrites to expand out the R-expr out (e.g. $(X=1) *(Y=1) *(V a l B=3) *((Y=1) *(V a l C=11)+(Y=9) *(V a l C=23)+(Y=5) *$ (ValC=0) $+\cdots$ ) $+\cdots$ ) which would create a large $\mathbf{R}$-expr and be inefficient, requiring hundreds of rewrites to access the relevant elements from the $b(X, Y)$ and $c(Y)$ relations.

Conversely, let us consider how line 473 would be written in a procedural language: ${ }^{159}$

[^98]```
function MatrixVectorProduct(b[0..N, 0..M], c[0..M])
    ret \(\leftarrow\) MakeZeroVector \((N)\)
    for \(X=0\) to \(N\) :
        for \(Y=0\) to \(M\) :
            \(\operatorname{ret}[\mathrm{X}] \leftarrow \operatorname{ret}[X]+\mathrm{b}[X, Y]{ }^{*} \mathrm{c}[Y]\)
    return ret
```

Algorithm 5. Example procedural implementation of matrix-vector product

Here, MatrixVectorProduct is able to assume that the representation of ' $b$ ' and 'c' are dense arrays/matrices of numbers. This makes accessing elements in the array (line 5) efficient, taking $O(1)$ time. Furthermore, the MatrixVectorProduct function is able to know the size of the matrix and vector using the integers $N$ and $M$ and have an iterator that loops over the domain of the matrix and vector, being the integers between $[0, N)$ and $[0, M)$ (lines 3,4 ). Finally, the place for the final result is allocated before the computation is started (line 2). This means that the procedural code does not have to store all the intermediate multiplications and instead aggregates the value along the way into the final location (in memory).

Our goal, with $\mathbf{R}$-expr-based rewriting, is to maintain the representational power of $\mathbf{R}$-exprs, while having the execution of $\mathbf{R}$-exprs be comparable to algorithm 5, in that we can avoid the overhead of naively rewriting $\mathbf{R}$-exprs.

### 11.2 Implementation Overview

Given the requirements that I listed above for a realistic implementation, I evaluated a number of different possibilities for the implementation of Dyna. I make no claim that this design is the "best" approach. However, the design presented in this chapter is the third implementation of Dyna that I have completed during my Ph.D., and it has been designed with the insight gathered from the prior two implementations. I believe that the design presented here is worthy of consideration for anyone looking to build a term rewriting system for a programming language.

First, for the choice of programming language, I choose to use Clojure [92]. Clojure is a LISP-like programming language that targets the Java Virtual Machine (JVM). Clojure is a reasonably fast language, running a little slower than Java but


Figure 11-2. A high-level figure representing the major components in the Dyna implementation. The user's program, queries, and updates start at the top and then are converted into the Dyna AST. The AST uses the same structured term class as section §2.1.1, allowing Dyna programs to manipulate the ASTs using macros (section §2.10.2). ASTs are converted into $\mathbf{R}$-exprs as in chapter $\S 7$. Memoization (chapter §10) is added to the program by rewriting the user's $\mathbf{R}$-exprs with the memoRead $\mathbf{R}$-exprs and constructing the relevant memo tables. The memoization policy (\$memo (•)) is represented as an $\mathbf{R}$-expr, which is processed specially by the memoization controller. Queries and updates made against the Dyna program are converted into $\mathbf{R}$-exprs and rewritten using Simplifyormalize (algorithm 7) until there are no more rewrites to apply, and there are no further updates to memos. The resulting $\mathbf{R}$-expr is then returned to the user. JIT compilation (chapter §12) attempts to make execution with $\mathbf{R}$-exprs faster.
faster than many other high-level languages. Clojure can also interface with Java libraries, which is convenient for interfacing Dyna with Java (and Python through the use of Java-Python interfaces). The main reason for choosing Clojure was that there are a few features that are essential for implementing Dyna. These are macros, the ability to modify ASTs of the host language, and a runtime evaluate function. Macros and AST manipulation are essential for making the implementation of the many different rewrite rules and $\mathbf{R}$-expr kinds efficient, and the runtime evaluation function is used to generate $\mathbf{R}$-expr-dependent procedural code. Admittedly, any LISP-like language satisfies these requirements.

Dyna's internal state is represented using $\mathbf{R}$-exprs. All the major components inside Dyna pass $\mathbf{R}$-exprs between them. The flow of data/R-exprs is shown in figure 11-2.

### 11.3 Dyna's Front-end

Dyna provides a Read-Print-Eval loop (REPL) as a front-end, as well as Java and Python API (section §2.3.1). The API, and REPL accept strings of Dyna code that represent queries and updates (as previously shown in section §2.3.1). The Dyna source is converted into $\mathbf{R}$-exprs and then rewritten completely.

The conversion process from a string of Dyna source code into an $\mathbf{R}$-expr is done with the help of two special R-exprs. The first is string_to_ast(S,Ast), which converts a string of Dyna code into the Dyna AST. The implementation of the string_to_ast(S,Ast) rewrite rules uses Antlr4 [110] ${ }^{160}$ to parse the string of Dyna code and assigns the AST to the AST variable. The Dyna AST is represented using the same term class used to represent structured terms (section §2.1.1). This is similar to Prolog where rules are represented using terms. This allows macros defined in Dyna to modify the Dyna AST before it is converted into R-exprs (section §2.10.2).

The second $\mathbf{R}$-expr is ast_to_rexpr(Ast, ResultVariable, VariableNameMap). ast_to_rexpr takes three arguments. The first variable is the AST itself. This can be a top-level AST, which comes from parsing an entire file or a line of code from the REPL, or this can be a smaller unit of the AST, which corresponds to an expression inside of the Dyna source code (as defined in chapter §2). The second variable

[^99]ResultVariable is the result from evaluating an expression in Dyna. In Dyna, all expressions return some value. For example, the AST corresponding to ' $1+2$ ' is well formed and will return the value 3 once evaluated. The reason we need the return variable as an argument is that $\mathbf{R}$-exprs do not have a return value but instead return the multiplicity, which indicates that the assignment is "true". Therefore, we must assign the return value of the Dyna expression to a given variable when converted into R-exprs. ${ }^{161}$ Asts that are not expressions, such as top-level asts and declaring rules, are made into expressions by defining them to return the dummy value of true. The third argument is a map from string variable names, as used in the source code, to the value object types: VariableNameMap: String $\rightarrow \mathcal{V}$. The R-expr is not required to use the same identifier as the source code; hence, the map tracks the association between names and variables. For example, $\{$ " $X$ " $\mapsto \operatorname{Var} \times 72$, " $Y$ " $\mapsto 7\}$.

The string_to_ast R-expr can be considered a "proper" R-expr, in that it can be completely defined as a relation on its arguments. On the other hand, ast_to_rexpr is not proper. When ast_to_rexpr is rewritten, it modifies the global ${ }^{162}$ state of the Dyna system. This includes changing the definition of user-defined $\mathbf{R}$-exprs (section §5.2.2.11).

Once the Dyna system, which is controlled by the API (section §2.3.1), is finished evaluating all queries and updates provided as a string, the results from queries are returned to the user. If the resulting $\mathbf{R}$-expr from a query is "sufficiently simple", then Dyna automatically extracts the value from the $\mathbf{R}$-expr representation so it is more easily interpretable by the driver program. This includes ground values such as integers and strings, as well as arrays and hash maps. If the result from a query is not a simple $\mathbf{R}$-expr, then the $\mathbf{R}$-expr itself will be returned to the driver program. Users are expected to write programs that result in simple $\mathbf{R}$-exprs and values, therefore an R-expr being return generally indicates that "something has gone wrong ${ }^{163 "}$ when evaluating a query. Hence, users are not expected to build driver programs that understand R-exprs. The Dyna REPL, which is built upon the

[^100]Dyna API, will attempt to pretty-print R-exprs.
A high-level overview of the front-end is shown in algorithm 6.

```
function DynAFrontEnd(Source)
    R\leftarrowproj(A,string_to_ast(Source, A)*ast_to_rexpr(A,true, {}))
    repeat
        ProcessUpdates() D Call pending updates on the update queue, algorithm 4
        global }\mp@subsup{}{}{162}\mathrm{ QueryResult }\leftarrow
        R2 \leftarrowSIMPLIFYNORMALIZE (R)}\triangleright Evaluate the program, modifying global as needed
        assert R2 == 1 \triangleright The top-level AST rewrites as multiplicity 1 upon success
    until isQuery(Source) and isEmpty(UpdateQueue) D If pending updates, run
    updates and redo queries, algorithm 4
    D Query results are saved into the QueryResult buffer by ast_to_rexpr
    if QueryResult matches (ResultVariable=Value):
        return Value }\triangleright\mathrm{ Avoid returning the }\mathbf{R}\mathrm{ -expr if the }\mathbf{R}\mathrm{ -expr is "sufficiently simple"
    else
        return QueryResult }\triangleright\mathrm{ Return the entire R-expr produced by the query
```

Algorithm 6. The front-end for Dyna is called on a single update or query at a time. An update/query is wrapped in a standard $\mathbf{R}$-expr (line 2) and is passed to SimplifyNormalize (line 6) where all execution happens via rewrites. If a query is performed that causes guessing (which can cause pending updates on the update queue), then those updates will be processed, and the queries will be redone (line 8 ). ${ }^{164}$ Once the query has converged, the result is returned to the user, with sufficiently simple values being cast into a usable format.

### 11.4 Realistic Rewriting, Part 1-Redesigning Simplify(Normalize)

In this chapter, we are trying to make our implementation of rewriting faster. As such, we are going to make some changes to the approach described in chapter §8 in hopes of making execution run faster.

[^101]Previously in section §8.2.1, we had the functions Simplify (section §8.A) and SimplifyNormalize (algorithm 1). The function Simplify recursively invokes itself on the R-expr performing any rewrites it can apply. To find relevant rewrites, Simplify matches the current R-expr and uses the context $\mathcal{C}$ to identify conjuncts without rearranging the R-expr. SimplifyNormalize invokes Simplify in a loop until no more rewrites can be applied.

The first problem with this design we will address is that Simplify chooses among all possible rewrites equally. However, not all rewrites are equally useful or can be matched quickly. To manage this, we separate our rewrites into three separate categories, depending on what information is needed to match the rewrite and what priority the rewrite should run at.

1. The first category we call standard rewrites. These rewrites only need a local view of the $\mathbf{R}$-expr and knowledge about bindings to variables, which is available in the context $\mathcal{C}$. These rewrites correspond to common operations such as performing addition between values (e.g. rewrite rule 28, plus $(1,2, X) \rightarrow(X=3))$. These rewrites are classified as standard as they are the most common kind of "useful" rewrites that we have. Furthermore, variable bindings can be efficiently tracked using an associative map (hash map) inside the context $\mathcal{C}$. This simplifies the context as it does not require complete knowledge about all conjunctive $\mathbf{R}$-exprs, which means that maintaining the context is efficient.
2. The second category of rewrites are inference rewrites. These rewrites require the complete context with all conjunctive $\mathbf{R}$-exprs. Because all conjunctive $\mathbf{R}$-exprs are tracked in the context, this is more "heavy" in that $\mathbf{R}$-exprs and indexes on $\mathbf{R}$-exprs are required. Furthermore, matching against the context is much more expensive. When the system only checks variable assignments, it can perform an $O(1)$ time lookup in a map. However, when matching with $\mathbf{R}$-expr in the context, it must scan through many potential matches.

Because these rewrites are expensive to match, they are run with the lowest priority.

The reason these rewrites have been named inference rewrites is that they usually infer the existence of new constraints using propagation. For example, rewrite rule 34 combines two lessthan constraints, e.g. lessthan(A,B)*
lessthan $(B, C) \xrightarrow{34}$ lessthan $(A, B) *$ lessthan $(B, C) *$ lessthan $(A, C)$. In fact, other rewrite systems sometimes call these kinds of rewrites propagators [72].
3. The third category of rewrites are constructor rewrites. These rewrites are run with the highest priority. They are run immediately when an $\mathbf{R}$-expr is constructed. These rewrites are mostly used to keep the $\mathbf{R}$-expr tidy. For example, rewrite rules 13 to 15 are the identities involving 0 and 1 , and will be removed from a conjunction immediately rather than being kept around in the $\mathbf{R}$-expr representation (e.g. $0 \star$ R $\rightarrow 0$ ).

With this division of rewrite methods, we have two different versions of the Simplify function: SimplifyOnlyFast and SimplifyAllRewrites. These new functions behave the same as the SIMPLIFY function from chapter $\S 8$, in that they take a context and an R-expr and rewrite it as another R-expr, but SimplifyOnlyFast only invoke the faster-to-match standard rewrites and SimplifyAllRewrites invokes both standard and inference rewrites. The constructor rewrites are invoked immediately upon the construction of an R-expr, and therefore are "hidden" and do not require a "visually obvious" call to Simplify in the code. The function Simplify refers to whichever version of SIMPLIFY is currently "active".

The SimplifyNormalize function is redesigned to first complete all of the fast rewrites by fixed-pointing the SimplifyOnlyFast function and then calling the SimplifyAllRewrites function once to check if there are any inference rewrites that can be performed. It will then go back to running only the fast rewrites until all of those are completed again. This modified version of SimplifyNormalize is shown in algorithm 7.

### 11.5 Declaration of R-exprs and Rewrites

As hinted at in the overview (section §11.2), we are going to make heavy use of Clojure macros in the implementation of Dyna. The reason for this is that Dyna, built on $\mathbf{R}$-exprs, is made up of hundreds of different $\mathbf{R}$-expr kinds and rewrite rules. Hence, we want to minimize the amount of code we need to write for each $\mathbf{R}$-expr and rewrite rule.

```
function SimplifyNormalize(R)
    \(\mathcal{C} \leftarrow\langle\}: \operatorname{Var} \mapsto \operatorname{val}\},\{\mathrm{R}\}\rangle \triangleright\) Initialize empty context, \(\langle\) variables to values, \(\mathbf{R}\)-exprs \(\rangle\)
    repeat
        repeat
            \(\mathrm{R}_{\text {old }} \leftarrow \mathrm{R}\)
            \(\mathrm{R} \leftarrow\) SimplifyOnlyFast \((\mathrm{R}, \mathcal{C}) \quad \triangleright\) Perform only rewrites that are "fast"
        until \(\mathrm{R}_{\text {old }}=\mathrm{R} \quad \triangleright\) Reached a fixed-point, no more fast rewrites
        \(R \leftarrow \operatorname{SimplifyAlLReWRites}(\mathrm{R}, \mathcal{C}) \quad \triangleright\) Perform all rewrites
        until \(R_{\text {old }}=R \quad \triangleright\) Reached a fixed-point, no more rewrites
        return \(R\)
```

Algorithm 7. Simplify Normalize with different priorities for fast rewrites.

### 11.5.1 Declaration of R-expr Kinds

To define R-expr kind in Clojure, the implementation provides the def-base-rexpr macro. This macro creates a class ${ }^{165}$ that is specialized for the implementation of a particular R-expr kind. It also sets up internal data structures associated with the $\mathbf{R}$-expr kind that are used for relevant meta-data, registering the $\mathbf{R}$-expr kind, and registering rewrites for the $\mathbf{R}$-expr kind.

Using def-base-rexpr we can define some core R-expr kinds as follows:

```
474 (def-base-rexpr conjunct [:rexpr-list args])
(def-base-rexpr unify [:var a :var b])
(def-base-rexpr proj [:hidden-var var :rexpr body])
```

The first argument is the name of the $\mathbf{R}$-expr, and the second argument is the names of fields associated with each R-expr, along with types for each field denoted using a Clojure keyword. The type annotation on the fields automatically configures how different methods on the R-expr behave. For example, : var represents a field that holds value types, which can be variables or constant values. This automatically configures the "rename variables" ${ }^{166}$ and "list all variable" ${ }^{167}$ functions to include the : var annotated fields. Fields marked : hidden-var contain variables that are projected out, and these variables are removed from the set of variables returned by "list all variables". Similarly, : rexpr represents a single R-expr, and :rexpr-list

[^102]represents an array of $\mathbf{R}$-exprs.
This annotation using keywords handles around $90 \%$ of all cases that we need when defining R-exprs. We can also override the definition of any method in the case that we need to change something that only applies to one $\mathbf{R}$-expr kind.

For example, we override the is-constraint? ${ }^{168}$ method to check that all of the sub-R-exprs contained in a conjunction are constraints, in which case their conjunction is also a constraint (line 479):

```
(def-base-rexpr conjunct [:rexpr-list args]
    (is-constraint? [this]
        (every? is-constraint? args)))
(def-base-rexpr unify [:var a :var b]
    (is-constraint? [this]
        true))
```


### 11.5.2 Declaration of Built-Ins

Most of the $\mathbf{R}$-expr kinds and rewrite rules that we have to define correspond to built-in R-exprs (sections § 5.2.2.3 and 6.3). Built-ins define numerical operations (plus, times, cosine, etc.), logical operations (lessthan, lessthaneq, etc.), and other primitive operations such as string concatenation.

To help define all of these R-expr kinds, we have another special macro: def-builtin-rexpr, which calls def-base-rexpr and def-rewrite, which will be defined shortly.

As an example, the complete definition of the plus $(\cdot, \cdot, \cdot) \mathbf{R}$-expr is as follows:

```
(def-builtin-rexpr plus 3
    (:allground (= v2 (+ v0 v1)))
    (v2 (+ v0 v1))
    (v1 (- v2 v0))
    (v0 (- v2 v1)))
(def-user-term "+" 2 (make-plus v0 v1 v2))
(def-user-term "-" 2 (make-plus v2 v1 v0))
```

The def-builtin-rexpr macro will define a plus R-expr which contains 3 variable fields which are automatically named v0, v1 and v2. Lines 484 to 487 define

[^103]the different rewrite rules we have for plus. Line 484 corresponds with rewrite rules 26 and 27 where all the arguments are ground (e.g. plus $(1,2,3) \rightarrow 1$ ). Similarly, lines 485 to 487 corresponds with rewrite rules 28 to 30 which define the rewrites for performing computation using plus. The variable that appears first in the parentheses is the variable that is assigned by the computation. The second expression in the parentheses defines how the computation is performed. For example, when the system rewrites plus $(1,2, x)$, it will have $v 0=1$ and $v 1=2$. The variable $v 2$ corresponds with the variable $x$ and is assigned the value returned from (+ 12 ). Hence, this is equivalent to rewriting plus $(1,2, X) \rightarrow(X=3)$.

Lines 489 and 490 define the plus R-expr under the names accessible from the Dyna source code. The ' + ' function (in the Dyna source) takes two arguments and returns the third argument. Represented as v0, v1 and v2 respectively. Recall from section §7.1.4, I mentioned that we do not have a subtraction $\mathbf{R}$-expr, but instead use the plus $(\cdot, \cdot, \cdot) \mathbf{R}$-expr with its arguments rearranged. This is what is happening on line 490 , with the $v 0$ and $v 2$ arguments switched.

The function make-plus is the constructor for the plus $\mathbf{R}$-expr. It is defined by the def-base-rexpr macro. The make-NAME functions internally run any constructor rewrites that are marked to run when an $\mathbf{R}$-expr is constructed.

### 11.5.3 Declaration of Rewrite Rules

One thing I attempted to make easy is the declaration of rewrite rules. We have literally hundreds of rewrite rules that need to be implemented. In previous rewritebased prototypes of Dyna that I developed, rewrites with large implementations frequently resulted in difficult-to-find bugs. ${ }^{169}$ Hence, having a short implementation for each rewrite makes it easier to visually inspect ${ }^{170}$ that a rewrite is correct.

The core mechanism for defining rewrites is the def-rewrite macro. This macro takes a number of different keyword arguments and a Clojure expression, which computes the result of a rewrite. For the remainder of this section, I will illustrate how the def-rewrite macro is used, to give an idea of what features

[^104]were necessary to concisely define $\mathbf{R}$-expr rewrite rules.

### 11.5.3.1 First Declaration of a Rewrite Rule

We will start by looking at the def-rewrite definition for rewrite rule 28, which performs the computation for built-in plus (e.g. plus ( $1,2, X) \rightarrow(X=3)$ ). Note: these rewrites on plus are automatically generated by the def-builtin-rexpr defined above on line 485 and do not appear directly in the implementation. I have chosen to present rewrite rule 28 because it is a simple rewrite which is useful for illustration purposes.

```
491 (def-rewrite
4 9 2 ~ : m a t c h ~ ( p l u s ~ ( : g r o u n d ~ A ) ~ ( : g r o u n d ~ B ) ~ ( : f r e e ~ C ) ) ~
493 (make-unify C (make-constant (+ (get-value A) (get-value B)))))
```

Figure 11-3. Basic version of def-rewrite for rewrite rule 28, e.g. plus $(1,2, X) \rightarrow(X=3)$. Prolog would notate this as plus $(+A,+B,-C)$.

In figure 11-3, the :match keyword argument specifies what R-expr this rewrite rules will match against. The arguments $A, B$ and $C$ are matched to the variables on the plus $(A, B, C)$ R-expr. The notation used by : match is positional, in that it is written in the order in which fields on the R-expr were defined. Internally, this generates code that directly accesses the relevant field on the plus R-expr class. ${ }^{171}$ The keywords : ground and : free are annotations that are used to match against the variable. : ground and : free cause the $\mathbf{R}$-expr matcher to check the context $\mathcal{C}$ to see if there is an assignment to a variable. We have more keywords defined to match more complex scenarios.

Rewrites return an R-expr that will replace the R-expr in the newly generated R-expr data structure returned by Simplify. This is illustrated by line 493, which returns the equality constraint (e.g. $(X=Y)$ ), written as unify in the Dyna implementation (section $\S 5.2 .2 .1$ ). The variable $C$ already is a variable ${ }^{172}$ so it can go

[^105]into the first argument of unify, which expects a value type. For the second value, the system will compute $A+B$, which will be a numerical value. First, it needs to read the value of $A$ and $B$. This is done using get-value, which will either get the value directly out in the case that it is a constant or read from the context $\mathcal{C}$ in the case that it is variable. The context $\mathcal{C}$ is globally ${ }^{162}$ accessible, hence usable by get-value. The make-constant constructor returns a value type, which is a constant. Hence, the value assigned to $C$ is embedded directly inside the unify R-expr.

### 11.5.3.2 Assignment Rewrites

The above definition of rewrite rules is very general in that it matches $\mathbf{R}$-exprs and returns an R-expr. However, we have certain rewrite patterns that are sufficiently frequent that we provide special handling. For example, we can assign a value to a variable directly as in figure 11-4.

```
(def-rewrite
    :match (plus (:ground A) (:ground B) C)
    :assigns-variable C
    (+ (get-value A) (get-value B)))
```

Figure 11-4. Rewrite rule defined which returns a value to be assigned to the variable $C$, instead of returning an $\mathbf{R}$-expr.

On line 497, we return the value to assign to the variable $C$ instead of an $\mathbf{R}$-expr which represents the assignment.

The definition in figure 11-4 is conceptually equivalent to figure 11-3. However, it is preferable to use figure 11-4. The reason is that figure 11-4 provides the system with additional metadata on what the rewrite does. In this case, the system can analyze the rewrite in figure $11-4$ to identify that $C$ is assigned a value after this rewrite runs. Trying to extract this information from figure 11-3 would require analysis of the Clojure source code that defines the rewrite. This information will be used when compiling a sequence of rewrites in chapter $\S 12$.

Furthermore, the rewrite in figure 11-4 handles two rewrites at once. Observe that on line 495 , I have removed the : free annotation for the variable C. If the
variable $C$ has already been assigned some value, the : assigns-variable behaves as an equality check (rewrite rules 26 and 27). If not, it assigns a value to C (rewrite rule 28).

### 11.5.3.3 Handling Invalid Inputs

In the above definition given in figure 11-4, we no longer return an $\mathbf{R}$-expr. This means that we cannot represent a failure or an invalid input by returning the $\mathbf{R}$-expr $0 .{ }^{173}$

To handle this, we have two ways to represent failure. The first is returning the 0 R-expr, and the second is throwing a UnificationFailure ${ }^{174}$ exception ${ }^{175}$ as in figure 11-5.

```
(def-rewrite
    :match (plus (:ground A) (:ground B) C)
    :assigns-variable C
    (let [val-A (get-value A)
            val-B (get-value B)]
        (when (or (not (number? val-A))
                            (not (number? val-B)))
            (throw (UnificationFailure.)))
        (+ val-A val-B)))
```

Figure 11-5. Rewrite rule checking inputs and throwing a UnificationFailure on line 505. Admittedly, this rewrite is checking the type of the value assigned to the variables A and B (lines 503 and 504) before it throws the exception, but recall that Dyna handles type errors by rewriting as 0 , and it is up to the system to detect this and report the 0 to the user as a type error (rewrite rule 26 and section §2.8.3).

Throwing a UnificationFailure and returning the 0 multiplicity are used inter-

[^106]changeably. However, there are some advantages to throwing the UnificationFailure exception. The exception stops rewriting the current R-expr and parent R-expr that would also be rewritten as 0 as a result. For example, suppose that the system is rewriting an $\mathbf{R}$-expr like $\mathrm{S}+\operatorname{proj}(\mathrm{X}, \mathrm{R} 1 * R 2 * R 3 * \cdots * R n)$ ), if $R 1$ is rewritten as $0(R 1 \rightarrow 0)$, then we would like the system to immediately stop rewriting the conjunction (R1*R2* $\cdots$ ) and the projection, and returns control flow to rewriting the disjunction.

### 11.5.3.4 Modifying the Context

Rewrite rules can access the context $\mathcal{C}$ as it is tracked with a global variable. ${ }^{176}$ We have already seen get-value, which can read the value of a variable from the context. We can also modify the context in place using methods like set-value!, which tracks the assignment of a variable directly into the context.

Figure 11-6 shows the value assigned to the variable $C$ being set directly into the context. The rewrite rule can therefore return the multiplicity 1 , instead of assignment to the variable $C$. This is fine, as we do not require a representation for C's value inside of the R-expr itself. Using the context in this way enables us to track the assignments to variables and avoid having to shuffle around equality constraints inside of the R-expr.

```
(def-rewrite
    :match (plus (:ground A) (:ground B) (:free C))
    (let [val-A (get-value A)
            val-B (get-value B)]
        (when (or (not (number? val-A))
                    (not (number? val-B)))
            (throw (UnificationFailure.)))
        (set-value! C (+ val-A val-B))
        (make-multiplicity 1)))
```

Figure 11-6. The variable $C$ is set directly into the context $\mathcal{C}$ (line 514) without a corresponding representation in the $\mathbf{R}$-expr, as was done in figure 11-3.

[^107]
### 11.5.3.5 Rewrite Priorities

As stated in section §11.4, we have three different priorities for which rewrites can run at: standard, construction, and inference. By default, a rewrite is created with the standard priority. A rewrite's priority is set using the : run-at keyword argument as in figure 11-7.

```
(def-rewrite
    :match {:rexpr (unify (:free A) (:ground B))
            :check (has-context?)}
    :run-at [:construction :standard]
    (do
        (set-value! A (get-value B))
        (make-multiplicity 1)))
```

Figure 11-7. A rewrite on the unify/equality constraint, which is run when it is first constructed. This rewrite directly adds the assignment of a variable in the context using set-value!, as described above. With this rewrite included and running when unify is constructed, this makes the rewrites defined in figure 11-3 and figure 11-6 equivalent. Note that construction rewrites run when the R-expr is constructed, which means there might be an active context that can track the assignment to a variable. Hence, line 518 defines a side condition as executable Clojure code that checks that there exists an active context before this rewrite matches.

Construction rewrites like figure 11-7 are used to keep the R-expr tidy. In this case, we do not want to have equality $\mathbf{R}$-exprs that assigns a value to a variable represented in the $\mathbf{R}$-expr. Instead, we prefer to track assignments using the context. Hence, this rewrite "eliminates" the equality constraint from the R-expr and instead tracks the assignments to variables using the context.

### 11.5.3.6 Inference Rewrites

Inference rewrites are the lowest-priority rewrites as they take the most effort to match. These rewrites have access to the complete context that tracks conjunctive R-exprs in addition to assignments to variables. Figure 11-8 shows rewrite rule 34, which combines two lessthan constraints to infer a third lessthan constraint.

```
(def-rewrite
    :match {:rexpr (lessthan A B (is-true? _)) 177
    :context (lessthan B C (is-true? _))}
    :run-at :inference
    :infers (make-lessthan A C (make-constant true)))
```

Figure 11-8. Implementation for rewrite rule 34, which combines two lessthan constraints together to infer a third. The : run-at (line 526) specifies that this runs during inference with the complete context. The :context expression in the : match checks in the context for a matching $\mathbf{R}$-expr. The variable $B$ used on lines 524 and 525 is required to be the same variable.

Note that the match expression on lines 524 and 525 only lists one $\mathbf{R}$-expr under the : rexpr matcher, and the second under the : context matcher. The reason is that Simplify recurses through the R-expr and performs matches quickly against the $\mathbf{R}$-exprs it encounters. In other words, once the first R-expr on line 524 is matched, Simplify then checks the context for a matching lessthan constraint. As such, a more accurate way to write rewrite rule 34 would be:

$$
\begin{aligned}
& \text { lessthan }(A, B) \xrightarrow{34} \text { lessthan }(A, B) * \text { lessthan }(A, C) \quad \text { if lessthan }(B, C) \in \mathcal{C} \\
& \text { and lessthan }(A, C) \not C \mathcal{C}
\end{aligned}
$$

where lessthan $(B, C)$ is checked if it is contained in the context as a side condition.
Additionally, observe that rewrite rule 34 checks that the newly inferred lessthan (A, C) does not already exist. This is done to avoid inferring the same constraint multiple times. This check is handled by the :infers keyword on line 527.

### 11.5.3.7 Combining R-exprs for Inference Rewrites

Often times when running inference rewrites that check the context, we do not care about which R-expr appears in the context and which R-expr was directly matched against. To handle this, we have the : match-combines keyword argument, which

[^108]```
(def-rewrite
    :match-combines [(lessthan A B (is-true? _))
                                    (lessthan B C (is-true? _))]
    :run-at :inference
    :infers (make-lessthan A C (make-constant true))
```

Figure 11-9. : match-combines expanded this rewrite rule into multiple rewrite rules of the form shown in figure 11-8. For each defined rewrite, one $\mathbf{R}$-expr will be matched in the local context, and the other $\mathbf{R}$-exprs will be matched using the context.
takes a list of $\mathbf{R}$-exprs that must be matched. The rewrite is expanded into several simpler rewrites of the form in figure 11-8.

### 11.5.3.8 Recursive Rewrites

As stated in section §8.2.2, we have special "rewrite rules" for recursive R-exprs (such as conjunction) that recursively invoke Simplify. To this end, we can use the function simplify as in figure 11-10, which automatically references the current simplify function (either SimplifyOnlyFast or SimplifyAllRewrites).

```
(def-rewrite
    :match (conjunct (:rexpr-list Rs))
    (make-conjunct (map simplify Rs)))}\mp@subsup{}{}{178
```

Figure 11-10. Recursive rewrite rules call simplify recursively applies rewrites to sub-R-exprs.

Recall that in section §11.5.3.3, I said that we throw a UnificationFailure exception instead of returning a 0 multiplicity $\mathbf{R}$-expr. Throwing an exception immediately aborts the loop on line 535 without any special handling.

[^109]
### 11.5.4 Conclusion of R-exprs and Rewrites Declarations

This section has served both as documentation for our Dyna implementation and as inspiration for anyone developing their own rewrite-based system. Given the number of rewrite rules required to implement a rewrite-based language, such as Dyna, I believe that it is critical to have a concise way to define rewrite rules. From experience, I can say that more than $90 \%$ of the rewrite rules can be handled succinctly, given the right abstraction. Allowing the remaining $10 \%$ of rewrite rules to fall back on the host language works well to ensure that the rewrite system is "sufficiently powerful".

### 11.6 Efficient R-exprs Kinds

At the beginning of this chapter, I gave a definition for an "efficient" program (section §11.1.1). So far, we have only seen the same "inefficient" R-exprs we have had since chapter $\S 5$. To make $\mathbf{R}$-exprs efficient, we are going to replace some $\mathbf{R}$-expr kinds with $\mathbf{R}$-exprs that are conceptually equivalent but have more efficient implementations. The way we accomplish this is by combining two or more $\mathbf{R}$-expr kinds into a single larger $\mathbf{R}$-expr. This will allow us to handle frequent cases with specialized code.

The two efficient R-expr kinds we currently have are efficient disjunction, which merge disjunctions and equality constraints, and efficient aggregation, which merge aggregation and projection. Future work should consider the implementation of additional efficient R-expr kinds. ${ }^{179}$

### 11.6.1 Efficient Disjunctions

The first efficient R-expr kind that we will look at is the efficient disjunction. We care a lot about the efficiency of disjunctions as they occur very frequently when translating a Dyna program into R-exprs, e.g., figure 11-11.

To start, the standard "inefficient" disjunction is defined as follows:

[^110]\[

$$
\begin{array}{c|cc}
{ }_{536} & a(0,0)=1 . & a(X, Y, V a l) \rightarrow(V a l=o n l y(\text { Inp }, \\
{ }^{537} & a(0,1)=2 . & (X=0) *(Y=0) *(\operatorname{Inp}=1)+ \\
{ }_{538} & a(1,0)=3 . & (X=0) *(Y=1) *(\operatorname{Inp}=2)+ \\
{ }^{539} & a(1,1)=4 . & (X=1) *(Y=0) *(\operatorname{Inp}=3)+ \\
& (X=1) *(Y=1) *(\operatorname{Inp}=4))) \\
& & \text { (b) Dyna } \text { R-expr }
\end{array}
$$
\]

Figure 11-11. Example Dyna program translated into an R-expr showing how ground values can be a frequent R-expr we must handle. Each rule in the Dyna program is converted into a branch of the disjunction, with the ground values present in the rules present in the R-expr as ground assignments to variables.

```
(def-base-rexpr disjunct [:rexpr-list args]
    ;; overridden functions on disjunction omitted
    )
```

Figure 11-12. Standard "inefficient" implementation of disjunction.

The declaration in figure 11-12 matches the disjunction we have used since chapter $\S 5$, with a disjunction of the form R1+R2+R3+R4 being represented as a list of $[R 1, R 2, R 3, R 4]$ inside of a disjunction R-expr kind. Unfortunately, this design has a fatal flaw. Finding the right disjunct requires scanning the entire list of disjunctions and applying Simplify to each sub-R-expr. For example, if we have the disjunct $(X=1) * R 1+(X=2) * R 2+(X=3) * R 3+(X=4) * R 4$, then finding a particular value of $X$ requires us to perform a linear time scan through all of the disjunctive $\mathbf{R}$-exprs.

To make the disjunction more efficient, we create an index for the assignment to the variable $X$. When $x$ has a known value, ${ }^{180}$ we can access the relevant sub-R-expr of the disjunct in $O(1)$ time by finding it using a data structure such as a hash table.

### 11.6.1.1 Requirements on the Efficient Disjunction Data Structure

To implement an efficient disjunction, which merges disjunction and ground assignments to variables, there are a few requirements that must be met. Currently, the

[^111]implementation only has hash-table-based tries as the only data structure backing the efficient disjunction (section §11.6.1.2). Future work should consider adding alternate implementations for disjunctions (section §16.1).

The data structure backing disjunctions combined with ground equality constraints needs to support the following:

1. In tracking the value assigned to a variable, the data structure needs to allow any ground value. For example, suppose we have an array containing disjunctive $\mathbf{R}$-exprs, where the index in the array corresponds with the assignment to a variable $X$. This would be able to represent the cases where $X$ is assigned a small integer, such as $(X=0) * R 1+(X=2) * R 2+\cdots$. However, the data structure also needs to handle the cases where $X$ is not a small integer, such as ( $X=$ "hello")*R3. This can be done by combining the array with some kind of auxiliary data structure to handle the cases where $X$ is not assigned an integer.
2. There needs to be some way to represent a "FREE VALUE". The reason is that we can have a disjunct like ( $X=1$ ) *R1+R2 where there is no known value of $X$ on the second branch of the disjunct. One solution is to represent the "FREE VALUE" internally using a special value that does not appear in $\mathcal{G}$ and, therefore, can be distinctly recognized. I will represent the "FREE VALUE" case using a question mark "?". This means that the previous $\mathbf{R}$-expr can be "normalized" into the disjunct ( $\mathrm{X}=1$ ) *R1+(X=?)*R2. ${ }^{181}$
3. There needs to be some way to efficiently make queries against the data structure that include the "FREE VALUE". This means queries of the form:

$$
\left(\left(X_{1}=1\right) \text { or }\left(X_{1}=?\right)\right) \text { and }\left(\left(X_{2}=2\right) \text { or }\left(X_{2}=?\right)\right) \text { and } \cdots \text { and }\left(\left(X_{n}=" N "\right) \text { or }\left(X_{n}=?\right)\right)
$$

The reason is that given an assignment $\left(X_{1}=1\right) *\left(X_{2}=2\right) * \cdots\left(X_{n}=" N "\right)$, we need to match all disjunctive branches that unify with this assignment. This includes branches where the assignment to a variable is known (e.g. $\left(X_{1}=1\right) *\left(X_{1}=1\right)$ ), and branches where there is no assignment to a variable (e.g. $\left(X_{1}=\right.$ ?) ). The only case that can be eliminated is when the assignment to a variable differs from the value we are querying for (e.g., $\left.\left(X_{1}=1\right) *\left(X_{1}=" n o t 1 "\right) \rightarrow 0\right)$.

[^112]
(b) Efficient Disjunctive Trie R-expr

Figure 11-13. A disjunctive $\mathbf{R}$-expr in (a) is converted into a trie in (b). Each level of the trie is associated with one of the variables (either $X$ or $Y$ in this example). Each level of the trie is an immutable hash map. This allows for efficient $O(1)$ retrievals in the case where a variable's value is known. A special value is associated with the case where the value of the variable is unknown, shown here as a question mark '?'. The leaf of the trie is a list that contains one or more disjunctive R-exprs.
4. There needs to be a way to "store" the R-expr in the data structure. In our case, the data structure is in memory, so we can store an opaque pointer to the R-expr. Future work could implement disjunctions with a disk-backed data structure or SQL database-backed disjunctions, in which case, they will have to implement serialization of R-exprs.

### 11.6.1.2 Efficient Disjunction using a Trie

The only data structure that we currently have for efficient disjunction is trie. Each level of the trie corresponds to a different variable. ${ }^{182}$ The leaf of the trie represents a disjunction of one or more sub-R-exprs. The $\mathbf{R}$-exprs in the leaves are treated as opaque pointers to $\mathbf{R}$-exprs. This is shown in figure 11-13.

By having lists at the leaf of the trie instead of an arbitrary $\mathbf{R}$-expr we can completely replace the standard disjunction with the trie-based disjunction. This is

[^113]useful as it allows us to integrate the trie disjunction into the R-expr rewrite with a single rewrite rule that runs at the construction of the standard disjunction. This automatically convert from the standard disjunction to the trie-backed disjunction, as shown in figure 11-14.

```
(def-base-rexpr disjunct-efficient [:var-list trie-variables
    :prefix-trie trie])
(def-rewrite
    :match (disjunct (:rexpr-list Rs))
    :run-at :construction
    (let [variable-order (vec (exposed-variables }\mp@subsup{}{}{183}\mathrm{ rexpr))]
        (make-disjunct-efficient variable-order
                        (convert-to-prefix-trie variable-order Rs))))
```

Figure 11-14. Declaration of the efficient disjunct, and a rewrite rule which converts from the standard disjunct to the efficient disjunct upon constructions. The variables on the trie are all exposed variables from the sub-R-exprs. The variable order is chosen arbitrarily (line 548). Future work should consider developing a variable ordering heuristic.

The path from the root of the trie to the leaves will have one level for each exposed variable in vars $(\cdot)$. This allows $\mathbf{R}$-exprs to be placed onto the most informative path of the trie without having secondary variable assignments contained in the $\mathbf{R}$-exprs in the leaves. To determine the assignments to the variables, recall that all the assignments to variables are in the context $\mathcal{C}$ and eliminated from the $\mathbf{R}$-expr itself (figure 11-7). This means that after invoking Simplify on each R-expr in the leaves, we are guaranteed that there are no assignments $\mathbf{R}$-exprs contained in the resulting $\mathbf{R}$-expr, and that all assignments to variables can be found using context $\mathcal{C}$. Any variable without a known assignment, as tracked by the context, is assumed to not have a value, and is placed under the "FREE VALUE", '?' branch.

### 11.6.2 Efficient Memoization uses Tries

In chapter $\S 10$, I introduced memoization using only abstract R-exprs. In chapter §10, I said that when an upstream dependency is changed, the system recomputes

[^114]the $\mathbf{R}$-expr using its original definition. If the system recomputes the entire $\mathbf{R}$-expr, this can be quite inefficient. The memoized $\mathbf{R}$-expr frequently represents hundreds of different disjunctive R-exprs (e.g. ( $X=1$ )*R1+(X=2)*R2++(X=n)*Rn).

To make memoization realistic, we will only recompute some subset of the memoized $\mathbf{R}$-exprs. In other words, if the currently memoized $\mathbf{R}$-expr is $(X=1) * R 1+(X=2) * R 2+$ $\cdots+(X=n) * R n$, then we might say that values which are memoized for $(X=2)$ are invalid and need to be recomputed. In which case, the system will modify the memoized $\mathbf{R}$-expr so that it gets $(X=1) * R 1+(X=2) * R 2$ new $+\cdots+(X=n) * R n$, where $R 2$ is replaced with R2new.

To implement this, we no longer treat the memoized $\mathbf{R}$-expr as a generic $\mathbf{R}$-expr. Instead, we explicitly use the efficient disjunctive trie (section §11.6.1) as the memoized R-expr. The trie supports "special rewrites" that implement efficient modification of the trie. These rewrites are exemplified by rewrite 86 and 87 , and allow us to delete entries matching a particular key ${ }^{184}$ from the trie, getting a new trie, or add $\mathbf{R}$-exprs to the trie.
$R_{\text {trie }}$ * $\operatorname{if}((X=1) *(Y=2), 0,1) \xrightarrow{86} R_{\text {new trie }} \quad \triangleright$ Delete from trie
$\mathrm{R}_{\text {trie }}+(\mathrm{X}=2) *(\mathrm{Y}=3) * \mathrm{~S} \xrightarrow{87} \mathrm{R}_{\text {new trie }} \quad \triangleright$ Add to trie

These "rewrite rules" make use of the trie implementation to share the internal structure with the existing trie as much as possible. These rewrites are invoked explicitly through special methods that are currently only supported by the trie.

### 11.6.3 Efficient Aggregation \& Projection

The second efficient R-expr kind we have is for aggregation and projection. So far, our rewrite strategy for projection and aggregation requires many steps of rewriting to rearrange them into a form that can be matched by the appropriate rewrite rules (rewrite rules 49 and 55). When considering aggregation R-exprs, there are often nested disjunctions and projections that need to be rewritten away before we can obtain the result of aggregation. For example, consider the program in figure 11-15.
${ }^{184}$ The key does not have to specify all of the variables.

```
a(Val) -> (Val=sum(Inp,
    proj(X, b(X, Inp))+
    proj(X,proj(Y,proj(Z, proj(Tmp1,proj(Tmp2,
        c(X,Y,Tmp1)*d(Y,Z,Tmp2)*
        times(Tmp1,Tmp2,Inp))))))))
```

551 a $+=\mathrm{b}(\mathrm{X})$.
$552 a+=c(X, Y) * d(Y, Z)$.
(b) R-expr

Figure 11-15. Dyna program with two contributions to the rule ' $a$ '.

This program has two contributions on lines 551 and 552 with each line having variables that are projected out. With the rewrite rules we have so far, we have to expand the projections using all possible assignments to the variables to solve this R-expr. (E.g. $\operatorname{proj}(X,((X=1)+(X=2)) * R) \rightarrow \operatorname{proj}(X,(X=1) * R)+\operatorname{proj}(X,(X=2) * R) \rightarrow$ $(R\{X \mapsto 1\})+(R\{X \mapsto 2\}))$

Ideally, the system would execute the $\mathbf{R}$-expr from figure 11-15 in a manner similar to the procedural implementation of the matrix-vector product in algorithm 5. Conceptually, the nested loops for figure 11-15 look something like the code in figure 11-16.

```
[ 1: }a\leftarrow0\quad\triangleright Initialize a container to hold the result of aggregatio
i- 2: for }x\in\operatorname{DomAIN}(\textrm{X},\textrm{b}(\textrm{X})):\triangleright Loop over the domain of projected variables from line 55
    (Inp=i)\leftarrow"SimpLIFY}((X=x)*\textrm{b}(\textrm{X}, Inp))" \triangleright Evaluate called user-defined rule
    a\leftarrowa+i D Record the contribution directly into the result container
for }x\in\operatorname{Domain}(\textrm{X},\textrm{c}(\textrm{X},\cdot)) :\triangleright eval line 552, start enumerating proj-ed var X's domain
        for y\in\operatorname{DomAIn}(Y,c(X,Y)): }\quad\triangleright\mathrm{ Enumerate projected var Y's domain
        \triangleright ~ E v a l u a t e ~ a ~ \mathbf { R } \text { -expr once enough arguments are known to trigger a rewrite}
        (Tmp1=t 1) \leftarrow"Simplify ((X=x)*(Y=y)*c(X,Y, Tmp1))"
        for z\in DomAin(Z, d(Y,Z)): }\triangleright\mathrm{ Enumerate projected var Z's domain
            (Tmp2= t2) \leftarrow"SIMPLIFY}((Y=y)*(Z=z)*d(Y,Z, Tmp2))"
            i\leftarrowt\mp@subsup{t}{1}{**" tr }2, \trianglerightEvaluate times(Tmp1,Tmp2,Inp)
            a\leftarrowa+i }\quad\triangleright\mathrm{ Once all R-exprs are evaluated, record the contribution
return }
```

Figure 11-16. An informal pseudocode example of the idealized evaluation of the rule 'a' from figure 11-15. First, a container is initialized to hold the result of aggregation on line 1. Then, the system sequentially runs the sub-R-exprs of the disjunction. For the first disjunct, the system loops over the entire domain of the projected variable $x$, line 2 . How the system figures out the domain of the variable will be discussed in section §11.7. During each iteration, the R-expr that represents a rule is Simplified. The resulting value is directly saved directly into the aggregator's accumulator variable $a$, line 4 .

For the second disjunct, we have multiple variables projected out, $\mathrm{X}, \mathrm{Y}$ and Z. The system loops through the domains of the first two variables on lines 5 and 6. Once it has enough information to evaluate an $\mathbf{R}$-expr, it is immediately evaluated, as on line 8. The remaining variables are still looped over (line 9) with the remaining R-exprs also being evaluated (line 10). Just like before, the contribution to aggregation is saved directly in the accumulator variable $a$ (line 12).

Finally, after evaluating the entire body of the aggregator, the result of the aggregation is returned (line 13).

To make the execution with $\mathbf{R}$-exprs like what is shown in figure 11-16, we will introduce two different efficient $\mathbf{R}$-expr kinds. The first $\mathbf{R}$-expr kind will be the outer aggregator, which initializes the container for aggregation and returns the final result of the aggregation. The outer aggregator corresponds with a term, and will combines the result from multiple rules. This corresponds with the solid red lines in figure 11-16 and is represented by lines 1 and 13. The second R-expr kind

```
a(Val) -> (Val=sum_aggregator_outer(
    sum_aggregator_inner([X], Inp, \trianglerightList of projected variables and input variable
        b(X, Inp) )+ \trianglerightBody of first inner aggregator
    sum_aggregator_inner([X,Y,Z,Tmp1,Tmp2], Inp,
        c(X,Y,Tmp1)*d(Y,Z,Tmp2)* }\triangleright\mathrm{ Body of second inner aggregator
        times(Tmp1,Tmp2,Inp)) ))
```

Figure 11-17. The R-expr from figure 11-15 with its aggregator and projections replaced with the efficient aggregator.
will be the inner aggregator, which simultaneously handles projection and nested loops over variables and will also compute the contribution to the aggregator. A single outer aggregator can have multiple inner aggregators internally, as in this example. The inner aggregator corresponds with a single rule, and terms can be defined using multiple rules. The inner aggregators correspond with the dashed blue lines in figure 11-16 with lines 2-4 and lines 5-12.

The R-expr representation for figure 11-15 with an efficient aggregator is shown in figure 11-17.

Note that in figure 11-16, the aggregator accumulator variable $a$ is initialized by the outer aggregator but used by the inner aggregator. The accumulator variable is passed between the outer and inner aggregator through a "globally" ${ }^{162}$ accessible pointer. Hence, the inner aggregator does not have well-defined R-expr semantics on its own. Instead, the inner aggregator is only well defined when it is nested inside an outer aggregator as in figure 11-17.

The implementation of the efficient aggregator is similar to the efficient disjunction. We define new R-expr kinds that hold the necessary details for the aggregator. A rewrite rule that runs at construction matches the standard aggregator and converts it into the efficient representation (figures 11-18 and 11-19).

```
(def-base-rexpr aggregator-outer [:string operator }\mp@subsup{}{}{185
    :var result
    :rexpr body])
(def-base-rexpr aggregator-inner [:hidden-var incoming
    :hidden-var-list projected
    :rexpr body])
(def-rewrite
    :match (aggregator operator }\mp@subsup{}{}{185}\mathrm{ (:var result) (:var incoming)
                            (is-true? is-conjunctive) ;; Recall §6.5.1
            (:rexpr body))
    :run-at :construction
    (make-aggregator-outer operator result
        (make-aggregator-inner incoming [] body)))
```

Figure 11-18. Declaration of the outer and inner efficient aggregator $\mathbf{R}$-expr, and rewrite rule which converts from the standard aggregator into the efficient aggregator.

The inner aggregator also uses several rewrite rules that match the body of the inner aggregator to move the inner aggregator closer to the computation performed by the $\mathbf{R}$-expr. The inner and outer aggregators can be separated by disjunctions and if-expressions (memoization). An inner aggregator nested under a disjunction corresponds to a single Dyna rule that partially defines a user-defined term. An example of rewrites that embed the inner aggregator is shown in figure 11-19.

When running the rewrites for the inner and outer aggregators, the outer ag-

[^115]```
(def-rewrite
    :match (aggregator-inner (:var incoming) (:var-list projected)
                            (proj (:var projv) (:rexpr body)))
    :run-at :construction
    (make-aggregator-inner incoming (cons projv projected) body))
(def-rewrite
    :match (aggregator-inner (:var incoming) (:var-list projected)
        (disjunct Rs))}\mp@subsup{)}{}{186
    :run-at :construction
    (make-disjunct
        (map (fn [R] (make-aggregator-inner incoming projected R))
            Rs)))
```

Figure 11-19. The inner aggregator rewrites its body to embeds itself in the $\mathbf{R}$-expr as close as possible to the $\mathbf{R}$-expr body which corresponds with a single Dyna rule. It matches against projection (line 567) and merges those variables into the inner aggregator's own projected variable list. For disjunction, it moves towards the leaves of the disjunct (line 572).
gregator sets up the aggregation accumulator where the intermediate result of aggregation is stored. The inner aggregator is only responsible for handling a single rule body. When the inner aggregator is completely rewritten its body and added its contribution to the outer aggregator's accumulator, it will be rewritten as 0 . Rewriting as 0 indicates that no delayed $\mathbf{R}$-exprs still need to be evaluated. ${ }^{189} \mathrm{~A}$ conceptual example of the rewrite rules that handle the inner and outer aggregator is shown in figure 11-20.

In the next section, I will define iterators, which is how the "Domain $(\mathrm{X}, \mathrm{b}(\mathrm{X})$ )" calls in figure 11-16 are implemented.

[^116]```
(def-rewrite
    :match (aggregator-outer operator result body)
    (let [;; stash parent frame aggregation accumulator container
                parent-accumulator (get-global *aggregator-accumulator*)]
            ;; initialize aggregation container into global
        (set-global *aggregator-accumulator* nil)
        (let [new-body (simplify body) ;; run rewrites on body
                    accumulator (get-global *aggregator-accumulator*)]
            ;; restore the parent frame aggregation accumulator into
        global
            (set-global *aggregator-accumulator* parent-accumulator)
            (if (= new-body (make-multiplicity 0))
                (do ;; done running the inner aggregators, return the value
                    (make-unify result (make-constant accumulator)))
                (do ;; inner aggregators not finished, return delayed
                    (make-aggregator-outer }\mp@subsup{}{}{187}\mathrm{ operator result
                (if (= nil accumulator)
                                    new-body
                                    (make-disjunct [new-body
                                    (make-aggregator-inner (make-constant accumulator)
                                    [] (make-multiplicity 1))]))))))))
(def-rewrite
    ;; match R-expr like `sum_aggregator_inner([], Inp, (Inp=7))`
    :match (aggregator-inner incoming []
                        (unify incoming (:ground value)))
    (do
        (set-global *aggregator-accumulator*
            (+ }\mp@subsup{}{}{188}\mathrm{ (get-global *aggregator-accumulator*)
                (get-value value)))
        (make-multiplicity 0)))
```

Figure 11-20. These example rewrites are only intended to be illustrative of how the inner and outer aggregators work together. Many details have been omitted. There is a single global *aggregator-accumulator* variable. This works fine, as the inner aggregator is always associated with the nearest outer aggregator. The aggregator accumulator from a higher scope is stashed in the local variable parent-accumulator as to not interfere with the parent's partially computed aggregation (line 581). Simplify is recursively called on the body R-expr (line 584). When an inner aggregator has been completely rewritten by SIMPLIFY, the resulting contribution is added to the global accumulator variable (line 603). The inner aggregator will rewrite itself as 0 (line 606). Here, rewriting as 0 indicates that there are no delayed $\mathbf{R}$-exprs, and that the result of the aggregation has been handled. This is conceptually similar to when the unify $\mathbf{R}$-expr rewrites as 1 once the assignment to a variable is tracked via the context (figure 11-7). ${ }^{189}$

### 11.7 Iterators

In chapter $\S 9$, I defined the optional constraint R-expr, which lifts useful conjunctions out of nested R-exprs. However, recall that optional constraints are only used to license various rewrites that are useful for rearranging the R-expr into something that can be rewritten. Iterators, as described in this section, are an implementation of optional constraints that work with variables that have a known finite domain. For example, iterators correspond with optional constraints of the form opt $((X=1)+(X=2)+\cdots+(X=1000))$, where this is a disjunction that represents 1000 different assignments to the variable X . Iterators are more limited than the optional constraints from chapter §9. However, by limiting iterators to only focus on ground assignments to variables, we can build efficient class abstractions and implementations. For example, an iterator can be backed by the efficient disjunctive trie (section §11.6.1.2), in which case it will stream the values assigned to a variable, as known by the trie. This is akin to having an iterator over the keys in a hash-table.

In addition to iterators having efficient implementation, iterators allow us to generate JIT-compiled code that is conceptually similar to the example code in figure 11-16. This is tractable as iterators and assignments to variables can be held in local variables of generated code at runtime, in the same way that is done when writing a procedural program. Compilation of $\mathbf{R}$-expr and will be discussed extensively in chapter $\S 12$.

### 11.7.1 Iterator Interface

Every R-expr defines its own getlterables method. The getlterables returns a set of IterablesWithMetadata that are supported by the R-expr (figure 11-21). Calling gettterables on an R-expr $R$ is equivalent to rewriting $R$ as $R * o p t(\cdots) * o p t(\cdots)$, using the rewrites defined in chapter §9, where opt $(\cdots)$ will represent a single iterator. The getlterables method's definition is informed by the R-expr's behavior with respect to rewriting optional constraints. For example, the getlterAbLES on a conjunction simply returns the union of all iterables. This follows as $(R * o p t(R o p t)) *(S * o p t(S o p t)) \rightarrow(R * S) * o p t(R o p t) * o p t(S o p t)$ is a valid rewrite which simply rearrange optional constraints inside of the conjunction.

The iterator interface is designed to be roughly comparable to iterating over

```
interface IterableWithMetadata {
    List<Variables> VariableOrder(); // order variables are assigned
    Set<Variables> VariablesCanEnumerate(); // variables with finite domains
    Iterator iterator(); // The iterator
}
interface Iterator {
    LazySeq<IterInstance> Run(); // Run without duplication
    LazySeq<IterInstance> RunNotFinite(); // Allow duplication
    Iterator BindValue(Object value); // directly assign value, no iteration
    int estimateCardinality(); // estimate of number of values returned
}
interface IterInstance {
    Object value(); // Value assigned to the iterated variable
    Iterator continuation(); // Iterator for next variable
}
interface LazySeq<IterInstance> { // LazySeq is provided by Clojure
    IterInstance head(); // Return the current value
    LazySeq<IterInstance> tail(); // Return tail, which contains next value
}
```

Figure 11-21. Java Interfaces for iterables and iterators.
tries, ${ }^{190}$ (figure 11-21). The IterableWithMetadata interface provides information about which variables the iterable supports. When the iterable is started, using the iterator method on line 610, the Iterator returned only iterates over the domain of a single variable (line 612). This corresponds to iterating over all of the value bindings to a variable at the first level of the trie. If we use the trie from figure 1113 b as an example, the first iterator will iterate over the bindings for the variable x , which are $\{?, 1,3\}$. To obtain an iterator for a second variable, the continuation method on line 620 is used. The second iterator is conditioned on the current value returned by the first iterator. To continue using the trie from figure 11-13b as an example, when the first iterator has $X=3$, the continuation iterator over $Y$ will loop over the values $\{4,9\}$, which correspond to the set of possible values that $Y$ can be assigned condition on ( $X=3$ ).

The order in which iterables loop over variables is indicated by the VariableOrder

[^117]method on line 608. The variable order typically corresponds to the order in which variables appear on the underlying data structure (which is usually a trie). In the case that the variable order supported by an IterableWithMetadata is not the desired variable order, temporary materialized tables can be created to rearrange iteration order of variables.

When using the iterator, we ideally want to only iterate over variables that have a known finite domain. Recall that this corresponds to an optional constraint of the form $\operatorname{opt}((X=1)+(X=2)+(X=3))$. The VariablesCanEnumerate method (line 609) returns the set of variables that have known finite domains. When a variable does not have a finite domain, the corresponding optional constraint has the form opt $((X=1)+(X=2)+1)$, which is not useful for splitting an $\mathbf{R}$-expr into smaller non-overlapping $\mathbf{R}$-exprs (previously discussed in section §9.2.2). When a variable does not have a finite domain, the iterator interface still allows the variable to be iterated. However, special care is required to handle the "FREE VALUE", which will be returned (indicating that any value is allowed to be assigned to the variable). For the purposes of implementation, and in an attempt to limit potential bugs, the Run method on line 613 cannot be used when a variable does not have a finite domain. Instead, a second method RunNotFinite on line 614 is used instead. Furthermore, the Run method guarantees that a value will not be repeated, whereas no guarantee is provided by the RunNotFinite method.

This design of iterating over a single variable at a time was chosen so that the system can interleave rewriting with Simplify with iteration. This was hinted at in figure 11-16, where some simplification on line 8 was performed before the variable Z was iterated.

Finally, an iterator can also be used as a filters. This is done using the BindValue method on line 615. The BindValue function allows us to avoid iterating through the domain of a variable at a particular level of the trie and directly go to the next variable. For example, if we have the $\mathbf{R}-\operatorname{expr}(X=3) * R_{\text {trie }}$, we do not need to use $R_{\text {trie }}$ to iterate over values of $X$. Furthermore, if $R_{\text {trie }}$ does not contain a branch for $(X=3)$ (meaning $(X=3) * R_{\text {trie }} \rightarrow 0$ ), then BindValue will return null, which can be used for filtering impossible values. Calling BindValue is generally more light weight than rewriting the R-expr with Simplify. Hence, when we have a conjunction of multiple iterators, the BindValue function can be used to filter out impossible values before rewriting is even attempted.

### 11.7.2 Different Kinds of Iterators

Inside of the Dyna implementation, there are many different classes that implement iterators (figure 11-21 only shows interface). Some of these iterators come from built-ins. For example, the integer range built-in's iterator enumerates integers. In this case, the iterator has its own class implementation, which efficiently represents the current integer as well as the min/max integer it is iterating to.

Outside of the built-ins, some core $\mathbf{R}$-expr kinds also interact with the iterators. For example, projection and aggregation will project out variables from an iterator when the variable is projected out in the $\mathbf{R}$-expr representation. In other words, if we have an $\mathbf{R}$-expr like in figure 11-22:

```
proj(X, (X=1)*((Y=1)+(Y=2)+(Y=3))+
    (X=2)*((Y=7)+(Y=2))+
    (X=3)*((Y=9)+(Y=1)))
```

Figure 11-22. Example R-expr with projection of $x$

When the system calls getlterables on the above R-expr in figure 11-22, it will only get an iterable over the variable $Y$. This corresponds to the optional constraint opt $((Y=1)+(Y=2)+(Y=3)+(Y=7)+(Y=9))$ pulled through the projection to the top of the $\mathbf{R}$-expr. This is handled by the implementation of getlterables for projection and has wrapping iterable, and iterator classes that modify the iterables returned from the body of the projection. The wrapping iterator will attempt to stream the iterated values of $Y$, instead of materializing the values of $Y$ into a data structure.

### 11.7.3 Iterators Attempt to Stream Values

The intention of iterators is to be able to make use of the underlying data structures without having to materialize everything as an R-expr. Much like streaming APIs and libraries, the iterator module has a number of different kinds of iterator types that wrap an iterator and modify it as needed. As already discussed for figure 11-22, the Dyna implementation has an iterator that will project out a variable from a wrapped iterator. The wrapper modifies the return values from the VariableOrder and VariablesCanEnumerate method from the IterableWi thMetadata (figure 11-
21) such that the projected out variable is not visible outside of the projection. In the case that an iterator's variable order is "unfavorable" to the projection (such as the variable order is $[X, Y, Z]$ and we are projecting out the variable $X$ ), then the wrapper will consolidate the values of $Y, Z$ using a temporary data structure when the Run method is invoked (line 613). Alternately, if the RunNotFinite method is used, then no consolidation is required (line 614).

## Conjunction of iterators

In some cases, we might want to combine two or more iterators on the same variable. For example, with a conjunction of two iterators from different sources, we might use one of the iterators as a filter as in figure 11-23. Checking that an iterator can bind to a particular value can be much faster than using Simplify to rewrite the $\mathbf{R}$-expr.

```
: for }\langlex,\mathrm{ continuation_1 }\rangle\in\mathrm{ iterator_1.Run():
    if iterator_2.BindValue ( }x\mathrm{ ) }\not=\mathrm{ null :
        yield value }
```

Figure 11-23. Merging conjunctive iterators using iterator_2 to filter out values which will not work (would cause the R-expr to rewrite as 0 ).

We can also use a conjunction of iterators in the case that we do not have a single iterator that can iterate over all of the variables that we are interested in. In this case, a second iterator can be used as in figure 11-24.

```
for }\langlex,continuation_1\rangle\initerator_1.Run() :
    if continuation_2 = iterator_2.BindValue(x) :
            for }\langley,_\rangle\in\mathrm{ continuation_1.Run() :
                for }\langlez,_\rangle\in\mathrm{ continuation_2.Run() :
            yield values }x&y&
```

Figure 11-24. If we have iterator_1 over $X \& Y$, and iterator_2 over $Y$ \& $Z$, then we can iterate over the variables $X, Y \& Z$ by combining these two iterators.

## Disjunction of iterators

To handle a disjunction like

$$
((X=1)+(X=2)+(X=3))+((X=1)+(X=7)+(X=11))
$$

we can have iterators from each branch of the disjunction. These iterators can be merged by first iterating through the domain of the first iterator and then iterating through the domain of the second iterator, using the first iterator to filter out values already seen. Assuming that the iterator efficiently implements the BindValue method, then this will be more efficient than creating an intermediate data structure for the values of x . This is shown in figure 11-25.

```
for }x\in\mathrm{ iterator_1.Run():
    yield value }
for }x\in\mathrm{ iterator_2.Run() :
    if not iterator_1.BindValue(X): }\\mathrm{ : Filter out values from iterator_1
        yield value X
```

Figure 11-25. Iterating through a disjunction of iterator_1 and iterator_2. The first iterator is used to filter out values of $x$ that have already been seen on line 4.

Note that the filtering in figure 11-25 corresponds to the more generally applicable if-expression rewrite rule 72. In this case, given the constraint if (Q1+Q2,1,0) (which can come from an optional constraint), is rewritten as if $(Q 1+Q 2,1,0) \rightarrow i f(Q 1,1,0)$ $+\operatorname{if}(Q 2,1,0) * i f(Q 1,0,1)$, which shows that we can split the disjunctive constraint (conditional of the if-expression) if we use if $(Q 1,0,1)$ to filter out redundant values.

### 11.7.4 Using Iterators

The iterator subsystem provides a run-iterators macro as an internal API that automatically determines how to use the available iterables to subdivide the R-expr. The run-iterators macro takes a list of variables that the system wants grounded (such as the list of projected variables in the case of aggregation), as well as a set of iterables returned by the Getlterables method. The run-iterators macro automatically chooses which iterators to use. Run-iterators can either select a single iterator or combine multiple iterators in the case where no single iterator will work. When there are multiple iterators that can iterate the same variable, ${ }^{191}$ the

[^118]run-iterators macro makes use of the estimateCardinality method (line 616) to select the iterator that reports the smallest estimated cardinality. When there does not exist an iterables that guarantees that some requested variable is grounded, then run-iterators can attempt a best effort iteration binding other variables. For example, if run-iterators is requested to ground the variables $X, Y$, and $Z$, where there only exists an iterator over $Y$ and $Z$, the system is allowed to iterate over $Y$ and $Z$ in hopes there is a constraint like plus $(X, Y, Z)$ contained in the $\mathbf{R}$-expr that is capable of determining the value of $X$ from an assignment to $Y$ and $Z$.

### 11.7.5 Efficient Aggregation uses Iterators

In figure 11-16, the aggregator makes use of a function of the form $\operatorname{Domain}(X, b(X))$ to loop over the domain of a variable. This comes from the iterator subsystem and is the run-iterators macro in the implementation. When run-iterators chooses the order in which variables will be bound, it is informed by the iterators available. (The iterator's variable order is informed by the order in the data structure's store data, such as in the trie.)

When a variable $X$ 's value is assigned, the aggregator will rewrite the $\mathbf{R}$-expr using the information $(X=x)$. If the $\mathbf{R}$-expr is rewritten as 0 , then this means that the current assignments of variables are inconsistent with the $\mathbf{R}$-expr, and the aggregator will skip the value $x$ and try the next possible assignment to $X$ without having run more deeply nested loops over variables. Attempting to rewrite the $\mathbf{R}$-expr before all variables have been bound is akin to the optimization of lifting operations outside of loops or propagating new bindings to variables once available in constraint logic programming through the constraint store.

The inner aggregator is responsible for performing the aggregation inside of the loop once all variables have been bound and the $\mathbf{R}$-expr is completely rewritten. If the $\mathbf{R}$-expr body of the aggregator is not rewritten completely, then the partially rewritten $\mathbf{R}$-expr is preserved by the aggregator in the hope of being rewritten later.

## Chapter 12

## Compilation of R-expr Rewrite Strategies

One unfortunate thing about research sometimes is that despite making significant progress, we can still find ourselves with an unsatisfactory system. In this case, the realistic implementation described in chapter $\S 11$ is easily 100 to 1000 times faster ${ }^{192}$ than the pure, minimal implementation presented in chapter $\S 8$. However, despite that, the "realistic implementation" is still too slow for the use cases that we care about. ${ }^{193}$

In hindsight, this is not terribly surprising. We can think of the R-expr as akin to an internal representation inside a compiler. Our rewrites are, therefore, akin to optimization passes. For example, a rewrite like 28 , plus $(1,2, x) \rightarrow(X=3)$ is essentially constant propagation. Unsurprisingly, an optimizing compiler being used to evaluate the entire program is much slower than the equivalent procedural instructions.

[^119]|  | LOAD_GLOBAL | $0(\mathrm{~b})$ |
| :---: | :--- | :--- |
| ${ }^{626} \mid \mathrm{a}=\mathrm{b} * \mathrm{c}$ | LOAD_GLOBAL | $2(\mathrm{c})$ |
| (a) Python | BINARY_OP | $5(*)$ |
|  | STORE_FAST | $0(\mathrm{a})$ |

(b) Python Bytecode

Figure 12-1. Example Python with procedural instructions

As such, in this chapter, I will discuss my efforts to compile R-expr rewriting. When rewriting $\mathbf{R}$-exprs that come from Dyna programs, there will be many similar R-expr "shapes" and sequences of rewrites performed. Hence, it is reasonable to believe that we can speed up the Dyna implementation by compiling R-expr rewrites.

The goal of our implementation is to maintain the flexibility of $\mathbf{R}$-exprs while running faster. The work presented here focuses on removing the overhead of the rewrite system and is not a "final stage" compiler. Instead, our goal is to generate procedural code instead of individual rewrites that should be comparable to what a user would write in a procedural language. ${ }^{194}$ Before we start looking at how we generate code, let us briefly discuss what some of the overheads are that we are aiming to eliminate.

### 12.1 What is Overhead?

Let us start by defining what "overhead" means in the context of R-exprs and executing Dyna programs. Consider the line of code in a procedural programming language in figure 12-1a. To evaluate figure 12-1, we are reading from two variables and calling the multiplication operation. If this was converted into efficient assembly instructions, we might have two read-from-memory-into-register instructions, followed by a multiplication instruction and then a final store instruction. These four instructions in figure 12-1b represent the required operations to evaluate this line of code.

[^120]Now, suppose that all variables are stored in a hash-map. Instead of performing a single read to load the variable, the system indirects through multiple operations performed by the hash-map: reading the base pointer of the hash-map, computing the hash code of the string " $b$ ", reading the size of the hash-map, computing the hash("b") mod the size of the hash-map, reading an entry in an array, checking equality between the string "b" and the entry found in the array. In this case, every operation the hash-map performs can be considered overhead. Now, overhead does not mean wasted. In this case, using a hash-map makes Python's implementation simpler. However, it does not contribute to the "useful work" of multiplying numbers.

### 12.1.1 Overhead with R-exprs Rewriting

When running a Dyna program represented as $\mathbf{R}$-exprs, there are similarly some operations that correspond to manipulations of the data that we are interested in and other operations that correspond to the overhead of $\mathbf{R}$-exprs. For example, some overheads we have with $\mathbf{R}$-exprs are:

1. Manipulations and accesses of the context $\mathcal{C}$. The context contains both variable bindings and other conjunctive $\mathbf{R}$-exprs. The assignments to variables are stored in a hash-map. ${ }^{195}$ This means that matching the : free and : ground preconditions requires accessing the context's hash-map. Similarly, when the context is used to find conjunctive $\mathbf{R}$-exprs, those $\mathbf{R}$-exprs are stored in sets. Finding the right conjunctive $\mathbf{R}$-expr requires searching through any relevant set. Additionally, conjunctive R-exprs are added to these sets when encountered.
2. To identify a rewrite, the system checks a list of potential rewrites. ${ }^{196}$ Rewrites are segmented by the kind of R-expr, but operations such as checking the ground/free state of variables use the context. Conjunctive $\mathbf{R}$-exprs require scanning through the context.

[^121]

Figure 12-2. This is a flame graph for CKY parsing using memoization and forward chaining of a small sentence without using the compiler. From this graph, we can observe which functions take the most time to run. We can observe two operations used by memoization (chapter §10): sending messages for updates and recomputation of values. Both of these operations use SIMPLIFY to rewrite specially crafted R-exprs to perform the bulk of the work. We can observe that the R-expr rewriting occupies $98 \%$ of the runtime.
3. The results of rewrites are either R-exprs or assignments to variables that are tracked by modifying the context (such as in figure 11-6). In both cases, this means creating many new Java objects.
4. Variable renaming and other similar manipulations of $\mathbf{R}$-exprs require scanning the entire $\mathbf{R}$-expr to perform the necessary rewrites. A new $\mathbf{R}$-expr is created and returned.
5. Finding iterators is a recursive function that returns a set of iterators. This creates many intermediate sets and iterator types. This happens every time we evaluate any disjunction under an aggregation.
6. Memoization and update messages are handled via a priority queue. ${ }^{197}$ Therefore, the push and pop operations take $O(\log n)$ time where $n$ is the number of updates messages pending. Comparably, a dynamic program with a fixed execution order does not require an agenda at all and entirely avoids this overhead. ${ }^{198}$

This is not intended as an exhaustive list of the sources of overhead. Furthermore, not all cases of overhead are "bad" or something we need to deal with. For example, if we allocate and deallocate a Java object within the same compilation unit, then the Java compiler can eliminate the allocation altogether. Unfortunately, given the design of Simplify from chapter $\S 11$, this optimization is unlikely to happen, as compilation units are likely only able to get as large as a single rewrite rule.

### 12.2 Compilation Overview

In the remainder of this chapter, I will detail how we compile R-exprs and their rewrites.

Our goal in compiling is to make the rewriting $\mathbf{R}$-exprs run faster. To make compilation work, we need compilable units and reentrancy so that the compiled units are reused. The most obvious approach would be to compile the Simplify

[^122]function, which is similar to a bytecode interpreter where R-exprs are the "bytecode". However, this has the same problems we had with memoization (section §10.2). It requires matching against a large part of the $\mathbf{R}$-expr to figure out when there is reentrancy. Instead, we will take an approach similar to the one taken with memoization. We will replace a sub-R-expr with a JIT-generated $\mathbf{R}$-expr kind that represents a compiled unit, such as in figure 12-3.

Generated $\mathbf{R}$-exprs and generated $\mathbf{R}$-expr rewrites will be generated using a just-in-time compiler (JIT) and are made up of multiple primitive $\mathbf{R}$-exprs and $\mathbf{R}$-expr rewrites (the $\mathbf{R}$-exprs and rewrites defined previously in chapters 5 and 6).

As we will see, our approach can be broken down into these two major components, which work in tandem:

1. Generating new JIT-generated R-expr kinds, which represent larger R-expr expressions state. These JIT-generated $\mathbf{R}$-exprs correspond to compilable units of code. This allows us to match with a larger $\mathbf{R}$-expr unit all at once. Therefore, we can save time by not having to match each R-expr individually. When a particular R-expr state is reencountered, we will reuse the generated $\mathbf{R}$-expr kind, enabling reentrancy (section §12.4.2).
2. Generating rewrites for the JIT-generated $\mathbf{R}$-expr kinds. The $\mathbf{R}$-expr kind on its own does not represent execution but rather a state of the program. Hence, we need to generate rewrites for the new $\mathbf{R}$-expr kinds. These rewrites correspond with multiple primitive rewrites being performed against the underlying R-expr state. Performing a single match allows us to avoid the overhead of performing each match individually and serializing the state as an $\mathbf{R}$-expr between rewrites.

### 12.3 Generating New JITted R-expr Kinds

The mechanism for generating a new $\mathbf{R}$-expr kind is conceptually straightforward. Every JIT-generated $\mathbf{R}$-expr kind is given a name like state1234, ${ }^{199}$ and corresponds to some primitive $\mathbf{R}$-expr state, which is tracked via external metadata. Primitive $\mathbf{R}$-expr kinds are those previously defined in chapter $\S 5$. The generated $\mathbf{R}$-expr

[^123]will have holes corresponding to variables and R-exprs not statically knowable. For example, exposed variables that are referenced by other $\mathbf{R}$-exprs may have different names in different contexts. Therefore, the generated $\mathbf{R}$-expr will have a hole representing a variable. Variables that are only used internally and projected out do not require a hole in the JITted $\mathbf{R}$-expr because they only represent internal data. For example, the variable $X$ in $\operatorname{proj}(X, \operatorname{times}(X, Y, Z) * \cdots)$ is only used locally, but the variables $Y$ and $Z$ are in the externally visible vars $(\cdot)$ of this expression, therefore, will be represented with holes.

Holes are typed as we saw in section §11.5.1, and can include sub-R-exprs, value types, and hidden variables. ${ }^{200}$ This is very important as it allows a generated $\mathbf{R}$-expr to represent part of a larger $\mathbf{R}$-expr and does not require that the entire $\mathbf{R}$-expr is represented at the same time. In other words, if we think of an $\mathbf{R}$-expr as a recursive tree data structure, then the generated $\mathbf{R}$-expr is allowed to represent a sub-tree all the way down to the leaves, and it is also allowed to represent a sub-tree that stops before it gets all the way to a leaf. This allows us to 1 ) create smaller JIT-generated units that are more likely to be reusable and 2) we can exclude R-exprs that we do not want to support in the JIT (I will revisit this point in section §12.5).

An example of creating a JIT-generated $\mathbf{R}$-expr is shown in figure 12-3.

### 12.4 Generating New Rewrites

The second part of JIT-compiling an $\mathbf{R}$-expr is to generate the corresponding rewrites. The high-level approach is inspired by tracing JIT compilers ([31, 74, 75], section $\S 3.6)$. A tracing JIT works by compiling a trace of operations that are performed rather than generating using the definition of a particular method. In the case of rewriting, a trace will correspond to a sequence of one or more rewrites applied to the state represented by the JIT-generated $\mathbf{R}$-expr.

To identify which rewrites can be applied, the JIT uses the same information that Simplify uses to match a rewrite. This includes the context $\mathcal{C}$, which contains information about conjunctive $\mathbf{R}$-exprs and the values assigned to the variables. To gain access to the context, we invoke the JIT-compiler from Simplify. In other

[^124]

Figure 12-3. This figure shows the first step in JIT compiling an $\mathbf{R}$-expr. The center-dot (•) represents a pointer to another $\mathbf{R}$-expr. We identify a subset of an $\mathbf{R}$-expr (represented by the red dashed box in the middle) that we want to compile into a single state. The selected $\mathbf{R}$-expr does not have to include all children $\mathbf{R}$-exprs, as represented by ' $g$ ( Y , Temp2)', which is not included in state687. This allows us to avoid including parts of the $\mathbf{R}$-expr that are unlikely to benefit from being compiled into the state. Variables such as X , Result are referenced by the external environment, and the variable Temp2 is hidden and referenced by the $\mathbf{R}$-expr in the hole. Variables that are only internal, such as Temp1, do not have variable slots to track their name.
words, when Simplify encounters a JIT R-expr kind that does not have a generated rewrite, ${ }^{201}$ it invokes the JIT compiler with the $\mathbf{R}$-expr and the current context. The process of creating JIT-generated rewrites therefore happens lazily-as is typically done when JIT compiling a program.

Any rewrite generated by the JIT compiler is immediately applied to the R-expr, in the usual way that Simplify applies rewrites to R-exprs. However, the newly generated rewrite will also be added to the collection of available rewrites that the system has. This means that the same rewrite can be applied in the future without requiring additional code generation. An example of this is shown in figure 12-4 with a rewrite being generated for the R-expr from figure 12-3 and applied immediately.

### 12.4.1 Combing Multiple Rewrites Into One

One of the core features of the JITted rewrites is that we can combine multiple steps of rewriting together into a single JIT-generated rewrite. This limits the overhead of matching and allows us to pass values between primitive rewrites using local variables.

The way the JIT compiler generates rewrites differs from what we have seen with the primitive rewrites from chapter §11. First, with the primitive rewrites, there is a single : match statement (section §11.5.3). The match statement is matched in full before the rewrite is applied. If we were to apply this same design to the JIT-generated rewrites, then we would end up with rewrite rules that are not widely applicable, as it would require matching all of the preconditions before anything can be rewritten. For example, suppose that we combine a sequence of 20 rewrite rules together. If we have to match against all 20 different conditions first, then it is likely that one of the rewrite's preconditions will not match, and the JIT-generated rewrite will get poor reuse.

To get around this issue, we alternate between checking preconditions and performing rewrites. A rewrite's preconditions are checked right before the rewrite

[^125]

Figure 12-4. A new JIT rewrite is generated for state687( $\cdots$ ) from figure 12-3. The JIT is invoked by the Simplify method, which recursively walks the $\mathbf{R}$-expr tree. When the JIT is invoked by Simplify, it has access to the current R-expr state687( $\cdots$ ) and the context $\mathcal{C}$. Hence, the JIT can see that $X$ is assigned the value ( $\mathrm{X}=7$ ). The JIT looks for rewrites that can be applied to the "primitive R-expr state" represented by state687( $\cdots$ ). In this case, the only rewrite that can be applied is to rewrite times( $3, X$, Temp1).

Observe that the result of this rewrite, state $865(\cdots)$, no longer has a reference to the variable $X$. The reason is that there are no $\mathbf{R}$-exprs in state $865(\cdots)$ that still refer to the variable $x$. Furthermore, we previously had that Temp1 was a hidden variable (due to the projection), but now that it has a known value that is not statically knowable, that value needs to be saved somewhere. Therefore, state865( $\cdots$ ) has an argument (hole) for the value assigned to Temp1, in this case that value is 21 .


Figure 12-5. This represents a single rewrite. The blue arrow in the middle is the trace of 4 different rewrites. The trace will include all possible rewrites from the time of compilation, starting from the initial state. The preconditions or match statements for a rewrite are represented as red lines and are checked before the rewrite is applied. If a precondition fails, then it takes branches to a generated $\mathbf{R}$-expr state. JIT-generated $\mathbf{R}$-expr states such as states $32,41,88$, and 72 are generated on demand at the same time that the rewrite was generated. If an existing state matches, then it will be reused rather than generating a new state. Preconditions that can be statically resolved to be true, either because they were already checked by another rewrite or because they consume statically known output from an earlier rewrite, do not generate any code in the generated rewrite (represented in gray).
is applied. If one of the rewrite's preconditions has been previously checked or can be statically determined, then it is not rechecked, and no code is generated for the precondition check, as is shown in figure 12-5.

### 12.4.1.1 What Happens If Only Some Rewrites Match?

Interlacing the checking of rewrite preconditions and applying rewrites means that we can end up in a scenario where only some of the rewrites have been applied. In this case, we need to generate code to handle the failed match. To accomplish this, we take inspiration from tracing. In tracing, when there is a conditional branch, a check is inserted to check that the branch goes the same way as when the JITted code was originally generated. If the conditional test branches the other way, the JITted code falls back to the new code and only then will generate the code under that branch [31, 74, 75]. An example of this was shown in the related work chapter with a small procedural program in section §3.6.

Adapting this idea to $\mathbf{R}$-exprs, we can observe that when a precondition fails to match, we need to represent the state of the primitive R-expr (that is being "conceptually" rewritten) at this point. Furthermore, we want to avoid generating all possible rewrites of what can happen upfront (e.g. generating a sequence of 20 rewrites without knowing that it can be used at least once is unproductive). First, to represent the state, we can use the generate-a-new-R-expr mechanism from section $\S 12.3$. An example of a single rewrite that interlaces checks and rewriting is shown in figure 12-5.

To complete the analogy to tracing, we need to generate code for branches that were previously not taken. Recall from section $\S 12.4$, I mentioned that when a JIT-generated R-expr is Simplified for the first time, we will attempt to generate a new rewrite. This new rewrite corresponds to generating the code for a branch that was not previously encountered when running. The end result of generating new $\mathbf{R}$-expr kinds and generating rewrites for those $\mathbf{R}$-exprs is a "flow graph" that intermixes $\mathbf{R}$-expr states as nodes with rewrites as transitions. This is shown in figure 12-6.


Figure 12-6. An example graph of JIT states and rewrites which go between the states. Rewrites are shown in blue and always start at an $\mathbf{R}$-expr state. A rewrite can go to different $\mathbf{R}$-expr states depending on how many primitive rewrites match. If a precondition fails, then it takes one of the red edges and immediately jumps to a JIT-generated $\mathbf{R}$-expr kind. These failure edges always start from a rewrite and go directly to a state that represents the $\mathbf{R}$-expr at this point. The JIT-generated $\mathbf{R}$-expr states are reentrant (e.g. states 32 and 22) in the case the same state is encountered multiple times.

### 12.4.2 How to Reenter JITted R-exprs

The generated $\mathbf{R}$-expr kinds represent the state of the program when we are rewriting. Hence, reusing a JIT state is how we get reentrancy into compiled code (as in figure 12-6).

To implement reuse of JIT-generated R-expr kinds, whenever we are tasked with generating a new JIT-generated $\mathbf{R}$-expr kind, we first check through the previously generated $\mathbf{R}$-expr kinds and see if there is anything applicable.

To check for equivalent JIT states, there are a few things that we need to handle. Namely, there are parts of the R-expr that we do not care about. First, different variable names can be used in equivalent $\mathbf{R}$-exprs. Because we are tracking the variable names using holes when a variable is renamed, this just results in a different
variable name being placed into the hole. Second, some R-exprs can be rearranged. For example, conjunctions are associative and commutative, so $R * S$ and $S * R$ are equivalent.

As a more concrete example, the following two R-exprs are equivalent:

```
    lessthan(A,B)*lessthan(A,C)*lessthan(C,D)*lessthan(A,E)
lessthan(X4,X0)*lessthan(X2,X4)*lessthan(X2,X1)*lessthan(X2,X7)
```

Figure 12-7. Equivalent R-exprs with different variables and ordering ${ }^{202}$

Checking for equivalent states uses a more relaxed definition of equality compared to what is used by Simplify from chapters 8 and 11. The first step is to identify a set of potential JIT states by computing a JIT state-specific hash-code. The hash-code ignores the value types (variables names and constants values) in the $\mathbf{R}$-expr and the order that sub- $\mathbf{R}$-exprs are in associative and commutative $\mathbf{R}$-exprs. The hash-code also ignores $\mathbf{R}$-exprs kinds that will likely be holes in the JIT-generated state (section §12.5). The implementation of this is done by having a second hashCode method call jitHashCode on all R-expr kinds. Using the JIT's hash-code, the system will quickly identifies a small set of potential JIT-generated R-expr kinds to consider.

Once we have identified a set of possible JIT R-expr kinds into which an R-expr can be converted, we need to identify if there is a way that an existing generated $\mathbf{R}$-expr kind can be equivalent to the $\mathbf{R}$-expr state that the rewrite is attempting to serialize (during the compilation of a rewrite). This is implemented using a Prolog-style backtracking and unification, which searches through different ways in which the holes on the JIT-generated $\mathbf{R}$-expr kind can be filled in such that the JIT-generated state is equivalent to the $\mathbf{R}$-expr being converted to a JIT state.

### 12.4.3 Abstract Evaluation of Primitive Rewrites

So far, I have only talked abstractly about "finding primitive rewrites that can be applied", and "checking preconditions". While this description is the high-level idea

[^126]of the JIT compiler, this level of detail does leave a large glaring hole that needs to be addressed, in how the rewrites are actually generated.

When the JIT compiler is called from Simplify, it has access to the current JITgenerated $\mathbf{R}$-expr state, the corresponding primitive $\mathbf{R}$-expr that state represents, context $\mathcal{C}$, and the Clojure source code for all primitive rewrites from chapter $\S 6$ that were defined using the def-rewrite macro (section §11.5.3). To merge rewrites and speed up the rewrites themselves, we use partial evaluation to convert the Clojure source code that defines a rewrite into a format that can be manipulated inside of the JIT compiler.

Partial evaluation has been well studied in the JIT compilation literature [42, $57,73,116,149$ ]. Partial evaluation is a process in which we evaluate the program before we know all the information, which will only happen when it is running. For example, if we have the expression $1+2+x$, we can partially evaluate this expression into $3+x$. The information about 1 and 2 was sufficient to determine the result of + . Conceptually, this is very similar to what we have already been doing with $\mathbf{R}$-exprs, in that we allow $\mathbf{R}$-expr to be partially rewritten when there is insufficient information to completely rewrite the $\mathbf{R}$-expr. The difference this time is that we are partially evaluating Clojure code instead of $\mathbf{R}$-exprs.

To implement partial evaluation rewrites defined as Clojure code, we have a partial-evaluate function that takes as arguments a Clojure expression represented as an AST (which is a recursive list data structure as is typical of LISP-like languages) and the information that is globally accessible to the rewrite, such as the context $\mathcal{C}$. The Clojure source code that partial-evaluate is passed as an argument is from a single primitive rewrite defined using the def-rewrite macro. ${ }^{203}$ The partial-evaluate function returns a data structure that includes any statically knowable values, the current value conditioned on the current context $\mathcal{C}$, and statically known type information. The exact information returned varies by the value's type that the Clojure expression returns, as will be described shortly. Currently, in our JIT compiler, we only distinguish between a few types that are useful for compiling R-expr rewrites. These types are: R-exprs, an array of R-exprs, value types (variables or constants), an array of value types, "primitive" value (any ground value such as integers or strings contained in $\mathcal{G}$, but also any Java value,

[^127]some of which may not be contained in $\mathcal{G}$ ), and function types that are callable. The partial-evaluate function does not distinguish the type of a "primitive" value, such as floating point vs integers. This would typically be necessary when doing low-level code generation. However, this is not a problem for us, as we are not directly generating a low-level representation but instead generating Clojure, which is sufficiently high-level and does not require this level of detail.

When partial-evaluate evaluates an expression that returns either a primitive value type (like an int or string), or a R-expr value type (a constant or variable), the partial-evaluate function returns the currently known value, with respect to the context $\mathcal{C}$ that the rewrite is being generated for, and an optional statically known value. It also returns a fragment of Clojure code that can be embedded in the generated rewrite to access the value at runtime.

Functions such as 'get-value' and 'is-ground?' receive special handling in the JIT-compiler. These functions cache their returned value in a local variable in the generated rewrite. This means that if a rewrite performs (get-value X) twice on the same variable, the second time, the generated code will return the local variable instead of calling (get-value X) in the generated rewrite. This is useful, as when generating the code for multiple rewrites sequentially, there are often multiple calls to get-value for the exact same variable.

The way partial-evaluate handles R-expr typed variables and functions that return $\mathbf{R}$-exprs is a bit more interesting. R-expr typed variables are tracked statically and have no runtime representation. ${ }^{204}$ For example, in figure 12-5, as the rewrite progresses from the initial state on the left to the final state on the right, there are no $\mathbf{R}$-exprs objects serialized inside of the rewrite. This means that if we write (make-lessthan A B (make-constant true)) ${ }^{205}$ in the def-rewrite for a rewrite rule, the $\mathbf{R}$-expr is only tracked statically and there is no corresponding code in the generated rewrite that needs to be executed at runtime. To make this work, we must slightly expand the $\mathbf{R}$-expr representations with additional classes. Namely, we define new value types ${ }^{206}$ to allow general Clojure expressions and

[^128]variable names that are local to the rewrite in addition to the named variables, which are looked up in the context $\mathcal{C}$, and constants. These new value types are only allowed within a single compilation unit that will generate a single rewrite. Therefore, when the $\mathbf{R}$-expr state is serialized into a JIT-generated $\mathbf{R}$-expr type, these new value types are saved into a hole for future use. For example, in figure 12-4, the number 21 on state865( $\cdots$ ) would have been held in a local variable that was referenced by the $\mathbf{R}$-expr state. Therefore, the value 21 is saved in the JIT-generated $\mathbf{R}$-expr kind so that the remaining $\mathbf{R}$-exprs can still refer to the value 21.

This design of not representing $\mathbf{R}$-exprs in the compiled code was inspired by the Truffle project [149] ${ }^{207}$ which is a framework for making JIT compilers by defining an AST interpreter. The AST is represented using standard Java objects. When the AST is compiled, the AST does not have a representation in the generated code outside of the program counter, just like we are doing with the R-expr typed variables.

### 12.4.4 Structure of Generated Rewrites

The way in which the generated code is structured is heavily influenced by the fact that we are generating Clojure code, which is translated into Java bytecode. ${ }^{209}$ This means that we cannot generate JUMP instructions in the code. Instead, we use Java exceptions and try/catch blocks to emulate JUMPS. These operations can be converted into JUMP instructions when translated into machine code by the Java Virtual Machine (JVM). A conceptual example of this is shown in figure 12-8 with a rewrite actually generated by the JIT compiler shown in figure 12-9.

### 12.4.5 Generating Aggregators in the JIT-Generated Rewrites

So far, the generated operations that we have talked about have only been about generating straight-line code that corresponds to a sequence of rewrites performed against an R-expr. As discussed previously in sections § 6.5 and 11.6.3, aggregation requires combining the results of multiple contributions. This requires splitting the

[^129]```
try:
    if not снеск(\cdots): }\triangleright\mathrm{ Check precondition 1 for first rewrite
        throw CheckFailure \triangleright throw/catch can be converted to JUMPs by the Java VM
    if not снеск(\cdots): 
        throw CheckFailure
        \vdots \vdots \triangleright More preconditions checked (omitted)
                                    Do Rewrite once all preconditions are checked
    local_Y \leftarrowGETVALUE(Y) \triangleright Values assigned to variables are cached into local variables
    local_z % GETVALUE(Z)
    local_x \leftarrow local_y + local_z \triangleright Only the "work" of the rewrite is generated
try: }\triangleright\mathrm{ Start Second rewrite
                                    \triangleright ~ S e c o n d ~ r e w r i t e ' s ~ b o d y ~ o m i t t e d ~
    catch CheckFailure:
                                    \triangleright JUMP target for failed matches of second rewrite
        \triangleright ~ L o c a l ~ v a r i a b l e s ~ u s e d ~ b y ~ s t a t e 7 3 1 ~ a r e ~ p a s s e d ~ a s ~ a r g u m e n t s
    return state731(.\cdots, local_x, local_z, ...)
catch CheckFailure: \triangleright \UMP target for failed matches of first rewrite
return null }\triangleright\mathrm{ Return null to indicate that no rewrite were performed
```

Figure 12-8. High-level example of how generated code nests try/catch blocks to represent each rewrite. The catch blocks are the JUMP targets for when the match fails. Subsequent rewrites become more deeply nested under the try blocks.

```
(fn [rexpr simplify] ; Rewrite function takes R-expr & active simplify func
    (StatusCounters/match_attempt)
    (let [**context** (ContextHandle/get)
            **threadvar** (ThreadVar/get)]
    (try
        (do ; rewrites that modify the R-expr may not generate operations
        (try
            (try
                (do ; status counter incremented after 1 rewrite is done. performing a
                    ; rewrite does not have "code" sometimes (e.g. rearranging R-expr)
                    (StatusCounters/jit_rewrite_performed)
                    (try
                    (do ; perform a test, otherwise throw JITRuntimeCheckFailed
                    (jit-precondition-to-check (is-bound-in-context? (. rexpr jv23463)
        **context**))
                    (do
                    (jit-precondition-to-check
                    (let* []
                        (jit-precondition-to-check (is-variable? (. rexpr jv23461)))
                    (not (is-bound-in-context? (. rexpr jv23461) **context**))))
                    (let* [local-cache23852 (get-value-in-context (. rexpr jv23463) **
        context**)]
                    (do (set-value-in-context! (. rexpr jv23461) **context**
        local-cache23852)
                    (try
                        (try
                            (try
                                    ; the final returned R-expr
                            (make-jit-rexpr24202 (. rexpr jv23462) (make-constant
        local-cache23852) (. rexpr jv23464))
                            (catch JITRuntimeCheckFailed
                            ; Failure branch, return a different R-expr
                (simplify (make-jit-rexpr24046 (. rexpr jv23462) (
        make-constant local-cache23852) (. rexpr jv23464)))))
                    (catch UnificationFailure
            ; due to disjunctions, a unification failure might not result in 0
            ; so we catch this exception is handle explicitly instead of bubbling
                    (make-multiplicity 0)))
                    (catch JITRuntimeCheckFailed 
                        ; Second failure branch, returns another different R-expr
                            (simplify (make-jit-rexpr23857 (. rexpr jv23464) (
    make-constant local-cache23852) (. rexpr jv23462)))))))))
                (catch UnificationFailure _
                (make-multiplicity 0))))
            (catch JITRuntimeCheckFailed _
            (simplify (make-jit-rexpr23624 (. rexpr jv23462) (. rexpr jv23461) (.
        rexpr jv23463) (. rexpr jv23464)))))
            (catch UnificationFailure _ (make-multiplicity 0))))
    ; the highest level runtime check returns nil to indicate no change/rewrites
    (catch JITRuntimeCheckFailed _ nil))))
```

Figure 12-9. Example generated Rewrite from JIT compiler. ${ }^{208}$

R-expr into smaller disjunctive R-exprs, which can each be processed individually. This is done through the use of iterators, which are designed to loop over the domain of a variable (section §11.7). By having iterators split the R-expr using assignments to variables, rather than returning arbitrary $\mathbf{R}$-exprs, we can easily represent this as a new JIT-generated state (line 8 of figure 12-10). The current value of a variable can be stored with a local variable at runtime. Conversely, an arbitrary $\mathbf{R}$-expr cannot be efficiently represented at runtime given the design of the JIT compiler.

When getting iterators from an R-expr, we create a precondition on finding an iterator that behaves the way that we need when considering the code that we have generated in the rewrite (lines 3 to 6 of figure 12-10). In the event that we cannot find an iterator that can support our compiled subroutine, the rewrite will be aborted with a precondition failure, just like in section §12.4.4 (line 6 of figure 1210). To implement this behavior inside of the JIT-compiler, there is special code for handing rewrites on the efficient inner and outer aggregator kinds (section §11.6.3), and the primitive rewrites for aggregation are not at all used by the JIT-compiler.

To accumulate the aggregated values, we use a globally accessible container that holds the current aggregated value. This is the same as was done in section §11.6.3 figure 11-20.

A conceptual example of an aggregator in a JIT-generated rewrite is shown in figure 12-10.

### 12.5 What is JITable?

Most of the primitive R-expr kinds can be included in a JIT-generated $\mathbf{R}$-expr state, however there are some $\mathbf{R}$-exprs that we exclude.

From section §12.4, we can observe that the JIT rewrite compiler requires that the $\mathbf{R}$-expr after a rewrite has been applied is predictable. If a rewrite does not result in a predictable $\mathbf{R}$-expr, then the JIT compiler would be required to generate many different branches for the different states that could result from a particular rewrite. If there are too many JIT-generated states, the JIT compilation will become

[^130]```
global aggregationResult \(\leftarrow\) agg_null \(\triangleright\) Where the resulting value of aggregation is
stored
aggregationResiduals \(\leftarrow 0 \quad \triangleright \mathbf{R}\)-expr of still un-evaluatable \(\mathbf{R}\)-exprs
allIterators \(\leftarrow\) Getlterables \((\cdots)^{210} \quad \triangleright\) Get iterators from \(\mathbf{R}\)-expr, section §11.7
iterator \(\leftarrow\) StartIteratorWithVarOrder(allIterators, \([\mathrm{X}, \mathrm{Y}, \mathrm{Z}])\)
if iterator \(==\) null \(: \triangleright\) If there is not the right iterator, the precondition for the rewrite fails
    throw CheckFailure \(\triangleright\) Abort rewrite, as described in section §12.4.4
for \(\langle x\), continuation \(\rangle \in\) iterator.Run(): \(\triangleright\) Run iterator
    \(\mathrm{R} \leftarrow\) state 941 ( \(x\), continuation, \(\cdots\) ) \(\triangleright\) JIT-generated \(\mathbf{R}\)-expr representing agg body
    \(\mathrm{R}^{\prime} \leftarrow \operatorname{SIMPLIFY}(\mathrm{R}, \mathcal{C}) \triangleright\) Evaluate the aggregator's body in its own \(\mathbf{R}\)-expr rewrite "unit"
    if \(R^{\prime} \neq 0\) : \(\quad \triangleright\) Check if there is an \(\mathbf{R}\)-expr residual
        aggregationResiduals \(\leftarrow\) aggregationResiduals \(+R^{\prime}\)
if aggregationResiduals \(==0\) : \(\quad \triangleright\) if no \(\mathbf{R}\)-expr residual
    \(\triangleright\) The aggregation is done, and the result is in aggregationResult
    if aggregationResult == agg_null :
        "return" 0
    else
        "return" (A=aggregationResult)
else
    \(\triangleright\) Return the aggregator with the delayed residual \(\mathbf{R}\)-expr
    "return" (A=sum(X, (X=aggregationResult)+aggregationResiduals))
```

Figure 12-10. Conceptual example of aggregation performed by a JIT-generated rewrite on the outer aggregator. The accumulator variable is initialized on line 1 and will be modified inside of Simplify, called on line 9. Like with efficient aggregators in section §11.6.3, the aggregation is done when the residual R-expr is rewritten as 0 (line 12).
inefficient, as it is difficult to benefit from reuse of existing JIT-generated states. Therefore, we exclude $\mathbf{R}$-exprs that are frequently rewritten in difficult-to-predict ways. To handle these R-expr kinds, they are represented using holes in the JITgenerated $\mathbf{R}$-expr; therefore, they are passed around and simplified as opaque R-expr pointers and not otherwise handled by the JIT.

The R-exprs kinds which we do not support are as follows:

1. An obvious case is the $\mathbf{R}$-expr that converts from the Dyna AST into $\mathbf{R}$-exprs (section §11.3). This R-expr can result in any possible R-expr depending on the shape of the program. Furthermore, these R-exprs are only used when translating the Dyna source. This rarely happens in the running of the program. Hence, we are unlikely to benefit from making it run faster.
2. Another case of unpredictable $\mathbf{R}$-exprs are disjunctions. For example, suppose that we have $R_{1}+R_{2}+R_{3}+\cdots+R_{n}$, if we consider that each $R_{i}$ could be independently rewritten as 0 , then there are potentially $2^{n}$ possible states in which this $\mathbf{R}$-expr could be rewritten. This means that we could potentially generate these $2^{n}$ states and the corresponding rewrites between them. To limit this, I only allow small ${ }^{211}$ disjunctions to be generated in the JIT. The idea here is that a small disjunction likely represents a function that would benefit from compilation. For example, the rectified linear function $f(x)=\max \{0, x\}$ would be represented as the $\mathbf{R}-\operatorname{expr}(A=\max (Y,(Y=0)+(Y=X))$ ) which includes a disjunction. In contrast, a large disjunction might represent a table of values or a memo. Hence, retrieving data from the hash table trie (section §11.6.1) would be more efficient than a chain of if-expressions.
3. Memo read $\mathbf{R}$-exprs are also excluded for the same reason as disjunctions.
4. We also do not include user-defined R-expr kinds in the JIT-generated R-expr kinds. The reason is that user-defined $\mathbf{R}$-exprs can be redefined at the REPL or by guessing with memoization. This means that any JIT state that included a user-defined $\mathbf{R}$-expr (and rewrites on those states) would have to be invalidated. Furthermore, user-defined $\mathbf{R}$-exprs often indirect through memo tables and
${ }^{211}$ Small disjunctions are currently set as a disjunction with fewer than 8 children. The limit of 8 was chosen without any data to inform this choice. In the future, one may wish to reconsider this limit and experimental with different configurations of the JIT compiler.
include a disjunction inside of their definition. Hence, a generated R-expr that includes a user-defined $\mathbf{R}$-expr would not be able to represent that much in the first place.

While limiting disjunctions, memos, and user-defined $\mathbf{R}$-exprs seems like a big limitation, we can still work around it. One way to work around this is by merging a JIT state with the $\mathbf{R}$-expr, which is held in as an opaque pointer inside of the JIT hole. This is accomplished by allowing JIT to create a rewrite that matches the $\mathbf{R}$-expr contained in the hole and add it to the current JIT state. This means that once the $\mathbf{R}$-expr contained in a hole has been sufficiently rewritten, it can be merged to create a larger JIT-generated $\mathbf{R}$-expr state.

### 12.6 Starting the JIT Compiler

So far, we have that we can take an R-expr and represent it as a JIT-generated $\mathbf{R}$-expr state and then create rewrites on those states. However, this leaves the question of how to create the initial seed JIT-generated states.

The way that I do this is by starting from 1) user-defined $\mathbf{R}$-exprs (as in figure 123) and 2) R-exprs that are nested under disjunctions and turning these into JITgenerated states. The reason to use user-defined $\mathbf{R}$-exprs is that these are likely to reappear a lot and be used in the same way. Hence, using a user-defined R-expr as a seed is akin to compiling a user-defined method. The second case for an $\mathbf{R}$-expr under a disjunction is that the disjunction itself does not get compiled. However, the leaves of the disjunction might still contain code.

This choice of seed is, of course, a heuristic and something that future work may wish to revisit.

### 12.7 Experiments: Benchmarks

Our ability to run real-world programs is still limited by the runtime performance of the implementation on the whole. In chapter $\S 16$, I discuss possible directions to further improve the runtime of Dyna. So, unfortunately, at this time, I do not have benchmarks on "real-world" problems. However, we can still benchmark a small

```
671 |f(X) += I for range(0, X, I). 212
(a) Dyna
\({ }^{672}\) def \(f(X)\) :
673
return \(\operatorname{sum}(\) range \((X))\)
```

(b) Python

| ${ }^{674}$ | $($ defn $f[X]$ |
| :--- | :--- |
| $($ reduce $+($ range $X)))$ |  |

(c) Clojure


Figure 12-12. The $y$-axis is the ratio between the JIT compiled version and the non-JIT compiled version (chapter $\S 11$ ). We can see that the JIT-compiled version runs roughly about five as fast as the non-JIT-compiled version (yay!). ${ }^{213}$
of the non-JIT-compiled version. The first run being slower is consistent with what our expectations as the first run is performs extra work generating the rewrites.

To try and understand why the JIT-compiled version is faster, we turn our attention to figure 12-13. Figure 12-13 shows the number of times that Simplify attempted to match a rewrite vs. the number of times a rewrite was applied. If you recall from section $\S 8$.A, the attempted matches come from the design of Simplify where it will dispatch against the $\mathbf{R}$-expr kind, and then sequentially check if any of the rewrites can be applied by checking preconditions. For the non JIT-compiled version, we first observe that SIMPLIFY attempts 60 matches per number summed. Without JIT-compilation, we observe that only $\frac{1}{3}$ of the rewrites match. This means

(a) This shows the number of rewrites attempted and the number of rewrites that were successfully matched and applied. Here, there are 60 rewrites attempted per number summed, with 130 rewrites as a constant overhead. There are 20 rewrites that are successfully applied per summed number and 50 that are a constant overhead.

(b) This shows the number of rewrites attempted and successfully applied when the JIT is used to generate $\mathbf{R}$-expr kinds and rewrites. Here we can observe that all rewrites attempted are successfully matched. Furthermore, there are only 10 additional rewrites per number summed (vs. 20 without the JIT), and there is a baseline of 40 rewrites of constant overhead. This shows that the JIT performs the same amount of work using fewer rewrites.

Figure 12-13. Creating JIT-generated $\mathbf{R}$-expr kinds significantly reduces the number of attempted rewrite matches, and the number of rewrites required to do productive work, in terms of rewrites matched.
$\frac{2}{3}$ of the time is wasted on checking if "useless" rewrites could apply. Conversely, after generating new $\mathbf{R}$-expr states and making JIT-generated rewrites, we see that $100 \%$ of the attempted matches match and run rewrites successfully. The JIT only generates rewrites for the scenarios that we have been previously encountered. Because this program is repeating the same operation multiple times (summing an integer), there are no unnecessary rewrites loaded into the system, and therefore no wasted effort in searching through matching rewrites. This is similar to the kinds of numerical programs that are typical of machine learning algorithms

In addition to observing that the JIT does help the runtime, we also compare it against other programming languages. In figure 12-14, I show the runtime per number summed for Dyna with and without JIT-compilation, Python, Clojure, and Java. We can observe that the JIT-compiled $\mathbf{R}$-exprs rewrites are 120 times slower than Python (6010 times slower than Java). Further delving into this in figure 12-15, we can profile where the JIT-compiled version is spending its time. It appears that much of the time spent is on non-R-expr rewriting activities, such as initializing hash-tables used for the context's internal data structures, computing hash-codes for $\mathbf{R}$-exprs, and accessing globally accessible objects such as the context. The good news from figure $12-15$ is that there is a lot of opportunity to make the system run faster through careful engineering to avoid these unnecessary overheads. The bad news from figure $12-14$ is that Dyna is 120 times slower than Python and likely still too slow for "real-world" problems.


Figure 12-14. Results of running the small benchmark programs from figure 12-11. The $y$-axis is the total time of running the benchmark 10 times and dividing it by the value of $x$, which is the number that is being summed to. Hence, this plot shows the scaling of time per summed integer. The non-JIT R-expr rewriting (chapter §11) is 31168 times slower than the fastest runtime, which is the Java program using a for-loop. The JIT compiled $\mathbf{R}$-exprs (this chapter) is 6010 times slower than Java.


Figure 12-15. This is a flame graph for the benchmark from figure 12-11a running with JIT-generated rewrites. We can observe that $85.19 \%$ of the runtime is taken up by a JIT-generated rewrite that represents the outer part of the aggregation (conceptually what is in figure 12-10). This rewrite internally calls simplify and performs other rewrites. We can observe that $23 \%$ of the overall time ( $27 \%$ of the rewrite) is spent rewriting the JIT-generated R-expr that represents the body of aggregation (line 9 in figure 12-10). The creation and manipulation of nested contexts account for $22 \%$ of the overall runtime ( $26 \%$ of the rewrite). $15 \%$ of the overall runtime is spent creating $\mathbf{R}$-exprs, more specially the profile seems to indicate that creating cached hash codes is somehow the expensive part of this. Approximately $10 \%$ of the overall runtime is spent reading and writing from global variables. This is, of course, subject to the profiler's ability to measure these typically fast functions, but accessing the global context does happen frequently, so it is conceivable that this is a correct measurement.

## Chapter 13

## Object Oriented Programming in a Pure Declarative Language

### 13.1 Dynabases

Object oriented design in Dyna with dynabases was originally proposed by Eisner and Filardo in [59] with an expanded discussion in Filardo's dissertation [66]. Unfortunately, the prior work did not include an implementation or complete details. This dissertation contributes for the first time working details of the syntax for dynabases (section §2.9) and a working implementation of dynabases. The design of dynabases in this dissertation has some differences from the one proposed in [59, 66], though these differences will be immaterial to most users.

Dynabases provide prototype-based inheritance [49, 59, 133, 136] for logic programming. A dynabase object can be extended by adding rules to a dynabase after it has been "created". Dynabases are open, in that additional rules can be added to the dynabase's "class definition" at any point, even after "instances" of the class have been instantiated. Allowing rules to be added to a dynabase is consistent with Dyna's out-of-order execution, where the program is a collection of rules, and the system is looking for a consistent assignment to all terms defined by the program, regardless of the order in which they were defined.

To implement this behavior, dynabases are not allocated objects, as is typical in a procedural programming language. Rather a dynabase is represented using an immutable dynabase term identifier, as will be described in section §13.2. The dynabase identifier can be used to call terms defined on the dynabase, much the same way that a structured-term (section §2.1.1) can be passed as an argument to a term. As such, we will see in section $\S 13.3$ that dynabases can be implemented as almost a purely syntactic transformation before the program is converted to $\mathbf{R}$-exprs (chapter §7). This is conceptually similar to what appeared in previous publications about Dyna's object-oriented design [59, 66] as well for prototype-based inheritance in Prolog with the LogTalk project [49].

### 13.2 Dynabase Object Representation

Dynabases allow closures capturing all variables present when they are instantiated. Dynabases also support single inheritance. Conceptually, we can think of a dynabase's identifier as an immutable list of dynabases types that the dynabase instance inherited from. If we are to use the Dyna notation from section §2.1.1, then a dynabase object would look something like:
${ }_{679}$ |[dynabase_123[CapturedVar1, CapturedVar2], dynabase_789[X, Y]]
Here, the numerical identifiers 123 and 789 on line 679 identify the dynabase class types that this object inherits from. The variables CapturedVar1, CapturedVar2, X, $Y$ are the captured variables that were present when the dynabase was instantiated. The presentation on line 679 is conceptually similar to what was presented previously in Eisner and Filardo [59].

While the array of identifiers on line 679 is conceptually correct, we instead going to use a slightly different design in the implementation for efficiency reasons. Instead, we are going to group the dynabases by dynabase class identifier into a map and then have a list of lists to track the captured variables. As a Java type, this looks like line 680:

680 | Map<DynabaseTypeID, List<List<G>>>
Here, DynabaseTypeID is an opaque class identifier. The identifier is generated to be unique when the Dyna program is converted from the Dyna AST into $\mathbf{R}$-exprs. The DynabaseTypeID identifier corresponds with the dynabase's "class" rather than
an instance of the dynabase. For the presentation in this chapter, I will use the line number on which a dynabase is defined on as the identifier.

The value associated with each dynabase class identifier is List<List<G$\gg$. The inner List< $\mathcal{G}>$ contains the values of the captured variables when the dynabase is created. The order in which the values are held in List $\langle\mathcal{G}\rangle$ is chosen arbitrarily, but the generated code will consistently reference the variables in the slots.

The second list of lists is to support the scenario where a dynabase type can appear more than once. This happens when self-inheritance is used. This will be discussed in more detail in section §13.5.

### 13.2.1 Why Capture All Variables?

Dynabases capture all local variables when created. The reason why it needs to capture all of the variables in the current rule is that a dynabase might be modified or referenced from places other than where it was constructed. For example, consider the dynabase on line 681.

```
f(X) = new {}.
db_a = f(1).
db_b = f(2).
db_a.a = 123.
db_b.a = 456.
assert db_a.a == 123.
assert db_b.a == 456.
assert f(1).a == 123.
```

Here the functor $f(X)$ has one variable $X$ and returns an "empty" dynabase. On lines 683 and 684, we construct two different dynabases. We then proceed to modify the dynabases by setting the "field" 'a' on lines 686 and 687. The values we have assigned in each of these dynabases are distinct from each other (lines 689 and 690). We, therefore, must have some way to track this distinction between dynabases.

If this were a procedural language, then we would use the fact that we are creating two different instances of an object with two different locations in memory as a result of the two different function calls on lines 683 and 684. However, by
the nature of being declarative and functional, we require that every time that we call a function with the same arguments (Dyna term), we get back the same result. Hence, we can get db_a again by calling f(1) a second time as on line 691.

Therefore, we would either need a design that is required to store any dynabase that is returned from a user-defined function or "recreate" the same dynabase multiple times. We have opted to recreate the same dynabase object multiple times. This is done by returning an immutable identifier from $f(1)$ rather than a mutable object. Furthermore, $f(\cdot)$ defined on line 681 must return different dynabase identifiers when $f(1)$ is called on line 683, and when $f(2)$ is called on line 684. To make this work, there needs to be a distinguishing feature between the call to $f(1)$ and $f(2)$. The only distinguishing feature is the assignment to variables present in the rule. (In this case, the variable x.) Hence, we capture the value of all variables when creating a dynabase.
Consequence of this design: An consequence of this design is that we cannot construct two (or more) dynabases at the same time in a single rule.

For example, the following program does not work:

```
692 does_not_work(X) = (A = new {}), (B = new {}), A.
```

The reason for this is that both the $A$ dynabase and the $B$ dynabase need to capture a ground value reference of the other (we do not allow for cycles in the ground values). In my opinion, this is not an issue, as it can easily be worked around by ensuring that there is at most one dynabase constructed per functor:

```
693 make_B(X) = new \(\}\).
694 works \((X)=(A=\) new \(\{ \}),(B=\) make_ \(B(X)), A\).
```

If a user writes a rule like line 692, this is detected and reported as an error to the user.

Future work may consider adding an automatic transformation to the AST-to-R-expr front-end to perform a transformation like line 693 to allow multiple dynabases to be defined in a single Dyna rule.

### 13.3 Desugared Dynabases

In converting a Dyna program with dynabases into an R-expr, there are two cases that we have to handle: constructing a dynabase and calling methods (ac-
cessing fields) on a dynabase. Both of these cases are handled by introducing new $\mathbf{R}$-expr types, which are called dynabaseCreate dynabaseTypeid $^{\text {(SuperDynabase, }}$ Var1, $\cdots, ~ V a r N, ~ R e s u l t) ~ a n d ~ d y n a b a s e A c c e s s d y n a b a s e T y p e I D(D y n a b a s e, ~ V a r 1, ~ \cdots, ~$ VarN).

The dynabaseCreate $\mathbf{R}$-expr is annotated with an opaque symbol dynabaseTypeID, ${ }^{214}$ which is picked to be unique when the program is translated from the Dyna source code. The variable SuperDynabase is the dynabase inherited from. The variables Var1, ..., VarN are the captured local variables (section §13.2.1).

To illustrate the translation process, consider the following Dyna program on lines 695 to 704 with dynabases. To make the translation easier to read, I will show the translation using "pseudo Dyna", which is an intermediate step in translating from Dyna into $\mathbf{R}$-exprs.

```
695 e = new {
    val = 123.
}.
f(X,Y) = new {
    val += X.
    val += Y.
    func(X) = $self.val * X.
}.
g = f(1,2).val.
```

With dynabases desugared:

| 705 | $e=$ dynabaseCreate_695(\$nil, Result), Result. | \% From line 695 |
| :--- | :--- | :--- | :--- |
| 706 | val(Dynabase) = dynabaseAccess_695(Dynabase), 123. | \% From line 696 |
| 707 | $f(X, Y)=$ dynabaseCreate_698(\$nil, X, Y, Result), Result. \% From line 698 |  |
| 708 | val(Dynabase) += dynabaseAccess_698(Dynabase, X, Y), X. \% From line 699 |  |
| 709 | val(Dynabase) += dynabaseAccess_698(Dynabase, X, Y), Y. \% From line 700 |  |
| 710 | func(DBase, X) = dynabaseAccess_698(DBase,_, Y), val(DBase)*X. \%From L701 |  |
| 711 |  |  |
| 712 | $g=\operatorname{val}(f(1,2))$. | \% From line 704 |

Note that the rule val appears on both dynabase_695 and dynabase_698. The distinguishing factor between the val rules are the dynabaseAccess $\mathbf{R}$-exprs. The

[^131]reason for this is that dynabases inherit from each other, and as such, there are rules that are defined in different dynabases that need to be combined. This means that a "compile-time" function dispatch only considers the functor name (val in this example). A secondary "runtime" dispatch selects the relevant rules that define the val terms from lines 706, 708 and 709. This is done by using the value assigned to the Dynabase variable to rewrite the dynabaseAccess R-expr.

The approach of grouping by functor name and using dynabaseAccess_698 is necessary to allow modifications to the dynabase after it has been initially defined. For example, consider line 715 where we define a new value $v$ on a dynabase.

```
713 db (X) = new {a=1.} for X > 0.
7 1 4 d b ( X ) = ~ n e w ~ \{ b = 2 . \} ~ f o r ~ X ~ < = ~ 0 . ~
715 db(X).v = 2 + X.
```

On line 715 we do not know the dynabase type we are modifying. In fact, we are actually defining ' $v$ ' on two different dynabase types at the same time. When we desugar line 715, we get something like line 716 below.

```
716 v(Dynabase) = (Dynabase = db (X)), 2 + X. % From line 715
```

Here, we have that the dynabase is represented using the variable Dynabase. We only need to find a value of $X$ such that the expression Dynabase $=d b(X)$, is true. This can be handled using rewrites on dynabaseCreate.

### 13.4 Dynabase Rewrite Rules

The dynabase rewrite rules are fairly straightforward. The inheritance relations between dynabase types are tracked via a global data structure and updated whenever a new dynabase instance is created by the dynabaseCreate rewrite rules. Some of the rewrite rules for dynabases are conditionally enabled and disabled depending on the inheritance between different dynabase types.

```
dynabaseCreate_698(SuperDB, X, Y, Result) \xrightarrow{}{88}}\mathrm{ (Result=dynabase object)
                                if SuperDB, X, Y \in\mathcal{G}
dynabaseCreate_698(SuperDB, X, Y, Result) \xrightarrow{}{89}\mathrm{ (SuperDB=dbase}
    object)*(X=x)*(Y=y)
                                    if Result }\in\mathcal{G}\mathrm{ and dynabase_698 does not self-inherit
dynabaseAccess_698(Dynabase, X, Y) \xrightarrow{}{90}((X=\mp@subsup{x}{1}{})*(Y=\mp@subsup{y}{1}{})+\quad\triangleright\mathrm{ allow self-inherit}
```

```
        (X=\mp@subsup{x}{2}{})*(Y=\mp@subsup{y}{2}{})+
        (X=\mp@subsup{x}{3}{})*(Y=\mp@subsup{y}{3}{})+\cdots) if Dynabase }\in\mathcal{G
dynabaseAccess_698(Dynabase, X, Y)*dynabaseAccess_695(Dynabase) 笙 0 if (
            dynabase_698 and dynabase_695 are incompatible (w.r.t. inheritance))
(dynabaseCreate_698(SuperDynabase, X, Y, Dynabase) *
dynabaseCreate_695(SuperDynabase2, Dynabase) ) \xrightarrow{}{92}0}\mathrm{ if (
    dynabase_698 and dynabase_695 are incompatible (w.r.t. inheritance))
(dynabaseCreate_698(SuperDynabase, X, Y, Dynabase)*
dynabaseAccess_695(Dynabase)) \xrightarrow{}{93}0\mathrm{ if (}
    dynabase_698 and dynabase_695 are incompatible (w.r.t. inheritance))
```

Rewrite rule 88 creates a dynabase object in the case that all of the arguments are ground. The SuperDB variable is the dynabase object from which the created dynabase inherits. In the case where there is no parent dynabase, then this variable is set to the value $\$ n i 1$.

Like with structured-terms, the dynabaseCreate $\mathbf{R}$-expr both creates dynabase objects and destructure dynabase objects, getting access to the variables captured by the dynabase closure. This is done using rewrite rule 89 , which requires that the Result variable is assigned dynabase that inherits from dynabase_698. Rewrite rule 89 is used when evaluating Dynabase $=\mathrm{db}(\mathrm{X})$ from line 715 . Further note that rewrite rule 89 is disabled in the case that a dynabase type self-inherits. I will discuss the reason further in section §13.5.

Rewrite rule 90 is used to match a dynabase object for rules that were defined inside of the dynabase's definition. This was used by lines 706 and 708 to 710 . Observe that rewrite rule 90 rewrites as a disjunction of bindings to the variables $X$ and $Y$. This is a disjunction over all of the times that a dynabase type has been inherited from in constructing a dynabase object. A single dynabase type appearing more than once in the case of self-inheritance, which will be discussed in section §13.5. Note, in the case that a dynabase type does not self-inherit, rewrite rule 90 will only rewrite a single assignment to the variables $X$ and $Y$ for a given assignment to the Dynabase variable.

Rewrite rules 91 to 93 implement "type checks" that cause an R-expr to be rewritten as 0 , thus eliminating dead code. Rewrite rules 91 to 93 do not require that the value of the Dynabase is known for these rewrites to apply. However, these rewrites depend on the information about how dynabases inherit from each other.

This inheritance information can be updated after the system has started running. As such, applications of rewrite rules 91 to 93 depend on the state of dynabase inheritance, and usage of these rewrite rules is tracked using the assumption mechanism previously discussed in section §10.4.1.

### 13.5 Self-Inheritance

As hinted at in the previous section, self-inheritance with dynabases allows for some strange interaction and causes rewrite rule 89 to become disabled in the case there is self-inheritance. So what is self-inheritance, why is it interesting, and why does it cause things to break? To answer this, let us work through a small example of self-inheritance to see what it is, how it works, and how we can build towards a logical inconsistency that prevents us from using rewrite rule 89 in the case of self-inheritance.

We define dynabase_717 on line 717, which inherits from the argument $x .{ }^{215}$ By passing in a dynabase that has already inherited from dynabase_717, we can self-inherit, as is done on lines 722 and 723.

```
f(X) = new (X) {
    z += 1.
}.
a = new {}.
af = f(a).
aff = f(af).
afff = f(aff).
assert af.z == 1
assert aff.z == 2.
assert afff.z == 3.
```

Observe that we have defined ' $z$ ' on line 718 to essentially count the number of times that dynabase_717 has been inherited from. The way this works is that ' $z$ ' is desugared to use the dynabaseAccess_717 R-expr:
${ }_{727}$ |z(Dynabase) += dynabaseAccess_717(Dynabase, X), 1. \% From line 718

[^132]And then the dynabaseAccess_717 R-expr will rewrite using rewrite rule 90 into a disjunction of the different values of $x$ that contributed to creating the 'afff' dynabase object:


Now, 'afff. z' returning the value 3 is exactly what we want to happen using the dynabaseAccess, which is used for rules defined inside of the dynabase's initial definition. Now, let us consider a scenario where dynabaseCreate, which is disabled, is used instead.

Before we do work through the dynabaseCreate scenario, let us build up to this case with a by defining 'w' as a standard, non-dynabase, Dyna term:

```
w(D) += 1.
assert w(a) == 1.
assert w(af) == 1.
assert w(aff) == 1.
assert w(afff) == 1.
```

There are no tricks here in defining ' $w$ ' in that we sum the value 1 for every unique term passed as an argument to ' $w$ '. Now, let us define w2 with a constraint on the variable ' D '.

```
w2(D) += (D = f(X)), 1.
assert is_null(w2(a)).
assert w2(af) == 1.
assert w2(aff) == 1.
assert w2(afff) == 1.
```

Here ' $(D=f(X)$ )' requires that there exists some value for $X$, such that the value held in the variable $D$ is returned. As already discussed in section $\S 13.3$, line 733 is equivalent to line 738.
${ }_{738} \mid f(X) \cdot w 2+=1$.
Now we can begin to see the problem. The interpretation of line 738 should be that the value of ' $w 2$ ' defined on the dynabase returned from the function $f(X)$ should be equivalent to the value of ' $z$ '. Hence, it should be counted multiple times. However, this is different from the interpretation of 'w2' as defined on line 733.

Rather than trying to reconcile the two ways in which 'w2' could be interpreted, we instead disable rewrite rule 89 . This means that if a user writes a rule like line 733
or line 738 they get back unrewritten R-expr instead of the value 1 or 3 . Admittedly, an unrewritten R-expr is rarely the desired result, but at least this does not give an incorrect answer. Furthermore, this issue can be avoided by defining rules inside of the dynabases definition, as with the definition of ' $z$ ' on line 718.

Finally, I want to reiterate that without self-inheritance, dynabases are conceptually very simple. The reason for the additional complexity of dynabases largely has to deal with self-inheritance, and for rewrites on dynabases to work as "compile time" rewrites when possible.

### 13.6 Comparison with 2011 Proposal for Dynabases

The design presented in this chapter is my own design for dynabases. However, it is not the only possible design for dynabases. When dealing with simple dynabases that do not use self-inheritance or "external modifications" of the dynabase (such as line 715), then all proposed implementations of dynabases exhibit similar behavior. However, how self-inheritance (section §13.5) and external modification (e.g. line 715) are handled does differ between the proposed implementations.

In Eisner and Filardo [59], the authors proposed "collect[ing] the relevant rules by recursively traversing [the dynabase's] parent [reference ${ }^{216}$ ]". This is, of course, the right high-level idea, but as we saw in section $\S 13.5$, there are many details that are needed to make this operational. Furthermore, I note that Eisner and Filardo [59] did not distinguish between dynabaseCreate and dynabaseAccess R-exprs as was done here. Instead, they only use Dyna-to-Dyna translations like (Dynabase=db(X)) (e.g. line 716), which depends on the interpretation of their "dynabaseCreate" operation. The issue of self-inheritance (the disjunction in rewrite rule 89) was not addressed in [59].

Filardo did expand on the [59] design in his dissertation [66] section 6.2. In section 6.2.2 "Rule-Collection Semantics With Recursive Owner Writes", Filardo also recognized the issue of self-inheritance. Filardo's section 6.2.2 attempted to fix the issue of self-inheritance by introducing a concept he called "focused dynabase" into the "collect-relevant-rules" function. Unfortunately, Filardo did not work out

[^133]the details of how the "focused dynabase" would integrate into the execution of the Dyna program that had dynabases, and how it works with calls to other terms defined on a dynabase. Although the design from [59] and [66] does not have an implementation and the description in [59, 66] is partially incomplete, I believe that regardless of how the design from $[59,66]$ is completed, it would run into the same kinds of issues that we have in the case of self-inheritance (section §13.5).

To illustrate some of the complications with dynabases, and how the way this chapter handles dynabases vs [59, 66], let us work through the following example program on lines 739 to 753 .

```
f(X) = new (X) {
    q += 1.
    r += $self.q.
    s += $self.u.
    t += $self.ss.s.
}.
a = new {
    ss = $self.
}.
af = f(a).
aff = f(af).
afff = f(aff).
afffb = new afff {
    u += 1.
}.
```

Here, I am defining dynabase_739 as a dynabase that self-inherits. Observe that the ' $q$ ' term on line 740 is the same as the ' $z$ ' from section $\S 13.5$ in that it counts the number of times that dynabase_739 is inherited from. However, this time, we additionally have the ' $r$ ' term that sums the value of ' $q$ ' every time it is added. The question we need to answer to understand how the system works is how \$self on line 741 is defined as rules are added by recursively traversing the parent references.

Before delving into the discussion about [59, 66], I will note that for this example running under the design presented in this chapter, we have afffb.r == 9 and afffb.t == 9, as \$self always reference the dynabase that is being called (in this case the value afffb). Hence, \$self.q on line 741 returns the value 3. Furthermore, the call to \$self. $u$ on line 742 will refer to the ' $u$ ' defined on line 752 , even though ' $u$ ' is defined in a descendant dynabase. This approach of making \$self reference the
dynabase object that and calling functions that have been overridden is consistent with the behavior that is typically used when doing object-oriented programming.

Now, we turn our attention to the design from [59, 66]. When evaluating afffb.r, I assume that the "focused dynabase" was intended to be afffb, which would have allowed the collect-relevant-rules function to identify that ' $r$ ' from line 741 needs to have 3 contributions. Next, we need to consider how \$self references the relevant dynabase.

A first possible interpretation of \$self would be to define \$self as only referencing the current dynabase (this would have allowed for values computed in parent dynabases to be cached using memoization and modified by the children), then this would mean that \$self is not able to reference functions on children dynabases. This is problematic for compositionality. This interpretation of \$self would mean that afffb.t is undefined as the \$self on line 742 would be unable to reference the ' $u$ ' defined on line 752. I note this design as [59] proposed caching the computation performed on parent dynabases. Filardo [66] rejects this idea and instead makes $\$$ self a dynamic variable that includes references to the rules defined on children dynabases.

Going to the second interpretation of \$self as a dynamic variable that references the "focused dynabase", this still leaves the question of how \$self and the ( $\$$ self $=\mathrm{db}(X)$ ) constraint from the Dyna-to-Dyna translation interact. Unfortunately, these details were glossed over. Presumably, had these details been added, it would have resulted in something like the dynabaseAccess $\mathbf{R}$-expr that can handle self-inheritance and match in the case where it is a parent of the current dynabase.

Finally, I will note that [59, 66] proposed an access control mechanism that was introduced through an additional hidden variable that tracked the owner dynabase. I have chosen to do away with the owner concept for dynabases. Instead, anyone who has a pointer to a dynabase can modify it. This simplifies the design of dynabases and the amount of information one has to learn to use Dyna.

## Chapter 14

## Folding and Speculative Rewrites for Recursive Programs

When rewriting calls to user-defined R-expr (section §5.2.2.11), our general approach is to expand the user definition (rewrite rule 74). This is sufficient for non-recursive programs, as, without recursion, we can expand the $\mathbf{R}$-expr to the maximum depth and have a finite-sized $\mathbf{R}$-expr where rewrites can be applied arbitrarily. Unfortunately, when there is recursion, we cannot expand the R-expr all of the way. ${ }^{217}$

One way we can handle recursive programs while still being able to leverage the power of $\mathbf{R}$-exprs is by using fold/unfold transformations. [22, 25, 58, 80, 97, 141] This essentially creates new automatically defined functions which are specialized to a specific case. This can allow us to make inferences for recursive programs, which may be useful in some cases.

The unfold transform is analogous to rewrite rule 74 that we have already seen. A user-defined $\mathbf{R}$-expr is expanded to its $\mathbf{R}$-expr definition, with variables renamed as needed.

The fold transform is the reverse of unfolding an R-expr. Folding generates a

[^134]new, automatically named R-expr. This can be written as rewrite rule 94, where ' $f$ ' is a newly chosen name.
$$
\mathrm{R}_{\mathrm{f}} \xrightarrow{94} \mathrm{f}\left(\mathrm{X}_{1}, \mathrm{x}_{2}, \mathrm{X}_{3}, \ldots, \mathrm{X}_{n}\right) \quad \text { where }\left\{\mathrm{X}_{1}, \mathrm{x}_{2}, \mathrm{X}_{3}, \ldots, \mathrm{X}_{n}\right\}=\operatorname{vars}(\mathrm{R})
$$

### 14.1 Why Fold a Program

The reason why one might want to fold a program is that, in the case of a recursive program, a folded definition can depend on itself. This allows us to perform rewrites that optimize recursive functions before it is unfolded in the context of the input to the Dyna program is processed. For example, if we have a higher-order function that performs an indirect function, we can create a specialized version of the function that does not perform an indirect call and can enable other rewrites the opportunity to optimize the R-expr. For example consider the map function below in figure 14-1 that applies a function to a list:
map(F, []) = [].
map(F, [Head|Tail]) =
[F(Head)|
map(F,Tail)].
(a) Dyna

```
```

```
    map(F, L, Res) }->\mathrm{ (Res=only(Inp,
```

```
    map(F, L, Res) }->\mathrm{ (Res=only(Inp,
        (L=[])*(Inp=[])+
        (L=[])*(Inp=[])+
```

    proj(Head,proj(Tail, proj(Tmp1, proj(Tmp2,
    ```
    proj(Head,proj(Tail, proj(Tmp1, proj(Tmp2,
        (L=[Head|Tail])*
        (L=[Head|Tail])*
        indirect_call(F, Head, Tmp1)*
        indirect_call(F, Head, Tmp1)*
        map(F, Tail, Tmp2)*
        map(F, Tail, Tmp2)*
    (Inp=[Tmp1|Tmp2])))))))
```

    (Inp=[Tmp1|Tmp2])))))))
    ```
(b) R-expr

Figure 14-1. A recursive function that makes an indirect call. Here the map function (lines 754 to 757 ) applies the function \(F\) to all elements in the list. The indirect call (line 756) is compiled into the indirect_call ( \(F, \cdots\) ) \(\mathbf{R}\)-expr.

The indirect_call R-expr makes the map function in figure 14-1 difficult to analyze and compile. Ideally, we want to eliminate all indirect_calls. In the case of the map function, this can accomplish by creating a specialized version as in figure 14-2. \({ }^{218}\)

\footnotetext{
\({ }^{218}\) In this example, folding away the indirect_call is similar to using a template in \(\mathrm{C}++\) which allows function calls to be static (rather than require a runtime indirection through a function pointer).
}
```

    b(L,Res) -> (Res=only(Inp,map_times_2(L, Inp)))
    proj(F, map(F,L,Res)*(F=times2[])) -> map_times_2
        (L, Res)
    times2(X) = X*2. map_times_2(L, Res) -> (Res=only(Inp,
b(L) = map( (L=[])*(Inp=[])+
times2[], L). proj(Head,proj(Tail,proj(Tmp1,proj(Tmp2,
assert b([1,2,3]) (L=[Head|Tail])*
== [2,4,6]. times(Head, 2, Tmp1)*
(a) Dyna
map_times_2(Tail, Tmp2)*
(Inp=[Tmp1|Tmp2])))))))

```
(b) R-expr with times \(_{2}\) folded into the map \(\mathbf{R}\)-expr.

Figure 14-2. Folding with map eliminates the indirect_call which can block other rewrites.

The remainder of this chapter is not intended to be a complete discussion about why one wants to fold programs or what kinds of programs folding can help solve. To find some discussion around why one wants to fold, how folding can solve certain kinds of programs, and how folding can be used to change the asymptotic runtime of a program, I suggest that the reader look into prior work such as Eisner and Blatz [58] and Vieira [141], both of which discuss folding on Dyna programs that have a single semiring. More generally, [22, 25, 80, 97] discuss folding on functional programming and logic programming.

\subsection*{14.2 How to Fold}

Folding a program represented as a collection of \(\mathbf{R}\)-exprs can be completed in a few steps. \({ }^{219}\)

First, observe that folding is only useful in the case of a recursive function. Nonrecursive functions can be expanded entirely. This means that we first identify if a user-defined \(\mathbf{R}\)-expr is recursive, or depends on a recursive \(\mathbf{R}\)-expr, and will target those user-defined \(\mathbf{R}\)-exprs only. These functions are identified by constructing

\footnotetext{
\({ }^{219}\) Folding on R-expr was implemented on the R-expr prototype https://github. com/argolab/ dyna-R, and has yet to be reimplemented on the Clojure implementation from chapter \(\S 11\).
}
a call graph and identifying which user-defined \(\mathbf{R}\)-exprs are depended on when expanding user-defined rewrites (rewrite rule 74). For example, if we have a userdefined \(\mathbf{R}\)-expr \(f(\cdots)\), which calls itself directly, then its rewrite rule will include a \(f(\cdots) \mathbf{R}\)-expr in the result of the rewrite:
\[
f(\cdots) \xrightarrow{74} \cdots * f(\cdots) * \cdots
\]

Or in the case of indirect calls such as with \(g(\cdots)\) and \(h(\cdots)\) that require computing the call graph:
\[
\begin{aligned}
& \mathrm{g}(\cdots) \xrightarrow{74} \cdots * \mathrm{~h}(\cdots) * \cdots \\
& \mathrm{~h}(\cdots) \xrightarrow{74} \cdots * \mathrm{~g}(\cdots) * \cdots
\end{aligned}
\]

Once the system has identified a user-defined \(\mathbf{R}\)-expr that is recursive, it unfolds the user-defined \(\mathbf{R}\)-expr so that it has a single large \(\mathbf{R}\)-expr that represents two or more steps of the recursive call. If the recursion happens indirectly (such as with \(g(\cdots)\) and \(h(\cdots)\) ), then the system unfolds the user-defined \(\mathbf{R}\)-exprs until it expands all the way to the recursive calls.

Once the recursive user-defined \(\mathbf{R}\)-exprs has been identified, the system will use folding to attempt to create new user-defined \(\mathbf{R}\)-exprs. These \(\mathbf{R}\)-exprs are identified as a conjunction of \(\mathbf{R}\)-exprs that are folded together. This is implemented by normalizing the variable names on the \(\mathbf{R}\)-expr and then using the fact that \(\mathbf{R}\)-exprs supports hashing and equality checks (chapter §8), and therefore can be used as keys in a hash-table.

If the newly folded \(\mathbf{R}\)-expr is different from the existing user-defined \(\mathbf{R}\)-exprs and previously folded \(\mathbf{R}\)-exprs, it is added to the set of automatically defined \(\mathbf{R}\)-exprs. If folded \(\mathbf{R}\)-expr already exists, then the existing \(\mathbf{R}\)-expr is referenced rather than creating another identical named \(\mathbf{R}\)-expr.

The newly created \(\mathbf{R}\)-expr will contain a call to a recursive user-defined \(\mathbf{R}\)-expr. Hence, it will be checked for additional folding opportunities within itself.

An example of this process is shown in the next section.

\subsection*{14.3 Example Using Folding to Solve a Recursive Program}

To see how folding a user-defined \(\mathbf{R}\)-expr works, let us work through an example of intersecting an even and odd length list in figure 14-3. Here, both even_list and odd_list are recursive functions. Independently, for these recursive functions, there is nothing that the system can do. However, once we take a conjunction of these on line 766, then there are rewrites that can identify that this term is impossible for all inputs provided we can work around the recursion using folding.
```

even_list([]).
even_list([X,Y|Tail]) :- even_list(Tail).
odd_list([X]).
odd_list([X,Y|Tail]) :- odd_list(Tail).
no_possible_list(List) :- even_list(List), odd_list(List).

```

Figure 14-3. Even and odd list length example. There is no list whose length is both even and odd, so the no_possible_list rule on line 766 as a null value for every list.

The \(\mathbf{R}\)-expr for figure 14 -3 is shown in figure 14-4:
```

even_list(L) $\rightarrow$ (true=exists(true,
(L=[]) +
proj(X, proj(Y, proj(Tail,
(L=[X,Y|Tail])*
even_list(Tail))))))
odd_list(L) $\rightarrow$ (true=exists(true,
$\operatorname{proj}(X,(L=[X]))+$
$\operatorname{proj}(X, \operatorname{proj}(Y, \operatorname{proj}(T a i l$,
( $L=[X, Y \mid$ Tail $]$ )*
odd_list(Tail))))))
no_possible_list(List) $\rightarrow$ (true=exists(true,
even_list(List)*odd_list(List)))

```

Figure 14-4. Even and Odd List translated into an R-expr. I am simplifying the presentation by omitting the Result variable as it would simply take on the value true at all times.

First, we can identify which user-defined \(\mathbf{R}\)-exprs are recursive in this program. We will identify both even_list, odd_list as recursive and determine that no_possible_list depends on them. Therefore, all of these R-exprs are candidates for folding.

When the system tries folding even_list and odd_list, no interesting/new \(\mathbf{R}\)-exprs are found. The resulting \(\mathbf{R}\)-exprs are equivalent to the existing even_list, and odd_list, so nothing is done. \({ }^{220}\)

However, when the system attempts to process no_possible_list, it unfolds both the even_list and odd_list and finds that it can use folding to create a new R-expr that represents the first step of even_list and odd_list, as shown in figure 14-5:

\footnotetext{
\({ }^{220}\) If a fold ends up getting created, then it would be equivalent to the existing \(\mathbf{R}\)-exprs, hence unproductive. An unproductive fold is not invalid, just not useful.
}
```

even_list(L)*odd_list(L) \xrightarrow{}{95} even_and_odd_list(L)
even_and_odd_list(L) \xrightarrow{}{96}
(true=exists(true,
(L=[])+
proj(X,proj(Y,proj(Tail,
(L=[X,Y|Tail])*
even_list(Tail))))))*
(true=exists(true,
proj(X, (L=[X]))+
proj(X,proj(Y,proj(Tail,
(L=[X,Y|Tail])*
odd_list(Tail))))))
no_possible_list(List) -> (true=exists(true,
even_and_odd_list(List)))

```

Figure 14-5. The no_possible_list user-defined \(\mathbf{R}\)-expr is defined in terms of the now folded even_and_odd_list. Observe that even_and_odd_list does not need to include aggregation at the top, like user-defined \(\mathbf{R}\)-exprs that come from Dyna programs, but rather is a conjunction of two aggregations that come from the definition of even_list and odd_list. Further rewrites can be performed against the definition of rewrite rule 96 .

The system can ten perform rewrites against rewrite rule 96, to determine that the ( \(\mathrm{L}=[\mathrm{l}\) ) and ( \(\mathrm{L}=[\mathrm{X}]\) ) are both impossible to satisfy. This is done by creating another \(\mathbf{R}\)-expr where ( \(L=[]\) ) is lifted to the of the \(\mathbf{R}\)-expr, and then observing that the resulting \(\mathbf{R}\)-expr rewrites as \(0 .{ }^{221}\) This results in the \(\mathbf{R}\)-exprs shown in figure 14-6.

\footnotetext{
\({ }^{221}\) E.g. \((A=\operatorname{sum}(X, R 1+R 2)) *(B=\operatorname{sum}(Y, S 1+S 2)) \xrightarrow{97}(A=\operatorname{sum}(X, R 2)) *(B=\operatorname{sum}(Y, S 1+S 2))\) if \(\forall_{E} \llbracket(A=\operatorname{sum}(X, R 1+R 2)) *(B=\operatorname{sum}(Y, S 1+S 2)) * \operatorname{proj}(X, \operatorname{proj}(A, R 1)) \rrbracket=0\).
}
```

even_list(L)*odd_list(L) \xrightarrow{}{95} even_and_odd_list(L)
even_and_odd_list(L) \xrightarrow{}{96}
(true=exists(true,
proj(X, proj(Y,proj(Tail,
(L=[X,Y|Tail])*
even_list(Tail))))))*
(true=exists(true,
proj(X,proj(Y,proj(Tail,
(L=[X,Y|Tail])*
odd_list(Tail))))))
no_possible_list(List) -> (true=exists(true,
even_and_odd_list(List)))

```

Figure 14-6. Analysis on rewrite rule 96 allows for ( \(L=[]\) ) and ( \(L=[X]\) ) to both be rewritten as 0 .

Now by rearranging rewrite rule 96, we can lift even_list and odd_list out of the aggregator as there are no other disjunctions present. This allows us to recognize that even_list and odd_list are conjuncts and will create an R-expr like figure 14-7.
```

even_list(L)*odd_list(L) \xrightarrow{}{95} even_and_odd_list(L)
even_and_odd_list(L) \xrightarrow{}{96}
(true=exists(true,
proj(X, proj(Y,proj(Tail,
(L=[X,Y|Tail])*
even_list(Tail)*
odd_list(Tail))))))
no_possible_list(List) -> (true=exists(true,
even_and_odd_list(List)))

```

Figure 14-7. even_list and odd_list are conjuncts inside of even_and_odd_list, rewrite rule 96 .

Using rewrite rule 95, we can apply the fold to rewrite rule 96's definition, which will allow us to identify that this has already been defined, and it recursively calls itself.
```

even_list(L)*odd_list(L) \xrightarrow{}{95} even_and_odd_list(L)
even_and_odd_list(L) \xrightarrow{}{96}
(true=exists(true,
proj(X, proj(Y,proj(Tail,
(L=[X,Y|Tail])*
even_and_odd_list(Tail))))))
no_possible_list(List) -> (true=exists(true,
even_and_odd_list(List)))

```

Figure 14-8. even_list and odd_list are folded together using rewrite rule 95, so we can now recognize that even_and_odd_list recurses to itself.

The same mechanism that identifies recursive functions also checks that all recursive functions have a base case. The reason is that recursive Dyna functions without a base case are equivalent to null for all values. Hence, we can rewrite rewrite rule 96 as 0 as shown in figure 14-9.
```

even_list(L)*odd_list(L) }\xrightarrow{}{95}\mathrm{ even_and_odd_list(L)
even_and_odd_list(L) \xrightarrow{}{96}0
no_possible_list(List) -> (true=exists(true,
even_and_odd_list(List)))

```

Figure 14-9. even_and_odd_list is found to be impossible due to the recursion without a base case. Hence, it is rewritten as 0

This sequence of rewrites and folding shows how the Dyna system can determine that there is no list that is simultaneously even and odd in length at the same time.

\subsection*{14.4 Folding Updatable User-Defined R-exprs}

Folded \(\mathbf{R}\)-expr depends on the original user-defined \(\mathbf{R}\)-expr. As such, if a userdefined \(\mathbf{R}\)-expr is modified (either by the API or by a user at the REPL), then all downstream folded \(\mathbf{R}\)-exprs have to be invalidated and refolded.

For user-defined R-exprs with memos, we completely disable folding for that
term. This is done by preventing user-defined R-exprs with memoization policies from unfolding when the system generates an \(\mathbf{R}\)-expr to check for folding opportunities (disabling rewrite rule 74). This is because memoized \(\mathbf{R}\)-exprs can include guesses, which cannot be bypassed (section §10.5). Furthermore, guesses are expected to change frequently during the running of the program as the system converges toward a final answer.

\section*{Chapter 15}

\section*{Properties of Simplify}

This chapter builds on the rewriting system defined in chapter §8. What we can prove about our rewriting system is unfortunately less interesting than what we might like-such is the reality of building a non-trivial programming language. Nevertheless, I claim the following properties about our rewrite system, and provide sketches for how these properties can be proven:
1. Soundness of Rewrites
2. Completeness on Datalog subset of Dyna
3. Incompleteness of Rewrites in general
4. Turing Completeness
5. Termination of SimplifyNormalize on bounded R-exprs

\subsection*{15.1 Is Simplify Sound and Complete?}

Being sound and complete are desirable properties of a logical system. A system is sound if it never proves a false statement. In other words, all statements that are proven true are actually true. A system is complete if there exists a proof for all true statements. In the case of term rewriting procedure (such as we have in this dissertation), completeness may instead be defined in terms of the proofs that are found through the systematic application of rewrite rules (e.g. Simplify given
in chapter §8), rather than claiming that there exist of a sequence of rewrites that can prove a statement.

When it comes to R-exprs, their semantics, and their rewrites, let us briefly discuss what it means for the rewrite system to prove a statement is true. Recall, the \(\mathbf{R}\)-exprs semantics is defined in terms of the multiplicity that it assigns to an environment \(E(\cdot)\) (chapter §5). As such, we will equate the idea of proving a statement as true with it having a non-zero multiplicity and is contained in the bag that the R-expr represents. Similarly, we will define false statements as those that have zero multiplicity and are not contained in the bag. To make this more concrete, we can say that for the rewrite system to prove an \(\mathbf{R}\)-expr R that does not contain free variables \((\operatorname{vars}(R)=\emptyset)^{222}\) true, there must exist a sequence of rewrites such that \(R \rightarrow * 1+S\), for some \(S .{ }^{223}\) Similarly, for the system to prove an R-expr \(R\) false, there must exist a sequence of rewrites \(R \rightarrow * \varnothing\) for it to be proven false.

\subsection*{15.1.1 Soundness of Simplify}

We claim that our rewrite system is sound in that all \(\mathbf{R}\)-exprs that are rewritten (and can be rewritten) into the form \(1+S\), for some \(S\), have a multiplicity greater than 1.

Proof: Our rewrite rules given in chapter §6 are semantics preserving with respect to the semantic interpretation given in chapter \(\S 5\). Hence, any R-expr \(R\) that is rewritten into the form \(1+S\) must be semantically equivalent and therefore have a multiplicity greater than 1 and be considered true. \({ }^{224}\)

\footnotetext{
\({ }^{222}\) Allowing free variables in the \(\mathbf{R}\)-expr requires more work to concisely define this. For example, the \(\mathbf{R}\)-expr \((X=1)\) has a free variable \(X\), but cannot rewrite as 1 or 0 . Hence, it is odd to think of this of \((X=1)\) as proving \(X\). Rather, by focusing on only the case without free variables, we do not have to define if \((X=1)\) is true or false. This restriction is not an issue though, as one can simply project out all free variables, in which case it can be rewritten as a multiplicity. E.g. proj \((X,(X=1)) \rightarrow 1\).
\({ }^{223}\) Writing the \(\mathbf{R}\)-expr as \(1+S\) allows for this to be more general, in that it can match all multiplicities \(\geq 1\), and matches the rewrite rule for if-expressions (rewrite rule 63). Alternately, we could write this as \(\operatorname{if}(R, 1,0)\), which would guarantee that this \(\mathbf{R}\)-expr is rewritten as 1 or 0 .
\({ }^{224}\) The rewrites in chapter \(\S 6\) have been checked by hand to match the semantic definitions in chapter \(\S 5\). Furthermore, the rewrites in the implementations in chapters 8 and 11 have also been checked by hand. Checking rewrites and their implementation by hand does not guarantee that the rewrites are bug-free. Any rewrite, or implementation of a rewrite, that is not semantic preserving is a bug.
}
```

fermat_was_right }
if(proj(A,proj(B,proj(C proj(N,proj(Aval,proj(BVal,proj(CVal,
power(A, N, Aval)*power(B, N, BVal)*power(C, N, CVal)*
add(AVal, BVal, CVal)*int(A)*int(B)*int(C)*int(N)*
lessthan(2, N)*lessthan(0, A)*lessthan(0, B)*lessthan(0, C)
)))))), 0, 1).

```

Figure 15-1. R-exprs are sufficiently powerful to express general mathematical expressions such as Fermat's last theorem shown here (I.e., there does not exist positive integers \(a, b, c, n\) where \(n \geq 3\) such that \(a^{n}+b^{n}=c^{n}\) ). While this Fermat's theorem was proven true [148], hence this \(\mathbf{R}\)-expr can be rewritten as 1 , it is unrealistic to include the necessary rewrites to solve this \(\mathbf{R}\)-expr.

\subsection*{15.1.2 R-exprs Rewrites are Incomplete}

Our rewrite system is not complete. Given the definitions given previously, for a rewrite system to be complete, any true statement that can be expressed must be rewritable into the form \(1+S\), for some \(S\). In other words, there exists a \(\mathbf{R}\)-expr \(R\) that is semantically equivalent to \(1+S\) for some \(s\), but there are no rewrites that can rewrite R into the form \(1+S\).

Proof: To prove that \(\mathbf{R}\)-exprs are not complete, we can simply demonstrate an \(\mathbf{R}\)-expr that is equivalent to 1, but cannot be rewritten as 1. Figure 15-1 shows an \(\mathbf{R}\)-expr that encodes Fermat's last theorem, which was proven true in [148]. Hence, the fermat_was_right \(\mathbf{R}\)-expr in figure \(15-1\) is semantically equivalent to the R-expr 1. However, our rewrite system does not include any rewrites that can rewrite the fermat_was_right \(\mathbf{R}\)-expr. (The proof of Fermat's last theorem requires mathematical properties that we currently do not have implemented for \(\mathbf{R}\)-exprs). The fact that there exists an \(\mathbf{R}\)-expr that cannot be rewritten shows that the rewrite system is incomplete.

Admittedly, users of the Dyna system are likely not expecting it to automatically prove Fermat's last theorem, hence this is not an informative result.

More generally, we can say that any sound rewriting system for \(\mathbf{R}\)-exprs cannot be complete in the sense that any two semantically equivalent \(\mathbf{R}\)-exprs can be non-deterministically rewritten into the same form. (In other words, for some pair of \(\mathbf{R}\)-exprs \(\mathrm{R}, \mathrm{R}^{\prime}\) with \(\llbracket \mathrm{R} \rrbracket_{E}=\llbracket \mathrm{R}^{\prime} \rrbracket_{E}\), there is no S such that \(\mathrm{R} \rightarrow{ }^{*} \mathrm{~S}\) and \(\mathrm{R}^{\prime} \rightarrow{ }^{*} \mathrm{~S}\).) The
reason is that \(\mathbf{R}\)-exprs are powerful enough to express any first-order statement about arithmetic. Hence, R-exprs are subject to Gödel's incompleteness theorem, which ensures that there exists an \(\mathbf{R}\)-expr that is a true, i.e., semantically equivalent to 1 , yet cannot be rewritten as 1 .

\subsection*{15.1.2.1 Completeness on Datalog Subset}

While R-expr and their rewrites are incomplete, we can prove that there are some subsets of the Dyna language that are complete.

First, I will prove that the Datalog subset of Dyna is complete when using the appropriate memoization policy.

Proof: First, let us briefly recall how Datalog works, which was previously discussed in section §3.1.2. Datalog memoizes all terms that have been proven true. Rules in a Datalog program combine terms that have been proven to deduce new terms that are also true. Furthermore, for a program to be a valid Datalog program, it must be stratified when aggregation or negation is used. This means that we do not have to worry about cyclic programs, which require the guessing mechanism from section §10.5. Finally, Datalog terms only contain ground values (e.g. the term foo \([1,2]\), but not bar \([\mathrm{X}, \mathrm{X}]\) ). This means that all of the terms can be easily enumerated by enumerating the ground assignments to variables in the term.

To emulate the memoize everything behavior, we can use the \$memo policy that returns "null" for all user-defined terms. \({ }^{225}\) This works because we are only concerned with the Datalog subset, which only proves ground terms. This means that the memo tables only consist of disjunctions of ground assignments to the variables present on a term.

For example, if we have \(\$\) memo \((f o o[X, Y])=" n u l l "\). as the memoization policy for a Datalog program. The foo/2 memo will look something like \((X=1) *(Y=2)\) * (Result=true) \(+(\mathrm{X}=7) *(\mathrm{Y}=3) *(\) Resul \(\mathrm{t}=\) true \()+\cdots+(\mathrm{X}=11) *(\mathrm{Y}=\) "hello" \() *(\) Result t true) ,

\footnotetext{
\({ }^{225}\) This would be equivalent to defining \(\$ m e m o(X)=" n u l l "\). However, due to limitations in the \$memo control mechanism, each user-defined term must be specified with its own \$memo (•) policy. Hence, this would require specifying multiple \$memo policies like:
\$memo(foo \([X, Y])=\) "null".
\$memo(bar \([X, Y, Z])=" n u l l "\).
}
where all of the values to \(X\) and \(Y\) are known ground values, and therefore enumerable. \({ }^{226}\)
Because Datalog rules combine the purely ground terms together by looping over the ground bindings to variables, we will exactly replicate this behavior using our standard execution procedure. Furthermore, Datalog programs are stratified \({ }^{227}\) when it comes to aggregation and negation, which ensures that the set of terms that are true and memoized will converge. Hence, Dyna's execution exactly matches what happens in a typical Datalog implementation.

Because our execution exactly replicates Datalog (when given an appropriate memoization policy), we can claim that our implementation is complete in the same way that Datalog inference is complete.

\subsection*{15.1.2.2 Emulating SLD Resolution}

Selective Linear Definite (SLD) resolution, as defined in Foundations of Logic Programming by Lloyd [101], is a technique to implement resolution on horn clauses. \({ }^{228}\) SLD resolution performs unification between non-recursive structural terms that can contain variables. For example, if the terms foo \([X, b a r[X, Y]]\) and foo \([A, \operatorname{bar}[B, B]]\) are unified together, then the variable \(X, Y\) and \(Z\) will all be unified together ( \(X=Y=Z\), hence foo[ \(X, \operatorname{bar}[X, X]]\) ). SLD resolution also includes an occurs check which ensures that terms recursive terms are identified as impossible. For example, the unification between \(f[X, X]\) and \(f[Y, g[Y]]\) would fail as this would result in \(Y=g[Y]\), and there is no value of \(Y\) that will make this expression true.

Horn clauses \({ }^{228}\) that are true can be forward chained to identify all terms that are true. This is the same sort of forward chaining that we have in a Datalog program without negation or aggregation. As such, SLD resolution is known to be sound and complete for horn clauses [101].

To see an example of SLD resolution, consider the following program.
\({ }_{767} \mid a(f[X])\).

\footnotetext{
\({ }^{226}\) In fact, memos of this form will be efficiently handled by the prefix-tries (section §11.6.1) and iterators (section §11.7).
\({ }^{227}\) There is nothing in Dyna that checks that a program is stratified and enforces stratification. Hence, I am assuming that the program is already in the Datalog subset and is correctly stratified, meaning it does not contain cycles that involve negation or aggregation.
\({ }^{228}\) Horn clauses are sometimes referred to as Pure-Prolog. The term Pure-Prolog is not used consistently across all publications, so I will avoid the term Pure-Prolog.
}
```

768 a(g[I,J]).
769 b(Y,Z) :- a(Y).
70 c(W) :- c(W).

```

First SLD resolution will deduce the terms \(a[f[X]]\) and \(a[g[I, J]]\) are true from lines 767 and 768 as there are no conditions for these horn clauses. SLD resolution will then use \(a[f[X]]\) to forward chain into the rule on line 769 , where the variable \(Y\) is unified with the structure \(f[X]\). This causes SLD resolution to deduce the term \(b[f[X], Z]\) is true. Next, SLD resolution forward chains \(a[g[I, J]]\), which causes \(\mathrm{b}[\mathrm{g}[\mathrm{I}, \mathrm{J}], \mathrm{Z}]\) to be deduced as true as well. Now the terms that have been proven true are \(a[f[X]], a[g[I, J]], b[f[X], Z]\), and \(b[g[I, J], Z]\). SLD resolution will not stop running as it has reached a fixed point where all possible terms have been deduced as true. Note that \(c[W]\) is never deduced as true, as there is no base case for \(c[W]\) to get it started.

Emulating SLD with R-exprs: SLD resolution can be emulated using R-exprs and our rewrites. However, it requires a slightly different control strategy from the one discussed in chapter \(\S 8\).

First, observe that we already have the necessary rewrites for unification and the occurs check (section §6.1.1). Second, our memo table is capable of memoizing structural terms with variables like foo \([\mathrm{X}, \mathrm{bar}[\mathrm{X}, \mathrm{Y}]]\) by memoizing an \(\mathbf{R}\)-expr containing structural unifications.

However, there is a slight difference from what SLD memoizes and uses to forward chain. Observe that in the previous SLD example, the terms a[f[X]] and \(a[g[I, J]]\) were each memoized individually and forward chained individually. Essentially, the disjunction between the different ways that \(X\) can be satisfied in \(a(X)\) is handled directly by the SLD algorithm. In the case of R-exprs, we instead would create a single memo for all of \(a(X)\), and then use the disjunction (section §5.2.2.5) to represent the different ways that \(X\) can be satisfied. In other words, when evaluating line 769 and expanding \(a(Y)\) using \(\mathbf{R}\)-exprs, we are going to have an \(\mathbf{R}\)-expr like \(b(Y, Z) \rightarrow\) (true=exist(true, \(\operatorname{proj}(X,(Y=f[X]))+\operatorname{proj}(I, \operatorname{proj}(J,(Y=g[I, J])))))\). In other words, because of our desire to make factored R-exprs to work around aggregation (section §8.2.3.1), we do not end up expanding out the terms entirely as SLD resolution does.

To emulate SLD resolution using \(\mathbf{R}\)-exprs. We could instead consider a system that uses the distributive rewrite to expand \(\mathbf{R}\)-exprs rather than factor them (e.g.
\(Q *(R+S) \xrightarrow{22} Q * R+Q * S)\). In this case, we will have that each \(R\)-expr would only represent a single conjunctive structural term with variables rather than the factored form that we currently have.

As a secondary issue, we also have aggregation in the program that needs to be removed before the distributive rewrite can be used to fully expand the R-expr. In principle, nested exists aggregators could be removed. \({ }^{229}\) E.g. (true=exists(true, \(\mathrm{Q} * \mathrm{R} *(\) true \(=e x i s t s(\) true, \(\mathrm{S} 1+\mathrm{S} 2)))) \xrightarrow{98}(\) true \(=e x i s t s(\) true, \(\mathrm{Q} * \mathrm{R} *(\mathrm{~S} 1+\mathrm{S} 2))) .{ }^{230}\)

In conclusion, if aggregators were removed from the R-expr, and the distributive rewrite was used to expand \(\mathbf{R}\)-exprs, rather than factor, then our memoization mechanism is sufficiently powerful to memoize representations of structured terms with variables, and our unification rewrites are the same as those used by SLD resolution. The system does not perform rewrites in this way because of our desire to work around aggregation (a frequent operation in Dyna programs).

\subsection*{15.2 Dyna is Turing Complete}

Dyna is Turing complete. This should not be surprising as Prolog is known to be Turing complete, and Dyna is a superset of Prolog. Proving Dyna is Turing complete is useful as it shows that all computable functions can be computed using Dyna using some program. However, being Turing complete does not mean that all ways in which a function can be represented will result in Dyna being able to compute using the result of the function. Hence, showing that Dyna is Turing complete does not contradict the previous proof that \(\mathbf{R}\)-exprs and rewrites on \(\mathbf{R}\)-exprs are incomplete.

Proof: Proving Turing completeness can be done by reducing \({ }^{231}\) from a known universal Turing machine into Dyna. I have chosen to use Rule- \(110^{232}\), as it has

\footnotetext{
\({ }^{229}\) The (true=exists (true, R)) aggregator essentially forces the multiplicity of \(R\) to 1 . Hence, as long as there is an exists aggregator at the top of the \(\mathbf{R}\)-expr, the entire \(\mathbf{R}\)-expr will have the same semantics.
\({ }^{230}\) Currently rewrite rule 98 is not included in the implementation.
\({ }^{231}\) This kind of reduction is the same kind of reductions that one sees with NP-complete proofs using a reduction from SAT.
\({ }^{232}\) https://en.wikipedia.org/wiki/Rule_110
}
been proven to be Turing complete [40] \({ }^{233}\) and is considered one of the simplest known Turing complete systems. The implementation of Rule-110 in Dyna is shown on lines 771 to 786:
```

77
772
773
774
7 7 5
776
7 7 7
778

```

```

780
781
782
783
784
84
turing_machine_step([1,1,1|X]) = [0|turing_machine_step([1,1|X])].
turing_machine_step([1,1,0|X]) = [1|turing_machine_step([1,0|X])].
turing_machine_step([1,0,1|X]) = [1|turing_machine_step([0,1|X])].
turing_machine_step([1,0,0|X]) = [0|turing_machine_step([0,0|X])].
turing_machine_step([0,1,1|X]) = [1|turing_machine_step([1,1|X])].
turing_machine_step([0,1,0|X]) = [1|turing_machine_step([1,0|X])].
turing_machine_step([0,0,1|X]) = [1|turing_machine_step([0,1|X])].
turing_machine_step([0,0,0|X]) = [0|turing_machine_step([0,0|X])].
turing_machine_step([A,B]) = [A,B].
turing_machine_step([A]) = [A].
turing_machine_step([]) = [].
turing_machine(X) := turing_machine(turing_machine_step(X)).
turing_machine(X) := X for X = turing_machine_step(X). % Turing
machine has reached a fixed point
78
786 final_output = turing_machine(start_state).

```

Rule-110 is a cellular automaton that uses a one-dimensional array consisting of 0 s and 1 s . The array is progressively rewritten using a pattern of 3 bits and then shifting over in the array by 1 position. Rule 110 is applied repeatedly using line 783 until it reaches a final state where the Turing machine stops changing, which is identified by line 784.

Observe that this program depends on recursion (line 783). Recursion is the only way in which a programmer can represent an unbounded computation (such as a Turing machine), which I will prove next.

\footnotetext{
\({ }^{233}\) The proof that Rule-110 is Turing complete requires several very dense math papers.
}

\subsection*{15.3 Termination of SimplifyNormalize on a Bounded Size R-expr}

A potential complication when designing a rewrite rule system is that the rewrite rules could oscillate unproductively, causing a terminating program to become nonterminating. We claim that all terminating programs will terminate when rewritten by SimplifyNormalize, and that the resulting R-expr will be "simple", according to some definition of "simple". This ensures that 1) we are doing something useful when rewriting (hence the result will be simple), and 2) that we are not accidentally causing non-termination. \({ }^{234}\)

First, the only way one can write a Dyna program that does not terminate is either through the use of a bad memoization policy (chapter \(\S 10\) and section §2.5), or through the use of recursion to write an unbounded loop. The function SimplifyNormalize is not directly involved with memoization; hence, memoization need not be considered in this proof. Therefore, for this proof to apply, we only need to ensure that there is no unbounded recursion, which means that the R-expr being rewritten by SimplifyNormalize is bounded in size.

\subsection*{15.3.1 Making an R-expr Bounded in Size}

In general, R-exprs are not bounded in size. Hence, to make the proof in the next section go through, we first need to ensure that the \(\mathbf{R}\)-expr that is being rewritten is bounded in size so that the proof will apply. To do this, we "modify" the \(\mathbf{R}\)-expr to ensure that the \(\mathbf{R}\)-expr is bounded in size.

The way this is done is by tracking the stack depth \({ }^{235}\) of user-defined rewrites. This is done by modifying the presentation from section \(\S 6.7\) so that there is a stack depth counter. Here, I denote the stack depth as a superscript on the user-defined \(\mathbf{R}\)-expr types and modify rewrite rule 74 so that the counter is incremented for all internal user-defined R-exprs.

\footnotetext{
\({ }^{234}\) This proof does not prove that this will terminate with the correct runtime, or that it will run "fast".
\({ }^{235}\) We do not really have a call stack in the traditional sense, as we are expanding the R-expr out using rewrite rule 74 , but we can think of the stack depth as corresponding to the depth at which a given call would appear if the program was executed under a traditional programming language.
}
```

proj(Tmp1,
times(100,Tmp1,Result)*
proj(Tmp2,
factorial(X) := times(99,Tmp2,Tmp1)*
factorial(X-1)*X.
factorial(0) := 1.
proj(Tmp3,
times(98,Tmp3,Tmp2)*
factorial4(97, Tmp3))))
(a) Dyna

```
(b) The factorial R-expr expanded up to \(K=4\) after some rewrites are applied.

Figure 15-2. Recursive Dyna program which requires many steps of recursion to get a "useful" result. If the program's expansion is cut short, with \(K=4\), then the resulting \(\mathbf{R}\)-expr is still correct, though not useful as in (b).
\(\mathrm{f}^{k}\left(\mathrm{X}_{1}, \ldots, \mathrm{X}_{n}\right) \xrightarrow{74} \cdots * \mathrm{f}^{k+1}(\cdots) * \cdots * \mathrm{~g}^{k+1}(\cdots) \star \cdots \quad\) if \(k<K\)

Now the rewrite rule 74 rule only runs if the stack depth is less than some max-stack-depth, denoted here as \(K\).

The approach of bounding the number of expansions of recursive user-defined \(\mathbf{R}\)-exprs is not without consequences. For example, if we set the stack limit too low, then we can end up with R-exprs that are not useful to the user. Figure 15-2 shows one such instance, where the factorial program stops expanding before reaching its base case. The R-expr that would be returned to the user's query would be partially expanded but does not return the desired numerical result of factorial.

Similarly, limiting the stack depth does not necessarily result in a small R-expr or a limited amount of memory consumed (as would commonly happen in a procedural stack-based programming language). For example, figure 15-3 shows a Dyna program that would create an exponential \(3^{K}\)-sized \(\mathbf{R}\)-expr when expanded to depth \(K\).

\footnotetext{
\({ }^{236}\) For example, \(f(X-1) * f(X-1) * f(X-1)\) could rewritten into \(Y=f(X-1), \quad Y * Y * Y\). In which case \(f(X-1)\) would only be called once, avoiding the exponential-sized expansion.
}
\(790 \mid f(X):=f(X-1) * f(X-1) * f(X-1)\).

Figure 15-3. Assuming that no rewrites are done to combine \(f(X-1)\) on line \(790,{ }^{236}\) then this program will expand to an exponential \(3^{K}\)-sized R-expr. Hence, the stack depth does not mean small program, or limited memory etc.

\subsection*{15.3.2 Outline: How to Prove Termination}

To prove that SimplifyNormalize terminates and Simplify reaches a fixed-point, we prove that there cannot be an infinite sequence of directional rewrites on a bounded size \(\mathbf{R}\)-expr. To do this, we define a set of energies \(\mathbb{E}\) such that there is no infinite descending sequence, and we define an energy function \(|\cdot|: \mathbf{R} \rightarrow \mathbb{E}\) such that
\[
|\operatorname{SimpLIFY}(R)| \leq|R|
\]
and
\[
(\operatorname{Simplify}(R) \neq R) \Longrightarrow(|\operatorname{SimpLify}(R)|<|R|) .
\]

Together, these properties mean that Simplify will either return the R-expr unmodified, meaning no rewrites were applied, or if it returns a different R-expr, where there were one or more rewrites applied to, it will have decreased the energy. Because there is no infinite descending sequence in \(\mathbb{E}\), this means that repeated applications of Simplify must reach a fixed point in a finite number of steps.

What is left to complete this proof is to define \(|\cdot|: \mathbf{R} \rightarrow \mathbb{E}\) and show that the rewrites defined in chapter \(\S 6\) will individually decrease the energy of the R-expr.

This proof is structured as follows. First, I will define \(|\cdot|_{\text {core }}: \mathbf{R} \rightarrow \mathbb{N}_{\geq 1}\) that is the energy for all R-expr types excluding built-in R-exprs (section §5.2.2.3). Second, in section \(\S 15.3 .4\), I will define a notion of energy specifically for built-ins. Finally in section \(\S 15.3 .5\), I will define \(|\cdot|: \mathbf{R} \rightarrow \mathbb{E}\) which applies to all \(\mathbf{R}\)-exprs.

\subsection*{15.3.3 Construction of "Core" Energy}

First, we define \(|\cdot|_{\text {core }}: \mathbf{R} \rightarrow \mathbb{N}_{\geq 1}\) as the energy for all \(\mathbf{R}\)-exprs except for built-in \(\mathbf{R}\)-expr types (section §5.2.2.3). The reason built-ins are excluded for now is that the rewrites on built-ins can be extended. This can be done by defining new rewrite
rules or by incorporating existing software libraries to solve conjunctions of built-in constraints (such as a linear programming solver or an SMT solver, section §16.8).
Non-recursive: \(\quad\) Non-recursive R-exprs have energies defined as follows:
1. \(|\mathrm{M}|_{\text {core }}=1\) - Multiplicities are a base case of the R-expr language as they cannot be rewritten further. They take on the smallest possible energy regardless of the value of \(M \in \mathcal{M}\).
2. \(|(X=G)|_{\text {core }}=2\) where \(G \in \mathcal{G}\) - Unification of a variable with a constant has energy 2 regardless of the value of \(G\). In the case where the unification with a constant is redundant, it is rewritten as a multiplicity of 1 , with an energy of 1 . Hence, this decreases the energy.
3. \(|(X=Y)|_{\text {core }}=3+\left\{\begin{array}{ll}0 & X \prec Y \\ 1 & X \succ Y\end{array}\right.\) where \(X, Y \in \tilde{\mathcal{V}}-\) With two variables in a unification \(\mathbf{R}\)-expr. One of the variables could is bound with a constant value, the energy will decrease. For example, \(|(X=Y)|_{\text {core }}>|(X=7)|_{\text {core }}\), where this rewrite could presumably happen as a result of some other \(\mathbf{R}\)-expr determining the value of \(Y\). E.g. \((Y=7) *(X=Y) \xrightarrow{5}(Y=7) *(X=7)\). To define the energy in the case that variables are reordered, there is an arbitrarily chosen order \(\prec\) which defines what variable should appear first in the unification R-expr.
4. \(\left|\left(X=f\left[U_{1}, \cdots, U_{n}\right]\right)\right|_{\text {core }}=3 m+3-\) Unification with a structured term that contains variables is assigned \(3 m+3\) where \(m\) is the number of free variables in the \(\mathbf{R}\)-expr. The reason we need to track the number of variables is that propagating a value into these \(\mathbf{R}\)-exprs must decrease the energy of the overall R-expr (rewrite rule 5). For example: \((X=9) *(Y=f[X, Z]) \rightarrow(X=9) *(Y=f[9, Z])\).

Recursive: Recursive R-exprs have energies defined as follows:
5. \(|R * S|_{\text {core }}=|R|_{\text {core }}+|S|_{\text {core }}-\) Conjunctive \(\mathbf{R}\)-exprs energies is defined as the sum of their sub-R-exprs. This means that any reduction in the energy of \(R\) or \(S\) will reduce the energy of the overall R-expr.
Furthermore, given that all R-exprs have a strictly positive energy, this means that removing a conjunct and unneeded constraints will reduce the energy. For example, rewrite rule 12 will multiply multiplicities, e.g. \(2 * 3 \xrightarrow{12} 6\), which takes this R-expr from an energy of 2 to 1 .
6. \(\left|R_{1}+R_{2}+\cdots+R_{n}\right|_{\text {core }}=2^{\left|R_{1}+R_{2}+\cdots+R_{n}\right|_{\text {core }}}\) - Disjunctive \(\mathbf{R}\)-exprs are raised to a power of 2. This ensures that nested \(\mathbf{R}\)-exprs are more expensive than \(\mathbf{R}\)-exprs which are not nested. Hence, the energy of an \(\mathbf{R}\)-expr is reduced when common sub-R-exprs are factored out.

Note that the presentation of disjunction presented here assumes a \(n\)-arity disjunction and not a recursively nested binary disjunction. However, the definition can be binarized by matching against the nested sub-R-exprs. \({ }^{237}\)
7. \(|\operatorname{proj}(X, R)|_{\text {core }}=2^{|R|_{\text {core }}}\) - Projected R-exprs are also nested. During simplification, we attempt to factor out expressions as much as possible; this means that more deeply nested sub-R-exprs will have higher energy. For example: \(|\operatorname{proj}(X, R * S)|_{\text {core }}>|R * \operatorname{proj}(X, S)|_{\text {core }}\).

Additionally, note that unlike structural unification (item 4), we do not have to check if \(x\) is a variable or a ground value. The reason is that if \(x \in \mathcal{G}\), then we can eliminate the projection.
8. \(\mid\) if \(\left.(Q, R, S)\right|_{\text {core }}=|Q|_{\text {core }}+|R|_{\text {core }}+|S|_{\text {core }}\)-if-expressions have an energy which is the sum of the nested expressions. The if-expressions are typically handled by first rewriting \(Q\) until it can be determined if it is rewritten as 0 or as a non-zero multiplicity. Once the if-expression is determined to be true or false, it is rewritten as either \(R\) or \(S\) (rewrite rules 63 and 64). Thus, the energy of this expression needs to be higher than \(R\) or \(S\). Additionally, rewrites can be performed directly on \(R\) or \(S\) for the purposes of making the \(\mathbf{R}\)-expr more efficient (e.g., as in the case of memoization chapter \(\S 10\) ). When this happens, the energy of the if-expression needs to still decrease, regardless if we rewrite R or \(\mathrm{S} .{ }^{238}\)
\({ }^{237}\) The binarised definition for disjunctions can be written as:
\[
|R+S|_{\text {core }}=2^{\wedge}\left(\left\{\begin{array}{ll}
\log _{2}|R|_{\text {core }} & \text { if } R \text { matches } A+B \\
|R|_{\text {core }} & \text { otherwise }
\end{array}\right\}+\left\{\begin{array}{ll}
\log _{2}|S|_{\text {core }} & \text { if } S \text { matches } A+B \\
|S|_{\text {core }} & \text { otherwise }
\end{array}\right\}\right)
\]

This definition works by matching if the nested expression is also a disjunction and undoing the \(\left.2^{\mid \mathrm{R}}\right|_{\text {core }}\) operation via \(\log _{2}\).
\({ }^{238}\) For the if-expression, an energy of \(|Q|_{\text {core }}+\max \left(|R|_{\text {core }},|S|_{\text {core }}\right)\) would not work, as if we have \(|R|_{\text {core }}>|S|_{\text {core }}\), and rewrites are performed on \(S\), then the energy of the if-expression would not decrease.

Bidirectional Rewrites: Observe that conjunctions, disjunctions, and projections have bidirectional R-exprs that are unproductive. By unproductive, I mean that a rewrite is only used to rearrange the \(\mathbf{R}\)-expr, and does not result in any meaningful change or decrease in energy. For example, the commutativity rewrites on conjunctions and disjunctions swap the order in which two sub-R-exprs are represented: e.g. \(R * S \stackrel{19}{\longleftrightarrow} S * R\). For these rewrite rules 18 to 21 and 40 , we do not actually use them and instead depend on the context \(\mathcal{C}\) to identify the necessary conjunctions between \(\mathbf{R}\)-exprs when rewriting; therefore, we do not rearrange the R-expr explicitly.

Some bidirectional rewrite rules are useful, though. For example, the distributivity rewrite can create factored \(\mathbf{R}\)-exprs from a disjunction ( \(R * S+R * Q \xrightarrow{22} R *(S+Q)\) ). These kinds of bidirectional rewrites are applied explicitly in a uni-directional manner to create more factored \(\mathbf{R}\)-exprs and used implicitly via the context \(\mathcal{C}\) to handle the "expanding out" direction of the rewrite (rewrite rules 22, 38 and 39, also recall section §8.2.3.1).
Aggregation: Aggregation requires a bit more care in how we define energy. Recall that we want to split the aggregation into smaller \(\mathbf{R}\)-exprs when it is an aggregation over a disjunction (e.g. \((A=\operatorname{sum}(X, R+S))\) ), so that each disjunct can be aggregated independently (rewrite rule 50). To accomplish this, we had to introduce aggregators that have a disjunction with ( \(\mathrm{X}=\) agg_null) to suppress the behavior of rewrite rule 48 which would cause it to be rewritten as 0 (e.g. ( \(A=\operatorname{sum}(X,(X=\) agg_null \()+R)\) ), see section \(\S 6.5 .1\) for the complete discussion). The reason we chose this representation for the aggregator \(\mathbf{R}\)-expr is that it makes it clear what the required behavior is from an aggregator when handing a disjunction. However, this ( \(\mathrm{X}=\) agg_null) does not behave like a disjunction but rather a different kind of behavior of the aggregator. Hence, for the purposes of defining the energy of an aggregator, we are going to include ( \(\mathrm{X}=\) agg_null) as part of the aggregator instead of as a nested disjunction. \({ }^{239}\)
9. \(|(A=\operatorname{sum}(X, R))|_{\text {core }}=\mid(A=\operatorname{sum}(X,(X=\) agg_null \()+R)) *\) not_equal \(\left.(A\), agg_null \()\right|_{\text {core }}\) +1 if ( \(R\) not match ( \(X=\) agg_null) \(+S\) ) - The "normal" aggregator is defined in terms of \((A=\operatorname{sum}(X,(X=\) agg_null \()+R))\), which contains the \((X=\) agg_null \()\) dis-

\footnotetext{
\({ }^{239}\) In the implementation, ( \(X=\) agg_null) is tracked as a boolean flag on the aggregator class instead of as a disjunction in the aggregator's body. See footnote 94 in section §6.5.1.
}
junction, which suppresses rewrite rule 48. The reason for this energy definition is that we can think of this as handling the first part of rewrite rule 50, where a constraint not_equal (A, agg_null) is added to the \(\mathbf{R}\)-expr to ensure that the case where \(R\) is 0 is handled correctly. This looks something like rewrite rule 99, and we can observe that this rewrite decreases the energy of the \(\mathbf{R}\)-expr.
\[
(A=\operatorname{sum}(X, R+S)) \xrightarrow{99}(A=\text { sum }(X,(X=\text { agg_null })+R+S)) * \text { not_equal }(X, \text { agg_null })
\]
10. \(\mid\left.(A=\operatorname{sum}(X,(X=\) agg_null \()+R))\right|_{\text {core }}=\left(2 \cdot|R|_{\text {core }}\right) 2\), where the left superscript denotes tetration \({ }^{240}\) - Aggregation with ( \(X=\) agg_null) follows the same basic idea of projection and disjunction in that more deeply nested sub-R-exprs will have a higher energy. As such, "solving" an aggregator, either by determining the value returned by an aggregator (rewrite rules 49, 51, 55 and 56) or splitting an aggregator into aggregations over small R-exprs (rewrite rules 50 and 53) will decrease the energy of the R-expr. In this case, we only have to concern ourselves with rewrite rule 53, as the other rewrites for aggregators either factor an R-expr out, which will obviously decrease the energy, or solve the aggregator without having to deal with any intermediate states.
\[
\begin{aligned}
& A=\operatorname{sum}\left(X, \quad\left(X=\operatorname{agg} \_n u l l\right)+R+S\right) \xrightarrow{53} \operatorname{proj}(B, \operatorname{proj}(C, \\
&\left(B=\operatorname{sum}\left(X,\left(X=\operatorname{agg\_ null)+R))}\right.\right.\right. \\
& *\left(C=\operatorname{sum}\left(X,\left(X=\operatorname{agg} \_n u l l\right)+S\right)\right) \\
&* \operatorname{plus}(B, C, A))) \\
& \text { if } R \neq(X=\text { agg_null }) \text { and } S \neq(X=\text { agg_null })
\end{aligned}
\]

Looking at rewrite rule 53, we can observe that there are two nested projections and that there are two nested aggregators. In other words, we need:
\[
\left\lvert\, \begin{array}{l|l}
\begin{array}{l}
(A=\operatorname{sum}(X, \\
\left(X=\operatorname{agg} \_n u l l\right) \\
+R+S))
\end{array} & \left.\right|_{\text {core }}
\end{array} \quad>2^{\wedge}\left(2^{\wedge}\left(\begin{array}{l}
\mid\left.(B=\text { sum }(X,(X=\text { agg_null })+R))\right|_{\text {core }}+ \\
\mid\left.(C=\text { sum }(X,(X=\text { agg_null })+S))\right|_{\text {core }}+ \\
\mid \text { plus }\left.(B, C, A)\right|_{\text {core }}
\end{array}\right)\right) .\right.
\]

This is what the energy will be once we expand rewrite rule 53 with two additional projections.

\footnotetext{
\({ }^{240}\) For example, \({ }^{3} a=a^{a^{a}}\). The notation and the term tetration are due to Rucker [115]:
\({ }^{n} a \stackrel{\text { def }}{=} \begin{cases}1 & \text { if } n=0 \\ \left.a^{(n-1} a\right) & \text { otherwise }\end{cases}\)
}

To make this work, first observe for a \(n\)-way splittable disjunction that
\[
\left|R_{1}+R_{2}+\cdots+R_{n}\right|_{\text {core }} \geq 2^{n}
\]
by the definition of the disjunction's energy. Therefore, the energy of a disjunction is an upper bound on the number of disjunctive branches that will need to be handled by the aggregator. In other words, the system could potentially introduce up to \(n\) projections and intermediate variables to handle aggregation over the disjuncts \(R_{1}\) to \(R_{n}\). Therefore, the energy of the aggregator's body is an upper bound \({ }^{241}\) on the number of times that rewrite rule 53 can be applied, and the number of projections that will be introduced. To handle this, we use tetration \({ }^{240}\) which takes a number to a power multiple times. In performing rewrite rule 53, other R-exprs are added in (such as plus(B, C, A) above), those are covered by multiplying \(|R|\) by 2 , which gives the whole energy for the aggregator as \(\left.{ }^{(2 \cdot|R| c o r e}\right) 2\).

\section*{User-defined R-exprs}

The user-defined R-expr that guarantees termination has a stack depth limit \(K\) (introduced in section §15.3.1) that limits the number of times that a user-defined R-expr will be expanded. This limited \(K\) allows us to define the energy of userdefined \(\mathbf{R}\)-exprs as a finite number.
11. \(\left|\mathrm{f}^{k}\left(\mathrm{X}_{1}, \ldots, \mathrm{X}_{n}\right)\right|_{\text {core }}=m+1+\left\{\begin{array}{ll}\left|\mathrm{R}_{f}^{k+1}\right|_{\text {core }} & \text { if } k<K \\ 0 & \text { otherwise }\end{array} \quad\right.\) - User-defined \(\mathbf{R}\)-exprs', energy is defined both in terms of the number of free variables \(m\) and in terms of the energy of the expanded \(\mathbf{R}\)-expr \(\mathrm{R}_{f}^{k+1}\). The reason for tracking the number of free variables is the same as with structured terms, if we propagate an assignment to a variable, we need to be able to record in the energy that some useful work was done in rewriting the R-expr.

The recursive part of the energy checks the stack depth \(k\) of the user's \(\mathbf{R}\)-expr. When \(k=K\), this means that the user's \(\mathbf{R}\)-expr is expanded as far as it will be expanded. Therefore, rewrite rule 74 will no longer run. We can now break the recursive definition of energy and replace that part with the value 0 .
\({ }^{241}\) This is very loose upper bound.

\subsection*{15.3.4 Energy for Built-in R-exprs}

To define the energy for built-ins, we construct a "pluggable" definition of energy that is dependent on the rewrites provided on the built-ins.

To do this, we introduce the function \(\mathcal{E}(\cdot): \eta R \int \rightarrow \mathbb{N}\) which is defined by the library of rewrites included in our system, implying that the rewrites on builtins will also terminate. \({ }^{242}\) The function \(\mathcal{E}(\cdot)\) maps from a bag-not an R-expr-of conjunctive, non-recursive \(\mathbf{R}\)-exprs to a finite natural number that is the upper bound on the number of rewriting steps which can be performed against that bag of constraints.

The way that we use \(\mathcal{E}(\cdot)\) is that we will gather built-ins from the \(\mathbf{R}\)-expr and use \(\mathcal{E}(\cdot)\) to determine what might happen. For example, if we start with the \(\mathbf{R}\)-expr \((X=1) *\) plus \((X, 2, Z)+l e s s t h a n(Z, W)\), then it is converted into the bag
\[
\begin{equation*}
\mathcal{R}(X=1) @ 1, \operatorname{plus}(X, 2, Z) @ 1, \text { lessthan }(Z, w) @ 1 \int, \tag{15.1}
\end{equation*}
\]
and passed to the function \(\mathcal{E}(\cdot)\). In this case, the function \(\mathcal{E}(\cdot)\) returns that the energy is 3 . This follows from the longest sequence of rewrites being:
\[
\begin{gather*}
\left.\begin{array}{c}
(X=1) * \\
(X=1) * \text { plus }(X, 2, Z) * \text { lessthan }(Z, W)
\end{array} \xrightarrow{5}, 2, Z\right) * \text { lessthan }(Z, W) \xrightarrow{28}  \tag{15.2}\\
(X=1) *(Z=3) * \text { lessthan }(Z, W) \xrightarrow{5}  \tag{15.3}\\
(X=1) *(Z=3) * \text { lessthan }(3, W) . \tag{15.4}
\end{gather*}
\]

Note that \(\mathcal{E}(\cdot)\) operates on the bag of constraints, not an R-expr (which contains disjunction, conjunctions, projections, and aggregations). As such, if we have an R-expr with a disjunction like ( \(X=1\) ) +plus \((X, 2, Z)+l e s s t h a n(Z, W)\), then the bag of constraints contained in the \(\mathbf{R}\)-expr is exactly the same as equation (15.1) above.

Additionally, we define that all built-ins will have an energy of 1 under the core energy definition. E.g. \(\mid\) lessthan \(\left.(X, Y)\right|_{\text {core }}=1\).

\section*{\(\underline{\text { Why we need } \mathcal{E}(\cdot)}\)}

As stated above, we are constructing a proof of termination independent of

\footnotetext{
\({ }^{242}\) As an example if we add an external linear programming solver as a library, there would a \(\mathcal{E}(\cdot)\) function that is associated with the linear programming library.
}
the built-in rewrites included. The reason for this is two-fold. First, we do not know the "complete set" of all built-in rewrites, and second, proving termination of built-in requires theories outside of what we are proving here. Hence, we represent the number of potential steps needed to rewrite a bag of built-ins using the \(\mathcal{E}(\cdot)\) function.

To see this, consider rewrite rule 34 which combines two lessthan constraints together to infer new lessthan constraints:
\[
\begin{align*}
& \text { lessthan }(A, B) * \text { lessthan }(B, C) * \text { lessthan (C, D) } \stackrel{34}{\longrightarrow}  \tag{15.6}\\
& \binom{\text { lessthan }(A, B) * \text { lessthan }(B, C) * \text { lessthan }(C, D) *}{\text { lessthan }(A, C)} \xrightarrow{34}  \tag{15.7}\\
& \binom{\text { lessthan }(A, B) * \text { lessthan }(B, C) * \text { lessthan }(C, D) *}{\text { lessthan }(A, C) * \text { lessthan }(B, D)} \xrightarrow{34}  \tag{15.8}\\
& \binom{\text { lessthan }(A, B) * \text { lessthan }(B, C) * \text { lessthan }(C, D) *}{\text { lessthan }(A, C) * \text { lessthan }(B, D) * \text { lessthan }(A, D)} \tag{15.9}
\end{align*}
\]

Proving termination of rewrite rule 34 requires a theory about lessthan and how many lessthan constraints can be inferred. Furthermore, we need to be able to "understand" the R-expr in that the energy of equation (15.6) must be greater than the energy of equations (15.7), (15.8) and (15.9) despite the fact that equation (15.6) has the smallest R-expr representation. Hence, energy is not a purely syntactic property of the \(\mathbf{R}\)-expr. This complexity for built-ins is punted into \(\mathcal{E}(\cdot)\).

\subsection*{15.3.5 Energy for all R-exprs}

Given \(|\cdot|_{\text {core }}\) as the energy for the core \(\mathbf{R}\)-exprs and \(\mathcal{E}(\cdot)\) as the energy for built-in \(\mathbf{R}\)-exprs, we need to construct a single definition of the energy \(|\cdot|\) that can be used for all \(\mathbf{R}\)-exprs at the same time.

To do this, we first need to convert from \(\mathcal{E}(\cdot)\), which operates on bags of nonrecursive \(\mathbf{R}\)-exprs, into a representation that operates over \(\mathbf{R}\)-exprs that includes disjunctions and aggregation.

To accomplish this, observe that in constructing the energy for an R-expr, we only need some upper bound-regardless of how loose the upper bound is. Furthermore, identifying which sub-R-exprs interact with each other is just as difficult as running

Simplify itself. Hence, it would be somewhat complex to determine the interaction between built-ins. Instead, we will construct an upper bound by considering all possible interactions which might happen between built-in. This can be done by summing the \(\mathcal{E}(\cdot)\) function over all power sets of the sub-R-exprs that are in a given R-expr.

A power set \(\mathcal{P}(x)\) is defined as a set of all subsets, or in this case sub-bags of \(x\). For example if we have the bag \(\{a @ 2, b @ 1\}\), the power set is:

Observe that the power set will contain the bags that correspond to the true interaction between the built-in R-exprs. Hence, we can be sure that \(\sum_{p \in \mathcal{P}(\mathrm{R})} \mathcal{E}(p)\) is an upper bound on the energy of the built-in \(\mathbf{R}\)-exprs.

All together, we can now define the energy of an \(\mathbf{R}\)-expr \(R_{k}\) as the ordered pair \(\mathbb{E}=\mathbb{N}_{\geq 1}^{2}:\)
\[
\left.\left.\left|\mathrm{R}_{k}\right| \stackrel{\text { def }}{=}\left\langle\min _{s=0,1,2, \ldots, k}\left(\left(\sum_{p \in \mathcal{P}\left(\mathrm{R}_{s}\right)} \mathcal{E}(p)\right)-(k-s)\right),\right| \mathrm{R}_{k}\right|_{\text {core }}\right\rangle
\]
where \(\mathbb{E}\) lexicographical ordered (equation 15.10 below).
The subscript \(k\) represents the number of times that built-ins have been rewritten. In other words, \(R_{0}\) is the \(\mathbf{R}\)-expr before any rewrites have been applied, and \(R_{1}\) is the \(\mathbf{R}\)-expr after one rewrite on the built-ins, and so on up to the "current" \(\mathbf{R}\)-expr \(\mathrm{R}_{k}\). The reason for this "dependency on the previous \(\mathbf{R}\)-exprs" energy is that new \(\mathbf{R}\)-exprs can be inferred, and the size of the power set will increase-as there are more \(\mathbf{R}\)-exprs contained in the power set. By using min on all power sets up to \(\mathrm{R}_{k}\), we ensure that the first value in the pair does not increase when the size of the power set increases.

The second number in this pair is \(|R|_{\text {core }}\) as previously defined (section \(\S 15.3 .3\) ). This value will decrease for all rewrites that do not modify any of the built-in \(\mathbf{R}\)-exprs, such as rearranging the \(\mathbf{R}\)-expr or creating a more factored \(\mathbf{R}\)-expr. The value of \(|R|_{\text {core }}\) may increase when there is a rewrite performed on a built-in, as previously shown with equations (15.6) to (15.9). However, every time \(|R|_{\text {core }}\) increases, it will correspond with a decrease in the first element of this pair.

We can therefore define an ordering on these order pairs as equation (15.10):
\[
\langle a, b\rangle<\langle c, d\rangle \stackrel{\text { def }}{=} \begin{cases}a<c & \text { if } a \neq c  \tag{15.10}\\ b<d & \text { otherwise }\end{cases}
\]

It can also be seen from the construction of this ordered pair that there does not exist an infinite descending sequence, as both the first and second values in the tuple are positive natural numbers \(\mathbb{N}_{\geq 1} .{ }^{243}\)

\subsection*{15.3.6 Checking Energy of Rewrites}

All of the directional rewrites in chapter \(\S 6\) have been checked to ensure that they decrease the energy. With the few notable exceptions that were previously mentioned in section §15.3.3, the reason a rewrite decreases the energy tends to be obvious. For example, consider rewrite rule 63 if \((1+M, R, S) \xrightarrow{63} R\). The energy of \(\mid\) if \(\left.(1+M, R, S)\right|_{\text {core }}=|1+M|_{\text {core }}+|R|_{\text {core }}+|S|_{\text {core }}\), which is clearly larger than \(|R|_{\text {core }}\). Furthermore, the definition of energy has that \(\mathbf{R}\)-exprs that are higher up in the \(\mathbf{R}\)-expr (more factored), have a lower energy. This means that rewrites that make the \(\mathbf{R}\)-expr more factored will reduce the energy. This matches with the principle from section \(\S 8.2 .3\) that we will create factored \(\mathbf{R}\)-exprs to get around the issues of aggregation.

This concludes the proof that Simplify will decrease the energy of the R-expr during every step under the assumptions that the R-expr is bounded in size and the built-ins and rewrites on the built-ins terminate. Therefore, SimplifyNormalize, which repeatedly invokes Simplify will eventually reach a stopping point when Simplify has reached a fixed point since the energy cannot decrease forever.

Admittedly, this proof is not useful for proving that SimplifyNormalize terminates in useful amount of time or achieves an optimal asymptotic runtime.

\footnotetext{
\({ }^{243}\) This could also be represented as an ordinal number \(\langle a, b\rangle=\omega^{a}+b\). Given that we do not need the "full power" of ordinal numbers, we have chosen to represent this as an ordered pair of two natural numbers instead of an ordinal number.
}

\section*{Chapter 16}

\section*{Future work}

This dissertation represents my work on the Dyna programming language and the development of the R-expr-based rewrite system. Over the years, our research group has discussed a number of potential directions for which additional work is required, and I have additionally thought of a number of potential project directions on my own. I think one could easily fill ten to twenty person-years adding to the Dyna system and turning it into a more practical tool for researchers. Here are some ideas that I think are worthy of consideration as well as brief comments about how one might go about achieving these ideas.

\subsection*{16.1 Additional Disjunctive R-exprs}

As noted in section \(\S 11.6 .1\), there is currently one efficient disjunctive R-expr kind. Realistically, there should be many different of efficient disjunctive kinds for different scenarios (e.g. [120]).

\subsection*{16.1.1 Improvements to the Trie}

The current trie implementation has a few limitations which should be addressed. First, there are no secondary indexes on the trie. This means that if we order the variables as \(X\) and then \(Y\), but want to enumerate the domain of \(Y\) using an iterator, we first need to enumerate all possible values of \(X\).

Second, the order in which variables are stored in the trie should somehow be
controllable. Currently, the system just picks the variable order arbitrarily-which is clearly suboptimal. The first step here would likely be to figure out some annotation that can be specified by the user about the order and indexes that a trie should maintain. Then any information collected from the user's annotation would have to be tracked and pushed all the way into the efficient disjunct. This is a little bit harder than it might sound at first. A disjunct does not have a one-to-one correspondence with a user-defined rule. Hence it can be tricky to track which disjunct should have what policy.

\subsection*{16.1.2 Dense Numerical Types}

Given that our target audience are researchers who write numerical algorithms, we should really have some way to represent matrices and dense numerical arrays. Ideally, this would be something that would not require any changes to the user's program to use dense array types, though that might be difficult in practice.

I think the first step towards this would be to define a dense array, matrix, and tensor types as R-exprs. From there, one could look into either custom syntax or just defining built-in functions that expose operations on the dense arrays. This would be akin to building a NumPy [90] like matrix library for Dyna.
```

792 (def-base-rexpr dense-array-type [:var array-index
:var output-value
:other pointer-to-dense-array])

```

Figure 16-1. Theoretical way in which a dense array could be exposed into an R-expr.

\subsection*{16.1.3 Backed by Something other than Memory}

Currently, all disjuncts are over R-exprs and held as structures in memory. There could realistically be a disjunctive type that represents other sources of data. For example, there could be an R-expr that reads data from a CSV file or a SQL database. There are no complications here to overcome in implementing this might be a good project to give to an undergraduate who wants to get started on the Dyna project.

\subsection*{16.2 User Studies}

The work in this dissertation focused on the implementation of the Dyna programming language, with the design completed before the start of my Ph.D. back in 2011 [59]. Now that we finally have a working version of Dyna, I think prioritizing user studies would be beneficial. I think this would be an opportunity to determine which future work should be prioritized.

As an example, I note that the landscape of ML and AI research and tooling has shifted significantly from when Dyna was initially conceived-with Dyna's design predating modern neural networks by at least five years.

\subsection*{16.2.1 Libraries Written in Dyna}

Related to the changing landscape of \(\mathrm{ML} / \mathrm{AI}\) research, most researchers are leveraging a large collection of existing software when developing ML models. In particular, neural networks are virtually never written from scratch and instead take advantage of a large collection of pre-existing modules from libraries such as PyTorch or Tensorflow [4, 23, 111]. If a user were to attempt the same in Dyna today, they only have core Dyna operations, which in this case are akin to low-level tensor manipulations.

I think that developing a useful library of neural techniques or automatic differentiation-which could either be done internally on R-exprs, or externally on source level Dyna. This would serve as a useful exercise in writing Dyna programs while simultaneously resulting in something that is useful to other users.

\subsection*{16.2.2 User-Friendliness}

The current system is at the level of a research prototype in that it works, but I think the system's user-friendliness can still be greatly improved. These projects are probably at the right level to get an undergraduate started on the Dyna project.
1. Better Syntax Errors-Syntax errors currently returned are from the Antlr4 [110] parser and could be greatly improved.
2. Better Semantic Errors-The only warning currently reported to the user is
attempting to use a user-defined term without any rules that define it (which is represented as multiplicity 0). Other issues, such as type incompatibility (e.g., \(\operatorname{int}(X) *\) string \((X) \rightarrow 0\) ), should probably also be reported, but it is currently difficult to determine if type incompatibility was an error or was intentional.
3. Better Display of \(\mathbf{R}\)-exprs-There is some code for printing \(\mathbf{R}\)-exprs to the terminal. I have attempted to make the printed representation as close to the representation presented throughout this dissertation, but it is not \(100 \%\) perfect. Furthermore, I do not believe that R-exprs are the best presentation for people who are trying to use the language (rather than understand its internals). As such, I think it would be worth reconsidering how \(\mathbf{R}\)-exprs are presented to the user, possibly designing some \(\mathbf{R}\)-expr to Dyna translation so that R-exprs can be presented as Dyna code.
4. The front-end parser currently reports an error if there is more than one dynabase defined per rule. This restriction could removed by adding additional transformations in the translation of dynabases to R-exprs, such as those described in section §13.2.1.
5. Debuggability and Visualization-There is currently no support for debugging the user's program. When I encounter issues, I have to drop into the Java debugger and step through \(\mathbf{R}\)-expr rewrites-which is not a pleasant experience.

One possible approach to debugging could be to go through the steps of rewriting that are performed. There is already support in the implementation for tracking how an R-expr is constructed (when the right command line flags are passed). So, this might be building a user interface to expose those details. A second approach could be to build a tool for visualizing the values of different user-defined rules in the program. This was previously done with Dyna 1.0 with the derivative Dynasty project [62].
6. Visualization and Interactions in Jupyter notebooks [99]-I have defined a \%dyna Jupyter notebook cell/line magic \({ }^{244}\), which allows for running simple Dyna programs from Jupyter notebooks. This user experience could also be improved with better visualizations.
\({ }^{244}\) https://ipython.readthedocs.io/en/stable/config/custommagics.html

\subsection*{16.3 More Rewrite Rules}

Our implementation currently contains a number of rewrite rules. However, this is by no means expected to be all possible rewrites. It is reasonable that one could invent new rewrites which can efficiently handle different scenarios. There could also be more rewrite rules that correspond to different mathematical operations. For example, there could be rewrite rules involving identities of trigonometric functions.

\subsection*{16.4 Automatic Configuration}

Currently, there are a number of different "decisions" that need to be made to configure how the program runs and can have a significant impact on the program's runtime. For example, we already have seen memoization, which is controllable by \$memo and \$priority. However, other things, such as variable order and indexing on disjunctions and variable loop order, should be considered.

Some of these proposed automatic configurations are akin to database query optimizers. However, controls like \$memo are much more complicated, as a bad policy can cause the system to become inoperable and have consequences outside of the scope of a single query (section §10.6.2).

I should also note that the automatic algorithm configuration was the central topic of Tim Vieria's dissertation [141], in that he explored different ways a program could be folded to make it more efficient (chapter §14). Many of the ideas in Tim’s dissertation should be adaptable to the system presented here.

\subsection*{16.4.1 Automatic Guessing on Cycles}

In the current implementation, all memoization decisions are manual. However, there are a few cases that could be automatically detected and would likely improve the user experience. One such example is the case of a cyclic program. The system will detect that a cycle happens and eventually stop expanding the \(\mathbf{R}\)-expr (section §15.3.1). However, it does not currently enable memoization automatically.

In the case that a cycle is detected that does not contain memoization, then
memoization could be automatically enabled, or a helpful error message could be reported to the user about how they can fix their program.

\subsection*{16.4.2 Automatic Prioritization of Updates}

Prioritization of updates is currently entirely dependent on the user's declarations. I believe there is an "easy to exploit" opportunity here to make this process automatic. The reason is that a bad update prioritization should not cause the system to accidentally not-terminate in most cases (like in the case of a bad \$memo). Hence, it is "mostly safe" to experiment with different prioritization when the system is running.

I believe that a possible approach would be to run tasks and track when "priority inversions" happens. Essentially, the system could detect priority inversions by tracking when a downstream value was previously computed has to be recomputed. A machine learning regression model could be trained to predict numerical priorities such that no priority inversions will happen. In the case that the program does not contain a cycle, then this would be akin to using features on the nodes in a graph to learn a topological ordering. Additional handling might be required in the case of cycles, though I believe that it might be possible to set up the machine learning algorithm such that cycles do not change the learned weights of the machine learning algorithm by having the notifications of priority inversion cancel each other out.

\subsection*{16.4.3 Automatic Variable Ordering}

The variable order used by disjunctive data structures, such as the trie, can have an impact on the efficiency of the system. A bad variable ordering should not cause any issues around non-termination, so it should be safe to explore different representations.

I am not sure if there is an "efficient" way in which runtime costs can be associated with a variable ordering. It is possible that the variable orderings up being zero-sum, in that improving the runtime of one operation could negatively impact the runtime of something else. I believe that doing this correctly will require a reinforcement learning approach, where different orders are experimented with
until something satisfactory is found [142].

\subsection*{16.5 Concurrency}

Currently, the system is single-threaded. I have tried to design the internal data structures with the intention of eventually being used in a concurrent, multithreaded environment, so hopefully, adding in concurrency is not too much of a chore. I think that the best approach would be to leave the rewriting done by Simplify and SimplifyNormalize as single-threaded and instead focus on doing parallel processing of the pending work on the UpdateQueue (section §10.7). The system could reasonably process updates in parallel, tracking if a conflict occurs much like a database. The update operations are not externally observable, so simple retry logic should be sufficient in most cases.

\subsection*{16.6 The Memoization Update Queue}

The UpdateQueve used in the implementation of memoization (section §10.7) is a priority queue that is controlled by \$priority (•) (section §10.8.5). Hence, this requires \(O(\log N)\) time to push and pop update operations from this queue, where \(N\) is the number of update messages on the queue. Comparably, a dynamic program with a predetermined execution order does not require a queue and will avoid this \(O(\log N)\) operations. When problems get sufficiently large, this extra \(O(\log N)\) overhead might become significant, so figuring out how to eliminate this overhead may be a good idea.

\subsection*{16.7 GPU Coprocessor}

GPUs are very important to modern ML algorithms and are \(100 \%\) essential to the implementation of large Neural Networks. I think that there are a few potential ways that GPU support could be added:
1. Expose existing GPU Kernels-There are a number of different GPU kernels that currently exist. Using existing kernels usually results in the best performance.

This could either be done by defining rules that match with the existing GPU kernels and calling those operations when the program matches an existing pattern. Alternately, one could simply define built-in operations for every kernel and require that the user manually reference the GPU kernel operations themselves.
2. Compiling Dyna to custom GPU Kernels-It should be possible to compile Dyna programs to a GPU. This might want to build on the work done for the JIT compiler (chapter §12), as it is already attempting to generate nested loops over variables' domains. Some challenges with this approach would be that the system still allows for unpredictable \(\mathbf{R}\)-exprs to be returned at various points while executing. This would mean that the system would either have to require some limits on what can be represented or somehow support R-exprs on the GPU.
3. GPU DSL-Dyna has support for implementing DSL in the language (section §2.10), one could create a GPU DSL which compiles into a custom R-expr type and defines its own rewrites so that it is integrated with the rest of the system.
4. Develop an explicit matrix type and library of matrix operations that use a GPU. This would be an extension of the dense numerical types proposed previously (section §16.1.2).

\subsection*{16.8 External Solvers}

Although we can continue to add many features to the \(\mathbf{R}\)-expr system to expand the ability of the Dyna programming language, we should also consider that there exist a number of powerful frameworks and software tools that might be worth leveraging in the Dyna project. For example, SMT solvers are able to efficiently solve SAT-like problems that require searching through possible assignments [47, 48, 113]. Being specially designed for a "more limited" problem formalism than Dyna, their implementation is much more specialized than we would be able to realistically emulate with R-exprs. Additionally, there are algebra systems designed to solve formulas, such as Mathematica and SymPy, both of which have a very
large collection of algebra rules [103, 150]. \({ }^{245}\)
Integrating external solvers into Dyna could be integrated as an external library (as suggested in section \(\S 2.10 .1\) with an example of a linear programming module) or as custom rewrite rules against R-exprs.

\subsection*{16.9 Advanced Update Propagation}

The updates of memoized values are currently handled by recomputation of the original \(\mathbf{R}\)-expr. While this is sufficient for the small programs with which we are experimenting with currently, in the future, it might be beneficial if updates could be aware of the values and kinds of updates that are being propagated through the system.

For example, suppose that we have that \(a(I, J)\) is representing a matrix, and there is a rank- 1 update to the matrix. A rank- 1 update could potentially modify all entries in the matrix \(a(I, J)\). However, knowing that \(a(I, J)\) received a rank1 update could allow for more asymptotically efficient computation to be used downstream.

\subsection*{16.10 Automatic Runtime Analysis and Folding of Programs}

This was part of Tim Vieira's dissertation [141]. However, that work was not done using the R-expr formalism, and the work to integrate these two bodies of work has not been started. Chapter \(\S 14\) discussed how folding can be done on R-exprs, which is critical for adapting Tim's dissertation work.

\subsection*{16.11 Improved Context \(\mathcal{C}\)}

The context \(\mathcal{C}\) currently uses a hash-map to track the current value assigned to variables. This means that every time that matcher preconditions such as :ground or : free are evaluated, or when the get-value function is called, the system

\footnotetext{
\({ }^{245}\) https://www.wolfram.com/engine/
}
consult the hash-map (section §11.5.3). Hence, the context and the hash-maps that are used in its implementation have the potential to become a major bottleneck. Currently, the hash-map used is the default Clojure hash-map. It is possible that simply replacing the hash-map with another hash-map implementation that is more efficient could improve the runtime.

As a more long term solution, redesigning the get-value-of-variable mechanism such that there are as few memory access operations as possible and no hard-topredict branches would probably be most beneficial. One possible solution to this might be to make a new class that implements the value type interface, which supports faster operations than the standard named variable. The context could then contain an array in addition to the hash-map that would enable retrieving the relevant value in a single memory access.

\section*{Chapter 17}

\section*{Conclusion}

This dissertation documents my work in implementing the Dyna language using R-expr based term rewriting. For the first time in the Dyna 2.0 project, we have an operational semantics for the language, an approach to implement the language, and an implementation at the level of the research prototype. A foundation has been laid for future work on the Dyna project. Although the material in this dissertation does work, I think there is still a lot of work to do to make it usable. I believe this will easily occupy the next several years of the Dyna project, and I hope to see this work continue, as I believe there is great promise.

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[^0]:    ${ }^{1}$ This is still true of large neural networks. However, neural networks research often builds on top of a large library such as PyTorch [111] or TensorFlow [4], which contain many existing neural net modules. Conceptually, similar software libraries could be implemented in Dyna.

[^1]:    ${ }^{2}$ Datalog will be explained further in section §3.1.2. In short, Datalog can be thought of as a programming language that operates over boolean values as a dynamic program. All statements in a Datalog program are materialized in database tables, which are queried (and joined) with a Prolog-like syntax.
    ${ }^{3}$ A semiring is an algebraic structure that has an additive and multiplicative operator, additive and multiplicative identity elements. Semirings do not have an additive inverse, like rings. For example, logic programming operates over the boolean semiring, 〈or, and, false, true $\rangle$. Dyna 1.0 semirings could include the real numbers $\langle+, *, 0,1\rangle$, or the min value (shortest path) $\langle\mathrm{min},+, \infty, 0\rangle$, but all rules had to use the same semiring.
    ${ }^{4}$ Ground terms means that all of the values in the expression are known. For example, the term foo( 1,2, "hello") is ground, as both integers 1 and 2 and the string "hello" are ground.
    ${ }^{5}$ A term with a free variable would be like $\operatorname{bar}(7, X)$ where the variable $X$ can represent any value.

[^2]:    ${ }^{6}$ The Dyna project has had many implementations and different dialects over the years. Details of these alternate implementations can be found on the project's website http://dyna.org.

[^3]:    ${ }^{7}$ At least more modern than when Logic Programming and Prolog first appeared in 1972 [38].

[^4]:    ${ }^{8}$ Modulo syntactic differences. A Prolog/Datalog program would require some changes to match Dyna's syntax exactly. The internal representation of Dyna, described in chapter $\S 5$ is capable of representing Prolog/Datalog programs.
    ${ }^{9}$ A bag can be thought of as a generalization of a set where an element can appear multiple times. A bag will be defined properly in chapter $\S 5$.

[^5]:    ${ }^{10}$ In the nomenclature of a SQL database, the declaration of a fact is equivalent to inserting a tuple into the database. E.g.: INSERT INTO parent VALUES ('charles', 'james');
    ${ }^{11}$ In SQL, a rule definition would be equivalent to defining a view that will be computed on demand from the defined facts. E.g.: CREATE VIEW grandparent AS SELECT a.parent, b.child FROM parent a INNER JOIN parent b ON a.child = b.parent;

[^6]:    ${ }^{12}$ Dyna also allows for an ampersand (\&) to be used as a "quote" on a term as an alternate way of writing this expression. E.g. \&termName (Argument1, Argument2, ...,ArgumentN) $\equiv$ termName[Argument1,Argument2, ...,ArgumentN].

[^7]:    ${ }^{13}$ The 'for' keyword will return the first sub-expression's value with the second sub-expression representing a side condition that must return the value true. The comma ',' is conceptually equivalent to 'for', but it returns the second sub-expression's value and the first sub-expression must return true. In other words, the expression 'A for $B$ ' is equivalent to the expression ' $B$, A'.
    ${ }^{14}$ Future work may consider adding in syntax for other types, such as sets. However, sets can be emulated using a map by using only the Key and setting a dummy value for all pairs, like: \{"X"->0, "Y"->0\}.
    ${ }^{15}$ The notation ${ }^{-}>$' was chosen to avoid conflicting with other syntax in the Dyna language.

[^8]:    ${ }^{16}$ Prolog and Datalog are "allowed" to stop early, but the exact circumstances under which this happens can differ. See the discussion in the related work chapter for more information. §3

[^9]:    ${ }^{17}$ The ' $:=$ ' aggregator accepts a special value of $\$$ null that can be returned by an expression to make the result of ': =' null, overriding other non-null contributions from earlier lines.
    ${ }^{18}$ Dyna is allowed to stop computation for a term's value early once it has determined that the value will not change further. This happens in the case of $\mid=, \&=,:-$ which have saturating values of true, false, true. Aggregators like := can stop the computation of earlier lines in the program if there is a value contributed from a line that occurs later. Similarly, ?= can stop all other computations once it has found something.
    ${ }^{19}$ Each term in the program can only use a single aggregator otherwise the value is defined to be an error. However, rules that define terms that share the same functor name can be defined using different aggregators. For example, we can write ' $a(X)+=2$ for $g(X)$.' and ' $a(X) \quad *=7$ for $q(X)$.' provided that $g(X)$ and $q(X)$ are non-overlapping.
    ${ }^{20}$ Previous Dyna papers have used the keyword with_key for this instead of the keyword arg. The keyword with_key is still supported by the Dyna front-end parser.
    ${ }^{21}$ Internally, Dyna tracks this using a pair of values, which includes the argument and the value returned. The Dyna implementation automatically references the value field by default or will reference the argument field using the macro $\$ \arg (\cdot)$ which suppresses the automatic access of the value field.

[^10]:    ${ }^{22} \mathrm{~A}$ Dyna program is declaritive in that defined rules can be rearranged and evaluated out of order. A Dyna script may contain procedural statements such as assert and print, which evaluate an expression with respect to the rules defined before the statement.
    ${ }^{23}$ There is a slight technicality here between "having" a value and "returning" a value. In Dyna, the term is an identifier, and the Dyna program is the function that takes the term as an argument and returns a value. In other words, in a functional language, you might have "return_value = named_function(Arg1, Arg2)" whereas in Dyna we have "return_value = dyna_program_function(named_term_identifier[Arg1, Arg2])" with the program itself being the function.

[^11]:    ${ }^{24}$ The $\$=$ is necessary in Prolog to represent a floating point addition constraint between the floats SS, CC, and the Result variable. If this was written with an equals sign only and without the dollar sign, it would end up assigning the term ' + ' $[S S, C C]$ (not the numerical value) to the variable Result.
    Alternately, Prolog has is which can be used to evaluate numerical expressions and could be used as Result is SS + CC, but is will only evaluate the expression on the right-hand side and assign it to the variable on the left-hand side. In other words, in Prolog the following is expression would result in an error: 5 is $2+X$.
    ${ }^{25}$ Prior work did include the idea that there would be a driver program querying and updating the Dyna program [61]. However, the design of the interface detailed here is a contribution of this dissertation.

[^12]:    ${ }^{26}$ Externally defined functions must have unique names and cannot be combined with terms that are defined by the Dyna program.

[^13]:    ${ }^{27}$ This is a contribution of this dissertation.

[^14]:    ${ }^{28}$ Business logic sometimes called domain-specific logic is the part of the program which encodes rules which are relevant to the real world problem being solved.

[^15]:    ${ }^{29}$ Note: the program on lines 158 to 161 will not execute under Prolog, as Prolog will get stuck performing a depth-first search on lines 158 and 159.

[^16]:    ${ }^{30}$ This is the halting problem, which states that there exists a program for which we can not prove or disprove that it terminates. This property is true for all Turing complete languages.

[^17]:    ${ }^{31}$ When there are multiple solutions, Dyna is only required to find a solution. We make no guarantees about which solution is returned. This is in contrast to Datalog, which guarantees that the minimal solution is returned (section §3.1.2), and answer set programming, which returns all solutions (section §3.1.6).

[^18]:    ${ }^{32}$ This example, and specifically line 173 , is discussed further in section $\S 4.2 .1$.

[^19]:    ${ }^{33}$ The \$memo control mechanism for memoization is a contribution of this dissertation.

[^20]:    ${ }^{34}$ The system internally uses the priority of $10^{16}$ for internal work, which should be run as soon as possible, and $-10^{16}$ for user updates when no \$priority is set. The system does not prevent you from setting a priority higher or lower than $\pm 10^{16}$, but be aware of these internal values as going higher or lower than $\pm 10^{16}$ can have unforeseen consequences.
    ${ }^{35}$ Future work may wish to add the ability for \$priority to handle updates.

[^21]:    ${ }^{36}$ The terminology of "null" vs. "unk" memos was adopted from Filardo and Eisner [67]-an earlier research paper published by our group.
    ${ }^{37}$ An "unk" memo is shorthand for specifying the case where all arguments must be ground. For example, the following policies are equivalent: \$memo(foo[X, Y: \$ground, Z: \$free]) = "unk". \$memo(foo[X:\$ground, $Y$ : \$ground, Z:\$ground]) = "null". This is discussed in greater detail in section §10.8.3.

[^22]:    ${ }^{38}$ Admittedly, Prolog included call/N, which supports higher-order functions, which have existed since early version of Prolog. However, earlier versions and proposals of Dyna were Datalog-inspired and did not include higher-order functions. Hence, the contribution of higher-order functions is specifically to Dyna and its syntax and not a contribution to logic programming in general.
    ${ }^{39}$ The modern PL constructions implementation and their syntax are all contributions of this dissertation work to the Dyna language.
    ${ }^{40}$ The indirect call can also use a expression, such as a call to another user-defined term. In this case, extra parenthesis are needed around the expression returning the function pointer. E.g. (foo $(1,2))(4,5)$

[^23]:    ${ }^{41}$ This is equivalent to defining a new int_or_string $(X)$ term using two rules to create a union type:
    int_or_string(X) :- int(X).
    int_or_string $(X)$ :- string $(X)$.

[^24]:    ${ }^{42}$ Warnings for dead code can only happen if Dyna can identify the dead code. In general, this is tricky as identifying dead code is equivalent to just running the program. Furthermore, some code might only be temporarily dead and become used when other rules are added to the program.
    ${ }^{43}$ There is currently disagreement among the Dyna team about the design of dynabases. It is possible that a future implementation of Dyna will have dynabases that behave differently. The design presented here is my own design and corresponds with how dynabases work in my implementation.

[^25]:    ${ }^{44} \$ \mathbf{s e l f}$ 's behavior is similar to that of self in Python. \$self is not required on the functor name of defined rules (bar on line 228), but it is required inside of the expression that defines the rule when referencing other terms defined on the dynabase.

[^26]:    ${ }^{45}$ There are have other proposals from Dyna team about how DSLs could be implemented in Dyna. The design here is my own design and a contribution of this dissertation.

[^27]:    ${ }^{46}$ It is helpful to observe the Dyna AST when developing macros by running in the REPL: print \$ast'\{ my_program = 1. \}.
    ${ }^{47}$ The macro auto_differentiate is not implemented at this time. I suggest in section §16.2.1 that future work should consider implementing libraries like this in Dyna.

[^28]:    ${ }^{48}$ Prolog systems will usually refer to rules using their name and arity (number of arguments). In this case, a/1 refers to the 'a' rule defined on lines 285 to 294.

[^29]:    ${ }^{49}$ I should note that unification is usually presented as working with structured terms (as in section §2.1.1) instead of delayed constraints on numerical values.

[^30]:    ${ }^{50}$ The notation $Y=g(X, X)$ is how Prolog does structural terms (section §2.1.1). Prolog does not have automatic evaluation, as we have in Dyna. Hence, Prolog uses parenthesis () for both calls and unification with structures.

[^31]:    ${ }^{51}$ See footnote 3 for an explanation Semirings.

[^32]:    ${ }^{52}$ This means that there are no infinite relations like in Prolog.

[^33]:    53 Documentation about Prolog's meta predicates that enable aggregation: https: //www.eclipseclp.org/doc/bips/kernel/allsols/bagof-3.html https://www. swi-prolog. org/pldoc/man?predicate=bagof/3 https://eclipseclp.org/doc/bips/lib/fd_global/ sumlist-2.html

    Documentation for aggregation in Datalog: https://souffle-lang.github.io/aggregates

[^34]:    ${ }^{54}$ On the contrary, in Prolog the unifications between variables are tracked globally and maintained using stack discipline when Prolog backtracks in the case of failure.
    ${ }^{55}$ When the expanded version of the program is greater than the max depth needed to represent all function calls in the program, then the solver is able to represent all intermediate values that the program would compute inside of the solver.
    ${ }^{56}$ As approaches that are built on SAT solving can take an exponential amount of time to solve.

[^35]:    ${ }^{57}$ It is conceptually possible that Dyna could add these features in the future as DSL in Dyna (e.g. section §2.10).
    ${ }^{58}$ SQL: Structured Query Language

[^36]:    ${ }^{59}$ Dyna can also reorder operations, like a database (section §2.4). At this time, Dyna does not include an optimizer that finds the best way to arrange computation, but this is proposed as future work (section §16.4).

[^37]:    ${ }^{60}$ Their "variable free" representation the integer index in a tuple to replace the variable name. We choose to use a name-based representation as we believe it is cleaner, section §5.1.1.
    ${ }^{61}$ The version of Dyna in this dissertation supports higher-order functions, lambdas, closures, lazy evaluation, immutable data structures such as hash maps, and the standard logic programming lists and structured terms. User-defined rules are side-effect-free and have a functional dependency between the arguments to the rule and the value returned from the aggregator.

[^38]:    ${ }^{62}$ Also called residuation in Antoy [6].

[^39]:    ${ }^{63}$ Prolog has also been compiled using method-based compilation with research projects such as the YAP compiler [43] and the Mercury project [39].

[^40]:    ${ }^{64}$ In a dynamic language like Javascript or Python, the trace may include a check that the MrPRINT function was not redefined.

[^41]:    ${ }^{65}$ There have been some syntactic changes to Dyna since 2011. None of those changes has made the language simpler.

[^42]:    ${ }^{66} \mathrm{~A}$ footgun is a bad feature which is easy to misuse and very likely to give surprising results in "common" use cases.

[^43]:    ${ }^{67}$ One approach for solving the shortest path would be to wait until the value of Start is known, and then start forward chaining the values for the paths. This corresponds to using the memoization policy of
    \$memo(path[Start:\$ground, X:\$free]) = "null".
    We can further implement Dijkstra's by ordering the shortest distances first using
    \$priority(path[Start,X]) = -path(Start,X).

[^44]:    ${ }^{68}$ Min multiplicity is sometimes used by other bag algebras for intersection.
    ${ }^{69}$ For example, in set builder notation, the set of even integers could be written as $\{x \in \mathbb{Z}: \exists y \in$ $\mathbb{Z}, y * 2=x\}$, where the boolean predicate $\exists y \in \mathbb{Z}, y * 2=x$ is only true for even integers. Note, when creating sets using an enumerable expression (iterable) in a program, this is called set comprehension. However, in mathematics (as we are doing here) this is called set-builder notation. The predicate that defines the set is not required to be computable. For example, we can use set-builder notation to define the set of all people you have met and will ever meet in your life. This set is well defined, but not something that we can express or construct in an executable program without also inventing a time machine.

[^45]:    ${ }^{70}$ Even if represented the set of integers between $\left[0,10^{100}\right.$ ) as a bag of $10^{100}$ integers, we still consider the bag of $10^{100}$ integers as simpler despite its enormous size, as it requires less "computation". This is explained in detail in section §15.3.

[^46]:    ${ }^{71}$ For this chapter, $E(\cdot)$ will only be used when it contains the necessary variables. Hence, we do not define what happens if a variable is not contained in $E(\cdot)$. In chapter $\S 8$, when discussing the implementation of $\mathbf{R}$-exprs and the rewrite rules, I will discuss what happens when a variable is not contained in $E(\cdot)$.

[^47]:    ${ }^{72}$ Note that the variables and ground terms are disjoint $(\mathcal{G} \cap \tilde{\mathcal{V}}=\emptyset)$.

[^48]:    ${ }^{73} \llbracket \cdot \rrbracket_{E}$ is undefined in the case that $\operatorname{vars}(\mathrm{R}) \nsubseteq$ domain $(E)$.
    ${ }^{74}$ Recall that a multiplicity is defined as $\mathcal{M}=\mathbb{N} \cup\{\infty\}$. Hence, 0 and 1 are the identity and annihilator elements of the multiplicities.

[^49]:    ${ }^{75}$ This is allowed as this does not change the multiplicity of the R-expr. e.g. $1 * 1 * 1=1$ or $0 * 0 * 0=0$
    ${ }^{76}$ If $R$ and $S$ are both constraints and non-overlapping, then the disjunction $R+S$ is also a constraint.

[^50]:    ${ }^{77}$ This fact will be used for some advanced rewrite rules discussed in chapter $\S 9$.
    ${ }^{78}$ If X is already contained in $E(\cdot)$, then it will be overridden, otherwise it is added as a new variable.
    ${ }^{79}$ Recall that $\mathcal{V}=\tilde{\mathcal{V}} \cup \mathcal{G}$, hence it is possible that $\mathrm{X} \in \mathcal{G}$ and a ground value. This case is "allowed" though not particularly interesting. We have $\llbracket \operatorname{proj}(\mathrm{g}, \mathrm{R}) \rrbracket_{E}=\llbracket \mathrm{R} \rrbracket_{E}$ as there is only one value $g \in \mathcal{G}$ that is equal to $g$.
    ${ }^{80}$ Usually $\llbracket \mathrm{R} \rrbracket_{E}[\mathrm{X}=x]$ will have a non-zero multiplicity for a finite number of values of $x \in \mathcal{G}$. Hence, this sum over $\mathcal{G}$ (which is infinite) can be computed by identifying the relevant subset of $\mathcal{G}$. This will be discussed extensively in the following chapters.
    ${ }^{81}$ For example, $\llbracket \operatorname{proj}(\mathrm{X}, 1) \rrbracket_{E}=\infty$, as this is equivalent to $\sum_{x \in \mathcal{G}} 1$, where $|\mathcal{G}|=\infty$.

[^51]:    ${ }^{82}$ Some aggregators such as min and max do not care about the number of copies, as long as it is more than one. Aggregators such as sum can sum an infinite number of zeros and still have the value zero, e.g. $0=\sum_{i=0}^{\infty} 0$.
    ${ }^{83}$ For example, the sum $\sum_{i=0}^{\infty}(-1)^{i}$ does not converge and therefore not well defined.

[^52]:    ${ }^{84}$ This is equivalent to $\alpha$-renaming in $\lambda$-calculus.
    ${ }^{85}$ This is similar to a let rec construction in a functional programming language.
    ${ }^{86}$ Dyna does not guarantee which assignment will be found. This is in contrast to other logic programming languages, such as Datalog (section §3.1.2), that guarantee a minimal assignment is found.

[^53]:    ${ }^{87}$ Some PDF viewers have a link preview feature when hovering over links. I strongly suggest that you turn that feature on.

[^54]:    ${ }^{88}$ We do not care which $\prec$ function is used, as long as it is consistent. An example would be variable string name comparisons.

[^55]:    ${ }^{89}$ This would have been equivalent to writing $f[X, Y, Z]=g[Q, H]$, and there is no assignment to $X, Y, Z, Q, H$ that makes this expression true.

[^56]:    ${ }^{90}$ We will see R-exprs represented as tries again in section §11.6.1.

[^57]:    ${ }^{91}$ Other non-recursive base-case ingredients of $\mathbf{R}$-exprs are structural equality constraints and user-defined $\mathbf{R}$-exprs for user-defined terms.

[^58]:    ${ }^{92}$ For example, with ( $I=2$ ) *R the rewrite sequence is as follows:
    $(I=2) \star \operatorname{proj}(J, \operatorname{plus}(I, 3, J) \star \operatorname{plus}(J, 4, K)) \xrightarrow{5}(I=2) \star \operatorname{proj}(J, \operatorname{plus}(2,3, J) \star p l u s(J, 4, K))$
    (Rewrite rule 5, equality propagation)
    $\xrightarrow{28}(\mathrm{I}=2) \star \operatorname{proj}(\mathrm{J},(\mathrm{J}=5) \star \operatorname{plus}(\mathrm{J}, 4, \mathrm{~K})) \quad$ (Rewrite rule 28 applied to pl us $(2,3, \mathrm{~J})$ )
    $\xrightarrow{5}(\mathrm{I}=2) * \operatorname{proj}(\mathrm{~J},(\mathrm{~J}=5) * \mathrm{plus}(5,4, \mathrm{~K})) \quad$ (Rewrite rule 5, equality propagation)
    $\xrightarrow{28}(\mathrm{I}=2) * \operatorname{proj}(\mathrm{~J},(\mathrm{~J}=5) *(\mathrm{~K}=9)) \quad$ (Rewrite rule 28 applied to $\mathrm{plus}(5,4, \mathrm{~K})$ )
    $\xrightarrow{39,41,13}(\mathrm{I}=2) *(\mathrm{~K}=9) \quad$ (Rewrite rules 13,39 and 41 used to eliminate projection)

[^59]:    ${ }^{93}$ The variable names are generated so to not conflict with vars $(R)$ or vars $(S)$. We are not literally using the variable names $B$ and $C$. Variables that appear on the right-hand side of a rewrite and inside a projection, such as B and C, are only chosen for the visual presentation of the rewrite, such as in the case of rewrite rule 50 .

[^60]:    ${ }^{94}$ The complexity here is fundamental and must be handled somehow. Essentially, we need to disable rewrite rule 48 so that when one disjunctive branch is rewritten as 0 , it does not cause the entire $\mathbf{R}$-expr to be rewritten as 0 (e.g. $A=\operatorname{sum}(X, 0) \rightarrow 0)$. Some alternative approaches to solving this problem could be to introduce a second aggregator $\mathbf{R}$-expr kind, like $A=s u m \_o f \_d i s j u n c t(X, R)$ or we could introduce a boolean flag onto the aggregator $\mathbf{R}$-expr kind to track if the ( $\mathrm{X}=$ agg_null) disjunct is present. In fact, in chapter $\S 11$, when discussing a realistic implementation of $\mathbf{R}$-exprs, the aggregation kind will use a boolean flag instead of the disjunction trick here. I have chosen to present the disjunction issue using the ( $X=a g g \_n u l l$ ) approach because I believe this presentation most clearly demonstrates the semantics of aggregation while still representing the issue of aggregating over a disjunction.

[^61]:    ${ }^{95}$ The rewrite rules for not_equal can be defined as: not_equal $(A, B) \rightarrow 1$ if $A, B \in \mathcal{G}$ and $A \neq B$ not_equal $(A, A) \rightarrow 0$

[^62]:    ${ }^{96}$ Admittedly, $A=\operatorname{sum}(X, 2 *(X=7))$ can be handled using rewrite rule 23 to first expand two copies of $(X=7)$ as $A=\operatorname{sum}(X,(X=7)+(X=7))$ and then use rewrite rule 50 to sum over the two disjuncts, however this will be less efficient than using rewrite rule 55 . For example, suppose that the value of $M$ is several hundred instead of 2 . In this case, using rewrite rules 50 and 53 would create a very large R-expr.

[^63]:    ${ }^{97}$ The condition $\forall_{E} \llbracket Q * Q 2 \rrbracket_{E}=0$ can be proven by checking if there exists a sequence of rewrites such that $\mathrm{Q} * \mathrm{Q} 2 \rightarrow^{*} 0$, meaning that regardless of the environment, the expression represents an empty bag relation.

[^64]:    ${ }^{98}$ Note that is standard in Prolog to represent functions as a constraint with the return value as the final argument. For example, append ([1,2], [3], $[1,2,3]$ ). This is conceptually equivalent to what we are doing here.

[^65]:    ${ }^{99}$ In other words, if we have an user-defined $\mathbf{R}$-expr kind like $a(X)$ with a rewrite rule like $a(X) \rightarrow a(X)+a(X)$, this will represent an infinitely large $\mathbf{R}$-expr, due to the recursive expansion. We will only have expanded this $\mathbf{R}$-expr up to some limited depth at any given moment. We can continue to perform more and more rewrites on this $\mathbf{R}$-expr to continue expanding it without end. As such, in the limit it is infinitely size, but for at any given time $t<\infty$ the $\mathbf{R}$-expr will be finite.
    ${ }^{100}$ Variable identifiers are allowed to be any Object that supports hashing and equality checks. In practice, the variable identifier is often a string.

[^66]:    ${ }^{101}$ Pseudocode for Simplify in section $\S 8 . A$
    ${ }^{102}$ Starved meaning that a rewrite is prevented from running because it does not get the "necessary resources" to run. In this case, starved resources mean that the rewrite was not matched (when it could have been) or that it was always not picked to run for some reason.
    ${ }^{103}$ There are bidirectional rewrites which can be used for rearranging the R-expr. Those will be addressed shortly in section §8.2.2.1.
    ${ }^{104} \mathrm{We}$ are done rewriting when there is nothing more to be done. We can identify this case by checking the argument to Simplify with its returned $\mathbf{R}$-expr: $\operatorname{Simplify}(R, \mathcal{C})==R$.
    ${ }^{105}$ A given $\mathbf{R}$-expr (with recursion represented as user-defined $\mathbf{R}$-exprs) is bounded in size, hence scanning through the entire $\mathbf{R}$-expr is guaranteed to be a bounded time operation.

[^67]:    ${ }^{106}$ SimplifyNormalize's non-termination corresponds with the fact that Dyna, and $\mathbf{R}$-expr rewriting, is Turing complete. Hence, programs that contain infinite loops, or rewrite forever without cycles, do not terminate. (section §15.2)
    ${ }^{107}$ SIMPLIFY is allowed to perform more than one rewrite during a single pass of rewriting.

[^68]:    ${ }^{108}$ This includes rewrite rules 18 to 21 as well as rewrites for permeability of projection and aggregation, rewrite rules 39 and 54 .

[^69]:    ${ }^{109}$ Our rewrites are semantics preserving, so a rewrite $R * S 1 \rightarrow R * S 2$ requires that $R * S 1$ is semantically equivalent to $R * S 2$, however it does not require that $S 1$ is semantically equivalent to S2.
    ${ }^{110}$ The context is a set, not a bag (like $\mathbf{R}$-exprs)-it is allowed to be implemented as a bag, but being a bag does not change the behavior of the context. The reason is that the context is used to identify if a constraint is conjunctive with the $\mathbf{R}$-expr currently being rewritten. We are allowed to duplicate constraints (section $\S 5.2 .2 .4$, rewrite rule 24), and no rewrites are performed on the context itself.
    ${ }^{111}$ The context can be seen as similar to the constraint store in constraint logic programming (section §3.1.5).
    ${ }^{112}$ Shallow copies of the context mean that internal structure is shared as much as possible. The internal data structures used by the context are immutable; hence, they can be shared between copies of the context.
    ${ }^{113}$ Prolog-based systems use an undo list rather than copies. Both making copies and undo lists solve the same problem. In my opinion, neither approach is strictly better than the other, as each has its own advantages and disadvantages.

[^70]:    ${ }^{114}$ In the visual presentation in this dissertation, the R-exprs which appear to the left are evaluated before the ones on the right. Hence, we can think of Simplify as evaluating in a left-to-right order.
    ${ }^{115}$ An oracle using only equality constraints and bidirectional rewrites on conjuncts and disjunctions is NP-hard in that it can be used to solve SAT formula. Therefore, it is unrealistic to require that we rearrange all R-exprs into an ideal form.

[^71]:    ${ }^{116}$ As an interesting side note, observe that the expansion in figure 8-6 is conceptually equivalent to what Prolog (and constraint logic programming built on Prolog) does. The reason is that each line of figure $8-6$ represents a conjunction of constraints. This would be that each conjunction is generated by Prolog when expanding the program using backward chaining (section §3.1.1).

[^72]:    ${ }^{117}$ By smaller bag, I mean that there are fewer elements in the bag. For example, the bag $\langle\langle X=$ $1\rangle @ 1 \int$ is smaller than the bag $\left\langle\langle X=1\rangle @ 2,\langle X=7\rangle @ 3 \int\right.$. A smaller bag can have more constraints and more conjunctive $\mathbf{R}$-exprs.

[^73]:    ${ }^{118}$ Note: The DPLL algorithm from [45] is no longer considered the state-of-the-art approach for solving SAT problems. However, current state-of-the-art methods for solving SAT (such as conflict driven clause learning, CDCL) still use DPLL at their "core".
    ${ }^{119}$ Branching is allowed to select any unassigned variable. The order of variables that are selected can have a big impact on the wall clock runtime of the DPLL algorithm, and there is a lot of research focused on developing better branching heuristics.

[^74]:    ${ }^{120}$ The SAT formula at this point is equivalent to $(A \vee B) \wedge(\bar{A} \vee \bar{B}) \wedge(A \vee \bar{B})$ which does not contain any unit clauses that can be used for propagation.

[^75]:    ${ }^{121}$ Admittedly, this is the distributive rewrite which, as already discussed in section §8.2.3.1, is used to factor the $\mathbf{R}$-expr rather than expanding out the $\mathbf{R}$-expr as done here. We are going to have a solution for this in section §11.7.

[^76]:    ${ }^{122}$ For example, $\operatorname{proj}(X, \operatorname{int}(X) * \operatorname{times}(3, X, Y)) *(X=" h e l l o ")$ cannot be rewritten further as the fact that $X$ is assigned to the string "hello" is not prevented by $\operatorname{proj}(X, \operatorname{int}(X) * \operatorname{times}(3, X, Y))$. However, if we have int $(X) * t i m e s(3, X, Y) *(X=" h e l l o ") \rightarrow 0$ this can be rewritten as 0 , as the int(•) constraint is conjunctive with ( $X=$ "hello").
    ${ }^{123}$ The set of values (named tuples of ground values) which are not mapped to zero.

[^77]:    ${ }^{124}$ Note: Rewrite rule 80 and rewrite rule 22 are usually performed simultaneously, in which case the rewrite looks like $R 1$ *opt (S1) +R2*opt (S2) $\xrightarrow{84} \mathrm{opt}(\mathrm{S} 1+\mathrm{S} 2) *(\mathrm{R} 1+\mathrm{R} 2)$. Rewrite rule 84 was not presented in the text as it is less general than rewrite rule 80.

[^78]:    ${ }^{125}$ Note, the $\mathbf{R}$-expr $(X=1)+(X=2)$ is a constraint as its multiplicity is $\leq 1$ in all cases. In general, our implementation assumes that all disjunctions are not constraints, but when the $\mathbf{R}$-expr is of a special form (such as a list of ground assignments) as we have here, then we can detect that it is a disjunctive constraint.

[^79]:    ${ }^{126}$ Or we could derive opt $((Y=1))$ from the first line, but when pulling the constraint on $Y$ through the aggregator we would get $\operatorname{opt}(\operatorname{proj}(Y,(Y=1))) \xrightarrow{41} \operatorname{opt}(1)$.

[^80]:    ${ }^{127}$ By this, I mean that Filardo allowed for modifications to the memoized values to be intermixed with when the message is sent about when change is about to happen or has already happened.

[^81]:    ${ }^{128}$ An internal update happens as a result of a cycle in the program. A cycle is where a value depends on itself. This is a specific kind of recursion that requires that the system solve an equation rather than just expand calls until a base case is reached. This was introduced previously in section $\S 2.5$ and will be reintroduced in section §10.4.
    ${ }^{129}$ Dyna allows for additional rules to be added to be added via the REPL. This can cause memos (cached computation) to become invalidated. Section §2.3
    ${ }^{130}$ Recall that \$memo is a user-definable rule which allows for controlling what and how something is memoized. Section $\S 2.7$

[^82]:    ${ }^{131}$ The signature should equivalence class the computation. Ideally, the signature will also allow for efficient look ups of memoized values (e.g. hash-table). This is not a strict requirement, as a system could (in theory) perform a linear scan over all memoized signatures.

[^83]:    ${ }^{132}$ This is actually $\Theta\left(\phi^{n}\right)$ where $\phi$ is the golden ratio.

[^84]:    ${ }^{133}$ The term pure function is a programming language term which would generally just be called a function in a mathematical setting. The returned result from a pure function is entirely determined by its arguments, which means that it returns the same value every time it is invoked and that the function is side-effect-free. Meaning that there is no other external observable behavior, such as modifying a global or class variable.

    134 An example of an impure function would be something that modifies global state. This can either be explicit, by mutating some global state, or implicitly by depending on the result of something like a random number generator (which maintains global state internally). E.g.
    $x=0$
    def not_pure():
    global x
    x += 1
    return x

[^85]:    ${ }^{135}$ Here, I have shown the explicit transformation of how memoization is added to a Python function. Python does provide a decorator that can perform this transformation automatically, using only one additional line of code. https://docs.python.org/3/library/functools.html\# functools.cache

[^86]:    ${ }^{136}$ The structure of Memo here is shown as a generic $\mathbf{R}$-expr. In practice, we might want to ensure that the $\mathbf{R}$-expr is represented as a hash-map so that it is efficient. Efficient disjunctive $\mathbf{R}$-expr kinds will be discussed in section §11.6.1.

[^87]:    ${ }^{137}$ The RMemo R-expr in this case might be better thought of as a "bulk" computed memo rather than an individual memo. This, of course, depends on the computation specified by S .
    ${ }^{138} \operatorname{Simplify} R$ in the context of $Q$ is equivalent to making the Simplify call $\operatorname{Simplify}(R, \mathcal{C}=\{Q\})$.
    ${ }^{139}$ In SImplify the system does not explicitly copy Q into the conditional. Instead, it depends on $Q$ by including it in the context. Hence, the system will instead check that $\operatorname{Simplify}(S, \mathcal{C}=\{Q\})$ returns an $\mathbf{R}$-expr that matches $1+\mathrm{T}$ for some T .

[^88]:    ${ }^{141} \mathrm{We}$ can check the signature by performing the rewrite sequence for $((N=0)+(N=1)) *(N=2) \rightarrow^{*} 0$.

[^89]:    ${ }^{142} \mathrm{We}$ have adopted the terminology of assumption from the JIT compiler literature [149].

[^90]:    ${ }^{143}$ Note: In a (future) parallel processing environment, it is possible that an invalid assumption to be "depended" on. This means that the computation that depends on an invalid assumption is already "stale" when it is computed. Most likely, it should be immediately thrown out and recomputed, as there is little use for a stale value.
    ${ }^{144}$ Future work has considered streaming queries, which are queries made against the Dyna program that are updated as the program changes rather than getting back a single value. Streaming queries can be implemented by subscribing to assumptions with a custom subscriber class that listens to invalidation messages.
    ${ }^{145}$ If the memoized $\mathbf{R}$-expr does not converge, e.g. $a=$ !a., then the memo not converging and running forever is consistent with the semantics of Dyna, section §2.5.

[^91]:    ${ }^{146}$ This can either be an R-expr, which was user-defined and changed via an external update, or a change resulting from a memo being recomputed.
    ${ }^{147}$ Proving two R-exprs are semantically equivalent is Turing-complete, hence can be difficult to check. We allow for one-sided error in that the system must prove that two $\mathbf{R}$-exprs are equivalent when it returns that two $\mathbf{R}$-exprs are equivalent. However, it is not required to prove that $\mathbf{R}$-exprs are not equivalent. Hence, it may overestimate that two R-exprs are not equivalent. We handled this by having a procedure that attempts to determine semantic equivalence in the case of minor reordering of the R-expr and renaming of variables. Section §12.4.2 also discusses checking R-exprs for semantic equivalence in the context of compilation.
    ${ }^{148}$ It is up to the user to write Dyna programs that converge. The Dyna system does not provide any automatic detection (or restrictions) that ensures that the programs converge (section §2.5).

[^92]:    ${ }^{149}$ This program will converge in the sequence $1,1.5,1.75,1.875, \ldots 2.0$. In theory, this is an infinitely long series converging to 2.0 , so one might think that this never stops/converges. In our case, however, we are using standard IEEE floating point numbers, which have limited numerical precision. Therefore, this sequence converges when run with floating point when it runs out of numerical precision and is rounded to 2.0 .
    ${ }^{150}$ This could have been computed using the closed form formula for a geometric series: $\frac{1}{1-r}$ where $r=\frac{1}{2}$ in this case.

[^93]:    ${ }^{151}$ In general, checking the semantic equivalence of $\mathbf{R}$-exprs is difficult unless we take special care to ensure that the structure of the memo is sufficiently "simple".

[^94]:    ${ }^{152}$ The queue has previously called an agenda in prior publications on Dyna. This was done previously to align the terminology with parsing algorithms [59].

[^95]:    ${ }^{153}$ For example, in the Fibonacci program figure 10-8, a memoization policy of \$memo(fib[X:\$free]) = "null". will cause the system to eagerly memoize all Fibonacci numbers, for which there are an infinite number of. Hence, this will not terminate.
    ${ }^{154}$ We could have used any ground value as the returned value from \$memo. The names "none", "null", and "unk" were chosen to maintain continuity with prior work on memoization on the Dyna project [66, 67].

[^96]:    ${ }^{155}$ In the case of a cycles, having a "null" memo can still be useful as the "null" memo will break the cycle, as seen in section $\S 10.5$.
    ${ }^{156}$ This is making use of the colon notation, which was presented as a Type Declaration section §2.8.3.

[^97]:    ${ }^{157}$ There are seven implementations of Dyna, including the one discussed in this chapter (depending on how you count). The major implementation attempts are linked at http://dyna.org/ \#downloads. The source discussed in this chapter is at https://github.com/argolab/dyna3. Previous Python prototypes that use rewriting can be found at https://github.com/argolab/dyna-R and https://github.com/argolab/dyna-R/blob/backend-v2/dyna_match_paper/rexprs.py.

[^98]:    ${ }^{158}$ At this point, we are not considering the constant factor overheads of implementation (such when choosing to write a program in slower Python vs faster C).
    ${ }^{159}$ Assuming that we are writing out the loops for the matrix-vector product and not using a library function.

[^99]:    ${ }^{160}$ Antlr4 is an off the shelf parser/lexer generator which generates a parser in Java.

[^100]:    ${ }^{161}$ This is similar to representing the return value as an argument in a Prolog program. For example, append([1, 2], [3], $[1,2,3]$ ).
    ${ }^{162}$ Throughout this chapter, I mention different global variables which are used to track the state of the system. More precisely, the global state is actually thread-local and specific to the current invocation of Simplify running. In this dissertation, the system is currently single-threaded, so there is no meaningful distinction.
    ${ }^{163}$ In the same way that a computer algebra system may return a large, complicated expression.

[^101]:    ${ }^{164}$ A guess will initially made with the 0 R-expr, which may cause the result from the query to be 0 also. Once the guess converges, the query may have a non 0 result. This was previously shown in section §10.5.

[^102]:    ${ }^{165}$ This is a Clojure deftype which gets compiled into a Java class file.
    166"Renaming variables" was previously denoted with the $R\{X \mapsto Y\}$ notation.
    ${ }^{167}$ The "list all variables" was previously denoted with vars( $R$ ).

[^103]:    ${ }^{168}$ Recall that constraints are $\mathbf{R}$-exprs whose multiplicity is at most 1 . (section $\S 5.2 .2 .4$ )

[^104]:    ${ }^{169}$ A bug might only be obvious after many subsequent steps of rewriting have been performed, which makes it very difficult to locate a bug
    ${ }^{170}$ At this time, there is no automatic checking that a rewrite is implemented correctly. This would require some executable representation of the semantic definitions (section §5.2) other than the rewrite rules, which we want to check.

[^105]:    ${ }^{171}$ This is the same as doing RexprObject. fieldName in Java. Coding this so that there is no indirection through method calls turned out to be very important for making this matching fast.
    ${ }^{172}$ Because $C$ is matched as : free, we know it must be a variable. In general, the $\mathbf{R}$-expr plus allows $C$ to be a value type that can be a constant or a variable.

[^106]:    ${ }^{173}$ Recall that an R-expr corresponds with a bag relation where 1 is returned when the assignment to the variables is contained in the bag (e.g. section §5.2.2.3). Hence, if we pass values to the plus $(\cdot, \cdot, \cdot)$ R-expr which are not supported, then this corresponds with multiplicity 0 (e.g. plus("hello", foo[1,2], X) $\rightarrow 0$ ).
    ${ }^{174}$ The name UnificationFailure was chosen as it aligns with the idea of unification failure in Prolog or CLP, which causes the system to backtrack and try another branch of a disjunction.
    ${ }^{175}$ Dyna runs on top of Java, and throwing an exception in Java is fast. This design may not work as well if Dyna is implemented with another programming language.

[^107]:    ${ }^{176}$ Previously in chapter $\S 8$ I denoted the context as being passed as an extra argument to all functions. This quickly becomes cumbersome, so it is easier to reference the context using a global (thread-local) variable.

[^108]:    ${ }^{177}$ Note, that (lessthan A B (is-true? _)) is the trinary version of lessthan discussed in section §7.1.4. The is-true? annotation is similar to : ground. However, it also checks value is true. The underscore _ is a placeholder variable.

[^109]:    ${ }^{178}$ The method map in Clojure returns a lazy sequence. We need to force the sequence to evaluate immediately using another Clojure function like vec or doall. I have omitted this in this example as it does not contribute to understanding the rewrite definition.

[^110]:    ${ }^{179}$ Some suggestions for efficient R-expr kinds would be dense matrix representations or GPU backed R-exprs and rewrites. (sections § 16.1 and 16.7)

[^111]:    ${ }^{180}$ Meaning that there is a sub-R-expr of the form $(X=7)$, where 7 is the known value.

[^112]:    ${ }^{181}$ The ( $X=$ ? ) is not an equality constraint, as we say that $(X=?) \rightarrow 1$ in all cases. This means that we can have something like $(X=7) *(X=?) \rightarrow(X=7)$. I have chosen to write this with an explicit value "?" rather than using an $\mathbf{R}$-expr like $(X=X)$ as this more closely matches the implementation.

[^113]:    ${ }^{182}$ The variable order is chosen arbitrarily. Future work should consider developing a better heuristic for this.

[^114]:    ${ }^{183}$ exposed-variables is the name in the implementation for $\operatorname{vars}(\cdot)$ from chapter $\S 5$.

[^115]:    ${ }^{185}$ In writing $\mathbf{R}$-exprs, we have been writing aggregators as ( $A=\operatorname{sum}(X, R)$ ) where each aggregator is its own $\mathbf{R}$-expr kind. There is a single aggregator implementation in the implementation, with different aggregators being distinguished by their name, represented here as operator.
    ${ }^{186}$ There is a similar rewrite rule for aggregator-inner which matches against the efficient disjunct. The efficient disjunct rule is a bit more complicated as it has to handle variables that are assigned, as well as the fact that some of those variable assignments might be projected out by the aggregator-inner.
    ${ }^{187}$ The delayed aggregator is an $\mathbf{R}$-expr like ResultVar=sum_aggregator_outer(R+sum_aggregator_inner([], Inp, ( $\operatorname{Inp}=$ accumulator))), where $R$ is the partially rewritten aggregator body as an $\mathbf{R}$-expr.
    ${ }^{188}$ The + on line 604 corresponds with running the addition between the current aggregated value and the value which has been computed from the inner aggregator. In the actual implementation, this would consult the aggregator operator rather than being hardcoded for $+=$ as done here to simplify this example.

[^116]:    ${ }^{189}$ The outer aggregator will have a disjunction of multiple inner aggregators (which represent each rule). The 0 is the identity element of disjunctions. This is conceptually similar to the unify $\mathbf{R}$-expr rewriting as 1 when its value has been saved in the context (figure 11-7). In that case, 1 is the identity element of the conjunction that contains the unify $\mathbf{R}$-expr.

[^117]:    ${ }^{190}$ Such as the trie in section §11.6.1.2 figure 11-13.

[^118]:    ${ }^{191}$ For example, if we have $((X=1)+(X=2)) *((X=1)+(X=2)+(X=3)+(X=4))$, then the first iterator iterates over $\{1,2\}$ and reports an estimated carnality of 2 , and the second iterator iterates over $\{1,2,3,4\}$ and reports an estimated carnality of 4 .

[^119]:    ${ }^{192}$ This observation was made on Python prototypes of term-rewriting. The code at https: //github.com/argolab/dyna-R/blob/backend-v2/dyna_match_paper/rexprs.py attempts to be a pure term rewriting-based system that closely matches chapter §8. https://github.com/ argolab/dyna-R was a prototype of the R-expr based system described in chapter §11, although it lacks several features added in the implementation for this dissertation: https://github.com/ argolab/dyna3.
    ${ }^{193}$ I note that the efficient disjunctions are asymptotically inefficient in some query modes. This asymptotic inefficiency was accounted for when making the judgment call to prioritize the work in this chapter over adding indices to the tries.

[^120]:    ${ }^{194}$ Our compilation targets Clojure, which is compiled into Java bytecode and eventually machine code by the Java virtual machine (JVM).

[^121]:    ${ }^{195}$ Variable identifiers are generic objects. Converting all variables to small integers that could be used as array indices is non-trivial because $\mathbf{R}$-expr objects are immutable and reused in multiple places.
    ${ }^{196}$ Recall that the pseudocode in section §8. A first matched against the type of the R-expr, and then checked each individual rewrite to see if any could apply.

[^122]:    ${ }^{197}$ \$priority returns a floating point value that is used to sort update messages (section §10.8.5).
    ${ }^{198}$ Admittedly, this is not addressed by JIT compilation in this dissertation, though future work should consider looking into this problem. Section §16.6

[^123]:    ${ }^{199}$ Names are generated randomly using the gensym method in Clojure, which generates a unique name using a global counter.

[^124]:    ${ }^{200}$ Hidden variables are how variables on projection are annotated, section §11.5.1.

[^125]:    ${ }^{201}$ This means that there are zero rewrites for a given R-expr kind. Not that there is no rewrite that can be applied. If there is a rewrite whose preconditions are not satisfied, then this will not invoke the JIT compiler. To allow JIT R-expr kinds to have more than one rewrite, the JIT compiler is still invoked with a low priority if there are no other rewrites that can be found to rewrite the top level R-expr.

[^126]:    ${ }^{202} A=X 2, B=X 7, E=X 1, C=X 4, D=X 0$.

[^127]:    ${ }^{203}$ Recall that last argument to def-rewrite is a Clojure expression that when evaluates returns an $\mathbf{R}$-expr or value depending on the keyword arguments passed to def-rewrite.

[^128]:    ${ }^{204} \mathbf{R}$-exprs that are contained in holes will have a variable which is treated as an opaque pointer to an R-expr.
    ${ }^{205}$ For example, the function make-lessthan is used by rewrite rule 34 and in the def-rewrite example figure 11-8.
    ${ }^{206}$ The "value type" is a Java interface that is implemented by both constants and variables. Here, we are simply defining new classes that also implement the value type interface.

[^129]:    ${ }^{207}$ The Truffle project is built on and closely integrated with the Graal compiler [56, 57].
    ${ }^{208}$ Some manual cleanup has been done to make the code presentable.
    ${ }^{209}$ Even if we generated Java bytecode directly, there are limitations on the JUMP instructions that can be generated, which is enforced by the Java Virtual Machine.

[^130]:    ${ }^{210}$ The JIT compiler also generates the necessary code to implement the Getlterables for all JIT-generated R-expr kinds.

[^131]:    ${ }^{214}$ There is one $\mathbf{R}$-expr class that implements the dynabaseCreate $\mathbf{R}$-expr, and dynabaseTypeID is held in a metadata field on the $\mathbf{R}$-expr class. This is the same as with structural-equality constraints that track the name of the structural-term through a metadata field, section §8.1.

[^132]:    ${ }^{215}$ An alternate way to get self-inheritance is by defining a term in terms of itself using recursion. Something like
    $f(N):=$ new $f(N-1)$ \{ $z+=N\}.$.
    $f(0):=$ new $\}$.

[^133]:    ${ }^{216}$ The dynabase's parent reference is akin to the SuperDynabase variable in section §13.3, which could be stored in a list data structure, as discussed at the beginning of section §13.2.

[^134]:    ${ }^{217}$ As the $\mathbf{R}$-expr will keep growing in size every time it is expanded. In other words, the program has a "data dependency" to ensure that the recursion eventually reaches a base case.

