1 Introduction

In the previous lecture we have seen backward chaining from a logical perspective, and how this can be seen as a foundation for backward-chaining logic programming languages like Prolog.

In this lecture we take a small step sideways: instead of considering all atoms to be negative we consider all atoms positive. This has a rather drastic impact on the operational behavior of proof search, leading to forward-chaining logic programming. This is also called bottom-up logic programming, although the direction is strangely reversed from the way we consider the proof construction process.

2 Reading Inference Rules from Premises to Conclusion

 Mostly over the last serious of lectures, we read inference rules by looking at the conclusion first and then the premises. This was so, because that is the direction of proof construction. In fact, the sequent calculus was specifically engineered by Gentzen to have this property!

Now we will read inference rules starting with the premises. For example, assume we would like to calculate the path relation in an undirected graph, where we say there is a path from vertex $x$ to $y$ if there a sequence of vertices $x = x_0, x_1, \ldots, x_n = y$ such that all $x_i$ and $x_{i+1}$ are connected by an edge. For simplicity, let us say that $n \geq 1$. 
We represent the vertices of a graph by constants, and the edge relation with a predicate $\text{edge}(x, y)$ if there is an edge from $x$ to $y$. Here is a specification of the path relation:

$$
\frac{\text{edge}(x, y)}{\text{edge}(y, x)} \quad \text{sym} \\
\frac{\text{edge}(x, y)}{\text{path}(x, y)} \quad \text{ep} \\
\frac{\text{path}(x, y)}{\text{path}(y, z)} \quad \text{trans}
$$

The first rule (sym) expresses we are working over an undirected graph. The second (ep) expresses that an edge represents a valid path, and the third that the path relation is transitive.

Read from the conclusion to the premises, backward logic programming search over this specification is useless. Even just the rule sym rule will lead to an infinite loop, and the trans rule has an unknown $y$ in the premise even if $x$ and $z$ are known in the conclusion.

Read from the premises to the conclusion, however, this is a decent program if we avoid re-deriving facts we already know. After while, this program must terminate because there are at most $O(n^2)$ facts of the form $\text{edge}(x, y)$ and $\text{path}(x, y)$ that could be derived. When inference reaches the point where any additional inference only infers facts we already know, we say the program has reached saturation and it halts. At this point we can answer any specific query simply by looking it up in the collection of derived facts, usually called the database.

### 3 Saturation

As another example, we first consider the usual bottom-up specifications of $\text{even}(n)$ and $\text{odd}(n)$ for unary numbers.

$$
\frac{\text{even}(z)}{\text{ev}_z} \\
\frac{\text{odd}(N)}{\text{ev}(s(N))} \quad \text{ev}_s \\
\frac{\text{even}(N)}{\text{odd}(s(N))} \quad \text{od}_s
$$

Reading these from the premise to the conclusion does not work: these rules would create an unbounded database with facts

$$
\text{even}(z), \text{odd}(s(z)), \text{even}(s(s(z))), \ldots
$$

But if we view these as introduction rules we can derive elimination rules that work in the other direction, using what we already know!

$$
\frac{\text{even}(s(N))}{\text{odd}(N)} \quad \text{ev}'_s \\
\frac{\text{odd}(s(N))}{\text{even}(N)} \quad \text{od}'_s \\
\frac{\text{odd}(z)}{\text{C} \quad \text{od}'_z}
$$
Note that there is no rule for \(\text{even}(z)\) because we cannot extract any information from that: there rule \(\text{ev}_z\) has no premises. In the last rule we have derived a contradiction, which is manifest in being able to conclude any proposition \(C\). If we want to use this as a program, we have to use a specific proposition. Unfortunately, \(\top\) is not part of the chaining fragment (for good reason), so we use a new atom \(\text{no}\), in the best tradition of the Prolog top level.

If we want to know if, say, the fact \(\text{even}(s(s(s(z))))\) is consistent with the definition of the predicate, we assert it in the database of facts and saturate the database using forward inference. Because of the simple nature of these rules, our hand is forced at each point of inference, and we obtain the following saturated database:

\[
\begin{align*}
\text{even}(s(s(z))))), \text{odd}(s(s(z))), \text{even}(s(z)), \text{odd}(z), \text{no}
\end{align*}
\]

This tells us that asserting that 3 is even is inconsistent with the definition of evenness. On the other hand, we we assert \(\text{even}(s(s(z)))\), we learn:

\[
\begin{align*}
\text{even}(s(s(z))), \text{odd}(s(z)), \text{even}(z)
\end{align*}
\]

This database is now saturated and there is no contradiction, so the assertion that 2 is even is consistent with the definition of evenness.

\section{Forward Chaining}

Our translation from rules to propositions leads us to the following propositions representing the downward-reading rules for even and odd numbers:

\[
\Gamma_{eo} = \forall n. \text{even}(s(n)) \supset \text{odd}(n),
\forall n. \text{odd}(s(n)) \supset \text{even}(n),
\text{odd}(z) \supset \text{no}
\]

We then ask, for example,

\[
\Gamma_{eo}, \text{even}(s(s(s(z)))) \rightarrow \text{no}
\]

to find out if the fact that 3 is even is consistent with the knowledge in \(\Gamma_{eo}\).

This search is now the exact opposite of goal directed, let’s call it database directed. We ignore the succedent (no) and saturate the database, which are the atoms in the antecedents. Only once we have saturated the database, do we even look at the succedent and see if it is a fact in the database (or,
more generally, can be proven from the database directly without further forward chaining).

For this intuition to work, we have to start by instantiating $n$ in the first proposition with \( s(s(z)) \) and then use the implication left rule to conclude \( \text{odd}(s(s(z))) \). It is this process we call forward chaining. To formalize this, we first recall the language and rules for backward chaining.

**Backward chaining fragment:** all atoms are negative, and the only polarity shift has the form $\downarrow P^-$. So far, we have only shown the rules for the connectives in red.

<table>
<thead>
<tr>
<th>Program formulas</th>
<th>$D^- ::= P^- \mid G^+ \supset D^- \mid \forall x. D^-(x) \mid D_1^- \land D_2^- \mid \top$</th>
</tr>
</thead>
<tbody>
<tr>
<td>Programs</td>
<td>$\Gamma^- ::= \cdot \mid \Gamma, D^-$</td>
</tr>
<tr>
<td>Goal formulas</td>
<td>$G^+ ::= \downarrow P^- \mid G_1^+ \land G_2^+ \mid \top \mid \exists x. G^+(x) \mid G_1^+ \lor G_2^+ \mid \bot$</td>
</tr>
</tbody>
</table>

**Forward chaining fragment:** all atoms are positive, and the only polarity shift has the form $\uparrow P^+$. 

<table>
<thead>
<tr>
<th>Program formulas</th>
<th>$D^- ::= \uparrow P^+ \mid G^+ \supset D^- \mid \forall x. D^-(x) \mid D_1^- \land D_2^- \mid \top$</th>
</tr>
</thead>
<tbody>
<tr>
<td>Database</td>
<td>$\Gamma ::= \cdot \mid \Gamma, D^- \mid \Gamma, P^+$</td>
</tr>
<tr>
<td>Goal formulas</td>
<td>$G^+ ::= P^+ \mid G_1^+ \land G_2^+ \mid \top \mid \exists x. G^+(x) \mid G_1^+ \lor G_2^+ \mid \bot$</td>
</tr>
</tbody>
</table>

The antecedents now mix the program formulas $D^-$ (sometimes called the IDB) and the database facts $P^+$ (sometimes called the EDB), while the succedent is always positive since negative atoms are not part of this fragment. We have the following three judgments

<table>
<thead>
<tr>
<th>Backward Chaining</th>
<th>Forward Chaining</th>
</tr>
</thead>
<tbody>
<tr>
<td>$\Gamma^- \xrightarrow{f} P^-$</td>
<td>stable sequent</td>
</tr>
<tr>
<td>$\Gamma, [D^-] \xrightarrow{f} P^-$</td>
<td>left focus</td>
</tr>
<tr>
<td>$\Gamma^- \xrightarrow{f} [G^+]$</td>
<td>right focus</td>
</tr>
<tr>
<td>$\Gamma \xrightarrow{f} C^+$</td>
<td></td>
</tr>
<tr>
<td>$\Gamma, [D^-] \xrightarrow{f} C^+$</td>
<td></td>
</tr>
<tr>
<td>$\Gamma \xrightarrow{f} [G^+]$</td>
<td></td>
</tr>
</tbody>
</table>

The rules for the connectives remain the same, with the exception of the order of premises in the $\supset$ rule.\(^1\) We also remove the rules concerned with negative atoms and add those for positive ones.

\(^1\)It is possible that in a combined forward/backward chaining language, there should be two forms of implication, presenting the premises in different order: $A \supset B$ proves $A$ before assuming $B$, and $B \subset A$ assumes $B$ before proving $A$. This first one would be employed in forward chaining, the second in backward chaining. Of course, logically the two are the same, but not operationally when viewed from the computation-as-proof-search perspective.
\[
\begin{align*}
D^{-} & \in \Gamma \quad \Gamma, [D^{-}] \overset{f}{\rightarrow} C^{+} \quad \text{focusL} \\
\Gamma & \overset{f}{\rightarrow} C^{+} \quad \Gamma, P^{+} \overset{f}{\rightarrow} C^{+} \quad \uparrow L \\
\Gamma, [D^{-}(X)] & \overset{f}{\rightarrow} C^{+} \quad \forall L^{*} \\
\Gamma & \overset{f}{\rightarrow} [G^{+}] \quad \Gamma, [D^{-}] \overset{f}{\rightarrow} C^{+} \\
\Gamma, [G^{+} \supset D^{-}] & \overset{f}{\rightarrow} C^{+} \\
\end{align*}
\]

Figure 1: Forward chaining fragment of Horn logic
Let’s observe these rules in action on our program and goal, where all atoms are positive. We have added shifts on the right-hand side of implications during polarization of the propositions.

\[
\Gamma_{eo} = \forall n. \text{even}(s(n)) \supset \uparrow \text{odd}(n), \\
\forall n. \text{odd}(s(n)) \supset \uparrow \text{even}(n), \\
\text{odd}(z) \supset \uparrow \text{no}
\]

\[
\Gamma_{eo}, \text{even}(s(s(z))) \xrightarrow{f} \text{no}
\]

Say we focus on the first proposition in \(\Gamma_{eo}\).

\[
\vdots \quad \vdots
\]

\[
\frac{\Gamma_{eo}, \text{even}(s(s(z))) \xrightarrow{f} [\text{even}(s(N))] \quad \Gamma_{eo}, \text{even}(s(s(z))), [\uparrow \text{odd}(N)] \xrightarrow{f} \text{no} \quad \forall L}{\Gamma_{eo}, \text{even}(s(s(z))), [\text{even}(s(N)) \supset \uparrow \text{odd}(n)] \xrightarrow{f} \text{no} \quad \forall L}
\]

Now we match \(\text{even}(s(s(z))))\) against \(\text{even}(s(N))\) which succeeds with substitution \(N = s(z)\), which is applied globally to the partial proof which then looks as follows:

\[
\vdots
\]

\[
\frac{\Gamma_{eo}, \text{even}(s(s(z))) \xrightarrow{f} [\text{even}(s(s(z))))] \quad \Gamma_{eo}, \text{even}(s(s(z))), [\uparrow \text{odd}(s(z))] \xrightarrow{f} \text{no} \quad \forall L}{\Gamma_{eo}, \text{even}(s(s(z))), [\text{even}(s(s(z)))) \supset \uparrow \text{odd}(n)] \xrightarrow{f} \text{no} \quad \forall L}
\]

In the right branch of the proof we now lose focus, and we have reached a stable sequent, adding another fact to the database.

\[
\vdots
\]

\[
\frac{\Gamma_{eo}, \text{even}(s(s(z))) \xrightarrow{f} [\text{even}(s(s(z))))] \quad \Gamma_{eo}, \text{even}(s(s(z))), [\uparrow \text{odd}(s(z))] \xrightarrow{f} \text{no} \quad \uparrow L}{\Gamma_{eo}, \text{even}(s(s(z))), \text{odd}(s(s(z))) \xrightarrow{f} \text{no} \quad \forall L}
\]
This will continue until we add \( \text{odd}(z) \) and then \( \text{no} \) to the database. At this point the goal looks like

\[
\vdash \\
\Gamma_{\text{eo}, \ldots, \text{no}} \rightarrow \text{no}
\]

Recall that all atoms are positive, so we can now focus on the succedent and complete the proof.

\[
\begin{array}{c}
\Gamma_{\text{eo}, \ldots, \text{no}} \xrightarrow{\text{id}^+} \\
\Gamma_{\text{eo}, \ldots, \text{no}} \xrightarrow{\text{[no]}} \Gamma_{\text{eo}, \ldots, \text{no}} \xrightarrow{\text{focus}_R} \\
\Gamma_{\text{eo}, \ldots, \text{no}} \xrightarrow{\text{no}}
\end{array}
\]

We can also see what happens if we focus on the “wrong” antecedent, that is, one that we cannot use with forward chaining.

\[
\begin{array}{c}
\Gamma_{\text{eo} \cdot \text{even}(s(s(z)))) \xrightarrow{\text{f}} [\text{odd}(s(N))], \Gamma_{\text{eo} \cdot \text{even}(s(s(z)))) \xrightarrow{[\text{even}(N)]} \xrightarrow{\supset} \Gamma_{\text{eo} \cdot \text{even}(s(s(z)))) \xrightarrow{\forall} \Gamma_{\text{eo} \cdot \text{even}(s(s(z)))) \xrightarrow{\forall} \Gamma_{\text{eo} \cdot \text{even}(s(s(z)))) \xrightarrow{\forall \text{even}(n)} \rightarrow L
\end{array}
\]

Experience shows that, generally, the reduction in the search space with forward chaining is not quite as drastic as with backward chaining. This is because there is only one goal (the succedent) but many clauses in the database (the antecedents). Nevertheless, there are many algorithms more easily described with forward chaining than with backward chaining. We will show one in Section 6 on unification.

5 Comparing Backward and Forward Chaining

We return to a simple example from last lecture

\[
a, a \supset b, b \supset c \rightarrow c
\]

If we polarize all atoms negatively, we obtain

\[
\bar{a}, \downarrow \bar{a} \supset \bar{b}, \downarrow \bar{b} \supset \bar{c} \rightarrow \bar{c}
\]
where the shift operators binds tightest.

Without backward chaining there are different proofs. In particular, we could apply \( \supset L \) to \( a \supset b \), or to \( b \supset c \). With backward chaining, there is only one possible proof: at each choice point, when we focus, only on possibility will succeed and the others will fail immediately. You should convince yourself that this is the case. We define \( \Gamma_0 = (a^-, \downarrow a^- \supset b^-, \downarrow b^- \supset c^-) \)

\[
\begin{align*}
\Gamma_0, [a^-] & \xrightarrow{f} a^- & \text{id}^- \\
\Gamma_0, [a^-] & \xrightarrow{f} a^- & \text{focus}_{L} \\\n\Gamma_0, [b^-] & \xrightarrow{f} b^- & \text{id}^- \\
\Gamma_0, [b^-] & \xrightarrow{f} b^- & \downarrow R \\\n\Gamma_0, [\downarrow a^- \supset b^-] & \xrightarrow{f} b^- & \supset L \\
\Gamma_0, [\downarrow a^- \supset b^-] & \xrightarrow{f} b^- & \text{focus}_{L} \\
\Gamma_0, [c^-] & \xrightarrow{f} c^- & \text{id}^- \\
\Gamma_0, [c^-] & \xrightarrow{f} c^- & \downarrow R \\\n\Gamma_0, [\downarrow b^- \supset c^-] & \xrightarrow{f} c^- & \supset L \\
\Gamma_0, [\downarrow b^- \supset c^-] & \xrightarrow{f} c^- & \text{focus}_{L} \\
\end{align*}
\]

Going back to our sequent

\[ a, a \supset b, b \supset c \rightarrow c \]

if we polarize all the atoms positively, in preparation for forward chaining, we get

\[ a^+, a^+ \supset \uparrow b^+, b^+ \supset \uparrow c^+, \uparrow c^+ \rightarrow c^+ \]

Now we must focus on \( a^+ \supset \uparrow b^+ \) first, then \( b^+ \supset \uparrow c^+ \), then \( c^+ \). All other attempts at focusing will either fail, or lead conclude a fact that is already
in the database. We abbreviate \(\Gamma_1 = (a^+, a^+ \supset \uparrow b^+, b^+ \supset \uparrow c^+)\)

\[
\begin{align*}
\frac{\Gamma_1, b^+ \xrightarrow{f} [c^+]}{\text{id}^+} & \quad \frac{\Gamma_1, b^+ \xrightarrow{f} [c^+]}{\text{focus}_R} \\
\frac{\Gamma_1, b^+ \xrightarrow{f} [b^+]}{\text{id}^+} & \quad \frac{\Gamma_1, b^+, [\uparrow c^+] \xrightarrow{f} c^+}{\uparrow L} \\
\frac{\Gamma_1, b^+ \xrightarrow{f} [c^+]}{\text{id}^+} & \quad \frac{\Gamma_1, b^+, [\uparrow c^+] \xrightarrow{f} c^+}{\supset L} \\
\frac{a^+ \in \Gamma}{\Gamma_1 \xrightarrow{f} [a^+]} & \quad \frac{\Gamma_1, b^+, [\uparrow b^+] \xrightarrow{f} c^+}{\text{focus}_L} \\
\frac{\Gamma_1, b^+ \xrightarrow{f} c^+}{\uparrow L} & \quad \frac{\Gamma_1, [\uparrow b^+] \xrightarrow{f} c^+}{\supset L} \\
\frac{\Gamma_1, [a^+ \supset \uparrow b^+] \xrightarrow{f} c^+}{\text{focus}_L} & \quad \frac{\Gamma_1 \xrightarrow{f} c^+}{\text{id}^+} \\
\end{align*}
\]

### 6 Unification

As a major and significant example of forward chaining, which is similar to many realistic applications of Datalog, we use unification itself. So far, we have just treated it informally, despite its complexity.

We describe the algorithm by a set of rules concerning a predicate \(t \equiv s\) for (first-order) terms \(t\) and \(s\). This set of rules can be translated to a collection of propositions \(\Gamma_u\) where all atoms are positive. We assert a new equality, adding it as an antecedent, and then saturate the database. If it produces no, then the new equality is inconsistent with all the information we already had. Otherwise, the new saturated database represents the “solution” and shows consistency.

We begin with two rules that compare the function symbol at the head of the two terms. We write \(\bar{t}\) for a sequence of terms.

\[
\begin{align*}
& f(\bar{t}) \equiv f(\bar{s}) & \text{con} = & \quad \frac{f(\bar{t}) \equiv g(\bar{s})}{\text{no}} \\
& \bar{t} \equiv \bar{s} & \quad \text{f \neq g} \\
\end{align*}
\]

The first rule expresses that if \(f(\bar{t})\) is equal to \(f(\bar{s})\) then the sequences of arguments must be equal. This means that function symbols are “uninterpreted”: they are used as data constructors, not to stand for arbitrary mathematical functions.
The second rule expresses that if the data constructors are different then
the terms are not equal. In other words, if we know they are equal, then
this is a contradiction.

These two rules also capture constants, since we think of a constant \( c \)
and \( c() \), with the empty sequence of arguments.

Now we need four rules for comparing sequences of arguments.

\[
\begin{align*}
(t, \bar{t}) & \equiv (s, \bar{s}) & \text{seq} = 1 \\
\bar{t} & \equiv \bar{s} \\
() & \equiv (s, \bar{s}) & \text{no} \quad \text{seq} \neq 1 \\
\bar{t} & \equiv () & \text{seq} \neq 2
\end{align*}
\]

Note that there is not rule for \( () \equiv () \), because such an equality contains no
information to extract.

At this point we have enough rules to decide equality, but not yet enough
to implement unification. Consider

\[
f(X, X) \equiv f(c(), d())
\]

This problem must fail, since \( X \) cannot be equal to \( c() \) and \( d() \) simulta-
neously, but the rules so far do not account for this. The simple device
of stating symmetry and transitivity of equality will solve this particular
problem.

\[
\begin{align*}
t & \equiv s \quad \text{sym} & t \equiv s \\
\text{sym} & \quad s \equiv r \\
\text{trans} & \quad t \equiv r
\end{align*}
\]

Now we deduce:

\[
\begin{align*}
f(X, X) & \equiv f(c(), d()) \\
(X, X) & \equiv (c(), d()) \\
X & \equiv c() \quad \text{given} \\
(X) & \equiv (d()) \\
X & \equiv d() \quad \text{by rule seq}_1 \\
c() & \equiv X \quad \text{by rule seq}_2 \\
c() & \equiv d() \quad \text{by rule sym} \\
\text{no} & \quad \text{by rule trans} \\
\text{no} & \quad \text{by rule con}_\neq
\end{align*}
\]

One could make these rules more efficient, for example, by restricting some
terms in symmetry and transitivity to be variables.
At this point we have arrived at Prolog-style unification. Unfortunately, we know that this is unsound, because an equation such as

\[ X = f(X) \]

is not recognized as inconsistent. We can incorporate this by adding some rules for the occurs-check that discover such inconsistencies.

\[
\begin{align*}
X & \notin f(t) & X & \notin X \\
X & \notin f(t) & X & \notin t \\
X & \notin f(t) & X & \notin (t, \bar{t}) \\
X & \notin Y & Y & \equiv t \\
X & \notin t \\
\end{align*}
\]

The last rule is necessary to obtain contradictions from problems such as

\[ X = f(Y), Y = g(X) \]

With a few optimizations, these rules can be seen to define Huet’s algorithm for (first-order) unification [Hue76]. This proceeds in two stages: in the first stage we saturate the equalities, and once they are saturated we perform the occurs-check. If implemented correctly, this will have complexity \(O(n \log(n))\), where \(n\) is the size of the input problem. Robinson’s original unification algorithm [Rob71] in contrast was exponential in the size of input, although in the context of his applications it performed quite well [CB83].

7 From Propositions to Rules of Inference

As we have seen in this lecture and also already in the last lecture, we can translate inference rules to propositions and then use either forward or backward chaining to specify and operational semantics for proof search.

Question: can we go the other way? That is, can we take proposition and turn them into inference rules? Another way to pose the question: can we take advantage of the chaining semantics to compile program propositions into “big-step” inference rules so we don’t have to play through focusing all the time?
Let’s try with the example of the type of the S combinator:

$$(a \supset (b \supset c)) \supset ((a \supset b) \supset (a \supset c))$$

In order to prove this, we first apply inversion as far as we can and arrive at the sequent

$$a, a \supset b, a \supset (b \supset c) \rightarrow c$$

Now we have to decide on a polarization. We’ll try all first all negative and then all positive.

Polarizing all atoms as negative we then have

$$a^-, \downarrow a^- \supset b^-, \downarrow a^- \supset (\downarrow b^- \supset c^-) \xrightarrow{f} c^-$$

Which propositions could we focus on? Not on $c^-$ in the succedent—there is no rule for that in the backward chaining fragment. But we can focus on each of the antecedents since they are all negative propositions. In each case we imagine what would happen if we focused on the proposition, not knowing the remaining antecedents $\Gamma^-$ or the conclusion $P^-$. But note that since antecedents are persistent, all propositions in $\Gamma_0 = (a^-, \downarrow a^- \supset b^-, \downarrow a^- \supset (\downarrow b^- \supset c^-))$ will always be present in $\Gamma$.

$$\begin{align*}
\Gamma^- & \xrightarrow{f} P^- \\
a^- & = P^- \\
\Gamma^- & \xrightarrow{\text{id}} P^- \\
\Gamma^- & \xrightarrow{\text{focusL}} P^-
\end{align*}$$

So for the focus on $a^-$ to succeed, the right-hand side $P^-$ must be equal to $a^-$. This gives use the derived rule

$$\begin{align*}
\Gamma^- & \xrightarrow{f} \Gamma^- \xrightarrow{R_1} a^-
\end{align*}$$

Focusing on the second proposition:

$$\begin{align*}
P^- & = b^- \\
\Gamma^- & \xrightarrow{\text{id}} P^- \\
\Gamma^- & \xrightarrow{\text{focusL}} P^-
\end{align*}$$
We see $P^- = b^-$ and we have one stable subgoal that will be the premise of the derived rule.

$$\Gamma^- \not\vdash a^- \quad R_2$$
$$\Gamma^- \not\vdash b^-$$

Finally, focusing on the third antecedent:

$$c^- = P^- \quad \text{id}^-$$
$$\Gamma^- , [c^-] \not\vdash P^-$$
$$\Gamma^- \not\vdash b^- \quad \downarrow R$$
$$\Gamma^- \not\vdash [\downarrow b^-]$$
$$\Gamma^- , [\downarrow b^- \supset c^-] \not\vdash P^- \quad \supset L$$
$$\Gamma^- \not\vdash a^- \quad \downarrow R$$
$$\Gamma^- \not\vdash [\downarrow a^-]$$
$$\Gamma^- , [\downarrow a^- \supset (\downarrow b^- \supset c^-)] \not\vdash P^- \quad \text{focusL}$$
$$\Gamma^- \not\vdash P^-$$

We read off $P^- = c^-$ and the two stable sequents that are the premises of the derived rules:

$$\Gamma^- \not\vdash b^- \quad \Gamma^- \not\vdash a^- \quad R_3$$
$$\Gamma^- \not\vdash c^-$$

Summarizing the three rules:

$$\Gamma^- \not\vdash a^- \quad R_1$$
$$\Gamma^- \not\vdash b^- \quad \Gamma^- \not\vdash a^- \quad R_2$$
$$\Gamma^- \not\vdash b^- \quad \Gamma^- \not\vdash a^- \quad R_3$$

To see how they prove our original sequent

$$a^- , \downarrow a^- \supset b^- , \downarrow a^- \supset (\downarrow b^- \supset c^-) \not\vdash c^-$$

we first take all the negative propositions away from the antecedents. That’s because instead of focusing on them, we should be using a derived rule. Here is the resulting proof.

$$\cdot \not\vdash a^- \quad R_1$$
$$\cdot \not\vdash b^- \quad \cdot \not\vdash a^- \quad R_3$$
$$\cdot \not\vdash c^-$$
Of course, this is much shorter and more efficient than the proof using the explicit focusing rules.

Circling back: let’s make all atoms positive

$$a^+, a^+ \supset \uparrow b^+, a^+ \supset (b^+ \supset \uparrow c^+) \overset{f}{\rightarrow} c^+$$

This time, we cannot focus on $a^+$ in the antecedent, but on the other two propositions and also on the succedent. Let’s do each in turn. Again, remember that $\Gamma_0 = (a^+ \supset \uparrow b^+, a^+ \supset (b^+ \supset \uparrow c^+))$ is a part of every sequent in a proof. The succedent can be any positive proposition $G^+$.

$$\frac{a^+ \in \Gamma}{\Gamma \overset{f}{\rightarrow} [a^+]} \quad \frac{\Gamma, b^+ \overset{f}{\rightarrow} G^+}{\uparrow L} \quad \frac{\Gamma, \uparrow b^+ \overset{f}{\rightarrow} G^+}{\supset L} \quad \frac{\Gamma, [a^+ \supset \uparrow b^+] \overset{f}{\rightarrow} G^+}{\text{focusL}}$$

Here, $a^+ \in \Gamma$ should be seen as a constraint, so we just write $\Gamma = (\Gamma', a^+)$ and obtain the rule

$$\frac{\Gamma', a^+, b^+ \overset{f}{\rightarrow} G^+}{\Gamma', a^+ \overset{f}{\rightarrow} G^+} S_1$$

For the second proposition:

$$\frac{a^+ \in \Gamma}{\Gamma \overset{f}{\rightarrow} [a^+]} \quad \frac{b^+ \in \Gamma}{\Gamma \overset{f}{\rightarrow} [b^+]} \quad \frac{\Gamma, c^+ \overset{f}{\rightarrow} G^+}{\uparrow L} \quad \frac{\Gamma, \uparrow c^+ \overset{f}{\rightarrow} G^+}{\supset L} \quad \frac{\Gamma, [a^+ \supset (b^+ \supset \uparrow c^+)] \overset{f}{\rightarrow} G^+}{\text{focusL}}$$

Again, collecting membership constraints we see $\Gamma = (\Gamma', a^+, b^+)$ and we get

$$\frac{\Gamma', a^+, b^+, c^+ \overset{f}{\rightarrow} G^+}{\Gamma', a^+ \overset{f}{\rightarrow} G^+} S_2$$
Finally, we can focus on the succedent:

\[
\frac{c^+ \in \Gamma}{\Gamma \xrightarrow{f} \{c^+\}} \quad \text{id}^+ \\
\frac{\Gamma \xrightarrow{f} \{c^+\}}{\Gamma \xrightarrow{f} c^+} \quad \text{focus}^R
\]

which gives us

\[
\frac{\Gamma', c^+ \xrightarrow{f} c^+}{S_3}
\]

In summary:

\[
\begin{align*}
\Gamma', a^+, b^+ \xrightarrow{f} G^+ & \quad \rightarrow S_1 & \Gamma', a^+, b^+, c^+ \xrightarrow{f} G^+ & \quad \rightarrow S_2 \\
\Gamma', a^+ \xrightarrow{f} G^+ & \quad \rightarrow S_3 & \Gamma', c^+ \xrightarrow{f} c^+ & \quad \rightarrow S_3
\end{align*}
\]

With these three rules we can drop $\Gamma_0$ since their only purpose would be to focus on them—and that has been replaced by the derived rules. Our big-step proof becomes:

\[
\begin{align*}
\frac{a^+, b^+, c^+ \xrightarrow{f} c^+}{S_3} & \quad \rightarrow S_3 \\
\frac{a^+, b^+ \xrightarrow{f} c^+}{S_2} & \quad \rightarrow S_2 \\
\frac{a^+ \xrightarrow{f} c^+}{S_1} & \quad \rightarrow S_1
\end{align*}
\]

Note that we have kept $a^+$ among the antecedents since we cannot focus on a positive atom (or any positive proposition, for that matter) in the antecedent.

8 Polarization: A Brief Roadmap

A critical component in understanding the various fragment and operational interpretations we have seen is polarization [Lau99]. We started from the inversion strategy.

Negative Propositions are those with invertible right rules, that is, if we see them in the succedent we can apply their right rule without considering any other choice and search remains complete.
**Positive Propositions** are those with invertible left rules, that is, if we see them in the antecedent we can apply their left rule without considering any other choice and search remains complete.

In order for every proposition to have a polarized version we need the so-called *shift* operators $\uparrow$ and $\downarrow$ to go between the two classes of propositions. We get:

\[
\begin{align*}
\text{Neg. props.} & \quad A^- ::= P^- \mid B^+ \supset A^- \mid \forall x. A^-(x) \mid A_1^- \land A_2^- \mid \top \mid \uparrow B^+ \\
\text{Pos. props.} & \quad B^+ ::= P^+ \mid B_1^+ \land B_2^+ \mid \top \mid \exists x. B^+(x) \mid B_1^+ \lor B_2^+ \mid \bot \mid \downarrow A^-
\end{align*}
\]

Atoms can be either negative ($P^-$) or positive ($P^+$). During polarization we can choose the polarization of each atom $p$ freely, but must assign the same polarity to each occurrence of $p$. Conjunction and truth have invertible left and right rules, so they appear in both rows.

*Chaining* is the opposite of inversion: we focus on a particular negative antecedent or positive succedent and continue to apply rules only to the single proposition in focus until the focusing phase is interrupted by a shift, changing the polarity of the proposition.

Chaining by itself is complete for proof search as long as the shifts are restricted such that we only have $B^- ::= \cdots \mid \downarrow P^-$ and $A^- ::= \cdots \mid \uparrow P^+$. The language so restricted is (an insignificant extension) of Horn logic.

In case all atoms are negative, chaining is called *backward chaining* (also called *top-down logic programming*), which is a goal-directed proof search strategy and the foundation of Prolog [Kow88].

In case all atoms are positive, chaining is called *forward chaining* (also called *bottom-up logic programming* [NR91]), which is a saturation-based proof search strategy and the foundation of Datalog.

There is the possibility of allowing both positive and negative atoms, but the resulting mixed chaining logic programming language has never been deeply investigated, as far as I am aware. However, chaining is complete for this language, so it is a plausible candidate for an interesting and expressive language.

If we allow arbitrary polarized propositions, then chaining alone is insufficient: positive (non-atomic) propositions can show up in the antecedents, and negative (non-atomic) propositions in the succedent. When such a proposition is encountered, we apply inversion until we once again reach a *stable sequent* which is characterized with only negative propositions and positive atoms as antecedents, and positive propositions and negative atoms as succedents.
Focusing = chaining + inversion was first discovered by Andreoli [And92], with two caveats: (1) his propositions were not implicitly polarized, and (2) focusing was defined for linear logic [Gir87], which we will only see later in this course. However, focusing (and also the chaining-only fragment based on Horn logic) has been remarkably robust in that it applies to a large number of reasonable substructural, modal, and other logics, both intuitionistic and classical.

We will discuss focusing in the next lecture.

References


