

2.2 Peephole Optimization

Idea:

- Slide a **small** window over the program.
- Optimize aggressively inside the window, i.e.,
 - Eliminate redundancies!
 - Replace expensive operations inside the window by cheaper ones!

Examples:

$$x = x + 1; \quad \Longrightarrow \quad x++;$$

// given that there is a specific increment instruction :-)

$$z = y - a + a; \quad \Longrightarrow \quad z = y;$$

// algebraic simplifications :-)

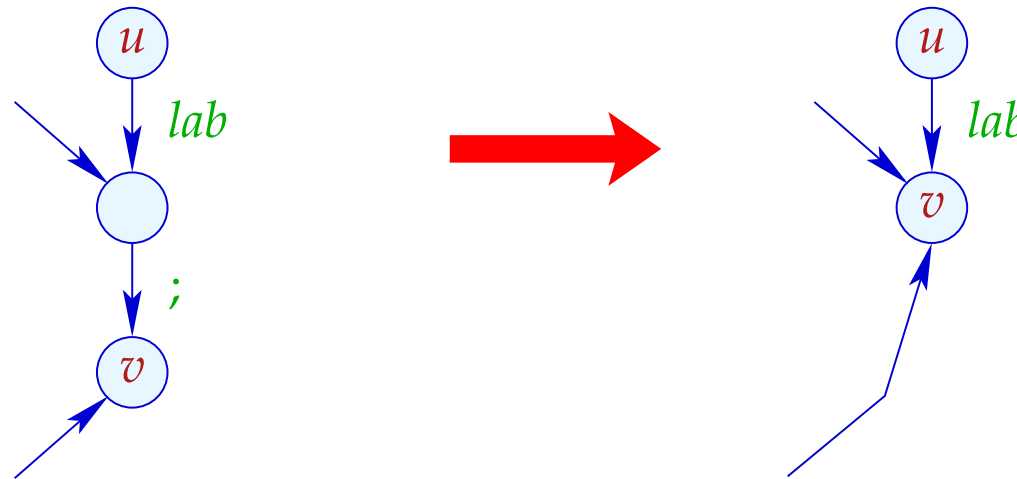
$$x = x; \quad \Longrightarrow \quad ;$$

$$x = 0; \quad \Longrightarrow \quad x = x \oplus x;$$

$$x = 2 \cdot x; \quad \Longrightarrow \quad x = x + x;$$

Important Subproblem:

nop-Optimization



- If $(v_1, ;, v)$ is an edge, v_1 has no further out-going edge.
- Consequently, we can identify v_1 and v :-)
- The ordering of the identifications does not matter :-))

Implementation:

- We construct a function $\text{next} : \text{Nodes} \rightarrow \text{Nodes}$ with:

$$\text{next } u = \begin{cases} \text{next } v & \text{if } (u, ;, v) \text{ edge} \\ u & \text{otherwise} \end{cases}$$

Warning: This definition is only recursive if there are
;-loops ???

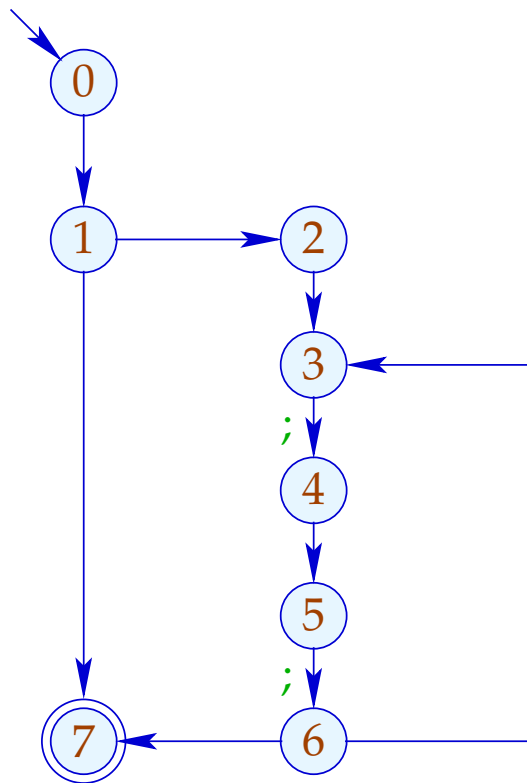
- We replace every edge:

$$(u, \text{lab}, v) \implies (u, \text{lab}, \text{next } v)$$

... whenever $\text{lab} \neq ;$

- All ;-edges are removed ;-)

Example:

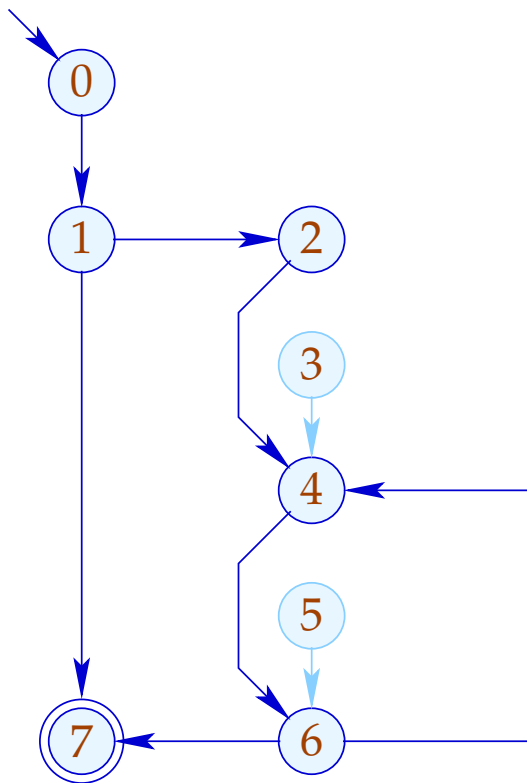


next 1 = 1

next 3 = 4

next 5 = 6

Example:



next 1 = 1

next 3 = 4

next 5 = 6

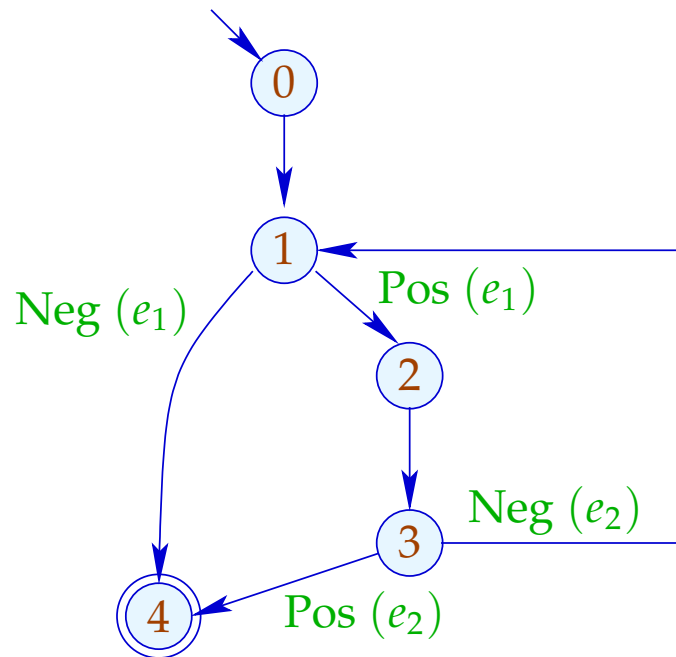
2. Subproblem: Linearization

After optimization, the CFG must again be brought into a **linearly arrangement** of instructions :-)

Warning:

Not every linearization is equally efficient !!!

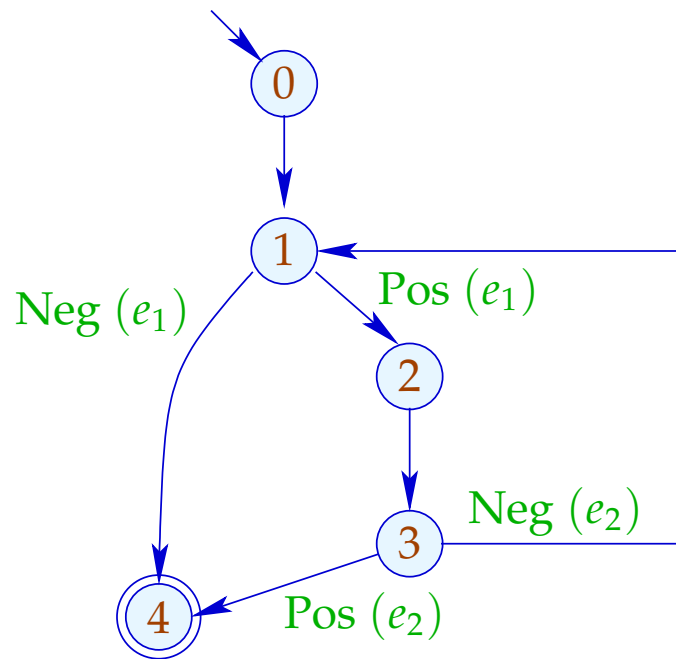
Example:



0:
1: if (e_1) goto 2;
4: halt
2: Rumpf
3: if (e_2) goto 4;
goto 1;

Bad: The loop body is jumped into :-(
:-(

Example:



0:
1: if (! e_1) goto 4;
2: Rumpf
3: if (! e_2) goto 1;
4: halt

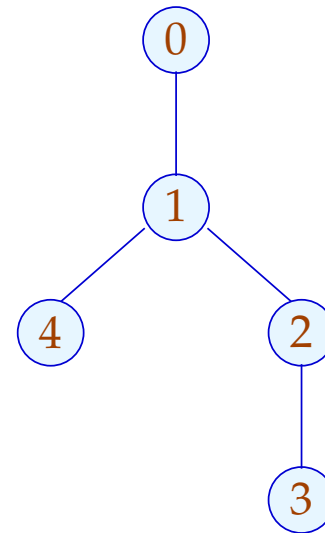
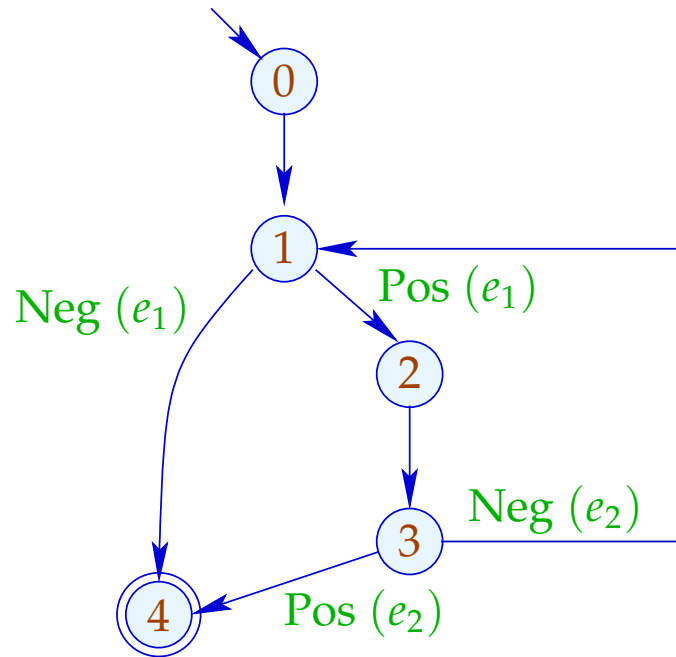
// better cache behavior :-)

Idea:

- Assign to each node a **temperature!**
- always jumps to
 - (1) nodes which have already been handled;
 - (2) **colder** nodes.
- **Temperature** \approx nesting-depth

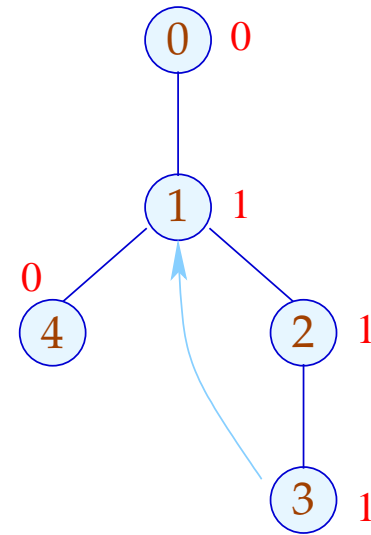
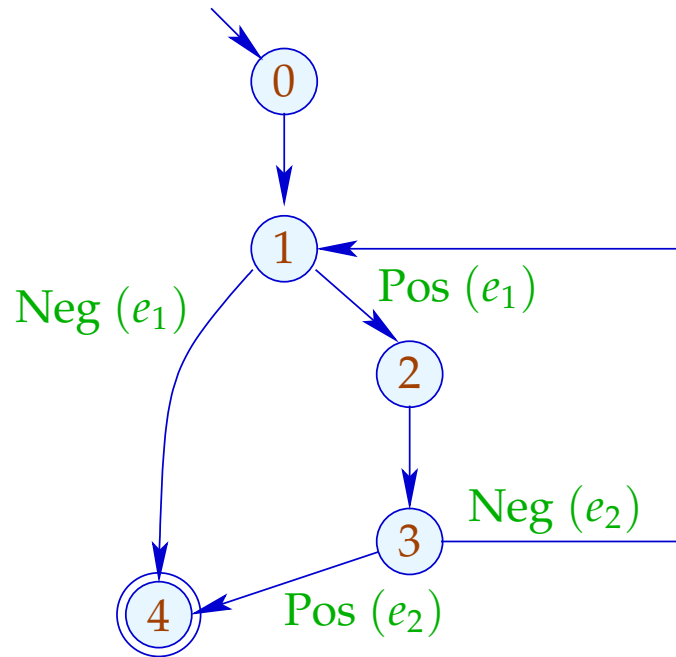
For the computation, we use the pre-dominator tree and strongly connected components ...

... in the Example:

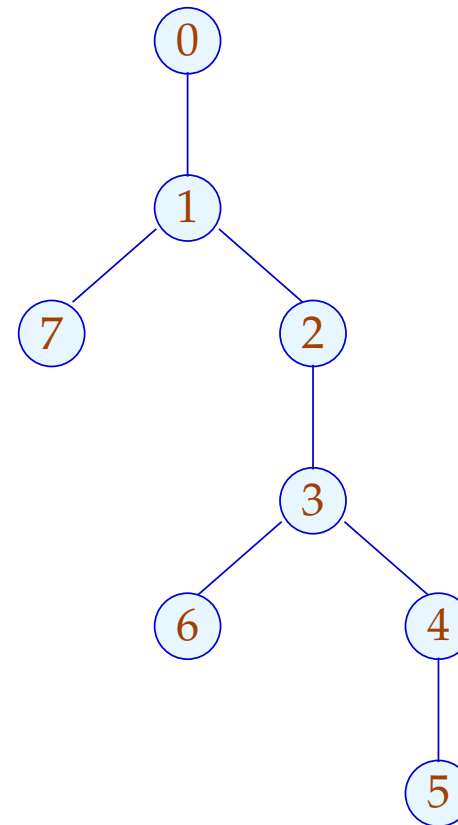
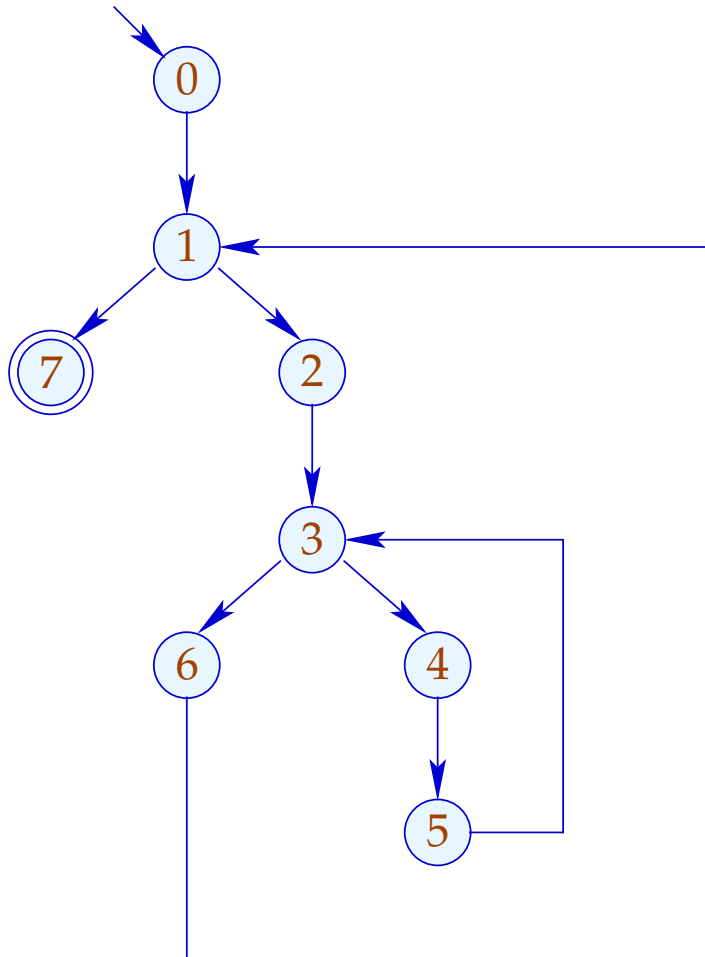


The sub-tree with back edge is **hotter** ...

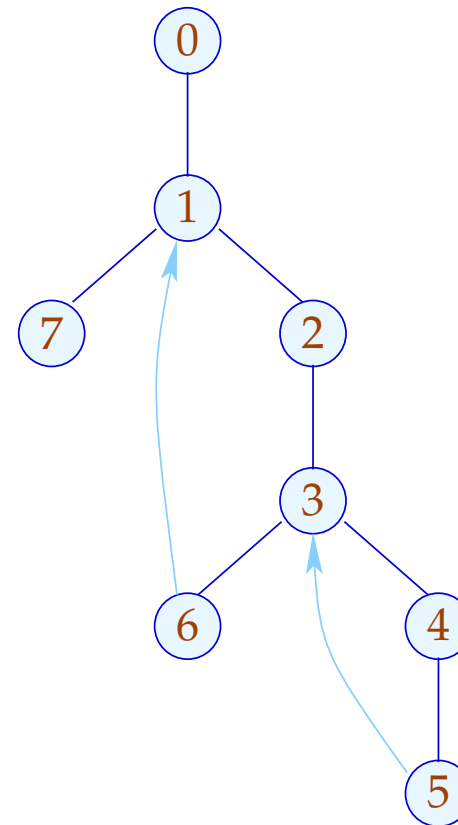
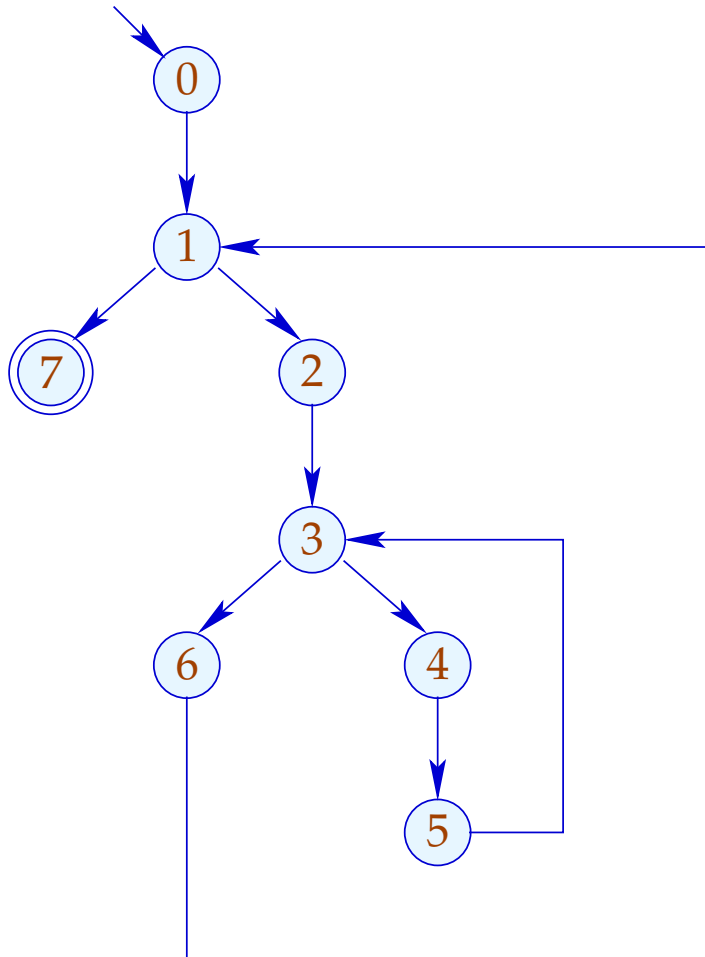
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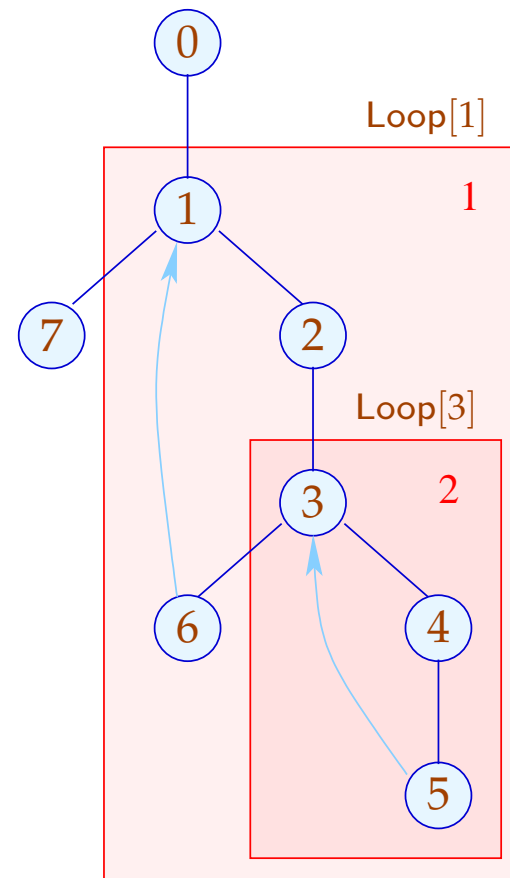
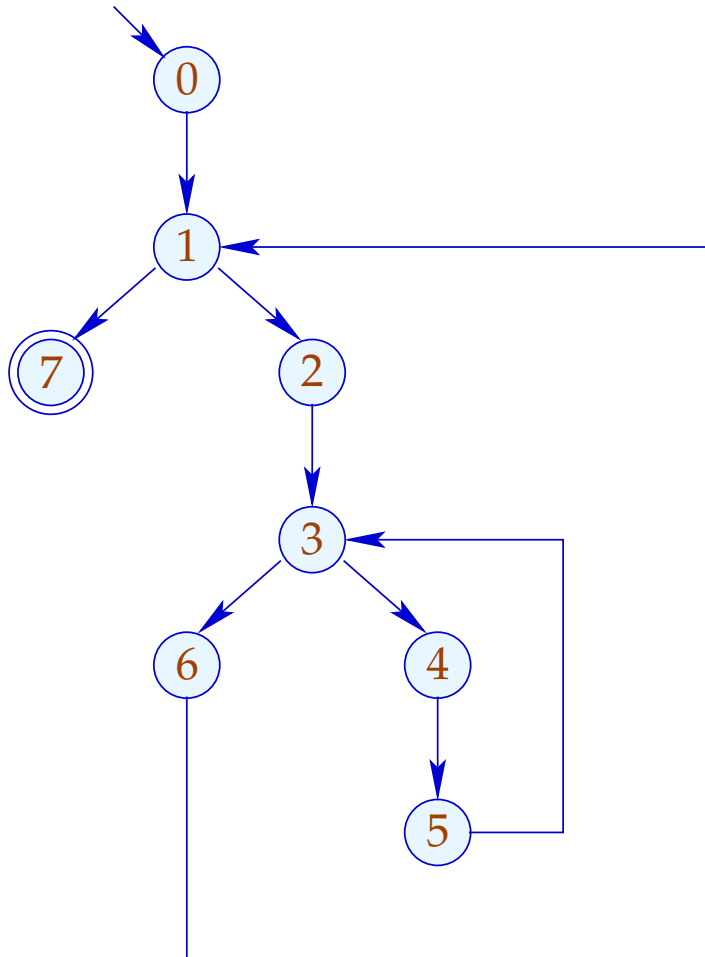
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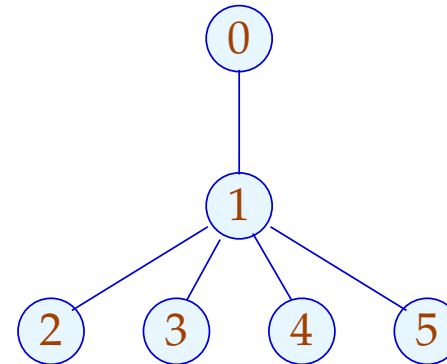
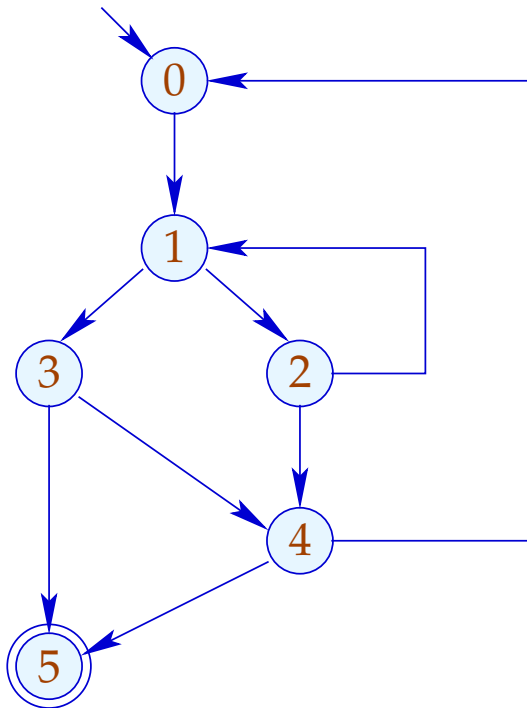


More Complicated Example:



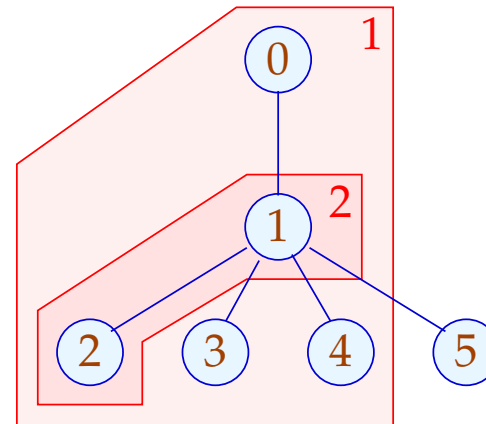
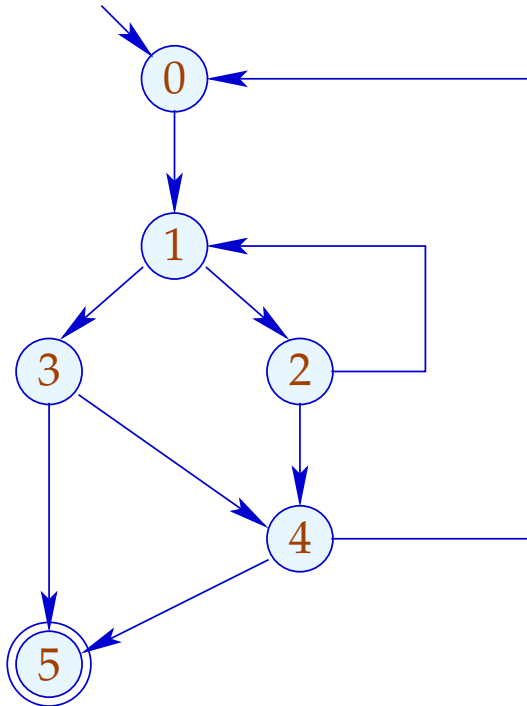
Our definition of `Loop` implies that (detected) loops are necessarily nested :-)

Is is also meaningful for do-while-loops with breaks ...



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Is is also meaningful for do-while-loops with breaks ...



Summary: The Approach

- (1) For every node, determine a temperature;
- (2) Pre-order-DFS over the CFG;
 - If an edge leads to a node we already have generated code for, then we insert a jump.
 - If a node has two successors with different temperature, then we insert a jump to the **colder** of the two.
 - If both successors are equally warm, then it does not matter ;-)

2.3 Procedures

We extend our mini-programming language by procedures without parameters and procedure calls.

For that, we introduce a new statement:

$$f();$$

Every procedure f has a definition:

$$f () \{ stmt^* \}$$

Additionally, we distinguish between **global** and **local** variables.

Program execution starts with the call of a procedure `main ()`.

Example:

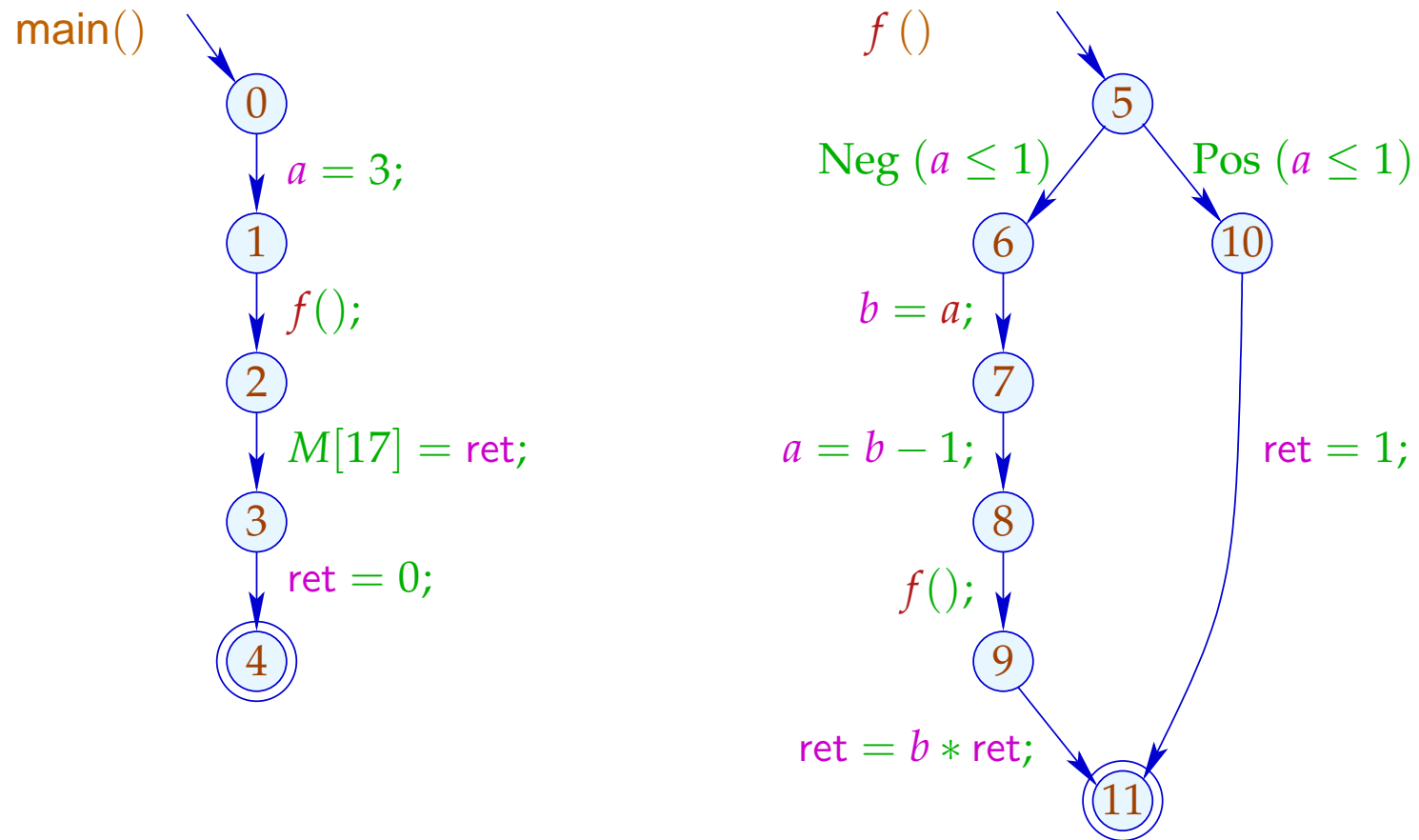
```
int a, ret;
main () {
    a = 3;
    f();
    M[17] = ret;
    ret = 0;
}

f () {
    int b;
    if (a ≤ 1) {ret = 1; goto exit;}
    b = a;
    a = b - 1;
    f();
    ret = b · ret;

    exit :
}
```

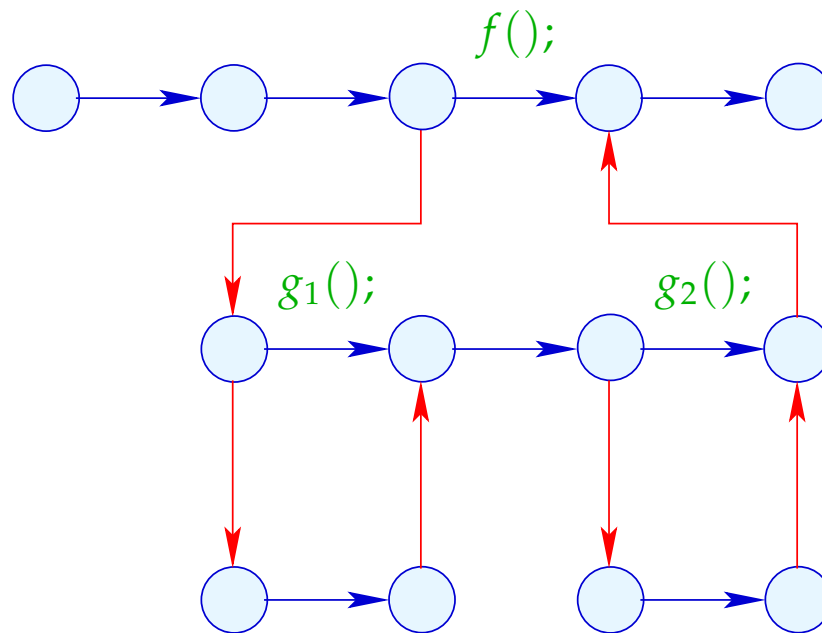
Such programs can be represented by a **set** of CFGs: one for each procedure ...

... in the Example:

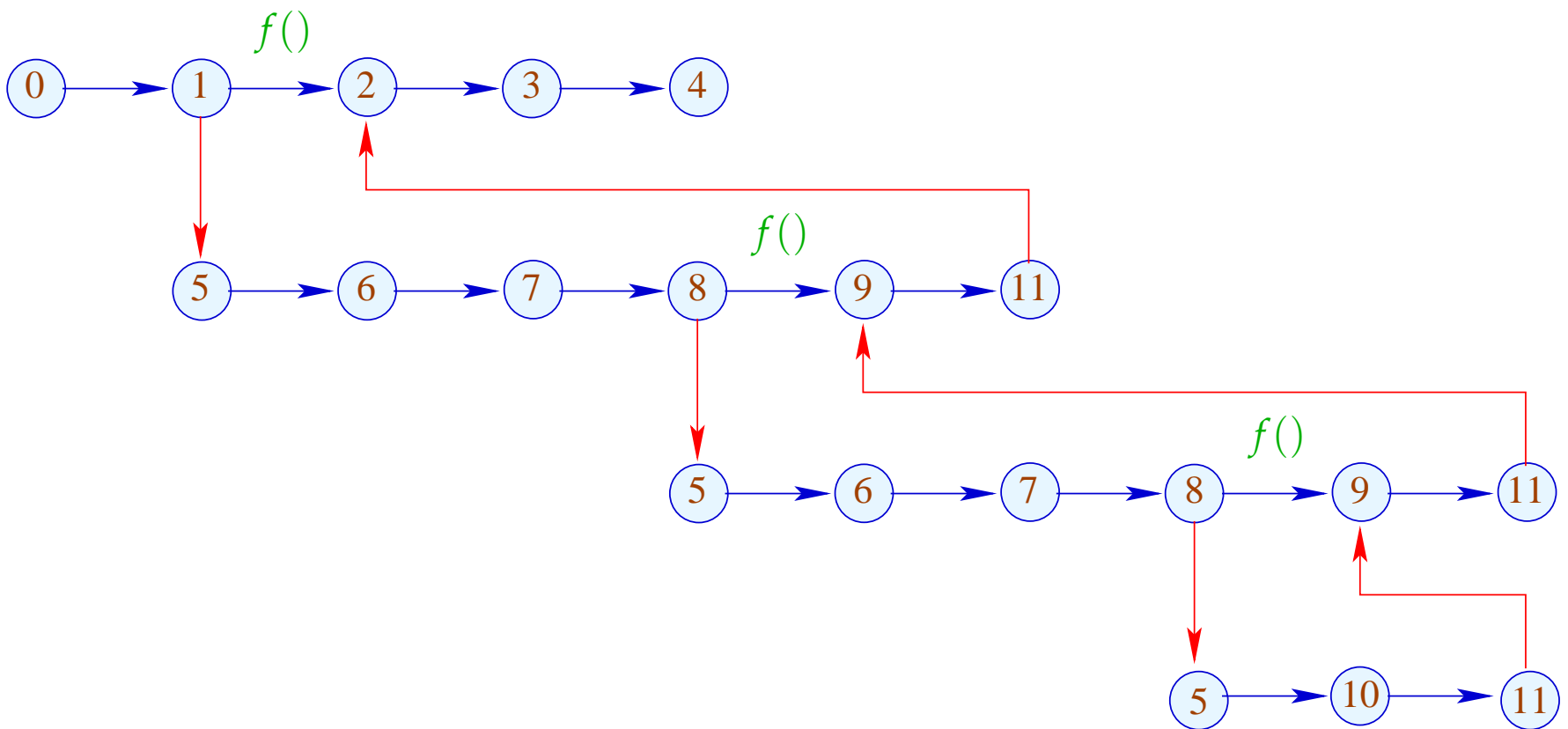


In order to optimize such programs, we require an extended operational semantics *;-)*

Program executions are no longer *paths*, but *forests*:



... in the Example:



The function $\llbracket \cdot \rrbracket$ is extended to computation forests: $w :$

$$\llbracket w \rrbracket : (\text{Vars} \rightarrow \mathbb{Z}) \times (\mathbb{N} \rightarrow \mathbb{Z}) \rightarrow (\text{Vars} \rightarrow \mathbb{Z}) \times (\mathbb{N} \rightarrow \mathbb{Z})$$

For a call $k = (u, f();, v)$ we must:

- determine the initial values for the locals:

$$\text{enter } \rho = \{x \mapsto 0 \mid x \in \text{Locals}\} \oplus (\rho|_{\text{Globals}})$$

- ... combine the new values for the globals with the old values for the locals:

$$\text{combine } (\rho_1, \rho_2) = (\rho_1|_{\text{Locals}}) \oplus (\rho_2|_{\text{Globals}})$$

- ... evaluate the computation forest inbetween:

$$\begin{aligned} \llbracket k \langle w \rangle \rrbracket (\rho, \mu) &= \text{let } (\rho_1, \mu_1) = \llbracket w \rrbracket (\text{enter } \rho, \mu) \\ &\text{in } (\text{combine } (\rho, \rho_1), \mu_1) \end{aligned}$$

Warning:

- In general, $\llbracket w \rrbracket$ is only partially defined :-)
- Dedicated global/local variables a_i, b_i, ret can be used to simulate specific calling conventions.
- The **standard** operational semantics relies on configurations which maintain a **call stack**.
- Computation forests are better suited for the construction of analyses and correctness proofs :-)
- It is an awkward (but useful) exercise to prove the equivalence of the two approaches ...

Configurations:

$$\begin{aligned} \text{configuration} & \quad \equiv \quad \text{stack} \times \text{store} \\ \text{store} & \quad \equiv \quad \text{globals} \times \mathbb{N} \rightarrow \mathbb{Z} \\ \text{locals} & \quad \equiv \quad (\text{Globals} \rightarrow \mathbb{Z}) \\ \text{stack} & \quad \equiv \quad \text{frame} \cdot \text{frame}^* \\ \text{frame} & \quad \equiv \quad \text{point} \times \text{locals} \\ \text{locals} & \quad \equiv \quad (\text{Locals} \rightarrow \mathbb{Z}) \end{aligned}$$

A *frame* specifies the local state of computation inside a procedure call *:-)*

The *leftmost* frame corresponds to the current call.

Computation steps refer to the current call :-)

The novel kinds of steps:

call $k = (u, f();, v) :$

$$\left((u, \rho) \cdot \sigma, \langle \gamma, \mu \rangle \right) \Longrightarrow \left((u_f, \{x \rightarrow 0 \mid x \in \text{Locals}\}) \cdot (v, \rho) \cdot \sigma, \langle \gamma, \mu \rangle \right)$$

u_f entry point of f

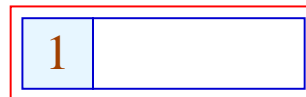
return:

$$\left((r_f, -) \cdot \sigma, \langle \gamma, \mu \rangle \right) \Longrightarrow \left(\sigma, \langle \gamma, \mu \rangle \right)$$

r_f return point of f

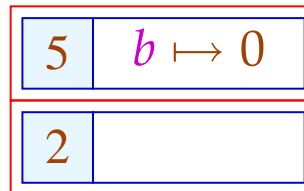
The call stack explicitly implements the DFS traversal through the computation forest :-)

... in the Example:



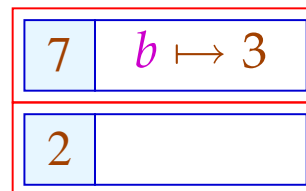
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5	$b \mapsto 0$
9	$b \mapsto 3$
2	

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... in the Example:

7	$b \mapsto 2$
9	$b \mapsto 3$
2	

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5	$b \mapsto 0$
9	$b \mapsto 2$
9	$b \mapsto 3$
2	

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11	$b \mapsto 0$
9	$b \mapsto 2$
9	$b \mapsto 3$
2	

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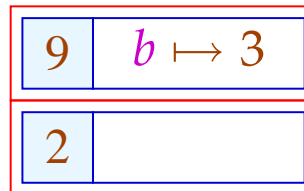
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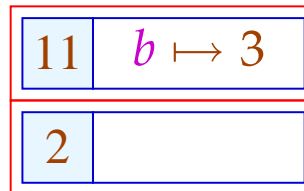
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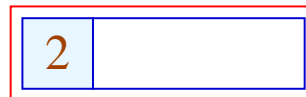
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This operational semantics is quite **realistic** :-)

Costs for a Procedure Call:

Before entering the body: ● Creating a stack frame;

- passing of the parameters;
- Saving the registers;
- Saving the return address;
- Jump to the body.

At procedure exit: ● Freeing the stack frame.

- Restoring the registers.
- Passing of the result.
- Return behind the call.

⇒ ... quite expensive !!!

1. Idea: Inlining

Copy the procedure body at every call site !!!

Example:

```
abs () {  
     $a_2 = -a_1$ ;  
    max ();  
}  
  
max () {  
    if ( $a_1 < a_2$ ) {  $ret = a_2$ ; goto _exit; }  
     $ret = a_1$ ;  
    _exit :  
}
```

... yields:

```
abs () {  
   $a_2 = -a_1$ ;  
  if ( $a_1 < a_2$ ) { ret =  $a_2$ ; goto _exit; }  
  ret =  $a_1$ ;  
  _exit :  
}
```

Problems:

- The copied block may modify the locals of the calling procedure ???
- More general: Multiple use of local variable names may lead to errors.
- Multiple calls of a procedure may lead to code duplication :-((
- How can we handle **recursion** ???

Detection of Recursion:

We construct the **call-graph** of the program.

In the Examples:

