



EECS 483: Compiler Construction

Lecture 18: Optimization and Dataflow Analysis

**March 23
Winter Semester 2025**

Slides adapted from Steve Zdancewic

Announcements

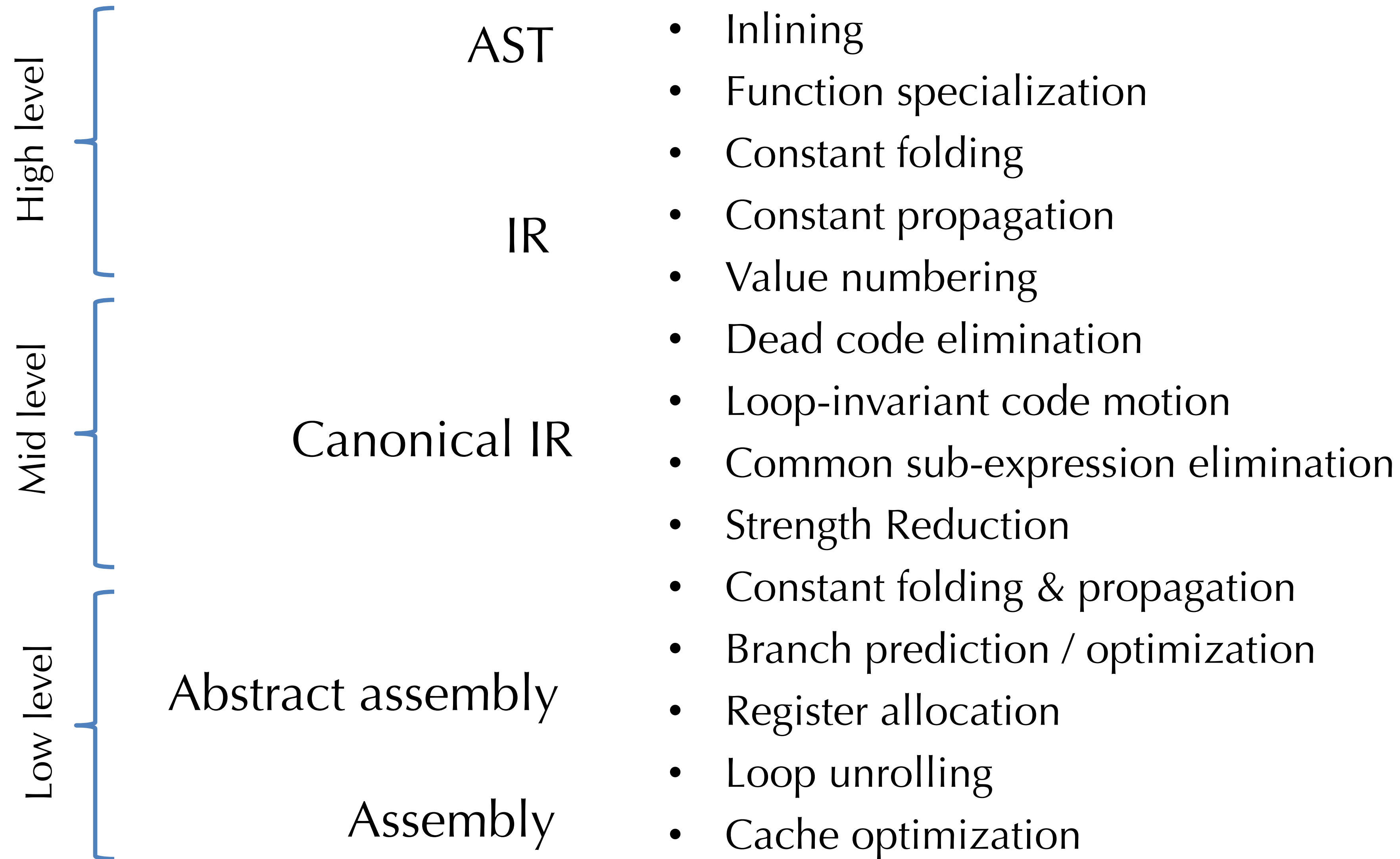
- Exam Grading almost done
- Assignment 4 due next Friday, April 4



Why optimize?

OPTIMIZATIONS, GENERALLY

When to apply optimization



Safety

- Whether an optimization is *safe* depends on the programming language semantics.
 - Languages that provide weaker guarantees to the programmer permit more optimizations but have more ambiguity in their behavior.
 - e.g., In C, loading from initialized memory is undefined, so the compiler can do anything if a program reads uninitialized data.
 - e.g., In Java tail-call optimization (which turns recursive function calls into loops) is not valid because of “stack inspection”.
- Example: *loop-invariant code motion*
 - Idea: hoist invariant code out of a loop

```
while (b) {  
  z = y/x;  
  ...  
}
```

// y, x not updated



```
z = y/x;  
while (b) {  
  ...  
}
```

// y, x not updated

- Is this more efficient?
- Is this safe?



A high-level tour of a variety of optimizations.

BASIC OPTIMIZATIONS

Constant Folding

- Idea: If operands are known at compile time, perform the operation statically.

`int x = (2 + 3) * y` → `int x = 5 * y`

`b & false` → `false`

- Performed at every stage of optimization...
- Why?
 - Constant expressions can be created by translation or earlier optimizations

Example: `A[2]` might be compiled to:

`MEM[MEM[A] + 2 * 4]` → `MEM[MEM[A] + 8]`

Constant Folding Conditionals

if (true) S \rightarrow S

if (false) S \rightarrow ;

if (true) S else S' \rightarrow S

if (false) S else S' \rightarrow S'

while (false) S \rightarrow ;

if (2 > 3) S \rightarrow

if (false) S \rightarrow ;

Algebraic Simplification

- More general form of constant folding
 - Take advantage of mathematically sound simplification rules
- *Mathematical identities:*
 - $a * 1 \rightarrow a$ $a * 0 \rightarrow 0$
 - $a + 0 \rightarrow a$ $a - 0 \rightarrow a$
 - $b \mid \text{false} \rightarrow b$ $b \& \text{true} \rightarrow b$
- *Reassociation & commutativity:*
 - $(a + 1) + 2 \rightarrow a + (1 + 2) \rightarrow a + 3$
 - $(2 + a) + 4 \rightarrow (a + 2) + 4 \rightarrow a + (2 + 4) \rightarrow a + 6$
- *Strength reduction:* (replace expensive op with cheaper op)
 - $a * 4 \rightarrow a \ll 2$
 - $a * 7 \rightarrow (a \ll 3) - a$
 - $a / 32767 \rightarrow (a \gg 15) + (a \gg 30)$
- *Note 1:* must be careful with floating point (due to rounding) and integer arithmetic (due to overflow/underflow)
- *Note 2:* iteration of these optimizations is useful... how much?
- *Note 3:* must be sure that rewrites terminate:
 - commutativity apply like: $(x + y) \rightarrow (y + x) \rightarrow (x + y) \rightarrow (y + x) \rightarrow \dots$

Constant Propagation

- If a variable is known to be a constant, replace the use of the variable by the constant
- Value of the variable must be propagated forward from the point of assignment
 - This is a *substitution* operation

Example:

```
int x = 5;  
int y = x * 2;  
int z = a[y];
```

 →

```
int y = 5 * 2;  
int z = a[y];
```

 →

```
int y = 10;  
int z = a[y];
```

 →

```
int z = a[10];
```

- To be most effective, constant propagation should be interleaved with constant folding

Copy Propagation

- If one variable is assigned to another, replace uses of the assigned variable with the copied variable.
- Need to know where copies of the variable propagate.
- Interacts with the scoping rules of the language.

- Example:

```
x = y;  
if (x > 1) {  
    x = x * f(x - 1);  
}
```

→

```
x = y;  
if (y > 1) {  
    x = y * f(y - 1);  
}
```

- Can make the first assignment to x **dead code** (that can be eliminated).

Dead Code Elimination

- If a side-effect free statement can never be observed, it is safe to eliminate the statement.

```
x = y * y // x is dead!  
...      // x never used  
x = z * z
```

```
→      ...  
x = z * z
```

- A variable is **dead** if it is never used after it is defined.
 - Computing such *definition* and *use* information is an important component of program analysis
- Dead variables can be created by other optimizations...

Unreachable/Dead Code

- Basic blocks not reachable by any trace leading from the starting basic block are *unreachable* and can be deleted.
 - Performed at the IR or assembly level
 - Improves cache, TLB performance
- Dead code: similar to unreachable blocks.
 - A value might be computed but never subsequently used.
- Code for computing the value can be dropped
- But only if it's **pure**, i.e., it has *no externally visible side effects*
 - Externally visible effects: raising an exception, modifying a global variable, going into an infinite loop, printing to standard output, sending a network packet, launching a rocket
 - Note: Pure functional languages (e.g., Haskell) make reasoning about the safety of optimizations (and code transformations in general) easier!

Inlining

- Replace a call to a function with the body of the function itself with arguments rewritten to be local variables:
- Example in C: inline `pow` into `g`

```
int g(int x) { return x + pow(x); }
int pow(int a) {
    var b = 1; var x = 0;
    while (x < a) {b = 2 * b; x = x + 1}
    return b;
}
```



```
int g(int x) {
    int a = x;
    int b = 1; int x2 = 0;
    while (x2 < a) {b = 2 * b; x2 = x2 + 1};
    tmp = b;
    return x + tmp;
}
```

note: renaming

- May need to rename variables to avoid *capture*
- Best done at the AST or relatively high-level IR.
- When is it profitable?
 - Eliminates the stack manipulation, jump, etc.
 - Can increase code size.
 - Enables further optimizations

Code Specialization

- Idea: create specialized versions of a function that is called from different places with different arguments.
- Example: specialize function `f` in:

```
class A implements I { int m() {...} }  
class B implements I { int m() {...} }  
int f(I x) { x.m(); }           // don't know which m  
A a = new A(); f(a);           // know it's A.m  
B b = new B(); f(b);           // know it's B.m
```

- `f_A` would have code specialized to dispatch to `A.m`
- `f_B` would have code specialized to dispatch to `B.m`
- You can also inline methods when the run-time type is known statically
 - Often just one class implements a method.

Common Subexpression Elimination

- *fold redundant computations together*
 - in some sense, it's the opposite of inlining
- Example:

```
a[i] = a[i] + 1
```

compiles to:

```
[a + i*4] = [a + i*4] + 1
```

Common subexpression elimination removes the redundant add and multiply:

```
t = a + i*4; [t] = [t] + 1
```

- For safety, you must be sure that the shared expression always has the same value in both places!

Unsafe Common Subexpression Elimination

- Example: consider this C function:

```
unit f(int[] a, int[] b, int[] c) {  
    var j = ...; var i = ...; var k = ...;  
    b[j] = a[i] + 1;  
    c[k] = a[i];  
    return;  
}
```

- The optimization that shares the expression `a[i]` is unsafe... why?

```
unit f(int[] a, int[] b, int[] c) {  
    var j = ...; var i = ...; var k = ...;  
    t = a[i];  
    b[j] = t + 1;  
    c[k] = t;  
    return;  
}
```



LOOP OPTIMIZATIONS

Loop Optimizations

- Program hot spots often occur in loops.
 - Especially inner loops
 - Not always: consider operating systems code or compilers vs. a computer game or word processor
- Most program execution time occurs in loops.
 - The 90/10 rule of thumb holds here too.
(90% of the execution time is spent in 10% of the code)
- Loop optimizations are very important, effective, and numerous
 - Also, concentrating effort to improve loop body code is usually a win

Loop Invariant Code Motion (revisited)

- Another form of redundancy elimination.
- If the result of a statement or expression does not change during the loop *and* it's pure, it can be hoisted outside the loop body.
- Often useful for array element addressing code
 - Invariant code not visible at the source level

```
for (i = 0; i < a.length; i++) {  
    /* a not modified in the body */  
}
```



```
t = a.length;  
for (i = 0; i < t; i++) {  
    /* same body as above */  
}
```

Hoisted loop-
invariant
expression

Strength Reduction (revisited)

- Strength reduction can work for loops too
- Idea: replace expensive operations (multiplies, divides) by cheap ones (adds and subtracts)
- For loops, create a *dependent induction variable*:
- Example:

```
for (int i = 0; i < n; i++) { a[i*3] = 1; } // stride by 3
```



```
int j = 0;  
for (int i = 0; i < n; i++) {  
    a[j] = 1;  
    j = j + 3; // replace multiply by add  
}
```

Loop Unrolling (revisited)

- Branches can be expensive, unroll loops to avoid them.

```
for (int i=0; i<n; i++) { S }
```



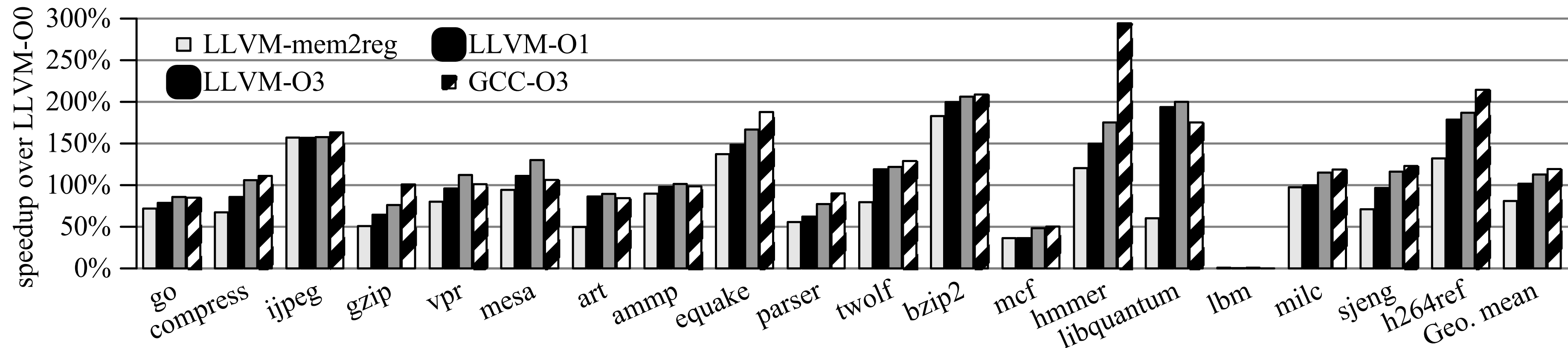
```
for (int i=0; i<n-3; i+=4) {S;S;S;S};  
for (      ; i<n; i++) { S } // left over iterations
```

- With k unrollings, eliminates $(k-1)/k$ conditional branches
 - So for the above program, it eliminates $3/4$ of the branches
- Space-time tradeoff:
 - Not a good idea for large S or small n
- Interacts with instruction caching, branch prediction



EFFECTIVENESS?

Optimization Effectiveness?



$$\% \text{speedup} = \left[\frac{\text{base time}}{\text{optimized time}} - 1 \right] \times 100\%$$

Example:

base time = 2s

optimized time = 1s

⇒ 100% speedup

Example:

base time = 1.2s

optimized time = 0.87s

⇒ 38% speedup

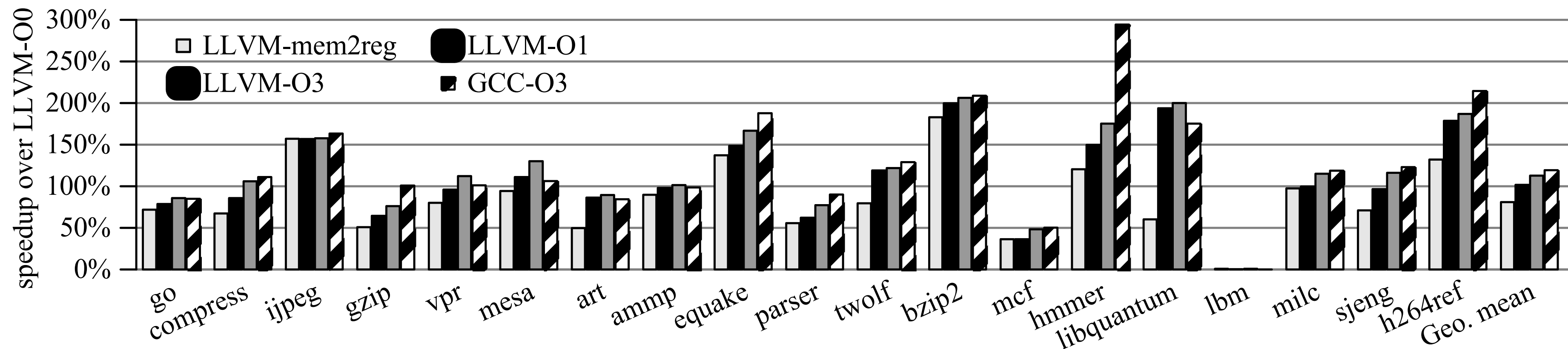
Graph taken from:

Jianzhou Zhao, Santosh Nagarakatte, Milo M. K. Martin, and Steve Zdancewic.

Formal Verification of SSA-Based Optimizations for LLVM.

In Proc. 2013 ACM SIGPLAN Conference on Programming Languages Design and Implementation (PLDI), 2013

Optimization Effectiveness?



- mem2reg: promotes alloca'd stack slots to temporaries to enable register allocation
- Analysis:
 - mem2reg alone (+ back-end optimizations like register allocation) yields ~78% speedup on average
 - -O1 yields ~100% speedup (so all the rest of the optimizations combined account for ~22%)
 - -O3 yields ~120% speedup
- Hypothetical program that takes 10 sec. (base time):
 - Mem2reg alone: expect ~5.6 sec
 - -O1: expect ~5 sec
 - -O3: expect ~4.5 sec



CODE ANALYSIS

Motivating Code Analyses

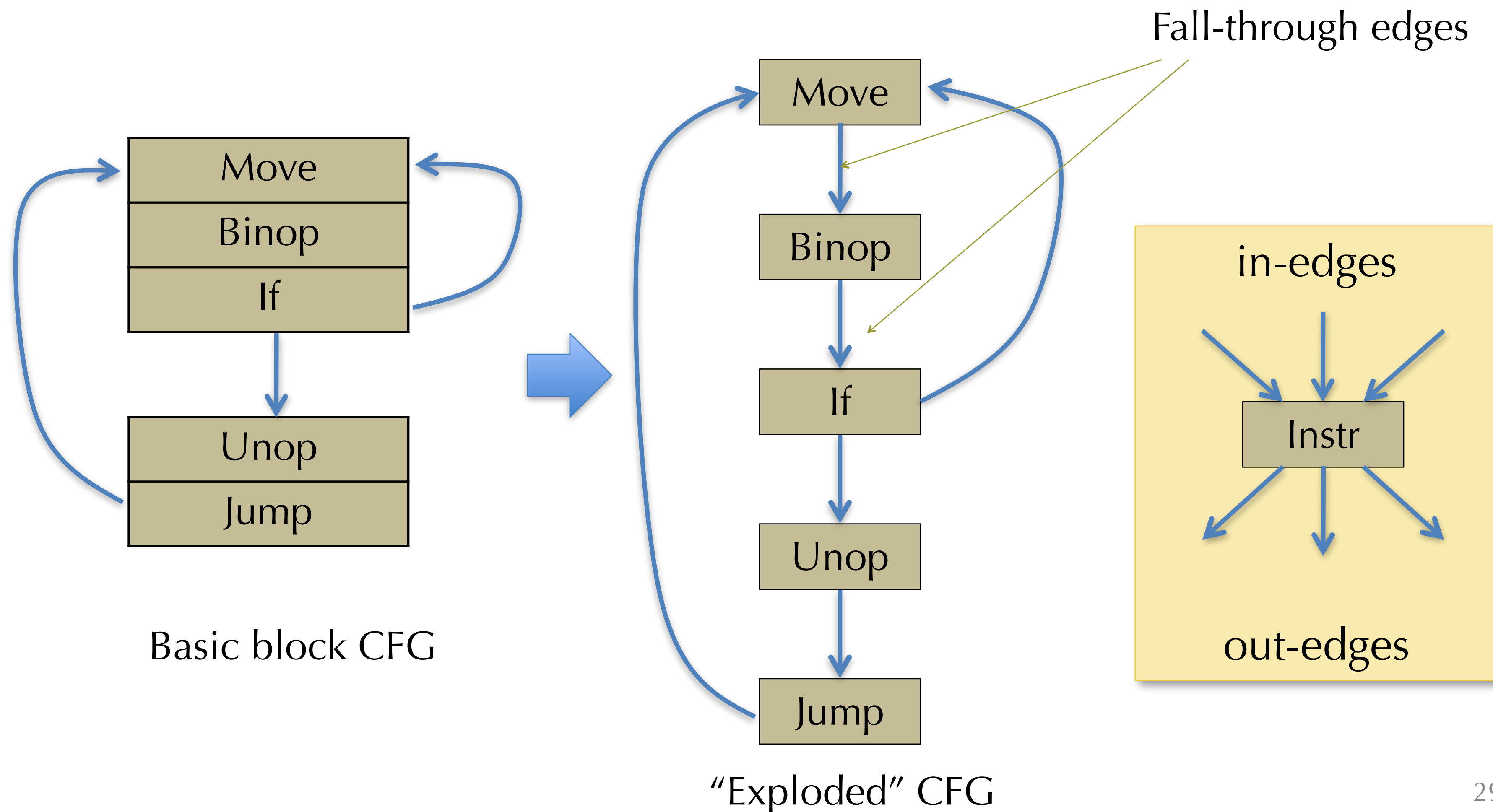
- There are lots of things that might influence the safety/applicability of an optimization
 - What algorithms and data structures can help?
- How do you know what is a loop?
- How do you know an expression is invariant?
- How do you know if an expression has no side effects?
- How do you keep track of where a variable is defined?
- How do you know where a variable is used?
- How do you know if two reference values may be aliases of one another?

Control-flow Graphs Revisited

- For the purposes of dataflow analysis, we use the *control-flow graph* (CFG) intermediate form.
- Recall that a basic block is a sequence of instructions such that:
 - There is a distinguished, labeled entry point (no jumps into the middle of a basic block)
 - There is a (possibly empty) sequence of non-control-flow instructions
 - The block ends with a single control-flow instruction (jump, conditional branch, return, etc.)
- A *control flow graph*
 - Nodes are blocks
 - There is an edge from B1 to B2 if the control-flow instruction of B1 might jump to the entry label of B2
 - There are no “dangling” edges – there is a block for every jump target.
- Note: the following slides are intentionally a bit ambiguous about the exact nature of the code in the control flow graphs:
 - at the x86 assembly level
 - an “imperative” C-like source level
 - at the LLVM IR level
 - Same general idea, but the exact details will differ
 - e.g. LLVM IR doesn’t have “imperative” update of %uid temporaries.
 - In fact, the SSA structure of the LLVM IR makes some of these analyses simpler.

Dataflow over CFGs

- For precision, it is helpful to think of the “fall through” between sequential instructions as an edge of the control-flow graph too.
 - Different implementation tradeoffs in practice...

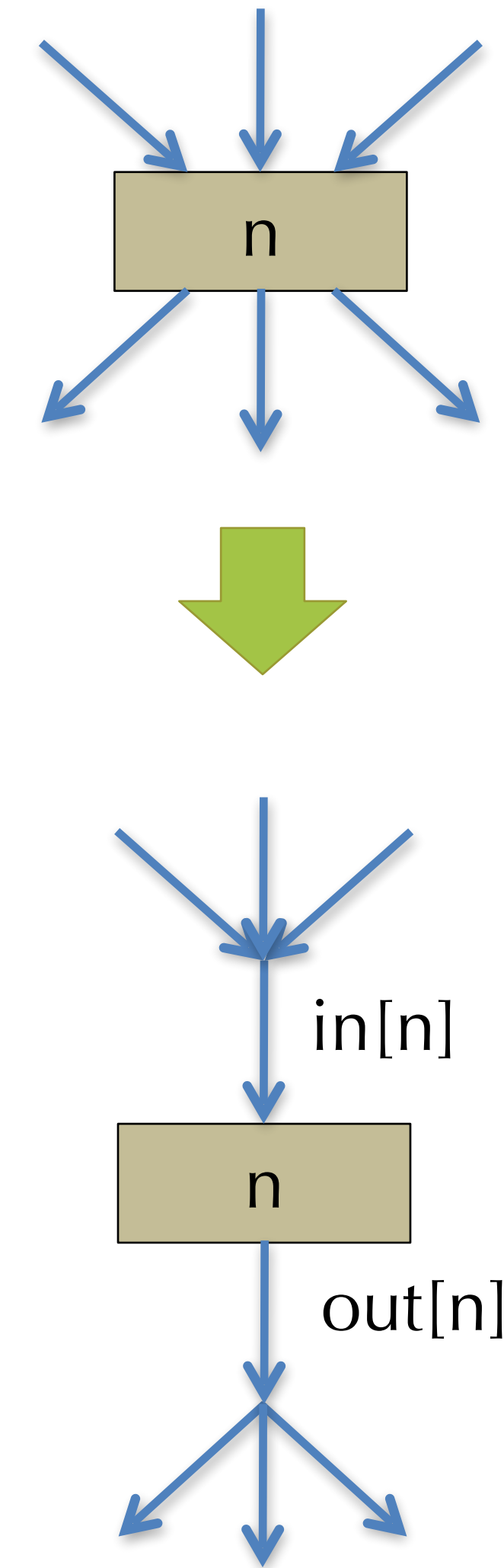


Dataflow Analysis

- *Idea*: compute liveness information for all variables simultaneously.
 - Keep track of sets of information about each node
- Approach: define *equations* that must be satisfied by any liveness determination.
 - Equations based on “obvious” constraints.
- Solve the equations by iteratively converging on a solution.
 - Start with a “rough” approximation to the answer
 - Refine the answer at each iteration
 - Keep going until no more refinement is possible: a *fixpoint* has been reached
- This is an instance of a general framework for computing program properties: dataflow analysis

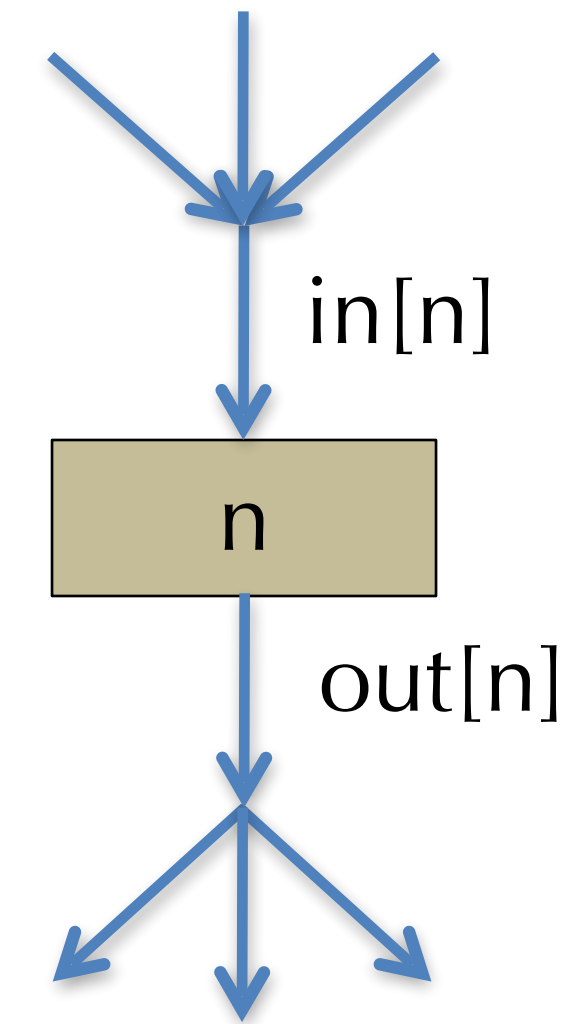
Dataflow Value Sets for Liveness

- Nodes are program statements, so:
- $use[n]$: set of variables used by n
- $def[n]$: set of variables defined by n
- $in[n]$: set of variables live on entry to n
- $out[n]$: set of variables live on exit from n
- Associate $in[n]$ and $out[n]$ with the “collected” information about incoming/outgoing edges
- For Liveness: what constraints are there among these sets?
- Clearly:
$$in[n] \supseteq use[n]$$
- What other constraints?



Other Dataflow Constraints

- We have: $\text{in}[n] \supseteq \text{use}[n]$
 - “A variable must be live on entry to n if it is used by n ”
- Also: $\text{in}[n] \supseteq \text{out}[n] - \text{def}[n]$
 - “If a variable is live on exit from n , and n doesn't define it, it is live on entry to n ”
 - Note: here '-' means “set difference”
- And: $\text{out}[n] \supseteq \text{in}[n']$ if $n' \in \text{succ}[n]$
 - “If a variable is live on entry to a successor node of n , it must be live on exit from n .”



Iterative Dataflow Analysis

- Find a solution to those constraints by starting from a rough guess.
 - Start with: $\text{in}[n] = \emptyset$ and $\text{out}[n] = \emptyset$
- The guesses don't satisfy the constraints:
 - $\text{in}[n] \supseteq \text{use}[n]$
 - $\text{in}[n] \supseteq \text{out}[n] - \text{def}[n]$
 - $\text{out}[n] \supseteq \text{in}[n']$ if $n' \in \text{succ}[n]$
- Idea: iteratively re-compute $\text{in}[n]$ and $\text{out}[n]$ where forced to by the constraints.
 - Each iteration will add variables to the sets $\text{in}[n]$ and $\text{out}[n]$ (i.e. the live variable sets will increase monotonically)
- We stop when $\text{in}[n]$ and $\text{out}[n]$ satisfy these equations: (which are derived from the constraints above)
 - $\text{in}[n] = \text{use}[n] \cup (\text{out}[n] - \text{def}[n])$
 - $\text{out}[n] = \bigcup_{n' \in \text{succ}[n]} \text{in}[n']$
 - Most SSA constructs: 1 successor
 - ret has 0 successors: its out set is empty
 - cbr has 2 successors: its out set is the **union** of the two out sets

Complete Liveness Analysis Algorithm

```
for all n, in[n] :=  $\emptyset$ , out[n] :=  $\emptyset$ 
repeat until no change in 'in' and 'out'
  for all n
    out[n] :=  $\bigcup_{n' \in \text{succ}[n]} \text{in}[n']$ 
    in[n] := use[n]  $\cup$  (out[n] - def[n])
  end
end
```

- Finds a *fixpoint* of the **in** and **out** equations.
 - The algorithm is guaranteed to terminate... Why?
- Why do we start with \emptyset ?

“Classic” Constant Propagation

- Constant propagation can be formulated as a dataflow analysis.

- Idea: propagate and fold integer constants in one pass:

x = 1;	→	x = 1;
y = 5 + x;		y = 6;
z = y * y;		z = 36;

- Information about a single variable:
 - Variable is never defined.
 - Variable has a single, constant value.
 - Variable is assigned multiple values.

Assertion Removal

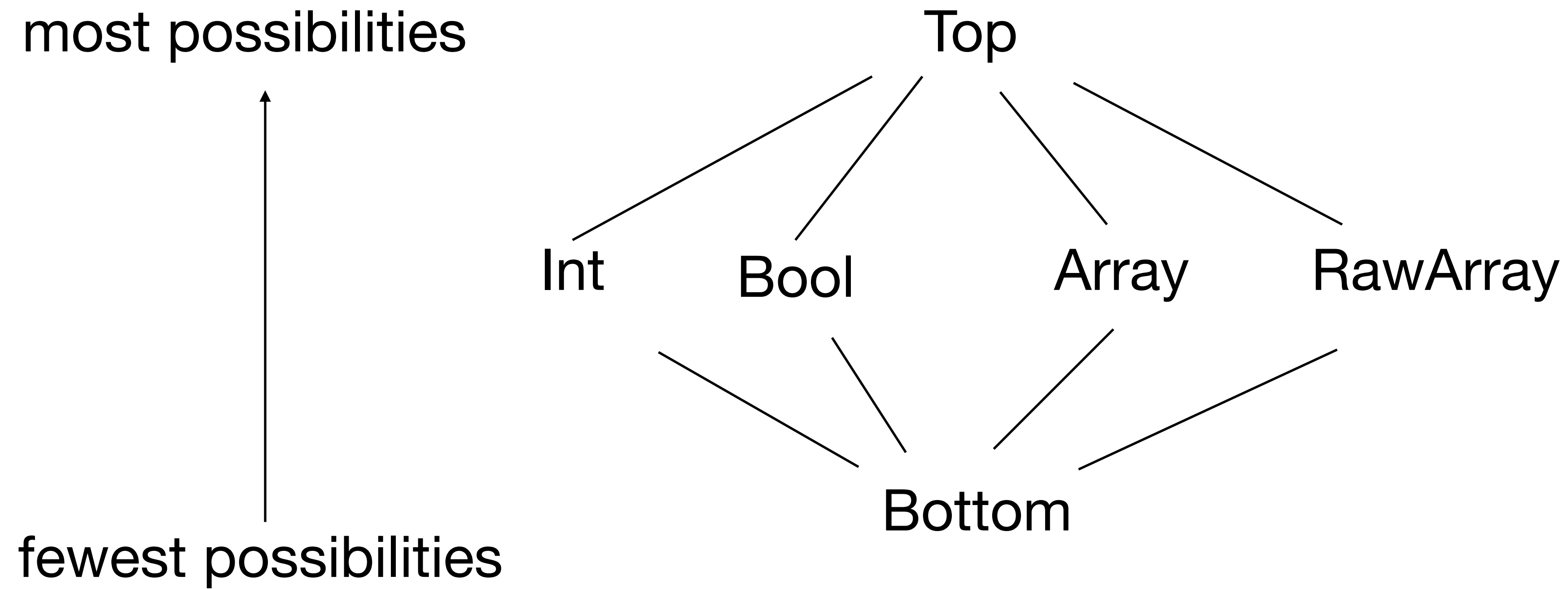
- Dynamic typing adds many runtime assertions into our program.
- ```
let x = g() in
let y = x + 2 in
let z = y * x in
...
```
- Current compilation always adds assertions that inputs are integers
- ```
x = f()
assertInt(x)
y = x + 2
assertInt(y)
assertInt(x)
y2 = y >> 1
z = y2 * x
...
```
- Which assertions can we remove?

Tag-checking Analysis

- At each program point, for each variable associate an approximation of what the possible values are:
 - Int: tagged integer, i.e., multiple of 2
 - Bool: tagged boolean, i.e., either 0b001 or 0b101
 - RawArray: untagged pointer to an array on the heap
 - Array: tagged array, i.e., a pointer tagged with 0b11
 - Top: any 64 bit value
 - Bottom: never assigned to, i.e., uninitialized
- Usage: If analysis determines x is an Int, then remove assertions `assertInt(x)`

similar for `assertArray`, `assertBool` etc.

Tag-checking Analysis



Straightline Code Example

```
x = f()  
assertInt(x)  
y = x + 2  
assertInt(y)  
assertInt(x)  
y2 = y >> 1  
z = y2 * x
```

Tag-checking Analysis

- For each operation in SSA, need to define "flow function" that says what possible tags are based on inputs.

Examples:

- $x = y + z$
 - if y and z are tagged Ints, then x is a tagged Int
 - otherwise x is Top
- $x = y * z$
 - if y or z is a tagged Int then x is a tagged Int
 - otherwise Top
- $x = y << n$
 - if n is at least 1 then x is tagged
 - if n is 0, then x is tagged if y is
- `assertInt(x)`
 - after this, x is always a tagged Int, because otherwise execution ended

Straightline Code Example

```
0 x = f()  
1 assertInt(x)  
2 y = x + 2  
3 assertInt(y)  
4 assertInt(x)  
5 y2 = y >> 1  
6 z = y2 * x  
7
```

```
0:  
1: {x: Top}  
2: {x: Int}  
3: {x: Int, y: Int}  
4: {x: Int, y: Int}  
5: {x: Int, y: Int}  
6: {x: Int, y: Int, y2: Top}  
7: {x: Int, y: Int, y2: Top, z: Int}
```

Straightline Code Example

```
0 x = f()  
1 assertInt(x)  
2 y = x + 2  
3 assertInt(y)  
4 assertInt(x)  
5 y2 = y >> 1  
6 z = y2 * x  
7
```

```
0:  
1: {x: Top}  
2: {x: Int}  
3: {x: Int, y: Int}  
4: {x: Int, y: Int}  
5: {x: Int, y: Int}  
6: {x: Int, y: Int, y2: Top}  
7: {x: Int, y: Int, y2: Top, z: Int}
```

Loop Example

```
extern g
def f(y,z):
    def loop(i,a):
        if i == 0:
            a - z
        else:
            loop(i - 1, a + g())
    in
    loop(y, 0)
```

```
f(y,z):
    loop(i,a):
        thn():
            assertInt(a)
            assertInt(z)
            r = a - z
            ret r
        els():
            assertInt(i)
            i' = i - 2
            x = g()
            assertInt(a)
            assertInt(x)
            a' = a + x
            br loop(i', a')
    b = i == 0
    cbr b thn() els()
    br loop(y, 0)
```

Loop Example

```
f(y,z):  
  loop(i,a):  
    thn():  
      assertInt(a)  
      assertInt(z)  
      r = a - z  
      ret r  
    els():  
      assertInt(i)  
      i' = i - 2  
      x = g()  
      assertInt(a)  
      assertInt(x)  
      a' = a + x  
      br loop(i', a')  
  b = i == 0  
  cbr b thn() els()  
br loop(y, 0)
```

What information do we use in internal blocks such as **loop**, **thn**, **els**?

These are **recursive**, same problem as liveness

Same solution: initialize their info to bottom, iteratively update them

Main difference: need to compute info based on **predecessors** of the block

What about top-level functions like **f**?

Initialize its inputs to **top** as they can be called with any snake values

In Diamondback, there is only one top-level function, **main(x)**, and its input is an **Array**.

Tag-checking Analysis

- When analysing a block $f(x_1, \dots)$: what information do we know about the parameters?
 - The value of x_1 could be the value of any a_1 for any $\text{br } f(a_1, \dots)$ in the program. These are the "predecessors" of f .
 - information for a parameter x_1 should be the best approximation of the information known for all instantiations a_1
 - roughly the "union" of all possible values it could be, i.e., the "least upper bound"
- Examples:
 - $f(x)$: ...
 - $\text{br } f(a)$ and $\text{br } f(b)$ where $a: \text{Int}$ and $b: \text{Int}$ then $x: \text{Int}$
 - $\text{br } f(a)$ and $\text{br } f(b)$ where $a: \text{Int}$ and $b: \text{Bool}$ then $x: \text{Top}$
 - generally: if all predecessors agree on the tag, then use that tag, if any disagree, use Top

```
f(y,z):  
  loop(i,a):  
    thn():  
      3 assertInt(a)  
      4 assertInt(z)  
      5 r = a - z  
      6 ret r  
    els():  
      7 assertInt(i)  
      8 i' = i - 2  
      9 x = g()  
     10 assertInt(a)  
     11 assertInt(x)  
     12 a' = a + x  
     13 br loop(i', a')  
  1 b = i == 0  
  2 cbr b thn() els()  
  0 br loop(y, 0)
```

f(y,z):

unmentioned vars: bottom

loop(i,a):

thn():

3 assertInt(a)

4 assertInt(z)

5 r = a - z

6 ret r

els():

7 assertInt(i)

8 i' = i - 2

9 x = g()

10 assertInt(a)

11 assertInt(x)

12 a' = a + x

13 br loop(i', a')

1 b = i == 0

2 cbr b thn() els()

0 br loop(y, 0)

0: {y:Top,z:Top}

1: bottom

2:

3: bottom

4:

5:

6:

7: bottom

8:

9:

10:

11:

12:

13:

f(y,z):

unmentioned vars: bottom

loop(i,a):

thn():

3 assertInt(a)

4 assertInt(z)

5 r = a - z

6 ret r

els():

7 assertInt(i)

8 i' = i - 2

9 x = g()

10 assertInt(a)

11 assertInt(x)

12 a' = a + x

13 br loop(i', a')

1 b = i == 0

2 cbr b thn() els()

0 br loop(y, 0)

0: {y:Top,z:Top}

1: bottom

2: {b:Top}

3: bottom

4: {a:Int}

5: {a:Int,z:Int}

6: {a:Int,z:Int,r:Int}

7: bottom

8: {i:Int}

9: {i:Int,i':Int}

10: {i:Int,i':Int,x:Top}

11: {i:Int,i':Int,x:Top,a:Int}

12: {i:Int,i':Int,x:Top,a:Int,x:Int}

13: {i:Int,i':Int,x:Top,a:Int,x:Int,a':Int}

f(y,z):

unmentioned vars: bottom

loop(i,a):

thn():

3 assertInt(a)

4 assertInt(z)

5 r = a - z

6 ret r

els():

7 assertInt(i)

8 i' = i - 2

9 x = g()

10 assertInt(a)

11 assertInt(x)

12 a' = a + x

13 br loop(i', a')

1 b = i == 0

2 cbr b thn() els()

0 br loop(y, 0)

0: {y:Top,z:Top}

1: bottom

2: {b:Top}

3: bottom

4: {a:Int}

5: {a:Int,z:Int}

6: {a:Int,z:Int,r:Int}

7: bottom

8: {i:Int}

9: {i:Int,i':Int}

10: {i:Int,i':Int,x:Top}

11: {i:Int,i':Int,x:Top,a:Int}

12: {i:Int,i':Int,x:Top,a:Int,x:Int}

13: {i:Int,i':Int,x:Top,a:Int,x:Int,a':Int}

How do we update blocks (1,3,7) based on the previous round?

```

f(y,z):
  loop(i,a):
    thn():
      3 assertInt(a)
      4 assertInt(z)
      5 r = a - z
      6 ret r
    els():
      7 assertInt(i)
      8 i' = i - 2
      9 x = g()
      10 assertInt(a)
      11 assertInt(x)
      12 a' = a + x
      13 br loop(i', a')
      1 b = i == 0
      2 cbr b thn() els()
      0 br loop(y, 0)

```

Previous iteration output

```

0: {y:Top,z:Top}
1: bottom
13: {i:Int,i':Int,x:Top,a:Int,x:Int,a':Int}

```

Next iteration, update loop based on its two predecessors

```

1: {y:Top,z:Top,i = y:Top,a = 0:Int}
  U {i = i':Int,a = a':Int}

```

simplification: remove variables that aren't in scope

```

f(y,z):
  loop(i,a):
    thn():
      3 assertInt(a)
      4 assertInt(z)
      5 r = a - z
      6 ret r
    els():
      7 assertInt(i)
      8 i' = i - 2
      9 x = g()
      10 assertInt(a)
      11 assertInt(x)
      12 a' = a + x
      13 br loop(i', a')
      1 b = i == 0
      2 cbr b thn() els()
      0 br loop(y, 0)

```

Previous iteration output

```

0: {y:Top,z:Top}
1: bottom
13: {i:Int,i':Int,x:Top,a:Int,x:Int,a':Int}

```

Next iteration, update loop based on its two predecessors

```

1: {y:Top,z:Top,i:Top,a:Int}

```

```

f(y,z):
  loop(i,a):
    thn():
      3 assertInt(a)
      4 assertInt(z)
      5 r = a - z
      6 ret r
    els():
      7 assertInt(i)
      8 i' = i - 2
      9 x = g()
      10 assertInt(a)
      11 assertInt(x)
      12 a' = a + x
      13 br loop(i', a')
  1 b = i == 0
  2 cbr b thn() els()
  0 br loop(y, 0)

```

Converges after one round

```

0: {y:Top,z:Top}
1: {y:Top,z:Top,i:Top,a:Int}
2: {y:Top,z:Top,i:Top,a:Int,b:Top}
3: {y:Top,z:Top,i:Top,a:Int,b:Top}
4: {y:Top,z:Top,i:Top,a:Int,b:Top}
5: {y:Top,z:Int,i:Top,a:Int,b:Top}
6: {y:Top,z:Top,i:Top,a:Int,b:Top,r:Int}
7: {y:Top,z:Top,i:Top,a:Int,b:Top}
8: {y:Top,z:Top,i:Int,a:Int,b:Top}
9: {y:Top,z:Top,i:Int,a:Int,b:Top,i':Int}
10: {y:Top,z:Top,i:Int,a:Int,b:Top,i':Int,x:Top}
11: {y:Top,z:Top,i:Int,a:Int,b:Top,i':Int,x:Top}
12: {y:Top,z:Top,i:Int,a:Int,b:Top,i':Int,x:Int}
13: {y:Top,z:Top,i:Int,a:Int,b:Top,i':Int,x:Int,a':Int}

```

```

f(y,z):
  loop(i,a):
    thn():
      3 assertInt(a)
      4 assertInt(z)
      5 r = a - z
      6 ret r
    els():
      7 assertInt(i)
      8 i' = i - 2
      9 x = g()
      10 assertInt(a)
      11 assertInt(x)
      12 a' = a + x
      13 br loop(i', a')
  1 b = i == 0
  2 cbr b thn() els()
  0 br loop(y, 0)

```

remove assertions based on inferred info

```

0: {y:Top,z:Top}
1: {y:Top,z:Top,i:Top,a:Int}
2: {y:Top,z:Top,i:Top,a:Int,b:Top}
3: {y:Top,z:Top,i:Top,a:Int,b:Top}
4: {y:Top,z:Top,i:Top,a:Int,b:Top}
5: {y:Top,z:Int,i:Top,a:Int,b:Top}
6: {y:Top,z:Top,i:Top,a:Int,b:Top,r:Int}
7: {y:Top,z:Top,i:Top,a:Int,b:Top}
8: {y:Top,z:Top,i:Int,a:Int,b:Top}
9: {y:Top,z:Top,i:Int,a:Int,b:Top,i':Int}
10: {y:Top,z:Top,i:Int,a:Int,b:Top,i':Int,x:Top}
11: {y:Top,z:Top,i:Int,a:Int,b:Top,i':Int,x:Top}
12: {y:Top,z:Top,i:Int,a:Int,b:Top,i':Int,x:Int}
13: {y:Top,z:Top,i:Int,a:Int,b:Top,i':Int,x:Int,a':Int}

```

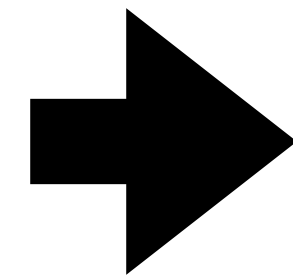
unfortunately, can't remove the assertInt(i) because we don't know initial value of your i is an Int

```

extern g
def f(y,z):
    def loop(i,a):
        if i == 0:
            a - z
        else:
            loop(i - 1, a + g())
    in
    loop(y, 0)

```

inline once



```

extern g
def f(y,z):
    def loop(i,a):
        if i == 0:
            a - z
        else:
            loop(i - 1, a + g())
    in
    if y == 0:
        0 - z
    else:
        loop(y - 1, 0 + g())

```

with this change, now `i` will be determined to always be an `Int`



GENERAL DATAFLOW ANALYSIS

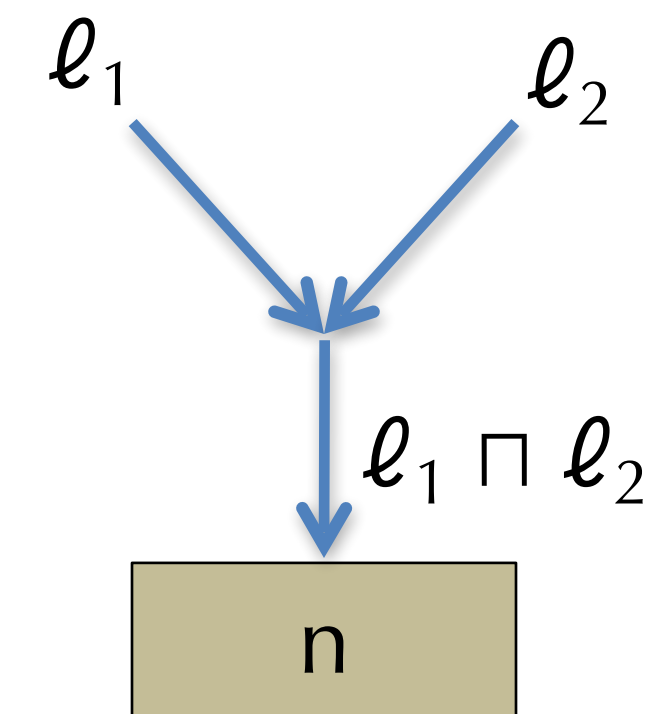
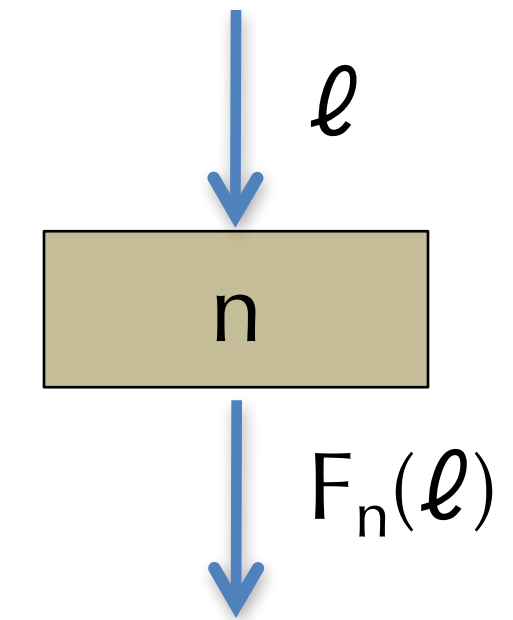
Common Features

- Analyses have a *domain* over which they solve constraints.
 - Liveness, the domain is sets of variables
 - Tag checking the domain is our tag sets
- Each analysis has a notion of **flow function**
 - Used to explain how information propagates across a node.
- Each analysis propagates information either *forward* or *backward*
 - Forward: defined in terms of predecessor nodes
 - Backward: defined in terms of successor nodes
- Each analysis has a way of aggregating information
 - Liveness uses union (**u**)
 - Tag analysis uses least upper bound

(Forward) Dataflow Analysis Framework

A forward dataflow analysis can be characterized by:

1. A domain of dataflow values \mathcal{L}
 - e.g. \mathcal{L} = the powerset of all variables
 - Think of $\ell \in \mathcal{L}$ as a property, then “ $x \in \ell$ ” means “ x has the property”
2. For each node n , a flow function $F_n : \mathcal{L} \rightarrow \mathcal{L}$
 - “If ℓ is a property that holds before the node n , then $F_n(\ell)$ holds after n ”
3. A combining operator \sqcap
 - “If we know *either* ℓ_1 *or* ℓ_2 holds on entry to node n , we know at most $\ell_1 \sqcap \ell_2$ ”
 - $\text{in}[n] := \sqcap_{n' \in \text{pred}[n]} \text{out}[n']$



Generic Iterative (Forward) Analysis

for all n , $\text{in}[n] := \top$, $\text{out}[n] := \top$

repeat until no change

 for all n

$\text{in}[n] := \bigwedge_{n' \in \text{pred}[n]} \text{out}[n']$

$\text{out}[n] := F_n(\text{in}[n])$

 end

end

- Here, $\top \in \mathcal{L}$ (“top”) represents having the “maximum” amount of information.
 - Having “more” information enables more optimizations
 - “Maximum” amount could be inconsistent with the constraints.
 - Iteration refines the answer, eliminating inconsistencies

Structure of \mathcal{L}

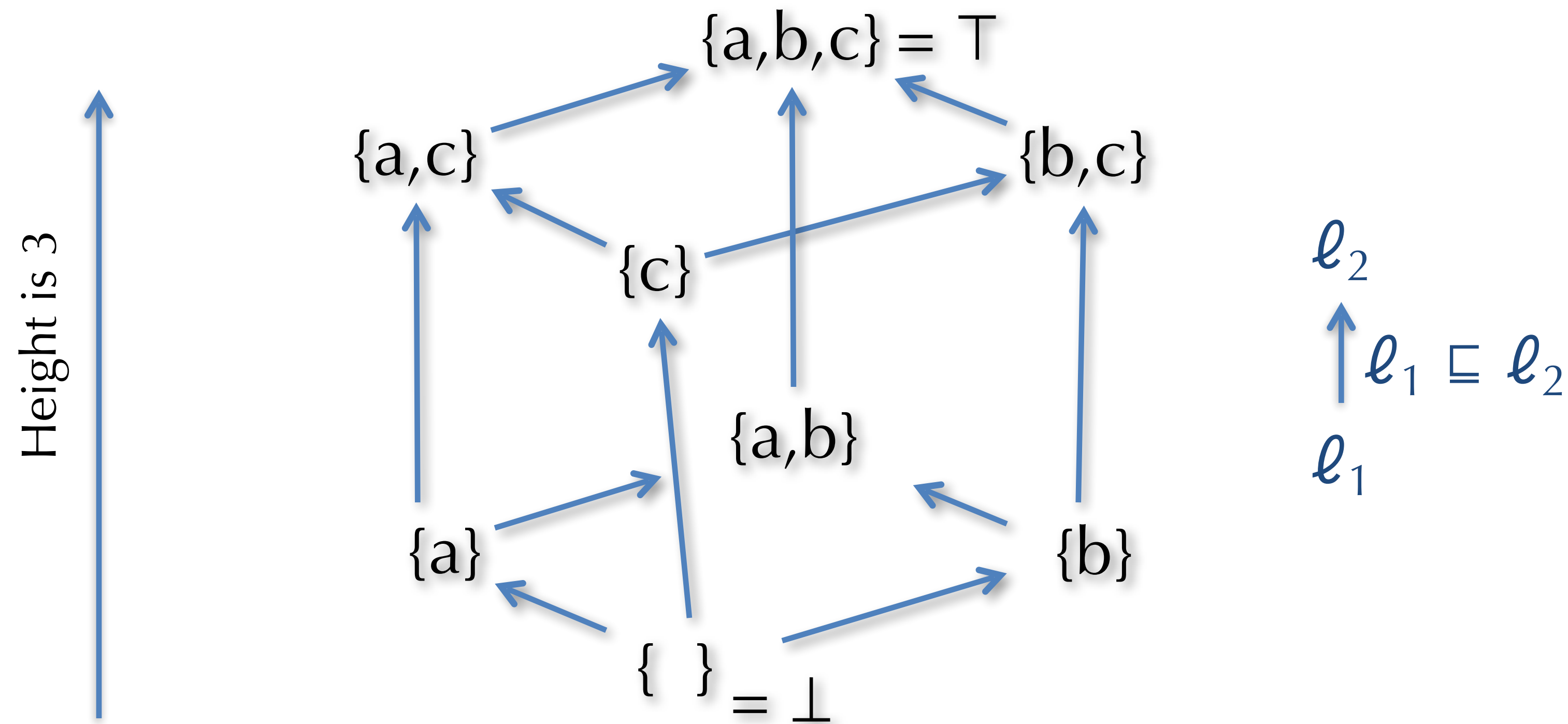
- The domain has structure that reflects the “amount” of information contained in each dataflow value.
- Some dataflow values are more informative than others:
 - Write $\ell_1 \sqsubseteq \ell_2$ whenever ℓ_2 provides at least as much information as ℓ_1 .
 - The dataflow value ℓ_2 is “better” for enabling optimizations.
- Example 1: for liveness analysis, *smaller* sets of variables are more informative.
 - Having smaller sets of variables live across an edge means that there are fewer conflicts for register allocation assignments.
 - So: $\ell_1 \sqsubseteq \ell_2$ if and only if $\ell_1 \supseteq \ell_2$
- Example 2: for available expressions analysis, larger sets of nodes are more informative.
 - Having a larger set of nodes (equivalently, expressions) available means that there is more opportunity for common subexpression elimination.
 - So: $\ell_1 \sqsubseteq \ell_2$ if and only if $\ell_1 \subseteq \ell_2$

\mathcal{L} as a Partial Order

- \mathcal{L} is a *partial order* defined by the ordering relation \sqsubseteq .
- A partial order is an ordered set.
- Some of the elements might be *incomparable*.
 - That is, there might be $\ell_1, \ell_2 \in \mathcal{L}$ such that neither $\ell_1 \sqsubseteq \ell_2$ nor $\ell_2 \sqsubseteq \ell_1$
- Properties of a partial order:
 - *Reflexivity*: $\ell \sqsubseteq \ell$
 - *Transitivity*: $\ell_1 \sqsubseteq \ell_2$ and $\ell_2 \sqsubseteq \ell_3$ implies $\ell_1 \sqsubseteq \ell_3$
 - *Anti-symmetry*: $\ell_1 \sqsubseteq \ell_2$ and $\ell_2 \sqsubseteq \ell_1$ implies $\ell_1 = \ell_2$
- Examples:
 - Integers ordered by \leq
 - Sets ordered by \subseteq or \supseteq

Subsets of $\{a,b,c\}$ ordered by \subseteq

Partial order presented as a Hasse diagram.



order \sqsubseteq is \subseteq

meet \sqcap is \cap

join \sqcup is \cup

Meets and Joins

- The combining operator \sqcap is called the “meet” operation.
- It constructs the *greatest lower bound*:
 - $\ell_1 \sqcap \ell_2 \sqsubseteq \ell_1$ and $\ell_1 \sqcap \ell_2 \sqsubseteq \ell_2$
“the meet is a lower bound”
 - If $\ell \sqsubseteq \ell_1$ and $\ell \sqsubseteq \ell_2$ then $\ell \sqsubseteq \ell_1 \sqcap \ell_2$
“there is no greater lower bound”
- Dually, the \sqcup operator is called the “join” operation.
- It constructs the *least upper bound*:
 - $\ell_1 \sqsubseteq \ell_1 \sqcup \ell_2$ and $\ell_2 \sqsubseteq \ell_1 \sqcup \ell_2$
“the join is an upper bound”
 - If $\ell_1 \sqsubseteq \ell$ and $\ell_2 \sqsubseteq \ell$ then $\ell_1 \sqcup \ell_2 \sqsubseteq \ell$
“there is no smaller upper bound”
- A partial order that has all meets and joins is called a *lattice*.
 - If it has just meets, it’s called a *meet semi-lattice*.

Another Way to Describe the Algorithm

- Algorithm repeatedly computes (for each node n):
- $\text{out}[n] := F_n(\text{in}[n])$
- Equivalently: $\text{out}[n] := F_n(\prod_{n' \in \text{pred}[n]} \text{out}[n'])$
 - By definition of $\text{in}[n]$
- We can write this as a simultaneous update of the vector of $\text{out}[n]$ values:
 - let $x_n = \text{out}[n]$
 - Let $\mathbf{X} = (x_1, x_2, \dots, x_n)$ it's a vector of points in \mathcal{L}
 - $\mathbf{F}(\mathbf{X}) = (F_1(\prod_{j \in \text{pred}[1]} \text{out}[j]), F_2(\prod_{j \in \text{pred}[2]} \text{out}[j]), \dots, F_n(\prod_{j \in \text{pred}[n]} \text{out}[j]))$
- Any solution to the constraints is a *fixpoint* \mathbf{X} of \mathbf{F}
 - i.e. $\mathbf{F}(\mathbf{X}) = \mathbf{X}$

Iteration Computes Fixpoints

- Let $\mathbf{X}_0 = (\top, \top, \dots, \top)$
- Each loop through the algorithm apply F to the old vector:
 $\mathbf{X}_1 = \mathbf{F}(\mathbf{X}_0)$
 $\mathbf{X}_2 = \mathbf{F}(\mathbf{X}_1)$
...
- $\mathbf{F}^{k+1}(\mathbf{X}) = \mathbf{F}(\mathbf{F}^k(\mathbf{X}))$
- A fixpoint is reached when $\mathbf{F}^k(\mathbf{X}) = \mathbf{F}^{k+1}(\mathbf{X})$
 - That's when the algorithm stops.
- Wanted: a maximal fixpoint
 - Because that one is more informative/useful for performing optimizations

Monotonicity & Termination

- Each flow function F_n maps lattice elements to lattice elements; to be sensible it should be *monotonic*:
- $F : \mathcal{L} \rightarrow \mathcal{L}$ is *monotonic* iff:
 $\ell_1 \sqsubseteq \ell_2$ implies that $F(\ell_1) \sqsubseteq F(\ell_2)$
 - Intuitively: “If you have more information entering a node, then you have more information leaving the node.”
- Monotonicity lifts point-wise to the function: $\mathbf{F} : \mathcal{L}^n \rightarrow \mathcal{L}^n$
 - vector $(x_1, x_2, \dots, x_n) \sqsubseteq (y_1, y_2, \dots, y_n)$ iff $x_i \sqsubseteq y_i$ for each i
- Note that \mathbf{F} is consistent: $\mathbf{F}(\mathbf{X}_0) \sqsubseteq \mathbf{X}_0$
 - So each iteration moves at least one step down the lattice (for some component of the vector)
 - $\dots \sqsubseteq \mathbf{F}(\mathbf{F}(\mathbf{X}_0)) \sqsubseteq \mathbf{F}(\mathbf{X}_0) \sqsubseteq \mathbf{X}_0$
- Therefore, # steps needed to reach a fixpoint is at most the height H of \mathcal{L} times the number of nodes: $O(Hn)$

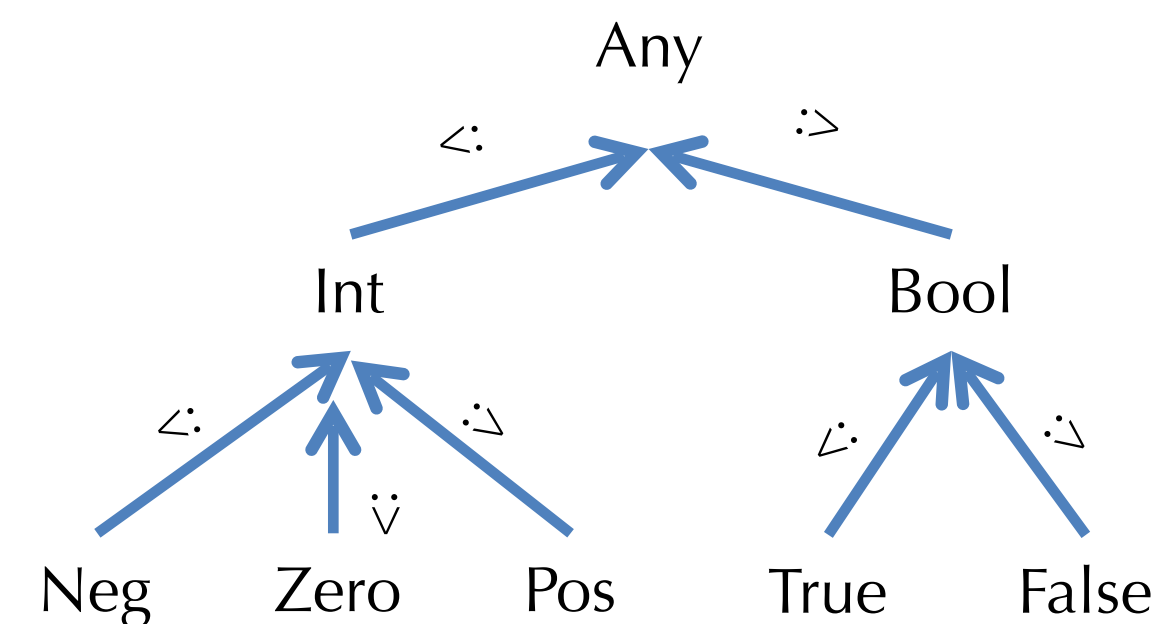
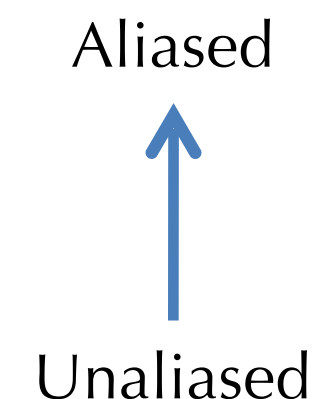
Building Lattices?

- Information about individual nodes or variables can be lifted *pointwise*:
 - If \mathcal{L} is a lattice, then so is $\{f : X \rightarrow \mathcal{L}\}$ where $f \sqsubseteq g$ if and only if $f(x) \sqsubseteq g(x)$ for all $x \in X$.

- Like *types*, the dataflow lattices are *static approximations* to the dynamic behavior:

- Could pick a lattice based on subtyping:

- Or other information:




- Points in the lattice are sometimes called dataflow “*facts*”

“Classic” Constant Propagation

- Constant propagation can be formulated as a dataflow analysis.

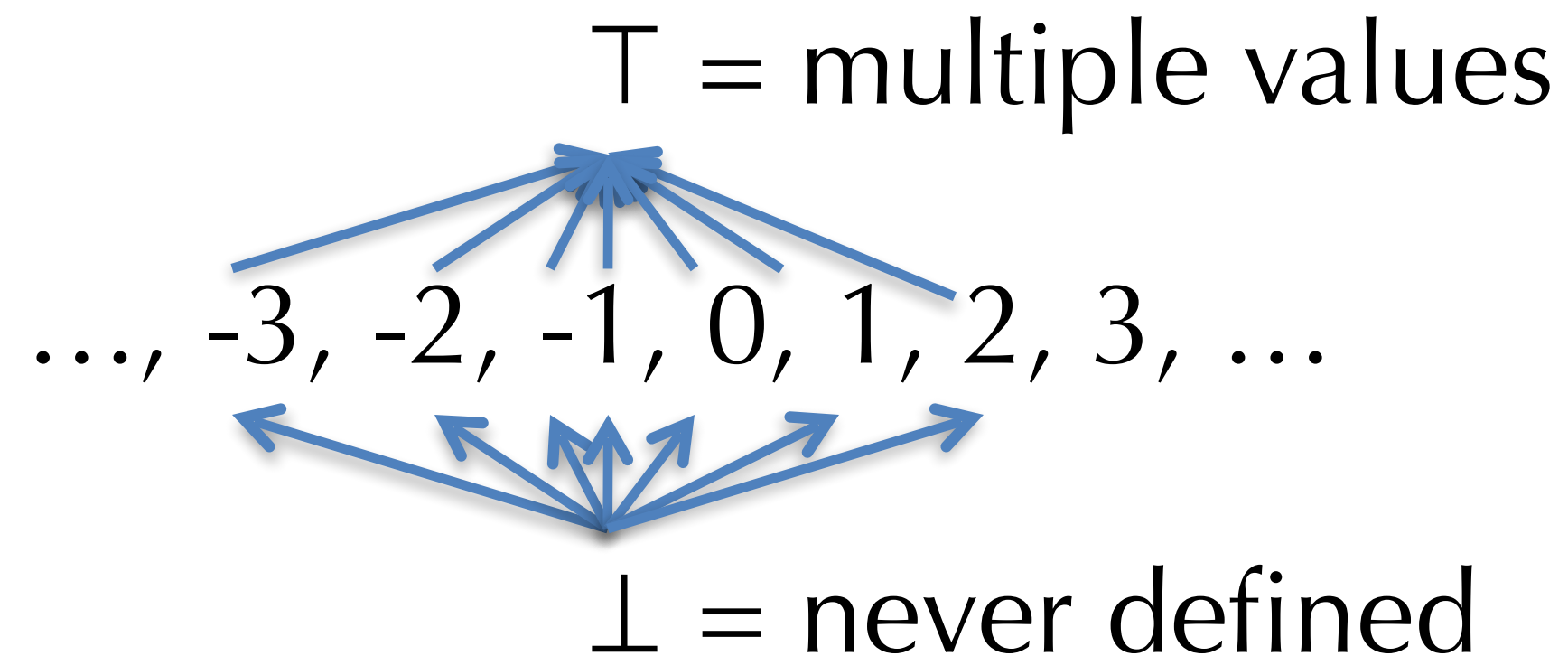
- Idea: propagate and fold integer constants in one pass:

$x = 1;$		$x = 1;$
$y = 5 + x;$		$y = 6;$
$z = y * y;$		$z = 36;$

- Information about a single variable:
 - Variable is never defined.
 - Variable has a single, constant value.
 - Variable is assigned multiple values.

Domains for Constant Propagation

- We can make a constant propagation lattice L for *one variable* like this:



- To accommodate multiple variables, we take the product lattice, with one element per variable.
 - Assuming there are three variables, x , y , and z , the elements of the product lattice are of the form (ℓ_x, ℓ_y, ℓ_z) .
 - Alternatively, think of the product domain as a context that maps variable names to their “*abstract interpretations*”
- What are “meet” and “join” in this product lattice?
- What is the height of the product lattice?

Flow Functions

- Consider the node $x = y \text{ op } z$
 - $F(\ell_x, \ell_y, \ell_z) = ?$
 - $F(\ell_x, \top, \ell_z) = (\top, \top, \ell_z)$
 - $F(\ell_x, \ell_y, \top) = (\top, \ell_y, \top)$
 - $F(\ell_x, \perp, \ell_z) = (\perp, \perp, \ell_z)$
 - $F(\ell_x, \ell_y, \perp) = (\perp, \ell_y, \perp)$
 - $F(\ell_x, i, j) = (i \text{ op } j, i, j)$
- “If either input might have multiple values the result of the operation might too.”
- “If either input is undefined the result of the operation is too.”
- “If the inputs are known constants, calculate the output statically.”
- Flow functions for the other nodes are easy...
 - Monotonic?

Dataflow Analysis: Summary

- Many dataflow analyses fit into a common framework.
- Key idea: *Iterative solution* of a system of equations over a *lattice* of constraints.
 - Iteration terminates if flow functions are monotonic.