

Interprocedural Analysis

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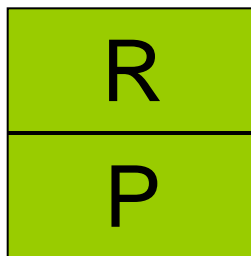
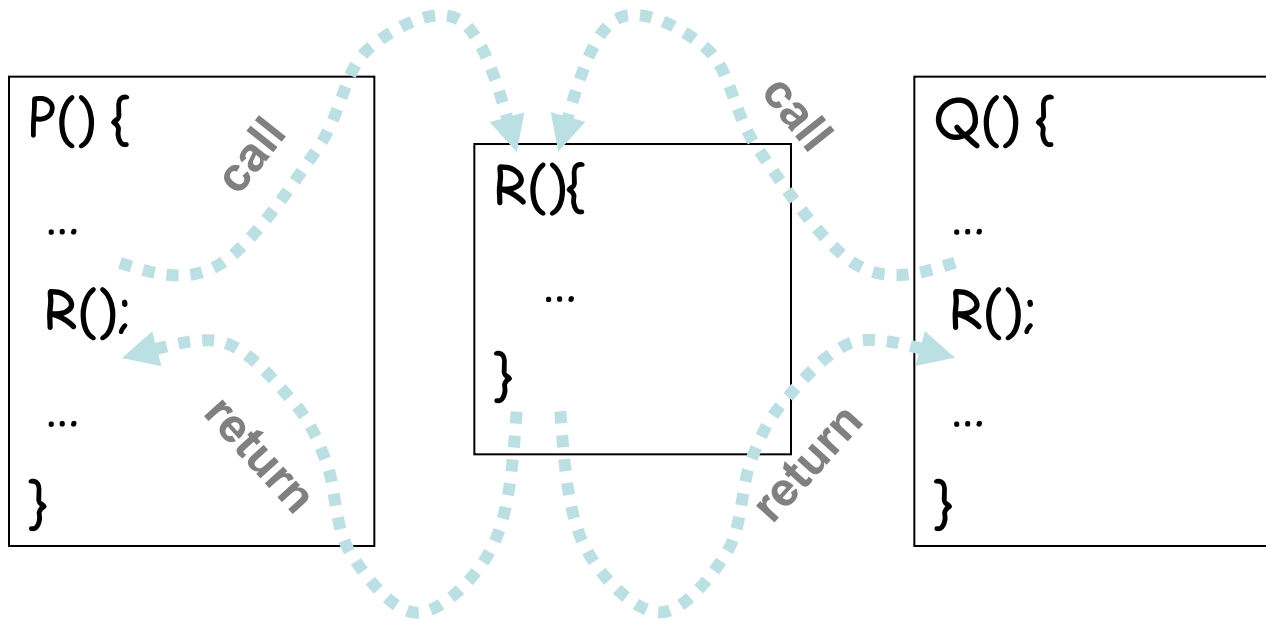
Bibliography

- **Textbook 2.5**
- Patrick Cousot & Radhia Cousot. Static determination of dynamic properties of recursive procedures In *IFIP Conference on Formal Description of Programming Concepts*, E.J. Neuhold, (Ed.), pages 237-277, St-Andrews, N.B., Canada, 1977. North-Holland Publishing Company (1978).
- **Two Approaches to interprocedural analysis by Micha Sharir and Amir Pnueli**
- **IDFS** Interprocedural Distributive Finite Subset Precise interprocedural dataflow analysis via graph reachability. *Reps, Horowitz, and Sagiv, POPL '95*
- **IDE** Interprocedural Distributive Environment Precise interprocedural dataflow analysis with applications to constant propagation. *Sagiv, Reps, Horowitz, and TCS '96*

Challenges in Interprocedural Analysis

- Respect call-return mechanism
- Handling recursion
- Local variables
- Parameter passing mechanisms
- The called procedure is not always known
- The source code of the called procedure is not always available
- Concurrency

Stack regime



Guiding light

- Exploit stack regime
 - Precision
 - Efficiency



Simplifying Assumptions

- ✗ Parameter passed by value
- ✗ No OO
- ✗ No nesting
- ✗ No concurrency
- ✓ Recursion is supported

Topics Covered

- The trivial approach
- Procedure Inlining
- The naive approach
- The call string approach
- Functional Approach
- Valid paths
- Interprocedural Analysis via Graph Reachability
- Beyond reachability
- [Demand analysis]
- [Advanced Topics]
 - Karr's analysis
 - Backward
 - MOPED

Undecidability Issues

- It is undecidable if a program point is reachable in some execution
- Some static analysis problems are undecidable even if the program conditions are ignored

The Constant Propagation Example

```
while (...) {  
    if (...) x_1 = x_1 + 1;  
    if (...) x_2 = x_2 + 1;  
    ...  
    if (...) x_n = x_n + 1;  
}  
y = truncate (1/ (1 + p2(x_1, x_2, ..., x_n))  
/* Is y=0 here? */
```

Conservative Solution

- Every detected constant is indeed constant
 - But may fail to identify some constants
- Every potential impact is identified
 - Superfluous impacts

The extra cost of procedures

- Call return complicates the analysis
- Unbounded stack memory
- Increases asymptotic complexity
- But can be tolerable in practice
- Sometimes even cheaper/more precise than intraprocedural analysis

A trivial treatment of procedure

- Analyze a single procedure
- After every call continue with conservative information
 - Global variables and local variables which “may be modified by the call” have unknown values
- Can be easily implemented
- Procedures can be written in different languages
- Procedure inline can help

Disadvantages of the trivial solution

- Modular (object oriented and functional) programming encourages small frequently called procedures
- Almost all information is lost

Procedure Inlining

- Inline called procedures
- Simple
- Does not handle recursion
- Exponential blow up

```
p1 {  
  call p2  
  ...  
  call p2  
}  
  
p2 {  
  call p3  
  ...  
  call p3  
}  
  
p3{  
  }  
}
```

A Naive Interprocedural solution

- Treat procedure calls as gotos
- Abstract call/return correlations
- Obtain a conservative solution
- Use chaotic iterations
- Usually fast

Simple Example

```
void main() {
```

```
    int x ;
```

```
    → x = p(7);
```

```
    x = p(9) ;
```

```
}
```

```
int p(int a) {
```

```
    return a + 1;
```

```
}
```


Simple Example

```
void main() {  
    int x ;  
    x = p(7);  
    x = p(9) ;  
}
```

```
→ int p(int a) {  
    [a ↦ 7]  
    return a + 1;  
}
```

Simple Example

```
void main() {  
    int x ;  
    x = p(7);  
    x = p(9) ;  
}
```

```
int p(int a) {  
    [a ↦ 7]  
    → return a + 1;  
    [a ↦ 7, $$ ↦ 8]  
}
```

Simple Example

```
void main() {  
    int x ;  
    x = p(7); ←  
    [x ↦ 8]  
    x = p(9) ; ←  
    [x ↦ 8]  
}
```

```
int p(int a) {  
    [a ↦ 7]  
    return a + 1;  
    [a ↦ 7, $$ ↦ 8]  
}
```

Simple Example

```
void main() {
```

```
  int x ;
```

```
  x = p(7);
```

```
    [x ↦ 8]
```

```
→ x = p(9) ;
```

```
    [x ↦ 8]
```

```
}
```

```
int p(int a) {
```

```
  [a ↦ 7]
```

```
  return a + 1;
```

```
  [a ↦ 7, $$ ↦ 8]
```

```
}
```

Simple Example

```
void main() {  
    int x ;  
    x = p(7);  
    [x ↦ 8]  
    x = p(9) ;  
    [x ↦ 8]  
}
```

```
⇒ int p(int a) {  
    [a ↦ τ]  
    return a + 1;  
    [a ↦ 7, $$ ↦ 8]  
}
```

Simple Example

```
void main() {  
    int x ;  
    x = p(7);  
    [x ↦ 8]  
    x = p(9);  
    [x ↦ 8]  
}
```

```
int p(int a) {  
    [a ↦ τ]  
    → return a + 1;  
    [a ↦ τ, $$ ↦ τ]  
}
```

Simple Example

```
void main() {  
    int x ;  
    x = p(7) ; ←  
    [x ↦ T]  
    x = p(9) ; ←  
    [x ↦ T]  
}
```

```
int p(int a) {  
    [a ↦ T]  
    return a + 1;  
    [a ↦ T, $$ ↦ T]  
}
```

The Call String Approach

- The data flow value is associated with sequences of calls (call string)
- Use Chaotic iterations
- To guarantee termination limit the size of call string (typically 1 or 2)
 - Represents tails of calls
- Abstract inline

Simple Example

```
void main() {  
    int x ;  
    → c1: x = p(7);  
    c2: x = p(9) ;  
  
}
```

```
int p(int a) {  
    return a + 1;  
}
```

Simple Example

```
void main() {  
    int x ;  
    c1: x = p(7);  
    c2: x = p(9) ;  
  
}
```

```
→ int p(int a) {  
    c1: [a ↦ 7]  
    return a + 1;  
}
```

Simple Example

```
void main() {  
    int x ;  
    c1: x = p(7);  
    c2: x = p(9) ;  
  
}
```

```
int p(int a) {  
    c1: [a ↦7]  
    → return a + 1;  
    c1:[a ↦7, $$ ↦8]  
}
```

Simple Example

```
void main() {  
    int x ;  
    c1: x = p(7); ←  
    ε: X ↦ 8  
    c2: x = p(9) ;  
  
}
```

```
int p(int a) {  
    c1: [a ↦ 7]  
    return a + 1;  
    c1:[a ↦ 7, $$ ↦ 8]  
}
```

Simple Example

```
void main() {  
    int x ;  
    c1: x = p(7);  
    ε: [x ↦ 8]  
    → c2: x = p(9) ;  
}
```

```
int p(int a) {  
    c1:[a ↦7]  
    return a + 1;  
    c1:[a ↦7, $$ ↦8]  
}
```

Simple Example

```
void main() {  
  int x ;  
  c1: x = p(7);  
  ε: [x ↦ 8]  
  c2: x = p(9) ;  
}
```

```
⇒ int p(int a) {  
  c1:[a ↦7]  
  c2:[a ↦9]  
  return a + 1;  
  c1:[a ↦7, $$ ↦8]  
}
```

Simple Example

```
void main() {  
  int x ;  
  c1: x = p(7);  
  ε: [x ↦ 8]  
  c2: x = p(9) ;  
}
```

```
int p(int a) {  
  c1:[a ↦7]  
  c2:[a ↦9]  
  → return a + 1;  
  c1:[a ↦7, $$ ↦8]  
  c2:[a ↦9, $$ ↦10]  
}
```

Simple Example

```
void main() {  
    int x ;  
    c1: x = p(7);  
    ε: [x ↦ 8]  
    c2: x = p(9) ; ←  
    ε: [x ↦ 10]  
}
```

```
int p(int a) {  
    c1:[a ↦7]  
    c2:[a ↦9]  
    return a + 1;  
    c1:[a ↦7, $$ ↦8]  
    c2:[a ↦9, $$ ↦10]  
}
```


Another Example

```
void main() {  
  int x ;  
  c1: x = p(7);  
  ε: [x ↦ 8]  
  c2: x = p(9) ;  
  ε: [x ↦ 10]  
}
```

```
int p(int a) {  
  c1:[a ↦7]  
  c2:[a ↦9]  
  return c3: p1(a + 1);  
  c1:[a ↦7, $$ ↦τ]  
  c2:[a ↦9, $$ ↦τ]  
}
```

```
int p1(int b) {  
  (c1|c2)c3:[b ↦τ]  
  return 2 * b;  
  (c1|c2)c3:[b ↦τ, $$ ↦τ]  
}
```

```

void main() {
    c1: p(7);
    ε: [X ↦ τ]
}

```

```

int p(int a) {
    c1: [a ↦ 7]  c1.c2+: [a ↦ τ]
    if (...) {
        c1: [a ↦ 7]  c1.c2+: [a ↦ τ]
        a = a - 1 ;
        c1: [a ↦ 6]  c1.c2+: [a ↦ τ]
        c2: p (a);
        c1.c2*: [a ↦ τ]

        a = a + 1;
        c1.c2*: [a ↦ τ]
    }
    c1.c2*: [a ↦ τ]

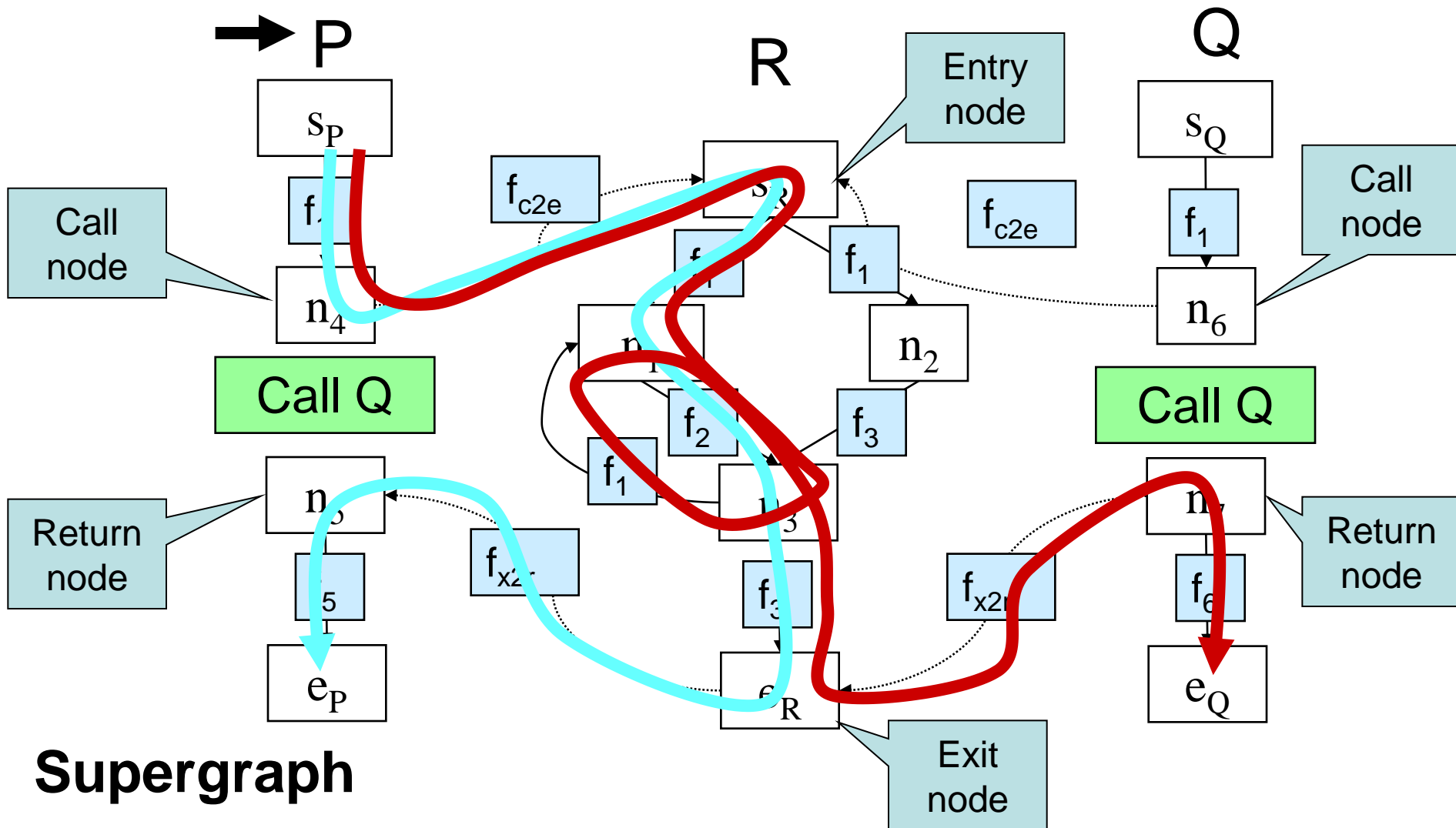
    x = -2*a + 5;
    c1.c2*: [a ↦ τ, x ↦ τ]
}

```

Summary Call String

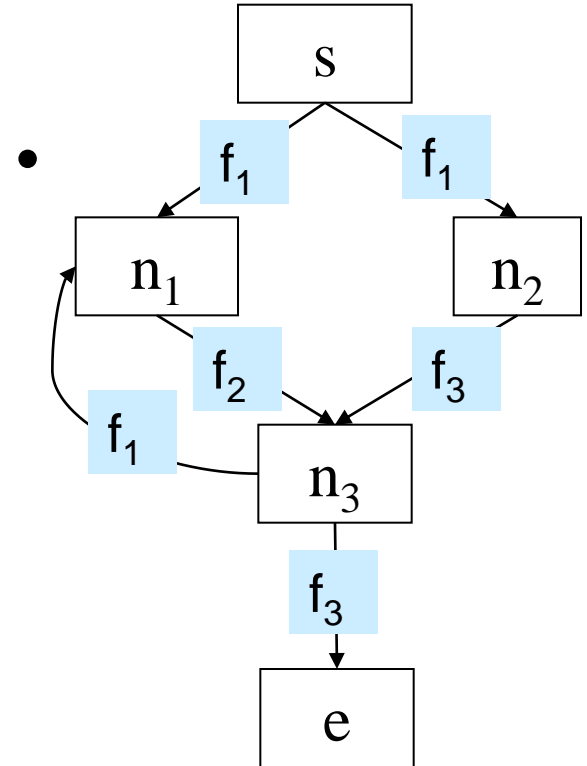
- Easy to implement
- Efficient for very small call strings
- Too expensive even for call strings of length 3
- For finite domains can be precise even with recursion
- Limited precision
- Order of calls can be abstracted
- Related method: procedure cloning

Interprocedural analysis

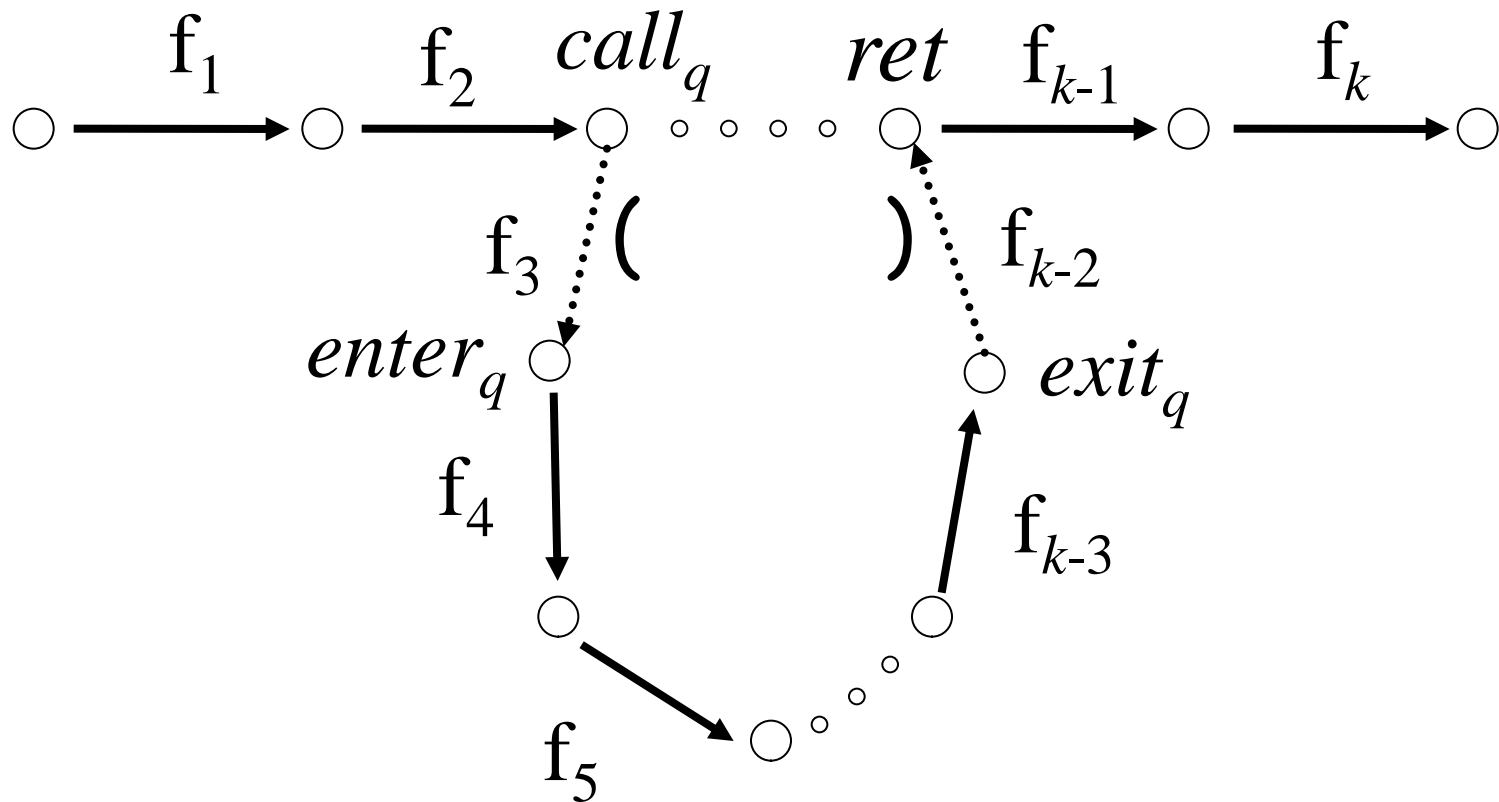


Paths

paths(n) the set of paths from s to n
((s,n₁), (n₁,n₃), (n₃,n₁)) –



Interprocedural Valid Paths

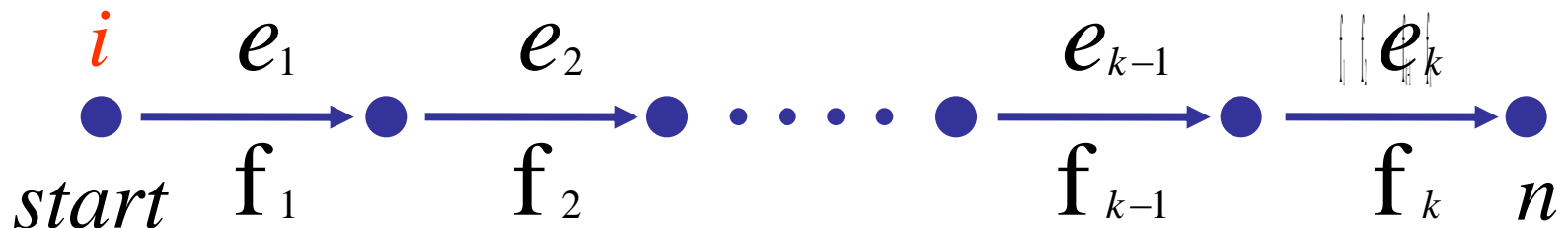


- IVP: all paths with matching calls and returns
 - And prefixes

Interprocedural Valid Paths

- **IVP** set of paths
 - Start at program entry
- Only considers matching calls and returns
 - aka, **valid**
- Can be defined via context free grammar
 - $\text{matched} ::= \text{matched } (_i \text{ matched }) _i \mid \varepsilon$
 - $\text{valid} ::= \text{valid } (_i \text{ matched } \mid \text{matched}$
 - *paths* can be defined by a regular expression

Join Over All Paths (JOP)



$$[[(e_1, \dots, e_k)]] = f_k \circ f_{k-1} \cdots \circ f_2 \circ f_1 \in L \rightarrow L$$

- $JOP[v] = \sqcup \{ [[(e_1, e_2, \dots, e_n)]] (v) \mid (e_1, \dots, e_n) \in \text{paths}(v) \}$
- $JOP \sqsubseteq LFP$
 - Sometimes $JOP = LFP$
 - precise up to “**symbolic execution**”
 - Distributive problem

The Join-Over-Valid-Paths (JVP)

- $vpaths(n)$ all valid paths from program start to n
- $JVP[n] = \sqcup \{ [[e_1, e_2, \dots, e]](1) \mid (e_1, e_2, \dots, e) \in vpaths(n) \}$
- $JVP \sqsubseteq JFP$
 - In some cases the JVP can be computed
 - (Distributive problem)

The Functional Approach

- The meaning of a procedure is mapping from states into states
- The abstract meaning of a procedure is function from an abstract state to abstract states
- Relation between input and output
- In certain cases can compute JVP

The Functional Approach

- Two phase algorithm
 - Compute the dataflow solution at the exit of a procedure as a function of the initial values at the procedure entry (functional values)
 - Compute the dataflow values at every point using the functional values

Phase 1

```
void main() {  
    p(7);  
}
```

```
→ int p(int a) {  
    [a ↦ a0, x ↦ x0]  
    if (...) {  
        a = a - 1 ;  
        p (a);  
        a = a + 1;  
    }  
    x = -2*a + 5;  
    [a ↦ a0, x ↦ -2a0 + 5]  
}
```

Phase 1

```
void main() {  
    p(7);  
}
```

```
int p(int a) {  
    [a ↦ a0, x ↦ x0]  
    if (...) {  
        a = a - 1 ;  
        p (a);  
        a = a + 1;  
    }  
    [a ↦ a0, x ↦ x0]  
    x = -2*a + 5;  
}
```

Phase 1

```
void main() {  
    p(7);  
}
```

```
int p(int a) {  
    [a ↦ a0, x ↦ x0]  
    if (...) {  
        a = a - 1 ;  
        p (a);  
        a = a + 1;  
    }
```

```
    [a ↦ a0, x ↦ x0]
```

```
    ⇒ x = -2*a + 5;
```

```
    [a ↦ a0, x ↦ -2a0 + 5]
```

```
}
```

Phase 1

```
void main() {  
    p(7);  
}
```

```
int p(int a) {  
    [a ↦ a0, x ↦ x0]  
    if (...) {  
        → a = a - 1 ;  
        [a ↦ a0-1, x ↦ x0]  
        p (a);  
        a = a + 1;  
    }  
    [a ↦ a0, x ↦ x0]  
    x = -2*a + 5;  
    [a ↦ a0, x ↦ -2a0 + 5]  
}
```

Phase 1

```
void main() {  
    p(7);  
}
```

```
int p(int a) {  
    [a ↦ a0, x ↦ x0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ a0-1, x ↦ x0]  
        → p (a);  
        [a ↦ a0-1, x ↦ -2a0+7]  
        a = a + 1;  
    }  
    [a ↦ a0, x ↦ x0]  
    x = -2*a + 5;  
    [a ↦ a0, x ↦ -2a0 + 5]  
}
```


Phase 1

```
void main() {  
    p(7);  
}
```

```
int p(int a) {  
    [a ↦ a0, x ↦ x0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ a0-1, x ↦ x0]  
        p (a);  
        [a ↦ a0-1, x ↦ -2a0+7]  
        → a = a + 1 ;  
        [a ↦ a0, x ↦ -2a0+7]  
    }  
    [a ↦ a0, x ↦ x0]  
    x = -2*a + 5;  
    [a ↦ a0, x ↦ -2a0 + 5]  
}
```

Phase 1

```
void main() {  
    p(7);  
}
```

```
int p(int a) {  
    [a ↦ a0, x ↦ x0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ a0-1, x ↦ x0]  
        p (a);  
        [a ↦ a0-1, x ↦ -2a0+7]  
        a = a + 1;  
        [a ↦ a0, x ↦ -2a0+7]  
    }  
    [a ↦ a0, x ↦ τ]  
    x = -2*a + 5;  
    [a ↦ a0, x ↦ -2a0 + 5]  
}
```

Phase 1

```
void main() {  
    p(7);  
}
```

```
int p(int a) {  
    [a ↦ a0, x ↦ x0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ a0-1, x ↦ x0]  
        p (a);  
        [a ↦ a0-1, x ↦ -2a0+7]  
        a = a + 1;  
        [a ↦ a0, x ↦ -2a0+7]  
    }  
}
```

[a ↦ a₀, x ↦ τ]

⇒ x = -2*a + 5;

[a ↦ a₀, x ↦ -2a₀ + 5]

}

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

```
→ int p(int a) {  
    [a ↦ 7, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        p (a);  
        a = a + 1;  
    }  
    x = -2*a + 5;  
}
```

$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

```
int p(int a) {  
    [a ↦ 7, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        p (a);  
        a = a + 1;  
    }  
    ↪ }
```

```
    [a ↦ 7, x ↦ 0]  
    x = -2*a + 5;
```

```
    }  
p(a0, x0) = [a ↦ a0, x ↦ -2a0 + 5]
```

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

```
int p(int a) {  
    [a ↦ 7, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        p (a);  
        a = a + 1;  
    }  
}
```


→ x = -2*a + 5;

```
[a ↦ 7, x ↦ -9]  
}
```

$$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$$

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

```
int p(int a) {  
    [a ↦ 7, x ↦ 0]  
    if (...) {  
         a = a - 1 ;  
        [a ↦ 6, x ↦ 0]  
        p (a);  
        a = a + 1;  
    }  
    [a ↦ 7, x ↦ 0]  
    x = -2*a + 5;  
    [a ↦ 7, x ↦ -9]  
}
```

$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

```
int p(int a) {
```

```
    [a ↦ 7, x ↦ 0]
```

```
    if (...) {
```

```
        a = a - 1 ;
```

```
        [a ↦ 6, x ↦ 0]
```

```
    → p (a);
```

```
        [a ↦ 6, x ↦ -9]
```

```
        a = a + 1;
```

```
    }
```

```
    [a ↦ 7, x ↦ 0]
```

```
    x = -2*a + 5;
```

```
    [a ↦ 7, x ↦ -9]
```

```
}
```

$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

$$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$$

```
int p(int a) {  
    [a ↦ 7, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ 6, x ↦ 0]  
        p (a);  
        [a ↦ 6, x ↦ -9]  
        → a = a + 1;  
        [a ↦ 7, x ↦ -9]  
    }  
    [a ↦ 7, x ↦ 0]  
    x = -2*a + 5;  
    [a ↦ 7, x ↦ -9]  
}
```

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

$$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$$

```
int p(int a) {  
    [a ↦ 7, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ 6, x ↦ 0]  
        p (a);  
        [a ↦ 6, x ↦ -9]  
        a = a + 1;  
        [a ↦ 7, x ↦ -9]  
    }  
    [a ↦ 7, x ↦ τ]  
    x = -2*a + 5;  
    [a ↦ 7, x ↦ -9]  
}
```

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

$$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$$

```
int p(int a) {  
    [a ↦ 7, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ 6, x ↦ 0]  
    }  
    → p (a);  
    [a ↦ 6, x ↦ -9]  
    a = a + 1;  
    [a ↦ 7, x ↦ -9]  
    }  
    [a ↦ 7, x ↦ τ]  
    x = -2*a + 5;  
    [a ↦ 7, x ↦ -9]  
}
```

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$

```
int p(int a) {  
    ⇒ [a ↦ 7, x ↦ 0] [a ↦ 6, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ 6, x ↦ 0]  
        p (a);  
        [a ↦ 6, x ↦ -9]  
        a = a + 1 ;  
        [a ↦ 7, x ↦ -9]  
    }  
    [a ↦ 7, x ↦ τ]  
    x = -2*a + 5;  
    [a ↦ 7, x ↦ -9]  
}
```

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$

```
→ int p(int a) {  
    [a ↦ τ, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ 6, x ↦ 0]  
        p (a);  
        [a ↦ 6, x ↦ -9]  
        a = a + 1;  
        [a ↦ 7, x ↦ -9]  
    }  
    [a ↦ 7, x ↦ τ]  
    x = -2*a + 5;  
    [a ↦ 7, x ↦ -9]  
}
```

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

$$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$$

```
int p(int a) {  
    [a ↦ τ, x ↦ 0]  
    if (...) {  
        ⇒ a = a - 1 ;  
        [a ↦ τ, x ↦ 0]  
        p (a);  
        [a ↦ 6, x ↦ -9]  
        a = a + 1;  
        [a ↦ 7, x ↦ -9]  
    }  
    [a ↦ 7, x ↦ τ]  
    x = -2*a + 5;  
    [a ↦ 7, x ↦ -9]  
}
```

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

$$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$$

```
int p(int a) {  
    [a ↦ τ, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ τ, x ↦ 0]  
        p (a);  
        [a ↦ τ, x ↦ τ]  
        a = a + 1;  
        [a ↦ 7, x ↦ -9]  
    }  
    [a ↦ 7, x ↦ τ]  
    x = -2*a + 5;  
    [a ↦ 7, x ↦ -9]  
}
```

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

$$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$$

```
int p(int a) {  
    [a ↦ τ, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ τ, x ↦ 0]  
        p (a);  
        [a ↦ τ, x ↦ τ]  
        a = a + 1;  
        [a ↦ τ, x ↦ τ]  
    }  
    [a ↦ 7, x ↦ τ]  
    x = -2*a + 5;  
    [a ↦ 7, x ↦ -9]  
}
```


Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$

```
int p(int a) {  
    [a ↦ τ, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ τ, x ↦ 0]  
        p (a);  
        [a ↦ τ, x ↦ τ]  
        a = a + 1;  
        [a ↦ τ, x ↦ τ]  
    }  
}
```



```
[a ↦ τ, x ↦ τ]  
x = -2*a + 5;  
[a ↦ 7, x ↦ -9]  
}
```

Phase 2

```
void main() {  
    p(7);  
    [x ↦ -9]  
}
```

$p(a_0, x_0) = [a \mapsto a_0, x \mapsto -2a_0 + 5]$

```
int p(int a) {  
    [a ↦ τ, x ↦ 0]  
    if (...) {  
        a = a - 1 ;  
        [a ↦ τ, x ↦ 0]  
        p (a);  
        [a ↦ τ, x ↦ τ]  
        a = a + 1;  
        [a ↦ τ, x ↦ τ]  
    }  
    [a ↦ τ, x ↦ τ]  
    x = -2*a + 5;  
    [a ↦ τ, x ↦ τ]  
}
```



Summary Functional approach

- Computes procedure abstraction
- Sharing between different contexts
- Rather precise
- Recursive procedures may be more precise/efficient than loops
- But requires more from the implementation
 - Representing relations
 - Composing relations

Issues in Functional Approach

- How to guarantee that finite height for functional lattice?
 - It may happen that L has finite height and yet the lattice of monotonic function from L to L do not
- Efficiently represent functions
 - Functional join
 - Functional composition
 - Testing equality

A Complicated Example

```
main() {  
    a = 0;  
    p();  
    print a;  
}
```

```
p() {  
    if (...) {  
        a = a + 2 + sgn(a - 100);  
        p();  
        a - 1 ;  
    }  
}
```

CFL-Graph reachability

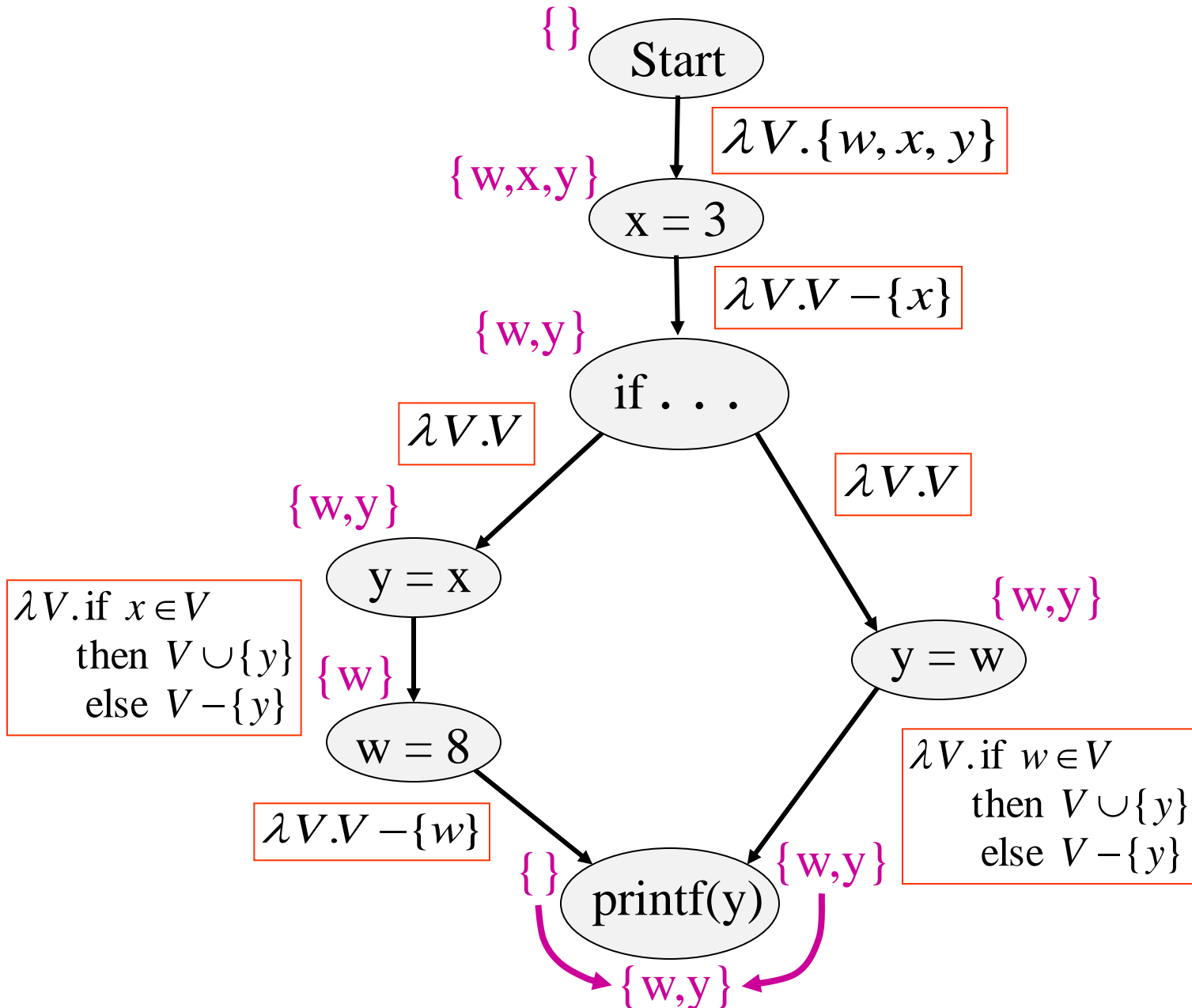
- Special cases of functional analysis
- Finite distributive lattices
- Provides more efficient analysis algorithms
- Reduce the interprocedural analysis problem to finding context free reachability



IDFS / IDE

- **IDFS** Interprocedural Distributive Finite Subset
Precise interprocedural dataflow analysis via graph reachability. *Reps, Horowitz, and Sagiv, POPL'95*
- **IDE** Interprocedural Distributive Environment
Precise interprocedural dataflow analysis with applications to constant propagation. *Reps, Horowitz, and Sagiv, FASE'95, TCS'96*
 - *More general solutions exist*

Possibly Uninitialized Variables



Efficiently Representing Functions

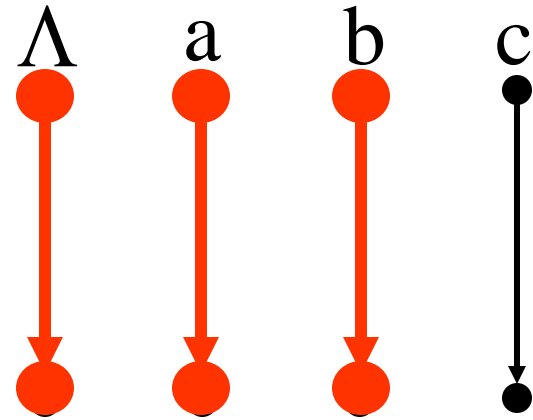
- Let $f:2^D \rightarrow 2^D$ be a distributive function
- Then: $f(X) = f(\emptyset) \cup \cup\{z \in X: f(\{z\})\}$

Representing Dataflow Functions

Identity Function

$$f = \lambda V.V$$

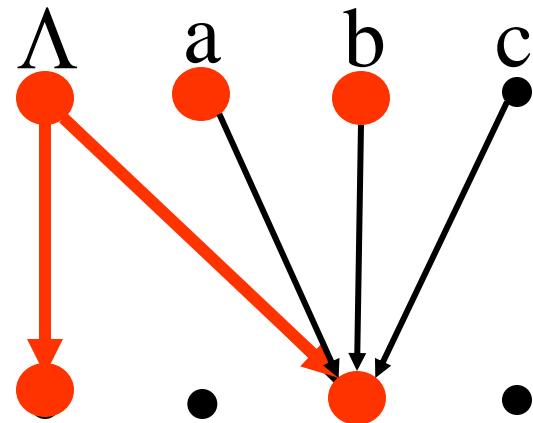
$$f(\{a, b\}) = \{a, b\}$$



Constant Function

$$f = \lambda V.\{b\}$$

$$f(\{a, b\}) = \{b\}$$

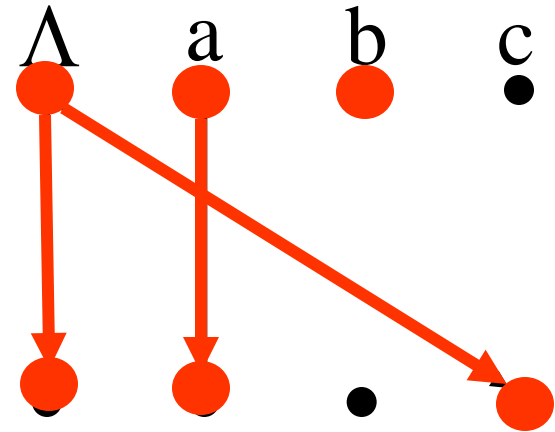


Representing Dataflow Functions

“Gen/Kill” Function

$$f = \lambda V. (V - \{b\}) \cup \{c\}$$

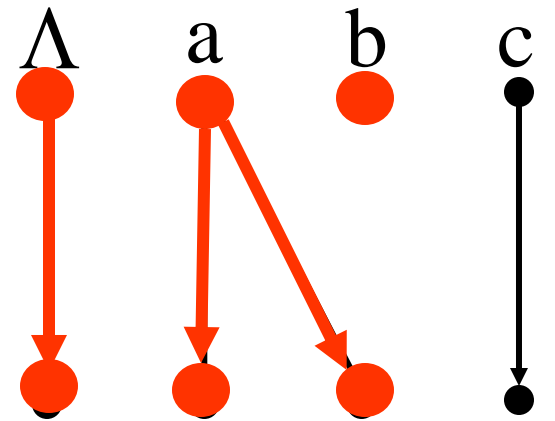
$$f(\{a, b\}) = \{a, c\}$$

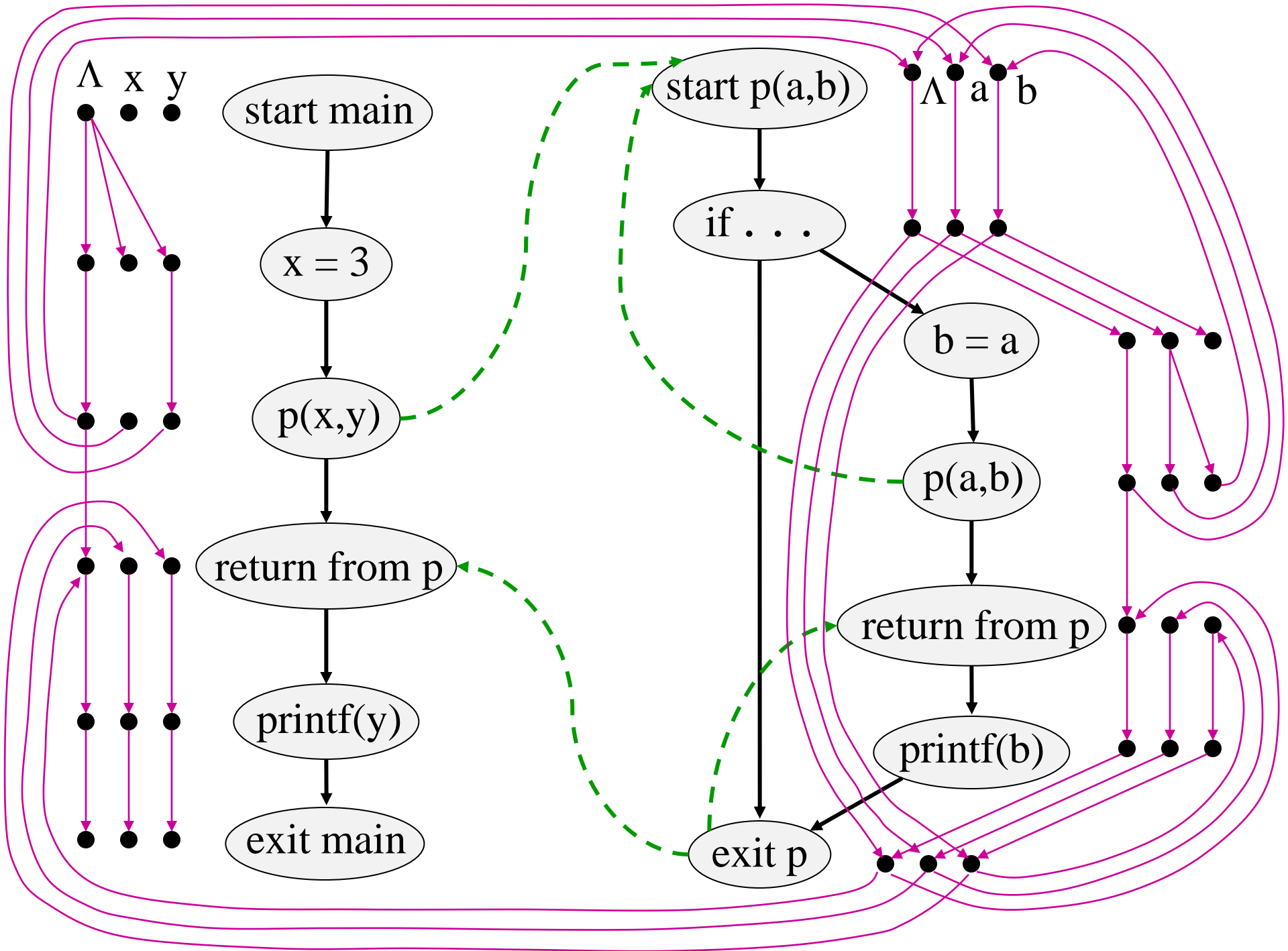


Non-“Gen/Kill” Function

$$f = \lambda V. \text{if } a \in V \\ \text{then } V \cup \{b\} \\ \text{else } V - \{b\}$$

$$f(\{a, b\}) = \{a, b\}$$

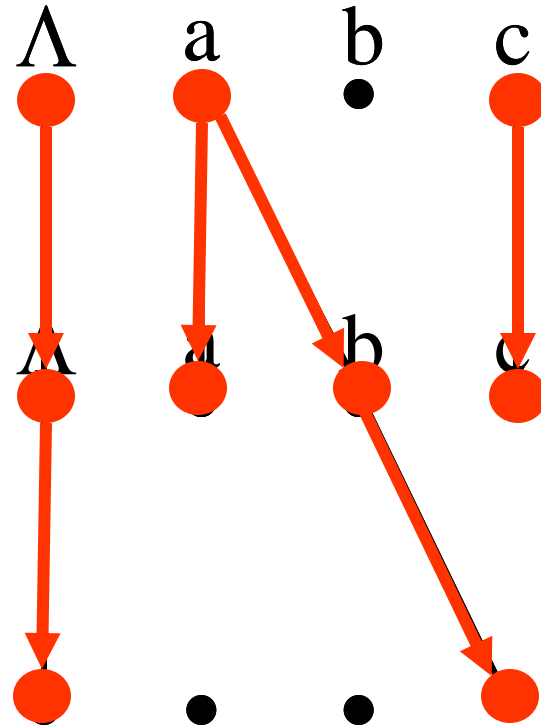




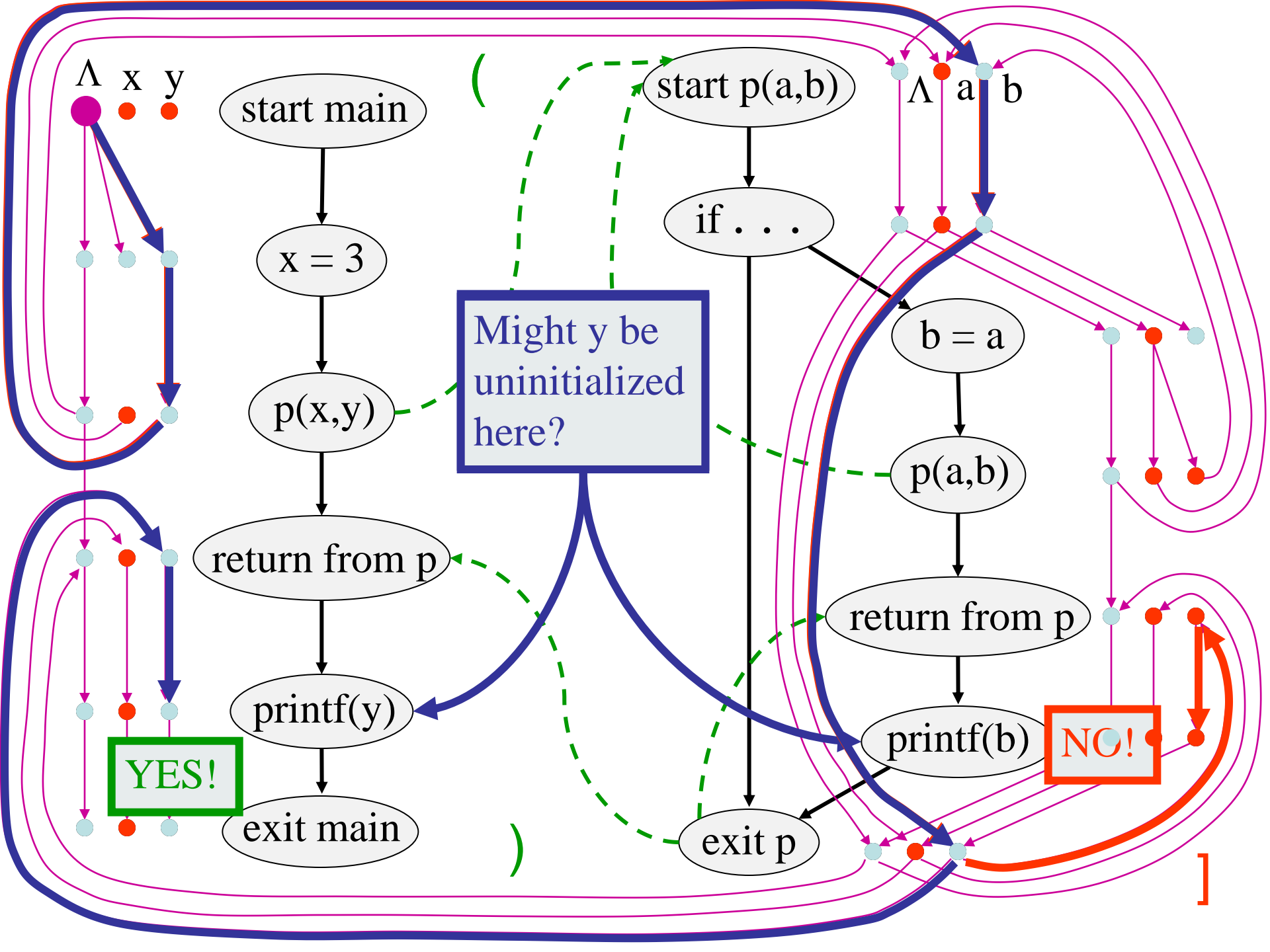
Composing Dataflow Functions

$f_1 = \lambda V. \text{if } a \in V$
 then $V \cup \{b\}$
 else $V - \{b\}$

$f_2 = \lambda V. \text{if } b \in V$
 then $\{c\}$
 else ϕ




$$f_2 \circ f_1(\{a, c\}) = \boxed{\{c\}}$$



Interprocedural Dataflow Analysis via CFL-Reachability

- Graph: Exploded control-flow graph
- L : $L(\text{unbalLeft})$
 - $\text{unbalLeft} = \text{valid}$
- Fact d holds at n iff there is an $L(\text{unbalLeft})$ -path from $\langle \text{start}_{\text{main}}, \Lambda \rangle$ to $\langle n, d \rangle$

Asymptotic Running Time

- CFL-reachability
 - Exploded control-flow graph: ND nodes
 - Running time: $O(N^3D^3)$
- Exploded control-flow graph  Special structure

Running time: $O(ED^3)$

Typically: $E \approx N$, hence $O(ED^3) \approx O(ND^3)$

“Gen/kill” problems: $O(ED)$

IDE

- Goes beyond IFDS problems
 - Can handle unbounded domains
- Requires special form of the domain
- Can be **much** more efficient than IFDS

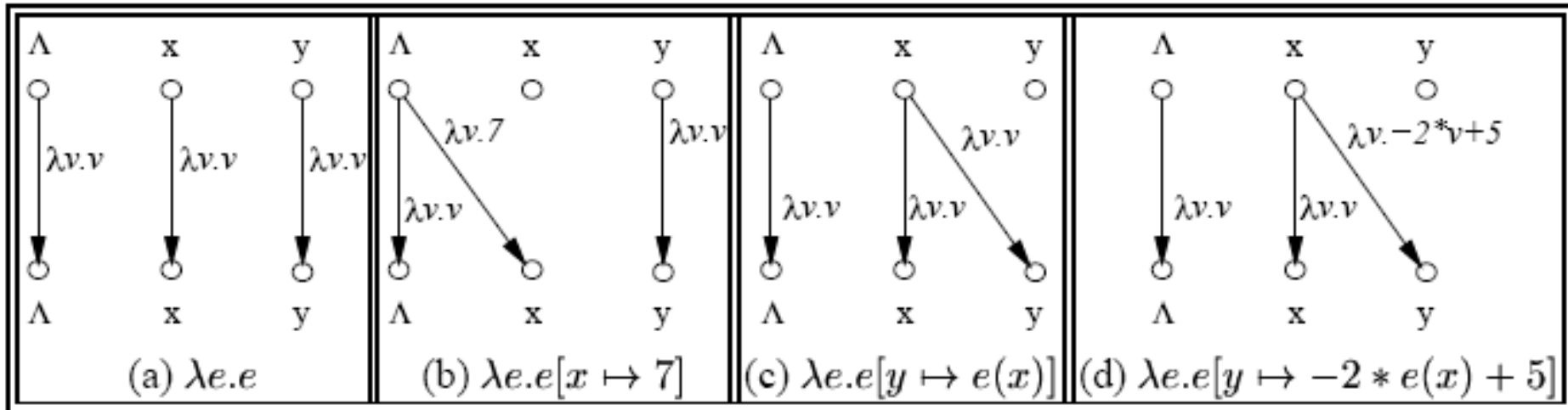
Example Linear Constant Propagation

- Consider the constant propagation lattice
- The value of every variable y at the program exit can be represented by:

$$y = \sqcup \{(a_x x + b_x) \mid x \in \text{Var}_*\} \sqcup c$$
$$a_x, c \in \mathbb{Z} \cup \{\perp, \top\} \quad b_x \in \mathbb{Z}$$

- Supports efficient composition and “functional” join
 - $[z := a * y + b]$
 - What about $[z:=x+y]$?

Linear constant propagation



Point-wise representation of environment transformers

IDE Analysis

- Point-wise representation closed under composition
- CFL-Reachability on the exploded graph
- Compose functions

```

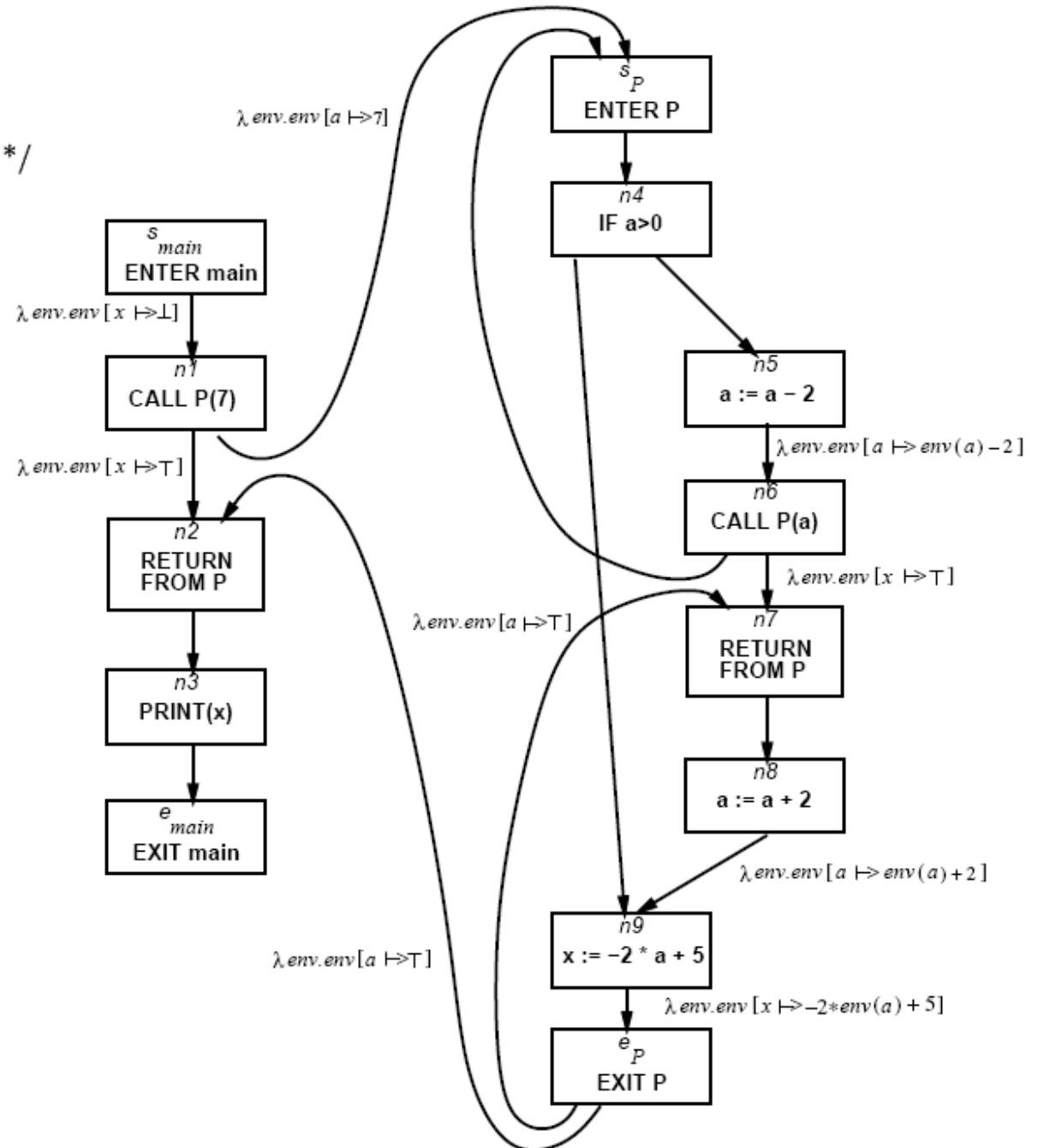
declare x: integer
program main
begin
    call P(7)
    print (x) /* x is a constant here */
end

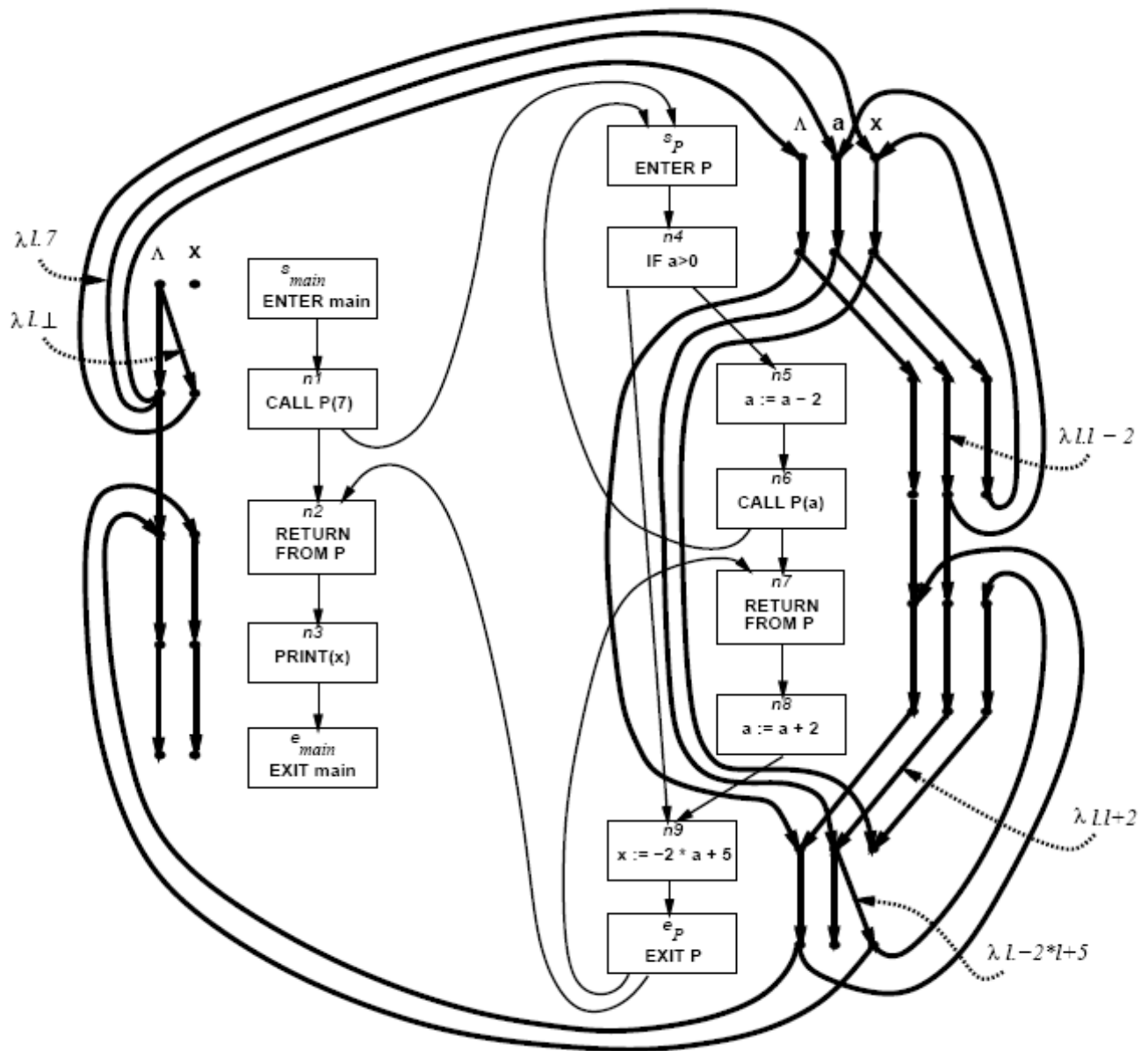
```

```

procedure P (value a : integer)
begin /* a is not a constant here */
    if a > 0 then
        a := a - 2
        call P (a)
        a := a + 2
    fi
    x := -2 * a + 5
    /* x is not a constant here */
end

```





Costs

- $O(ED^3)$
- Class of value transformers $F \subseteq L \rightarrow L$
 - $\text{id} \in F$
 - Finite height
- Representation scheme with (efficient)
 - Application
 - Composition
 - Join
 - Equality
 - Storage

Conclusion

- Handling functions is crucial for abstract interpretation
- Virtual functions and exceptions complicate things
- But scalability is an issue
 - Small call strings
 - Small functional domains
 - Demand analysis