

Abstract Interpretation IV

Semantics and Application to Program Verification

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Selected advanced topics:

- **reduced products** of abstract domains
- **disjunctive** abstract domains
- **inter-procedural** analysis

Practical session:

- implement a reduced product
- help with the project

Reduced products

Idea

Theory:

- the set of abstract domains is a **lattice**,
- ordered by abstraction, which is a partial order, i.e.:
 $(C, \leq) \xleftrightarrow[\alpha]{\gamma} (A, \sqsubseteq)$ denotes that C is more concrete than A .
 (every property of A can also be represented exactly in C)
- there is a least upper bound \sqcup for arbitrary sets of domains and a greatest lower bound \sqcap .

Application: reduced product

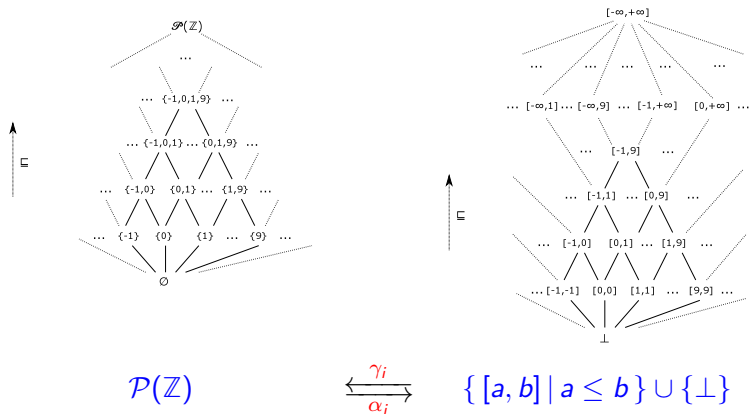
Effective construction for the **least upper bound** $A_1 \sqcup A_2$,
 able to represent properties expressible in either A_1 or A_2

Benefit

We can design **more precise** analyses
 by combining existing abstractions

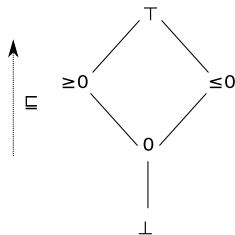
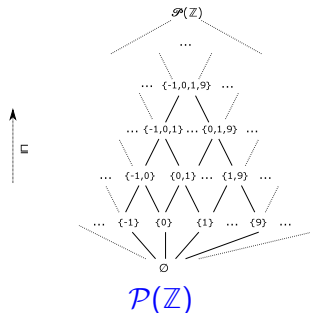
Abstract domain lattice

Reminder: interval abstraction



- $\alpha_i(S) \stackrel{\text{def}}{=} [\min S, \max S]$
- $\gamma_i([a, b]) \stackrel{\text{def}}{=} \{x \in \mathbb{Z} \mid a \leq x \leq b\}$

Reminder: sign abstraction

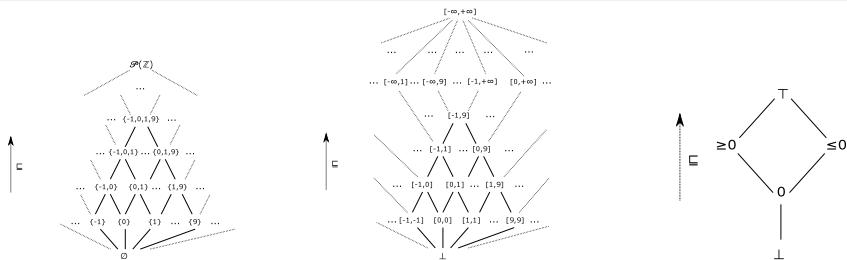


$$\begin{array}{c} \xleftarrow{\gamma_s} \\ \xrightarrow{\alpha_s} \end{array} \quad \{\perp, 0, \leq 0, \geq 0, \top\}$$

$$\begin{array}{lll} \gamma_s(\perp) & \stackrel{\text{def}}{=} & \emptyset \\ \gamma_s(0) & \stackrel{\text{def}}{=} & \{0\} \\ \gamma_s(\geq 0) & \stackrel{\text{def}}{=} & \mathbb{N} \\ \gamma_s(\leq 0) & \stackrel{\text{def}}{=} & -\mathbb{N} \\ \gamma_s(\top) & \stackrel{\text{def}}{=} & \mathbb{Z} \end{array}$$

$$\alpha_s(S) \stackrel{\text{def}}{=} \begin{cases} \perp & \text{if } S = \emptyset \\ 0 & \text{if } S = \{0\} \\ \geq 0 & \text{else, if } \forall s \in S, s \geq 0 \\ \leq 0 & \text{else, if } \forall s \in S, s \leq 0 \\ \top & \text{otherwise} \end{cases}$$

Composing abstractions



$$\mathcal{P}(\mathbb{Z}) \begin{array}{c} \xleftarrow{\gamma_i} \\ \xrightarrow{\alpha_i} \end{array} \{ [a, b] \mid a \leq b \} \cup \{ \perp \} \begin{array}{c} \xleftarrow{\gamma'_s} \\ \xrightarrow{\alpha'_s} \end{array} \{ \perp, 0, \leq 0, \geq 0, \top \}$$

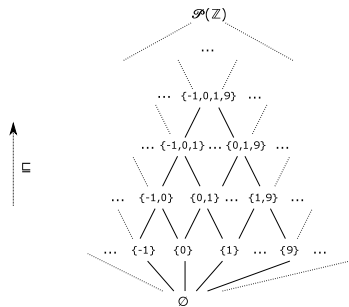
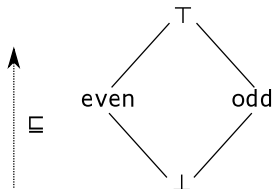
where: $\gamma'_s(\perp) \stackrel{\text{def}}{=} \perp$ $\gamma'_s(\top) \stackrel{\text{def}}{=} [-\infty, +\infty]$
 $\gamma'_s(\geq 0) \stackrel{\text{def}}{=} [0, +\infty]$ $\gamma'_s(\leq 0) \stackrel{\text{def}}{=} [-\infty, 0]$ $\gamma'_s(0) \stackrel{\text{def}}{=} [0, 0]$

We can **compose Galois connections**:

If $(X_1, \sqsubseteq_1) \begin{array}{c} \xleftarrow{\gamma_1} \\ \xrightarrow{\alpha_1} \end{array} (X_2, \sqsubseteq_2) \begin{array}{c} \xleftarrow{\gamma_2} \\ \xrightarrow{\alpha_2} \end{array} (X_3, \sqsubseteq_3)$, then
 $(X_1, \sqsubseteq_1) \begin{array}{c} \xleftarrow{\gamma_1 \circ \gamma_2} \\ \xrightarrow{\alpha_2 \circ \alpha_1} \end{array} (X_3, \sqsubseteq_3)$.

Proof: $(\alpha_2 \circ \alpha_1)(c) \sqsubseteq_3 a \iff \alpha_1(c) \sqsubseteq_2 \gamma_2(a) \iff c \sqsubseteq_1 (\gamma_1 \circ \gamma_2)(a)$

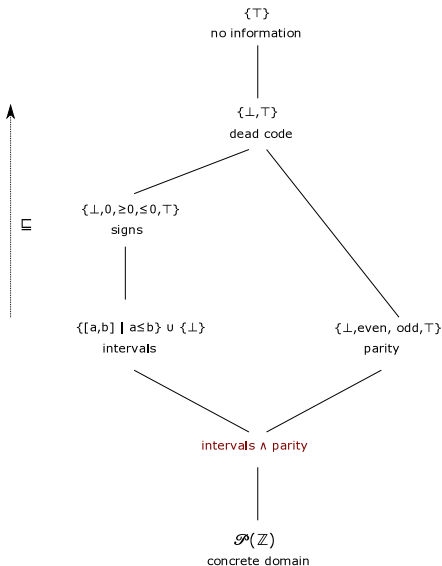
Parity domain

 $\mathcal{P}(\mathbb{Z})$ 

$$\begin{array}{c} \xrightarrow{\gamma_P} \\ \xleftarrow{\alpha_P} \end{array}$$
 $\{\perp, \top, \text{even}, \text{odd}\}$

$$\begin{array}{lll} \gamma_P(\perp) & \stackrel{\text{def}}{=} & \emptyset \\ \gamma_P(\text{even}) & \stackrel{\text{def}}{=} & 2\mathbb{Z} \\ \gamma_P(\text{odd}) & \stackrel{\text{def}}{=} & 2\mathbb{Z} + 1 \\ \gamma_P(\top) & \stackrel{\text{def}}{=} & \mathbb{Z} \end{array}$$

$$\alpha_P(S) \stackrel{\text{def}}{=} \begin{cases} \perp & \text{if } S = \emptyset \\ \text{even} & \text{else, if } S \subseteq 2\mathbb{Z} \\ \text{odd} & \text{else, if } S \subseteq 2\mathbb{Z} + 1 \\ \top & \text{otherwise} \end{cases}$$

Part of the abstraction lattice for $\mathcal{P}(\mathbb{Z})$ 

- signs are more abstract than intervals;
every sign information can be expressed as an interval
- parities and intervals are incomparable;
no common property, except \perp and \top
- $\mathcal{P}(\mathbb{Z})$ is the most concrete domain;
- $\{\top\}$ is the most abstract domain;
- **intervals \wedge parities** is the coarsest abstract domain more precise than intervals and parities.

Reduced product construction

Simple product

Algebraic structure:

Given two domains $(\mathcal{D}_1^\#, \sqsubseteq_1)$ and $(\mathcal{D}_2^\#, \sqsubseteq_2)$, we use **pairs** of abstract elements to represent **conjunctions** of properties.

- $\mathcal{D}_{1 \times 2}^\# \stackrel{\text{def}}{=} \mathcal{D}_1^\# \times \mathcal{D}_2^\#$
- $\gamma_{1 \times 2}(X_1^\#, X_2^\#) \stackrel{\text{def}}{=} \gamma_1(X_1^\#) \cap \gamma_2(X_2^\#)$
- $\alpha_{1 \times 2}(S) \stackrel{\text{def}}{=} (\alpha_1(S), \alpha_2(S))$
- $(X_1^\#, X_2^\#) \sqsubseteq_{1 \times 2} (Y_1^\#, Y_2^\#) \iff X_1^\# \sqsubseteq_1 Y_1^\# \text{ and } X_2^\# \sqsubseteq_2 Y_2^\#$

Abstract operators in $\mathcal{D}^\#$:

Applied in parallel (independently) in each abstract domain:

- $(X_1^\#, X_2^\#) \cup_{1 \times 2}^\# (Y_1^\#, Y_2^\#) \stackrel{\text{def}}{=} (X_1^\# \cup_1^\# Y_1^\#, X_2^\# \cup_2^\# Y_2^\#), ;$
- $(X_1^\#, X_2^\#) \nabla_{1 \times 2}^\# (Y_1^\#, Y_2^\#) \stackrel{\text{def}}{=} (X_1^\# \nabla_1^\# Y_1^\#, X_2^\# \nabla_2^\# Y_2^\#);$
- $S^\#[[s]]_{1 \times 2}(X_1^\#, X_2^\#) \stackrel{\text{def}}{=} (S^\#[[s]]_1(X_1^\#), S^\#[[s]]_2(X_2^\#)).$

Simple products: limitations

```

V ← 1;
while V ≤ 10 do V ← V + 2 done;
● if V ≥ 12 then ● V ← 0 ●;
  
```

Analysis in the product domain of intervals and parities:

	intervals	parities	product: intervals × parities
●	$V \in [11, 12]$	V odd	$(V \in [11, 12]) \wedge (V \text{ odd})$
●	$V = 12$	V odd	$(V = 12) \wedge (V \text{ odd})$
●	$V = 0$	V even	$(V = 0) \wedge (V \text{ even})$

Identical to two **separate** analyses:

- at ●, we get $(V = 12) \wedge (V \text{ odd})$, which represents \emptyset ;
- at ●, we apply $V \leftarrow 0$ **independently** on intervals and parities, which gives $(V = 0) \wedge (V \text{ even})$, instead of \emptyset !

⇒ huge loss of precision

Fully reduced product

Idea: propagate information between domains

Given Galois connections (α_1, γ_1) and (α_2, γ_2) over \mathcal{D}_1^\sharp and \mathcal{D}_2^\sharp , we define a **reduction operator** ρ as:

$$\rho : \mathcal{D}_{1 \times 2}^\sharp \rightarrow \mathcal{D}_{1 \times 2}^\sharp$$

$$\rho(X_1^\sharp, X_2^\sharp) \stackrel{\text{def}}{=} (\alpha_1(\gamma_1(X_1^\sharp) \cap \gamma_2(X_2^\sharp)), \alpha_2(\gamma_1(X_1^\sharp) \cap \gamma_2(X_2^\sharp)))$$

i.e., the best representation of $\gamma_{1 \times 2}(X_1^\sharp, X_2^\sharp)$ in both domains

Application:

use ρ to transfer information between domains **after abstract operations:**

- $(X_1^\sharp, X_2^\sharp) \cup_{1 \times 2}^\sharp (Y_1^\sharp, Y_2^\sharp) \stackrel{\text{def}}{=} \rho(X_1^\sharp \cup_1^\sharp Y_1^\sharp, X_2^\sharp \cup_2^\sharp Y_2^\sharp)$,
- $S^\sharp[[s]]_{1 \times 2}(X_1^\sharp, X_2^\sharp) \stackrel{\text{def}}{=} \rho(S^\sharp[[s]]_1(X_1^\sharp), S^\sharp[[s]]_2(X_2^\sharp))$.

Warning:

ρ should not be used on fixpoint iterates with widening $\nabla (X_{n+1} \stackrel{\text{def}}{=} \rho(X_n \nabla F(X_n)))$
 \implies this could prevent the convergence in $\mathcal{D}_1^\sharp \times \mathcal{D}_2^\sharp$!

Analysis with reduction: exemple

```

V ← 1;
while V ≤ 10 do V ← V + 2 done;
• if V ≥ 12 then V ← 0

```

Reduction ρ between intervals and parities $\rho([a, b], p)$:

First refine interval bounds $[a, b]$ using parity information p , then refine the parity information using the refined bounds:

- let $a' = a + 1$ if $a \notin \gamma_p(p)$, $a' = a$ otherwise;
- let $b' = b - 1$ if $b \notin \gamma_p(p)$, $b' = b$ otherwise;
- if $a' > b'$, return (\perp, \perp) ;
- if $a' = b'$, return $([a', b'], \alpha_p(a))$;
- otherwise, return $([a', b'], p)$.

Example:

At •, $\rho([11, 12], \text{odd}) = ([11, 11], \text{odd})$
 \implies the “then” branch is not reachable.

Partial reduction

The optimal reduction ρ is well-defined but:

- ρ assumes we have Galois connections;
 - there is no general effective algorithm to compute ρ .
- (similar to the case of optimal operators, defined as $F^\# \stackrel{\text{def}}{=} \alpha \circ F \circ \gamma$)

Partial reduction:

Practical definition, when the optimal reduction is not available:

- $\rho(X_1^\#, X_2^\#) = (Y_1^\#, Y_2^\#)$ is a **partial reduction** if:
- $Y_1^\# \sqsubseteq_1 X_1^\#$ and $Y_2^\# \sqsubseteq_2 X_2^\#$ (improvement)
- $\gamma_{1 \times 2}(Y_1^\#, Y_2^\#) = \gamma_{1 \times 2}(X_1^\#, X_2^\#)$ (soundness)

Example:

$$\rho(X_1^\#, X_2^\#) \stackrel{\text{def}}{=} \begin{cases} (\perp_1, \perp_2) & \text{if } X_1^\# = \perp_1 \text{ or } X_2^\# = \perp_2 \\ (X_1^\#, X_2^\#) & \text{otherwise} \end{cases}$$

In practice, an analyzer contains many abstract domains (for expressiveness) with limited reductions between them (for efficiency).

Disjunctive domains

Motivation

Remark: most domains abstract **convex sets** (conjunctions of constraints)

$\Rightarrow \cup^\#$ causes a loss of precision!

The need for non-convex invariants

```

X ← rand(10, 20);
Y ← rand(0, 1);
if Y > 0 then X ← -X;
• Z ← 100/X
  
```

Concrete semantics:

At •, $X \in [-20, -10] \cup [10, 20]$

\Rightarrow there is **no division by zero**

Abstract analysis:

Convex analyses (intervals, polyhedra) will find $X \in [-20, 20]$

(with intervals, $[-20, -10] \cup^\# [10, 20] = [-20, 20]$)

\Rightarrow **possible division by zero**

(false alarm)

Disjunctive domains

Principle:

generic constructions to **lift** any numeric abstract domain to a domain able to represent disjunctions exactly

Example constructions:

- **powerset** completion
unordered “soup” of abstract elements
- **state partitioning**
abstract elements keyed to selected subsets of environments
- **path-sensitive** analyses
partition with respect to the **history** of execution

each construction has its strength and weakness
they can be combined during an analysis to exploit the best of each

Powerset completion

Powerset completion

Given: $(\mathcal{E}^\#, \sqsubseteq, \gamma, \cup^\#, \cap^\#, \nabla, S^\#[\text{stat}])$

abstract domain $\mathcal{E}^\#$

ordered by \sqsubseteq , which also acts as a sound abstraction of \subseteq (i.e., $\subseteq^\# = \sqsubseteq$)

with concretization $\gamma : \mathcal{E}^\# \rightarrow \mathcal{P}(\mathcal{E})$

sound abstractions $\cup^\#, \cap^\#, S^\#[\text{stat}]$ of $\cup, \cap, S[\text{stat}]$, and a widening ∇

Construct: $(\hat{\mathcal{E}}^\#, \hat{\sqsubseteq}, \hat{\gamma}, \hat{\cup}^\#, \hat{\cap}^\#, \hat{\nabla}, \hat{S}^\#[\text{stat}])$

- $\hat{\mathcal{E}}^\# \stackrel{\text{def}}{=} \mathcal{P}_{\text{finite}}(\mathcal{E}^\#)$ (finite sets of abstract elements)
- $\hat{\gamma}(A^\#) \stackrel{\text{def}}{=} \cup \{ \gamma(X^\#) \mid X^\# \in A^\# \}$ (join of concretizations)

Example: using the interval domain as $\mathcal{E}^\#$

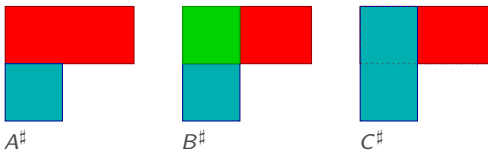
$$\hat{\gamma}(\{[-10, -5], [2, 4], [0, 0], [2, 3]\}) = [-10, -5] \cup \{0\} \cup [2, 4]$$

Ordering

Issue: how can we compare two elements of $\hat{\mathcal{E}}^\sharp$?

- $\hat{\gamma}$ is generally not injective
there is no canonical representation for $\hat{\gamma}(A^\sharp)$
- testing $\hat{\gamma}(A^\sharp) = \hat{\gamma}(B^\sharp)$ or $\hat{\gamma}(A^\sharp) \subseteq \hat{\gamma}(B^\sharp)$ is difficult

Example: powerset completion of the interval domain



$$A^\sharp = \{\{0\} \times \{0\}, [0, 1] \times \{1\}\}$$

$$B^\sharp = \{\{0\} \times \{0\}, \{0\} \times \{1\}, \{1\} \times \{1\}\}$$

$$C^\sharp = \{\{0\} \times [0, 1], [0, 1] \times \{1\}\}$$

$$\hat{\gamma}(A^\sharp) = \hat{\gamma}(B^\sharp) = \hat{\gamma}(C^\sharp)$$

B^\sharp is more costly to represent: it requires three abstract elements instead of two
 C^\sharp is a covering and not a partition (red \cap blue = $\{0\} \times \{1\} \neq \emptyset$)

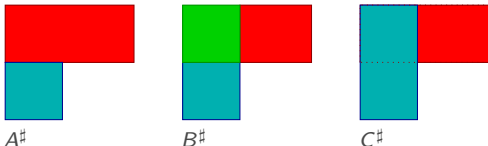
Ordering (cont.)

Solution: sound approximation of \sqsubseteq

$$A^\# \hat{\sqsubseteq} B^\# \stackrel{\text{def}}{\iff} \forall X^\# \in A^\#: \exists Y^\# \in B^\#: X^\# \sqsubseteq Y^\# \quad (\text{Hoare powerdomain order})$$

- $\hat{\sqsubseteq}$ is a partial order (when \sqsubseteq is)
- $\hat{\sqsubseteq}$ is a sound approximation of \sqsubseteq (when \sqsubseteq is)
 $A^\# \hat{\sqsubseteq} B^\# \implies \hat{\gamma}(A^\#) \sqsubseteq \hat{\gamma}(B^\#)$ but the converse may not hold
- testing $\hat{\sqsubseteq}$ reduces to testing \sqsubseteq finitely many times

Example: powerset completion of the interval domain



$$\hat{\gamma}(A^\#) = \hat{\gamma}(B^\#) = \hat{\gamma}(C^\#)$$

$$B^\# \hat{\sqsubseteq} A^\# \hat{\sqsubseteq} C^\#$$

Abstract operations

Abstract operators

- $\hat{S}^\# \llbracket stat \rrbracket A^\# \stackrel{\text{def}}{=} \{ S^\# \llbracket stat \rrbracket X^\# \mid X^\# \in A^\# \}$
 apply *stat* on each abstract element independently
- $A^\# \hat{\cup}^\# B^\# \stackrel{\text{def}}{=} A^\# \cup B^\#$
 keep elements from both arguments without applying any abstract operation
 $\hat{\cup}^\#$ is **exact**
- $A^\# \hat{\cap}^\# B^\# \stackrel{\text{def}}{=} \{ X^\# \cap^\# Y^\# \mid X^\# \in A^\#, Y^\# \in B^\# \}$
 $\hat{\cap}^\#$ is **exact** if $\cap^\#$ is (as \cup and \cap are distributive)

Galois connection:

in general, there is **no abstraction function** $\hat{\alpha}$ corresponding to $\hat{\gamma}$

Example: powerset completion $\hat{\mathcal{E}}^\#$ of the interval domain $\mathcal{E}^\#$

given the disc $S \stackrel{\text{def}}{=} \{ (x, y) \mid x^2 + y^2 \leq 1 \}$

$\alpha(S) = [-1, 1] \times [-1, 1]$ (optimal interval abstraction)

but there is no best abstraction in $\hat{\mathcal{E}}^\#$



$\alpha(S)$



not $\hat{\alpha}(S)$

Dynamic approximation

Issue: the size $|A^\#|$ of elements $A^\# \in \hat{\mathcal{E}}^\#$ is unbounded
 every application of $\hat{\cup}^\#$ adds some more elements
 \implies efficiency and convergence problems

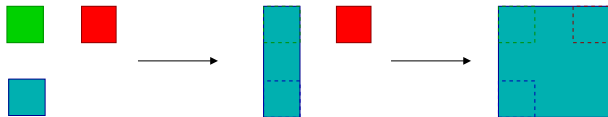
Solution: to reduce the size of elements

- redundancy removal

$simplify(A^\#) \stackrel{\text{def}}{=} \{X^\# \in A^\# \mid \forall Y^\# \neq X^\# \in A^\#: X^\# \not\sqsubseteq Y^\#\}$
 no loss of precision: $\hat{\gamma}(simplify(A^\#)) = \hat{\gamma}(A^\#)$

- collapse: join elements in $\mathcal{E}^\#$

$collapse(A^\#) \stackrel{\text{def}}{=} \{\cup^\# \{X^\# \in A^\#\}\}$



large loss of precision, but very effective: $|collapse(A^\#)| = 1$

- partial collapse: limit $|A^\#|$ to a fixed size k by $\cup^\#$
 but how to choose which elements to merge? no easy solution!

Widening

Issue: for loops, abstract iterations $(A_n^\#)_{n \in \mathbb{N}}$ may not converge

- the size of $A_n^\#$ may grow arbitrarily large
- even if $|A_n^\#|$ is stable, some elements in $A_n^\#$ may not converge if $\mathcal{E}^\#$ has infinite increasing sequences

\implies we need a **widening** $\hat{\vee}$

Widenings for powerset domains are **difficult to design**

Example widening: collapse after a fixed number N of iterations

$$A_{n+1}^\# \stackrel{\text{def}}{=} A_n^\# \hat{\vee} B_n^\# \stackrel{\text{def}}{=} \begin{cases} \text{simplify}(A_n^\# \hat{\cup}^\# B_n^\#) & \text{if } n < N \\ \text{collapse}(A_n^\#) \nabla \text{collapse}(B_n^\#) & \text{otherwise} \end{cases}$$

(this is very naïve, see Bagnara et al. STTT06 for more interesting widenings)

State partitioning

State partitioning

Principle:

- partition *a priori* \mathcal{E} into **finitely** many sets
- abstract each **partition of \mathcal{E} independently** using an element of \mathcal{E}^\sharp

Abstract domain:

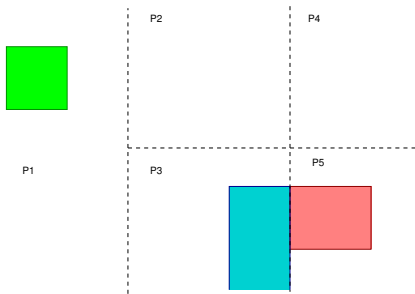
Given an abstract partition $P^\sharp \subseteq \mathcal{E}^\sharp$, i.e., a set such that:

- P^\sharp is finite
- $\bigcup \{ \gamma(X^\sharp) \mid X^\sharp \in P^\sharp \} = \mathcal{E}$
for generality, we have in fact a covering, not a partitioning of \mathcal{E}
i.e., we can have $X^\sharp \neq Y^\sharp \in P^\sharp$ with $\gamma(X^\sharp) \cap \gamma(Y^\sharp) \neq \emptyset$

We define $\tilde{\mathcal{E}}^\sharp \stackrel{\text{def}}{=} P^\sharp \rightarrow \mathcal{E}^\sharp$

representable in memory, as P^\sharp is finite

Ordering



Example: $\mathcal{E}^\#$ is the interval domain

$P^\# = \{P_1, P_2, P_3, P_4, P_5\}$ where

$$P_1 = [-\infty, 0] \times [-\infty, +\infty]$$

$$P_2 = [0, 10] \times [0, +\infty]$$

$$P_3 = [0, 10] \times [-\infty, 0]$$

$$P_4 = [10, +\infty] \times [0, +\infty]$$

$$P_5 = [10, +\infty] \times [-\infty, 0]$$

$$X^\# = [P_1 \mapsto [-6, -5] \times [5, 6],$$

$$P_2 \mapsto \perp,$$

$$P_3 \mapsto [9, 10] \times [-\infty, -1],$$

$$P_4 \mapsto \perp,$$

$$P_5 \mapsto [10, 12] \times [-3, -1]]$$

- $\tilde{\mathcal{E}}^\# \stackrel{\text{def}}{=} P^\# \rightarrow \mathcal{E}^\#$
- $\tilde{\gamma}(A^\#) \stackrel{\text{def}}{=} \cup \{ \gamma(A^\#(X^\#)) \cap \gamma(X^\#) \mid X^\# \in P^\# \}$
- $A^\# \stackrel{\sim}{\sqsubseteq} B^\# \stackrel{\text{def}}{\iff} \forall X^\# \in P^\#: A^\#(X^\#) \sqsubseteq B^\#(X^\#)$ (point-wise order)
- $\tilde{\alpha}(S) \stackrel{\text{def}}{=} \lambda X^\# \in P^\#. \alpha(S \cap \gamma(X^\#))$
if $\mathcal{E}^\#$ enjoys a **Galois connection**, so does $\tilde{\mathcal{E}}^\#$

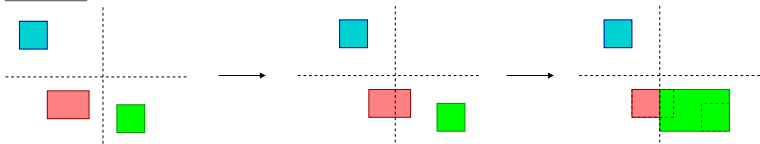
Abstract operators

Abstract operators: point-wise extension from $\mathcal{E}^\#$ to $P^\# \rightarrow \mathcal{E}^\#$

- $A \overset{\sim}{\cup} B \stackrel{\text{def}}{=} \lambda X^\# \in P^\#. A(X^\#) \cup^\# B(X^\#)$
- $A \overset{\sim}{\cap} B \stackrel{\text{def}}{=} \lambda X^\# \in P^\#. A(X^\#) \cap^\# B(X^\#)$
- $A \overset{\sim}{\vee} B \stackrel{\text{def}}{=} \lambda X^\# \in P^\#. A(X^\#) \vee^\# B(X^\#)$
- $\tilde{S}^\#[[e \leq 0?]] A^\# \stackrel{\text{def}}{=} \lambda X^\# \in P^\#. S^\#[[e \leq 0?]] A^\#(X^\#)$
- $\tilde{S}^\#[[V \leftarrow e]] A^\#$ is **more complex**

any $S^\#[[V \leftarrow e]] A^\#(X^\#)$ may escape its partition $X^\#$; we must cut them at partition borders and glue the pieces falling into the same partition

example: $X \leftarrow X + 2$



$$\tilde{S}^\#[[V \leftarrow e]] A^\# \stackrel{\text{def}}{=} \lambda X^\#. \cup^\# \{ X^\# \cap^\# S^\#[[V \leftarrow e]] A^\#(Y^\#) \mid Y^\# \in P^\# \}$$

Example analysis

Example

```

X ← rand(10, 20);
Y ← rand(0, 1);
if Y > 0 then X ← -X;
• Z ← 100/X

```

Analysis:

- $\mathcal{E}^\#$ is the interval domain
- partition with respect to the sign of X
 $P^\# \stackrel{\text{def}}{=} \{X^+, X^-\}$ where
 $X^+ \stackrel{\text{def}}{=} [0, +\infty] \times \mathbb{Z} \times \mathbb{Z}$ and $X^- \stackrel{\text{def}}{=} [-\infty, 0] \times \mathbb{Z} \times \mathbb{Z}$
- at • we find:
 $X^+ \mapsto [X \in [10, 20], Y \mapsto [0, 0], Z \mapsto [0, 0]]$
 $X^- \mapsto [X \in [-20, -10], Y \mapsto [1, 1], Z \mapsto [0, 0]]$
 \implies no division by zero

Path partitioning

Path sensitivity

Principle: partition wrt. the **history of computation**

- keep different abstract elements for different execution **paths**
e.g., different branches taken, different loop iterations
- **avoid** merging with $\cup^\#$ elements at control-flow **joins**
at the end of **if** \dots **then** \dots **else**, or at loop head

Intuition: as a program transformation

```
X ← rand(-50, 50);
if X ≥ 0 then
  Y ← X + 10
else
  Y ← X - 10;
assert Y ≠ 0
```

→

```
X ← rand(-50, 50);
if X ≥ 0 then
  Y ← X + 10;
  assert Y ≠ 0
else
  Y ← X - 10;
  assert Y ≠ 0
```

the **assert** is tested in the context of each branch
instead of after the control-flow join

the interval domain can prove the assertion on the right, but not on the left

Abstract domain

Formalization: we consider here only **if** \dots **then** \dots **else**

- \mathcal{L} denote **syntactic labels** of **if** \dots **then** \dots **else** instructions
- **history abstraction** $\mathbb{H} \stackrel{\text{def}}{=} \mathcal{L} \rightarrow \{\text{true}, \text{false}, \perp\}$

$H \in \mathbb{H}$ indicates the outcome of the last time we executed each test:

- $H(\ell) = \text{true}$: we took the **then** branch
- $H(\ell) = \text{false}$: we took the **else** branch
- $H(\ell) = \perp$: we never executed the test

Notes:

- \mathbb{H} can remember the outcome of several successive tests
 $\ell_1 : \text{if } \dots \text{ then } \dots \text{ else}; \ell_2 : \text{if } \dots \text{ then } \dots \text{ else}$
- for tests in loops, \mathbb{H} remembers only the last outcome
while \dots **do** $\ell : \text{if } \dots \text{ then } \dots \text{ else}$
- we could extend \mathbb{H} to longer histories with $\mathbb{H} = (\mathcal{L} \rightarrow \{\text{true}, \text{false}, \perp\})^*$
- we could extend \mathbb{H} to track loop iterations with $\mathbb{H} = \mathcal{L} \rightarrow \mathbb{N}$

- $\mathcal{E}^\# \stackrel{\text{def}}{=} \mathbb{H} \rightarrow \mathcal{E}^\#$

use a different abstract element for each abstract history

Abstract operators

- $\mathcal{E}^\# \stackrel{\text{def}}{=} \mathbb{H} \rightarrow \mathcal{E}^\#$
- $\gamma(A^\#) = \bigcup \{ \gamma(A^\#(H)) \mid H \in \mathbb{H} \}$
- $\underline{\llcorner}, \check{\cup}^\#, \check{\lrcorner}^\#, \check{\vee}$ are **point-wise**
- $\check{S}^\# \llbracket V \leftarrow e \rrbracket$ and $\check{S}^\# \llbracket e \leq 0? \rrbracket$ are **point-wise**
- $\check{S}^\# \llbracket \ell : \text{if } c \text{ then } s_1 \text{ else } s_2 \rrbracket A^\#$ is more complex
 - we **merge** all information about ℓ

$$C^\# = \lambda H. A^\#(H[\ell \mapsto \text{true}]) \cup^\# A^\#(H[\ell \mapsto \text{false}]) \cup^\# A^\#(H[\ell \mapsto \perp])$$
 - we compute the **then branch**, where $H(\ell) = \text{true}$

$$T'^\# = \check{S}^\# \llbracket s_1 \rrbracket (\check{S}^\# \llbracket c? \rrbracket T^\#)$$
 where

$$T^\# = \lambda H. C^\#(H) \text{ if } H(\ell) = \text{true}, \perp \text{ otherwise}$$
 - we compute the **else branch**, where $H(\ell) = \text{false}$

$$F'^\# = \check{S}^\# \llbracket s_2 \rrbracket (\check{S}^\# \llbracket \neg c? \rrbracket F^\#)$$
 where

$$F^\# = \lambda H. C^\#(H) \text{ if } H(\ell) = \text{false}, \perp \text{ otherwise}$$
 - we **join** both branches: $T'^\# \check{\cup}^\# F'^\#$
 the join is exact as $\forall H \in \mathbb{H}$: either $T'^\#(H) = \perp$ or $F'^\#(H) = \perp$

\implies we get a semantic by induction on the syntax of the original program

Complex example

Linear interpolation

```

X ← rand(TX[0], TX[N]);
I ← 0;
while I < N ∧ X > TX[I + 1] do
    I ← I + 1;
done;
Y ← TY[I] + (X - TX[I]) × TS[I]

```

Concrete semantics: table-based interpolation based on the value of X

- look-up index I in the interpolation table: $TX[I] \leq X \leq TX[I + 1]$
- interpolate from value $TY[I]$ when $X = TX[I]$ with slope $TS[I]$

Analysis: in the interval domain

- without partitioning:
 $Y \in [\min TY, \max TY] + (X - [\min TX, \max TX]) \times [\min TS, \max TS]$
- partitioning with respect to the **number of loop iterations**:
 $Y \in \bigcup_{I \in [0, M]} TY[I] + ([0, TX[I + 1] - TX[I]) \times TS[I]$
more precise as it keeps the relation between table indices

Inter-procedural analyses

Overview

- Analysis on the **control-flow graph**
reduce function calls and returns to *gotos*
useful for the project!
- **Inlining**
simple and precise
but not efficient and may not terminate
- **Call-site** and **call-stack abstraction**
terminates even for recursive programs
parametric cost-precision trade-off
- **Tabulated abstraction**
optimal reuse of analysis partial results
- We also mentioned summary-based abstractions last week,
leveraging relational domains for modular bottom-up analysis

in general, these different abstractions give incomparable results;
there is no clear winner

Analysis on the control-flow graph

Inter-procedural control-flow graphs

Extend control-flow graphs:

- one **subgraph** for each function
- additional **arcs** to denote function **calls** and **returns**

we get one big graph without procedures nor calls, only **gotos**

⇒ reduced to a **classic analysis** based on **equation systems**

but difficult to use in a denotational-style analysis by induction on the syntax

Note: to simplify, we assume here **no local variable and no function argument**:

- locals and arguments are transformed into globals
- only possible if there are no recursive calls

Example: Control-flow graph

Example

main :

```

R ← -1;
X ← rand(5, 10); f();
X ← 80; f()

```

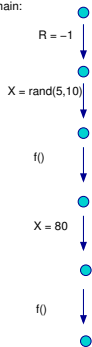
f :

```

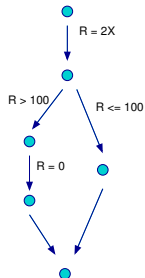
R ← 2 × X;
if R > 100 then R ← 0

```

main:



f:



create one control-flow graph for each function

Example: Control-flow graph

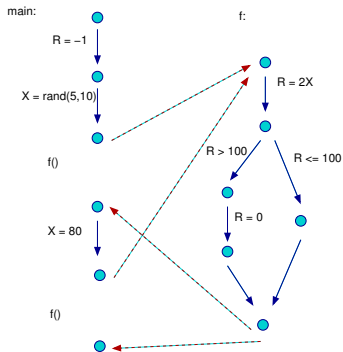
Example

main :

```
R ← -1;
X ← rand(5, 10); f();
X ← 80; f()
```

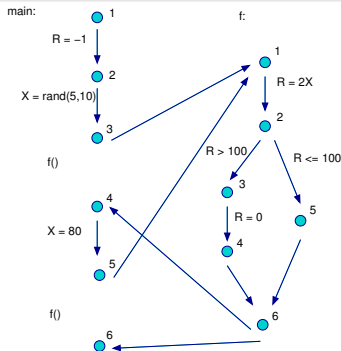
f :

```
R ← 2 × X;
if R > 100 then R ← 0
```



replace **call** instructions with **gotos**

Example: Equation system

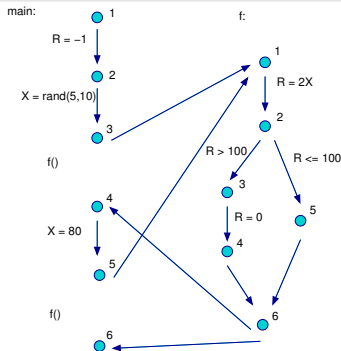


$$\begin{aligned} \mathcal{S}_{\text{main},1} &= \top \\ \mathcal{S}_{\text{main},2} &= \mathcal{S}[\![R \leftarrow 1]\!] \mathcal{S}_{\text{main},1} \\ \mathcal{S}_{\text{main},3} &= \mathcal{S}[\![X \leftarrow \text{rand}(5, 10)]\!] \mathcal{S}_{\text{main},2} \\ \mathcal{S}_{\text{main},4} &= \mathcal{S}_{f,6} \\ \mathcal{S}_{\text{main},5} &= \mathcal{S}[\![X \leftarrow 80]\!] \mathcal{S}_{\text{main},4} \\ \mathcal{S}_{\text{main},6} &= \mathcal{S}_{f,6} \end{aligned}$$

$$\begin{aligned} \mathcal{S}_{f,1} &= \mathcal{S}_{\text{main},3} \cup \mathcal{S}_{\text{main},5} \\ \mathcal{S}_{f,2} &= \mathcal{S}[\![R \leftarrow 2X]\!] \mathcal{S}_{f,1} \\ \mathcal{S}_{f,3} &= \mathcal{S}[\![R > 100]\!] \mathcal{S}_{f,2} \\ \mathcal{S}_{f,4} &= \mathcal{S}[\![R \leftarrow 0]\!] \mathcal{S}_{f,3} \\ \mathcal{S}_{f,5} &= \mathcal{S}[\![R \leq 100]\!] \mathcal{S}_{f,2} \\ \mathcal{S}_{f,6} &= \mathcal{S}_{f,4} \cup \mathcal{S}_{f,5} \end{aligned}$$

- each variable \mathcal{S}_i denotes a set of environments at a control location i
- we can derive an abstract version of the system
e.g.: $\mathcal{S}_{f,2}^\sharp = \mathcal{S}^\sharp[\![R \leftarrow 2X]\!] \mathcal{S}_{f,1}^\sharp$, $\mathcal{S}_{f,6}^\sharp = \mathcal{S}_{f,4}^\sharp \cup^\sharp \mathcal{S}_{f,5}^\sharp$, etc.
- we can solve the abstract system, using widenings to terminate
c.f. project

Example: Equation system



$$\begin{aligned}
 S_{\text{main},1} &= \top \\
 S_{\text{main},2} &= S[R \leftarrow 1] S_{\text{main},1} \\
 S_{\text{main},3} &= S[X \leftarrow \mathbf{rand}(5, 10)] S_{\text{main},2} \\
 S_{\text{main},4} &= S_{f,6} \\
 S_{\text{main},5} &= S[X \leftarrow 80] S_{\text{main},4} \\
 S_{\text{main},6} &= S_{f,6}
 \end{aligned}$$

$$\begin{aligned}
 S_{f,1} &= S_{\text{main},3} \cup S_{\text{main},5} \\
 S_{f,2} &= S[R \leftarrow 2X] S_{f,1} \\
 S_{f,3} &= S[R > 100] S_{f,2} \\
 S_{f,4} &= S[R \leftarrow 0] S_{f,3} \\
 S_{f,5} &= S[R \leq 100] S_{f,2} \\
 S_{f,6} &= S_{f,4} \cup S_{f,5}
 \end{aligned}$$

using intervals we get the following solution:

$$S_{\text{main},1}^{\#} : X, R \in \mathbb{Z}$$

$$S_{\text{main},2}^{\#} : X \in \mathbb{Z}, R = -1$$

$$S_{\text{main},3}^{\#} : X \in [5, 10], R = -1$$

$$S_{\text{main},4}^{\#} : X \in [5, 80], R \in [0, 100]$$

$$S_{\text{main},5}^{\#} : X = 80, R \in [0, 100]$$

$$S_{\text{main},6}^{\#} : X \in [5, 80], R \in [0, 100]$$

$$S_{f,1}^{\#} : X \in [5, 80], R \in [-1, 100]$$

$$S_{f,2}^{\#} : X \in [5, 80], R \in [10, 160]$$

$$S_{f,3}^{\#} : X \in [5, 80], R \in [101, 160]$$

$$S_{f,4}^{\#} : X \in [5, 80], R = 0$$

$$S_{f,5}^{\#} : X \in [5, 80], R \in [10, 100]$$

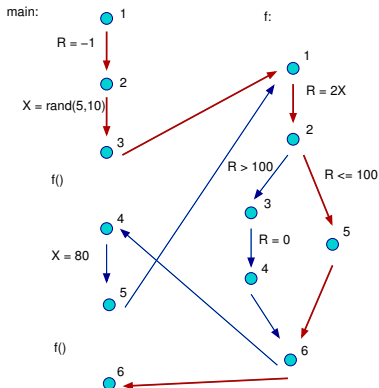
$$S_{f,6}^{\#} : X \in [5, 80], R \in [0, 100]$$

Imprecision

In fact, in our example, $R = 0$ holds at the end of the program
but we find $R \in [0, 100]$!

\implies the analysis is **imprecise**

Explanation: the control-flow graph adds **impossible executions paths**



General case: concrete semantics

Procedures

Syntax:

- \mathcal{F} finite set of procedure names
- $body : \mathcal{F} \rightarrow stat$: procedure bodies
- $main \in stat$: entry point body
- \mathbb{V}_G : set of **global** variables
- \mathbb{V}_f : set of **local** variables for procedure $f \in \mathcal{F}$
 procedure f can only access $\mathbb{V}_f \cup \mathbb{V}_G$
 $main$ has no local variable and can only access \mathbb{V}_G
- $stat ::= f(expr_1, \dots, expr_{|\mathbb{V}_f|}) \mid \dots$

procedure call, $f \in \mathcal{F}$, setting all its local variables

local variables double as **procedure arguments**

no special mechanism to return a value (a global variable can be used)

Concrete environments

Notes:

- when f calls g , we must **remember** the value of f 's locals \mathbb{V}_f in the semantics of g and **restore** them when returning
- several copies** of each $V \in \mathbb{V}_f$ may exist at a given time due to recursive calls, i.e.: cycles in the call graph

⇒ concrete environments use **per-variable stacks**

Stacks: $\mathcal{S} \stackrel{\text{def}}{=} \mathbb{Z}^*$ (finite sequences of integers)

- push**(v, s) $\stackrel{\text{def}}{=} v \cdot s$ ($v, v' \in \mathbb{Z}, s, s' \in \mathcal{S}$)
- pop**(s) $\stackrel{\text{def}}{=} s'$ when $\exists v: s = v \cdot s'$, undefined otherwise
- peek**(s) $\stackrel{\text{def}}{=} v$ when $\exists s': s = v \cdot s'$, undefined otherwise
- set**(v, s) $\stackrel{\text{def}}{=} v \cdot s'$ when $\exists v': s = v' \cdot s'$, undefined otherwise

Environments: $\mathcal{E} \stackrel{\text{def}}{=} (\cup_{f \in \mathcal{F}} \mathbb{V}_f \cup \mathbb{V}_G) \rightarrow \mathcal{S}$

for \mathbb{V}_G , stacks are not necessary but simplify the presentation

traditionally, there is a single global stack for all local variables

using per-variable stacks instead also makes the presentation simpler

Concrete semantics

Concrete semantics: on $\mathcal{E} \stackrel{\text{def}}{=} (\cup_{f \in \mathcal{F}} \mathbb{V}_f \cup \mathbb{V}_G) \rightarrow \mathcal{S}$

variable reads and updates only consider the **top of the stack**;

procedure calls **push** and **pop** local variables

- $E[V] \rho \stackrel{\text{def}}{=} \text{peek}(\rho(V))$
- $S[V \leftarrow e] R \stackrel{\text{def}}{=} \{ \rho[V \mapsto \text{set}(x, \rho(V))] \mid \rho \in R, x \in E[e] \rho \}$
- $S[f(e_{V_1}, \dots, e_{V_n})] R = R_3$, where:
 - $R_1 \stackrel{\text{def}}{=} \{ \rho[\forall V \in \mathbb{V}_f: V \mapsto \text{push}(x_V, \rho(V))] \mid \rho \in R, \forall V \in \mathbb{V}_f: x_V \in E[e_V] \rho \}$
(evaluate each argument e_V and push its value x_V on the stack $\rho(V)$)
 - $R_2 \stackrel{\text{def}}{=} S[\text{body}(f)] R_1$ (evaluate the procedure body)
 - $R_3 \stackrel{\text{def}}{=} \{ \rho[\forall V \in \mathbb{V}_f: V \mapsto \text{pop}(\rho(V))] \mid \rho \in R_2 \}$ (pop local variables)
- initial environment: $\rho_0 \stackrel{\text{def}}{=} \lambda V \in \mathbb{V}_G. 0$

other statements are unchanged

Semantic inlining

Semantic inlining

Naïve abstract procedure call: mimic the concrete semantics

- assign **abstract variables** to stack positions:

$$\mathbb{V}^\# \stackrel{\text{def}}{=} \mathbb{V}_G \cup (\cup_{f \in \mathcal{F}} \mathbb{V}_f \times \mathbb{N})$$

$\mathbb{V}^\#$ is **infinite**, but each abstract environment uses finitely many variables

- $\mathcal{E}_\mathbb{V}^\#$ abstracts $\mathcal{P}(\mathbb{V} \rightarrow \mathbb{Z})$, for any finite $\mathbb{V} \subseteq \mathbb{V}^\#$

$V \in \mathbb{V}_f$ denotes $(V, 0)$ in $\mathbb{V}^\#$

push V : shift variables, replacing (V, i) with $(V, i + 1)$, then add $(V, 0)$

pop V : remove $(V, 0)$ and shift each (V, i) to $(V, i - 1)$

- $S^\# \llbracket f(e_1, \dots, e_n) \rrbracket X^\#$ is then reduced to:

$$X_1^\# = S^\# \llbracket \mathbf{push} V_1; \dots; \mathbf{push} V_n \rrbracket X^\# \quad (\text{add fresh variables for } \mathbb{V}_f)$$

$$X_2^\# = S^\# \llbracket V_1 \leftarrow e_1; \dots; V_n \leftarrow e_n \rrbracket X_1^\# \quad (\text{bind arguments to locals})$$

$$X_3^\# = S^\# \llbracket \mathit{body}(f) \rrbracket X_2^\# \quad (\text{execute the procedure body})$$

$$X_4^\# = S^\# \llbracket \mathbf{pop} V_1; \dots; \mathbf{pop} V_n \rrbracket X_3^\# \quad (\text{delete local variables})$$

Limitations:

- does not terminate in case of **unbounded recursivity**
- requires **many abstract variables** to represent the stacks
- procedures must be **re-analyzed for every call**

full context-sensitivity: precise but costly

Example

Example

main :

```
R ← -1;  
f(rand(5, 10));  
f(80)
```

f(X) :

```
R ← 2 × X;  
if R > 100 then R ← 0
```

Analysis using intervals

- after the first call to f , we get $R \in [10, 20]$
- after the second call to f , we get $R = 0$

Call-site abstraction

Call-site abstraction

Abstracting stacks: into a **fixed, bounded** set $\mathbb{V}^\#$ of variables

- $\mathbb{V}^\# \stackrel{\text{def}}{=} \bigcup_{f \in \mathcal{F}} \{V, \hat{V} \mid V \in \mathbb{V}_f\} \cup \mathbb{V}_G$
 two copies of each local variable
 V abstracts the value at the top of the stack (current call)
 \hat{V} abstracts the rest of the stack
- $S^\#[\text{push } V] X^\# \stackrel{\text{def}}{=} X^\# \cup^\# S^\#[\hat{V} \leftarrow V] X^\#$
 $S^\#[\text{pop } V] X^\# \stackrel{\text{def}}{=} X^\# \cup^\# S^\#[V \leftarrow \hat{V}] X^\#$
 weak updates, similar to array manipulation
 no need to create and delete variables dynamically
- assignments and tests always access V , not \hat{V}
 \implies strong update (precise)

Note: when there is no recursivity, \hat{V} , **push** and **pop** can be omitted

Call-site abstraction

Principle: merge all the contexts in which each function is called

- we maintain two global maps $\mathcal{F} \rightarrow \mathcal{E}^\sharp$:
 - $C^\sharp(f)$: abstracts the environments when calling f
 - $R^\sharp(f)$: abstracts the environments when returning from f
gather environments from all possible calls to f , disregarding the call sites
- during the analysis, when encountering a call $S^\sharp[\![\text{body}(f)]\!] X^\sharp$:
 - we return $R^\sharp(f)$
 - but we also replace C^\sharp with $C^\sharp[f \mapsto C^\sharp(f) \cup^\sharp X^\sharp]$
- $R^\sharp(f)$ is computed from $C^\sharp(f)$ as

$$R^\sharp(f) = S^\sharp[\![\text{body}(f)]\!](C^\sharp(f))$$

Call-site abstraction

Fixpoint:

there may be **circular dependencies** between C^\sharp and R^\sharp

e.g., in $f(2); f(3)$, the input for $f(3)$ depends on the output from $f(2)$

\implies we compute a fixpoint for C^\sharp by iteration:

- initially, $\forall f: C^\sharp(f) = R^\sharp(f) = \perp$
- analyze *main*
- while $\exists f: C^\sharp(f)$ **not stable**
 - apply **widening** ∇ to the iterates of $C^\sharp(f)$
 - update** $R^\sharp(f) = S^\sharp \llbracket \text{body}(f) \rrbracket C^\sharp(f)$
 - analyze** *main* and all the procedures **again**
(this may modify some $C^\sharp(g)$)

\implies using ∇ , the analysis always terminates in finite time

we can be more efficient and avoid re-analyzing procedures when not needed

e.g., use a workset algorithm, track procedure dependencies, etc.

Example

Example

main :

```
R ← -1;
f(rand(5, 10));
f(80)
```

f(*X*) :

```
R ← 2 × X;
if R > 100 then R ← 0
```

Analysis: using intervals (without widening as there is no dependency)

- first analysis of *main*: we get \perp (as $R^\sharp(f) = \perp$)
but $C^\sharp(f) = [R \mapsto [-1, -1], X \mapsto [5, 10]]$
- first analysis of *f*: $R^\sharp(f) = [R \mapsto [10, 20], X \mapsto [5, 10]]$
- second analysis of *main*: we get
 $C^\sharp(f) = [R \mapsto [-1, 20], X \mapsto [5, 80]]$
- second analysis of *f*: $R^\sharp(f) = [R \mapsto [0, 100], X \mapsto [5, 80]]$
- final analysis of *main*, we find $R \in [0, 100]$ at the program end
less precise than $R = 0$ found by semantic inlining

Partial context-sensitivity

Variants: k -limiting, k is a constant

- **stack:**

assign a distinct variable for the k highest levels of V
 abstract the lower (unbounded) stack part with \hat{V}
 more precise than keeping only the top of the stack separately

- **context-sensitivity:**

each syntactic call has a unique **call-site** $\ell \in \mathcal{L}$
 a call stack is a sequence of nested call sites: $c \in \mathcal{L}^*$
 an **abstract call stack** remembers the last k call sites: $c^\# \in \mathcal{L}^k$
 the $C^\#$ and $R^\#$ maps now distinguish abstract call stacks
 $C^\#, R^\# : \mathcal{L}^k \rightarrow \mathcal{E}^\#$
 more precise than a partitioning by function only

larger k give more precision but less efficiency

Example: context-sensitivity

Example

main :

$R \leftarrow -1;$

$l_1 : f(\mathbf{rand}(5, 10));$

$l_2 : f(80)$

f(X) :

$R \leftarrow 2 \times X;$

if $R > 100$ **then** $R \leftarrow 0$

Analysis: using intervals and $k = 1$

- $C^\#(l_1) = [R \mapsto [-1, 1], X \mapsto [5, 10]]$
 $\implies R^\#(l_1) = [R \mapsto [10, 20], X \mapsto [5, 10]]$
- $C^\#(l_2) = [R \mapsto [10, 20], X \mapsto [80, 80]]$
 $\implies R^\#(l_2) = [R \mapsto [0, 0], X \mapsto [80, 80]]$
- at the end of the analysis, we get $R = 0$
 more precise than $R \in [0, 100]$ found without context-sensitivity

Tabulation abstraction

Cardinal power

Principle:

the semantic of a function is $S[\![\text{body}(f)]\!] : \mathcal{P}(\mathcal{E}) \rightarrow \mathcal{P}(\mathcal{E})$

\implies abstract it as an **abstract function in $\mathcal{E}^\# \rightarrow \mathcal{E}^\#$**

we use a partial function as the image of most abstract elements is not useful

Analysis: tabulated analysis

- use a global partial map $F^\# : \mathcal{F} \times \mathcal{E}^\# \rightarrow \mathcal{E}^\#$
- $F^\#$ is initially empty, and is filled **on-demand**
- when encountering $S^\#[\![\text{body}(f)]\!] X^\#$
 - return $F^\#(f, X^\#)$ if defined
 - else, compute $S^\#[\![\text{body}(f)]\!] X^\#$, store it in $F^\#(f, X^\#)$ and return it

Optimizations: trade precision for efficiency

- if $X^\# \sqsubseteq Y^\#$ and $F^\#(f, X^\#)$ is not defined, we can use $F^\#(f, Y^\#)$ instead
- if the size of $F^\#$ grows too large, use $F^\#(f, \top)$ instead
sound, and ensures that the analysis terminates in finite time

Example

Example

main :

```

R ← -1;
f(rand(5, 10));
f(80)

```

f(X) :

```

R ← 2 × X;
if R > 100 then R ← 0

```

Analysis using intervals

- $F^\# =$
 $[(f, [R \mapsto [-1, -1], X \mapsto [5, 10]]) \mapsto [R \mapsto [10, 20], X \mapsto [5, 10]],$
 $(f, [R \mapsto [10, 20], X \mapsto [80, 80]]) \mapsto [R \mapsto [0, 0], X \mapsto [80, 80]]]$
- at the end of the analysis, we get again $R = 0$

here, the function partitioning gives the same result as the call-site partitioning

Dynamic partitioning: complex example

Example: McCarthy's 91 function

main :

$Mc(\text{rand}(0, +\infty))$

$Mc(n)$:

if $n > 100$ **then** $r \leftarrow n - 10$
else $Mc(n + 11); Mc(r)$

- in the concrete, when terminating:
 $r = n - 10$ when $n > 101$, and $r = 91$ when $n \in [0, 101]$
- using a widening ∇ to choose tabulated abstract values $F^\sharp(f, X^\sharp)$
 we find:

$n \in [0, 72]$	\Rightarrow	$r = 91$
$n \in [73, 90]$	\Rightarrow	$r \in [91, 101]$
$n \in [91, 101]$	\Rightarrow	$r = 91$
$n \in [102, 111]$	\Rightarrow	$r \in [91, 101]$
$n \in [112, +\infty]$	\Rightarrow	$r \in [91, +\infty]$

(source: Bourdoncle, JFP 1992)