

Inter-Procedural Analysis and Points-to Analysis

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- Inter-Procedural analysis
- Call graph construction (fast)
- Call graph construction (precise)
- Call graph construction (fast and precise, not today – requires SSA)

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Inter-Procedural Analysis

- What is inter-procedural dataflow analysis
 - DFA that propagates dataflow values over procedure boundaries
 - Finds the impact of calls to caller and callee
- Tasks:
 - Determine a conservative approximation of the called procedures for all call sites
 - Referred to as Call Graph construction (more general: Points-to analysis)
 - Tricky in the presents of function pointers, polymorphism and procedure variables
 - Perform conservative dataflow analysis over basic-blocks of procedures involved
- Reason:
 - Allows new analysis questions (code inlining, removal of virtual calls)
 - For analysis questions with intra-procedural dataflow analyses, it is more precise (dead code, code parallelization)
- Precondition:
 - Complete program
 - No separate compilation
 - Hard for languages with dynamic code loading

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Call / Member Reference Graph

- A **Call Graph** is a rooted directed graph where the nodes represent methods and constructors, and the edges represent possible interactions (calls):
 - from a method/constructor (caller) to a method/constructor (callee).
 - root of the graph is the main method.
- Generalization: **Member Reference Graph** also including fields (nodes) and read and write accesses (edges).

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Proper Call Graphs

- A proper call graph is in addition
 - Conservative: Every call $A.m() \rightarrow B.n()$ that may occur in a run of the program is a part of the call graph
 - Connected: Every member that is a part of the graph is reachable from the main method
- Notice
 - We may have several entry points in cases where the program in question is not complete.
 - E.g., an implementation of the `ActionListener` interface will have the method `actionPerformed` as an additional entry point if we neglecting the `java.swing` classes.
 - Libraries miss a main method
 - In general, it is hard to compute, which classes/methods may belong to a program because of dynamic class loading.

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Techniques for Inter-Procedural Analysis

- Intra-procedural analysis on an inlined basic block graphs (textbook approach)
- Simulated execution

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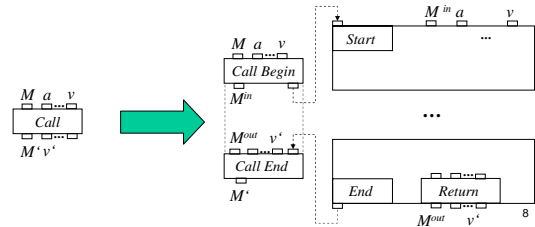
“Inlined” basic block graphs

- Given call graph and a bunch of procedures each with a basic block graph
- Inline basic block graphs
 - Split call nodes (and hence basic blocks) into callBegin and callEnd nodes
 - Connect callBegin with entry blocks of procedures called
 - Connect callEnd with exit blocks of procedures called
- Entry (exit) block of main method gets start node of forward (backwards) dataflow analysis
- Polymorphism is resolved by explicit dispatcher or by several targets
- Inter-procedural dataflow analysis now possible as before ab intra-procedural analysis

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“Inlining” of basic block graphs

- New node: begin and end of calls distinguished
- Edges: connection between caller and callees

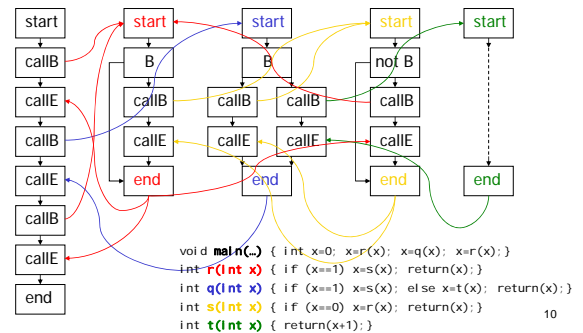


Example Program

```
public class One {
    public static void main(String[] args) {
        int x=0; x=r(x); x=q(x); x=r(x);
        System.out.println("Result: " + x);
    }
    static int r(int x) {
        if (x==1) x=s(x); return(x);
    }
    static int q(int x) {
        if (x==1) x=s(x); else x=t(x); return(x);
    }
    static int s(int x) {
        if (x==0) x=r(x); return(x);
    }
    static int t(int x) {
        return(x+1);
    }
}
```

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Example



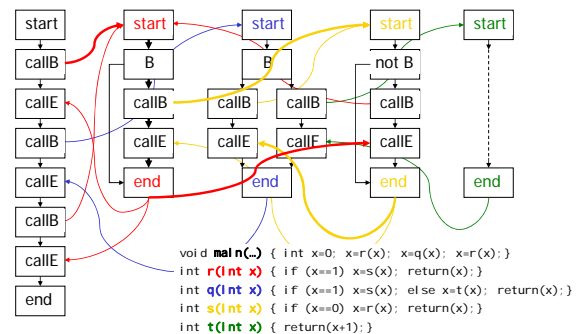
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Unrealizable Path

- Data gets propagated along path that never occur in any program run:
 - Calls to one **method** returning to another **method**
 - CallBegin** → Method Start → Method End → **CallEnd**
- Makes analysis conservative
- Still correct
- (And still more precise than corresponding intra-procedural analyses)

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Example: Unrealizable Path



Simulated Execution

- Starts with analyzing main
- Interleaving of [analyze method](#) and the [transfer function of calls](#)
- A method (intra-procedural analysis):
 - propagates data values analog the edges in basic-block graph
 - updates the analysis values in the nodes according to their transfer functions
 - If node type is a call then ...
- Calls' transfer function and only if the target method input changed:
 - Interrupts the processing of a caller method
 - Propagates arguments ($v_1 \dots v_n$) to the all callees
 - Processes the callees (one by one) completely
 - Iterate to local fixed point in case of recursive calls
 - Propagates back and merges (supremum) the results r of the callees
 - Continue processing the caller method ...

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Comparison

- Advantages of Simulated Execution
 - Fewer non realizable path, therefore:
 - More precise
 - Faster
- Disadvantages of Simulated Execution
 - Harder to implement
 - More complex handling of recursive calls
 - Leaves theory of monotone DFM and Abstract Interpretation

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Outline

- Inter-Procedural analysis
- [Call graph construction \(fast\)](#)
- Call graph construction (precise)
- Call graph construction (fast and precise, not today – requires SSA)

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Call Graph Construction in Reality

- The actual implementation of a call graph algorithm involves a lot of language specific considerations and exceptions to the basic rules. For example:
 - Field initialization and initialization blocks
 - Exceptions
 - Calls involving inner classes often need some special attention.
 - How to handle possible call back situations involving external classes.

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Why computing Call Graphs

- Elimination of dead code
 - classes never loaded, no objects created from, and
 - methods never called, and
 - branches never used in any program run
- Elimination of polymorphism (usage refers to a statically known class, not to its super-class nor siblings in the class hierarchy)
- Detection of [aliases](#) (usages refer to the same object) and [strangers](#) (usages are guaranteed not to refer to the same object)
- Detection of singletons (usage refers to a single object, not to a set of objects)
- ...

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Call Graphs: The Basic Problem

- The difficult task of any call graph construction algorithm is to approximate the set of members that can be targeted at different call sites.
- What is the target of call site $a.m()$?
- Depends on classes of objects potentially bound to designator expression a ?
- Not decidable, in general, because:
 - In general, we do not have exact control flow information.
 - In general, we can not resolve the polymorphic calls.
 - Dynamic class loading. This problem is in some sense more problematic since it is hard to make useful conservative approximations.

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Declared Target

- We say that the **declared target** of $a.m()$ occurring in a method $x.x()$ is the method $m()$ in the declared type of the variable a in the scope of $x.x()$.
- When using declared targets, **connectivity** can be achieved by ...
 - ... inserting (virtual) calls from super to subtype method declarations
 - ... keeping (potentially) dynamically loaded method nodes reachable from the main method (or as additional entry points).

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Discussion

- Class objects (static objects) are treated as objects
- Stack objects are considered part of *this*
 - Let a be a local variable or parameter, resp.
 - $a.m()$ is a usage of whatever a contains (target), i.e. $N(a)$, in whatever *this* contains (source), i.e. $N(this)$.

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Generalized Call Graphs

- A call graph is a directed graph $G=(V, E)$
 - vertices $V = \text{Class}.m$ are pairs of classes *Class* and methods / constructors / fields *m*
 - edges E represent usage: let a and b be two objects: a uses b (in a method / constructor execution x of a occurs a call / access to a method / constructor / field y of b) $\Leftrightarrow (\text{Class}(a).x, \text{Class}(b).y) \in E$
- An **generalized** call graph is a directed graph $G=(V, E)$
 - vertices $V = N(o).m$ are pairs of finite abstractions of runtime objects o using a so called name schema $N(o)$ and methods / constructors / fields m
 - edges E represent usage: let a and b be two objects: a uses b (in a method / constructor execution x of a occurs a call / access to a method / constructor / field y of b) $\Leftrightarrow (N(a).x, N(b).y) \in E$
- A name schema N is an abstraction function with finite co-domain
- $\text{Class}(o)$ is a special name schema and, hence, describes a special call graph

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Name Schemata

- One can abstract from objects by distinguishing:
 - Just heap and stack (decidable, not relevant)
 - Objects with same class (not decidable, relevant, efficient approximations)
 - Objects with same class but syntactic different creation program point (not decidable, relevant, expensive approximations)
 - Objects with same creation program point but with syntactic different path to that creation program point (not decidable, relevant, approximations exponential in execution context)
 - Different objects (not decidable)
 - ...

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Decidability

- **Not** decidable in general: reduction from termination problem
 - Add a new call (not used anywhere else before the program exit)
 - If I could compute the exact call graph, I know if the program terminates or not
- Decidable if name schema abstract enough (then not relevant in practice)

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Approximations

- Simple conservative approximation
 - from static semantic analysis
 - declared class references in a class A and their subtypes are **potentially** uses in A
 - $a.x$ really uses $b.y \Rightarrow (N(a).x, N(b).y) \in E$
- Simple optimistic approximation
 - from profiling
 - actually used class references in an execution of class A (a number of executions) are **guaranteed** uses in A
 - $a.x$ really uses $b.y \Leftarrow (N(a).x, N(b).y) \in E$

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Simplification

- For a first try, we consider only one name schema:
 - Distinguish objects of different classes / types
 - Formally, $N(o) = \text{Class}(o)$
- Consequently, a call graph is ...
 - a directed graph $G = (V, E)$
 - vertices V are pairs of classes and methods / constructors / fields
 - edges E represent usage: let A and B be two classes: $A.x$ uses $B.y$ (i.e. an instance of A executes x using an method / constructor / field y instance of B)

$$\Leftrightarrow (A.x, B.y) \in E$$
- Not decidable still, we discuss optimistic and conservative approximations

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Algorithms to discuss

All algorithms these are **conservative**:

- Reachability Analysis – RA
- Class Hierarchy Analysis – CHA
- Rapid Type Analysis – RTA
- ...
- (context-insensitive) Control Flow Analysis – 0-CFA
- (k -context-sensitive) Control Flow Analysis – k -CFA

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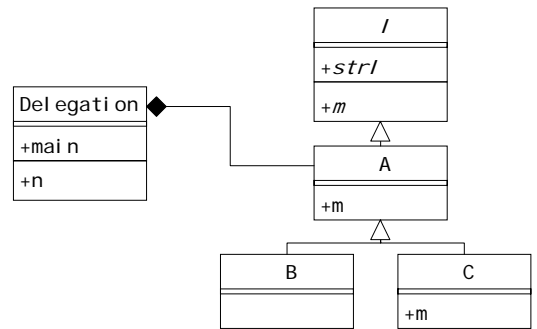
Reachability Analysis – RA

- Worklist algorithm maintaining reachable methods
 - initially *main* routine in the *Main* class is reachable
- For this and the following algorithms, we understand that
 - Member (field, method, constructor) names n stand for complete signatures
 - R denotes the worklist and finally reachable members
 - R may contain fields and methods/constructors. However, only methods/constructors may contain other field accesses/call sites for further processing.
- RA:
 - $\text{Main.main} \in R$ (maybe some other entry points too)
 - $M.m \in R$ and $e.n$ is a field access / call site in m

$$\Rightarrow \forall N \in \text{Program}: N.n \in R \wedge (M.m, N.n) \in E$$

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Example



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Example

```
public class Delegation {
    public static void main(String args[]) {
        A i = new B();
        i.m();
        Delegation.n();
    }
    public static void n() {
        new C().m();
    }
}

interface I {
    public String str! = "Printing I string";
    public void m();
}

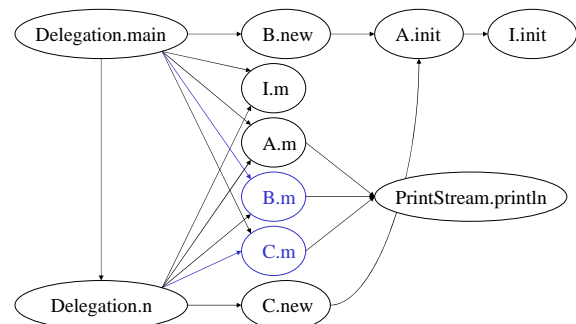
class A implements I {
    public void m() {System.out.println(str!);}
}

class B extends A {
    public B() {super();}
    public void m();
}

class C extends A {
    public void m() {System.out.println("Printing C string");}
}
```

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RA on Example



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Class Hierarchy Analysis – CHA

- Refinement of RA
- $Main.main \in R$
- $M.m \in R$
 - $e.n$ is a field access / call site in $M.m$
 - $type(e)$ is the static (declared) type of access path expression e
 - $subtype(type(e))$ is the set of (declared) sub-types of $type(e)$
- $\Rightarrow \forall N \in subtype(type(e)): N.n \in R \wedge (M.m, N.n) \in E$

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Example

```
public class Delegation {
    public static void main(String args[]) {
        A i = new B();
        i.m();
        Delegation.n();
    }
    public static void n() {
        new C().m();
    }
}

interface I {
    public String str1 = "Printing I string";
    public void m();
}

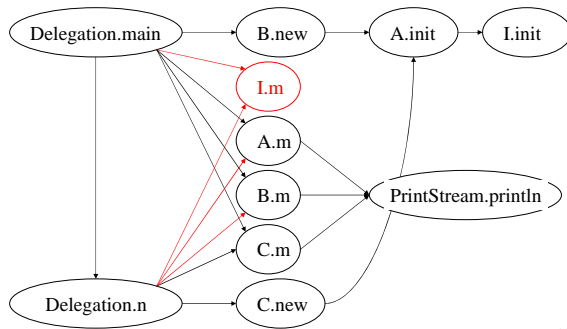
class A implements I {
    public void m() {System.out.println(str1);}
}

class B extends A {
    public B() {super();}
}

class C extends A {
    public void m() {System.out.println("Printing C string");}
}
```

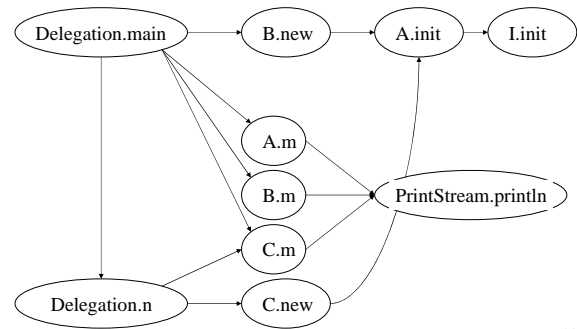
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CHA on Example



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CHA on Example



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Rapid Type Analysis – RTA

- Still simple and fast refinement of CHA
- Maintains reachable methods R and instantiated classes S
- Fixed point iteration: whenever S changes, we revisit the worklist R
- $Main.main \in R$
- For all class (static) methods s : $class(s) \in S$
- $M.m \in R$
 - $new N$ is a constructor call site in $M.m$
 - $\Rightarrow N \in S \wedge N.new \in R \wedge (M.m, N.new) \in E$
 - $e.n$ is a field access / call site in $M.m$
 - $\Rightarrow \forall N \in subtype(type(e)) \wedge N \in S: N.n \in R \wedge (M.m, N.n) \in E$

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Example

```
public class Delegation {
    public static void main(String args[]) {
        A i = new B();
        i.m();
        Delegation.n();
    }
    public static void n() {
        new C().m();
    }
}

interface I {
    public String str1 = "Printing I string";
    public void m();
}

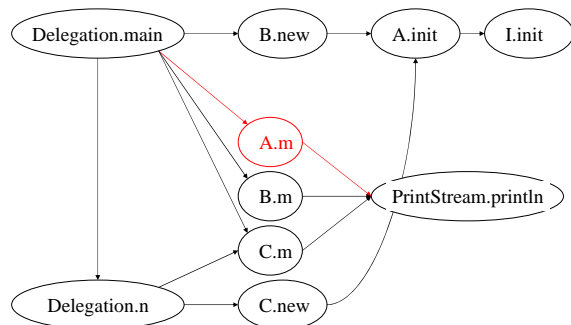
class A implements I {
    public void m() {System.out.println(str1);}
}

class B extends A {
    public B() {super();}
}

class C extends A {
    public void m() {System.out.println("Printing C string");}
}
```

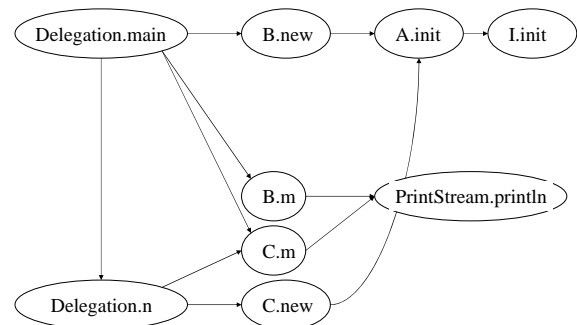
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RTA on Example



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RTA on Example



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Context-Insensitive Control Flow Analysis – 0-CFA

- RTA assumes that **any** constructed class object of a type can be bound to an access path expression of the same type
- Considering the control flow of the program, the set of reaching objects further reduces
- Example:

```
main() {
    A a = new A();
    a.n();
    sub();
}

sub() {
    A a = new B();
    a.n();
}

class A {
    public void n() {...}
}

class B extends A {
    public void n() {...}
}
```

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Context-Sensitive Control Flow Analysis – k -CFA

- 0-CFA merges objects that can reach an access path expression (designator) via different call paths
- One can do better when distinguishing the objects that can reach an access path expression via paths differing in the last k nodes of the call paths

```
main() {
    A a = new A();
    X.dispatch(a);
    sub();
}

sub() {
    A a = new B();
    X.dispatch(a);
}

class A {
    public void n() {...}
}

class B extends A {
    public void n() {...}
}

class X {
    public static void dispatch(A a) { a.n(); }
}
```

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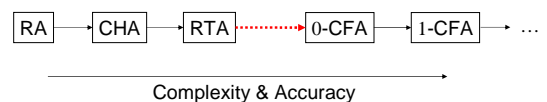
Control Flow Analysis

- Requires data flow analysis
- 0-CFA: has already high memory consumption in practice (still practical)
- k -CFA: is exponential in k
 - Requires a refined name schema (and, hence, even more memory)
 - Does not scale in practice (if extensively used)
 - Solutions discussed later today
 - One idea (current research):
 - Make k adaptive over the analysis
 - Focus on specific program parts
 - Reduce k to max 1

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Order on Algorithms

- Increasing complexity
- Increasing accuracy



- Analyses between RTA and 0-CFA?

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Analyses Between RTA and 0-CFA

- RTA uses **one** set S of instantiated classes
- Idea:
 - Distinguish **different** sets of instantiated classes reaching a specific field or method
 - Attach them to these fields, methods
 - Gives a more precise "local" view on object types possibly bound to the fields or methods
 - Regards the control flow between methods but
 - Disregards the control flow within methods
- Fixed point iteration

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Notations

- Subtypes of a **set** of types:
 $subtype(S) ::= \bigcup_{N \in S} subtype(N)$
- Set of parameter types $param(m)$ of a method m : all static (declared) argument types of m excluding $type(this)$
- Return type $return(m)$ of a method m : the static (declared) return type of m

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Separated Type Analysis – XTA

- Separate type sets S_m reaching methods m and fields x (treat fields x like methods pairs set_x, get_x)
- $Main.main \in R$
- $M.m \in R$
 - For all class (static) methods s : $class(s) \in S_{M,m}$
 - $new N$ is a **constructor** call site in $M.m$
 - $\Rightarrow N \in S_{M,m} \wedge N.new \in R \wedge (M.m, N.new) \in E$
 - $e.n$ is a field access / call site in $M.m$
 - $\Rightarrow \forall N \in subtype(type(e)) \wedge N \in S_{M,m} : N.n \in R \wedge$
 $subtype(param(N.n)) \cap S_{M,m} \subseteq S_{N,n} \quad \wedge$
 $subtype(result(N.n)) \cap S_{N,n} \subseteq S_{M,m} \quad \wedge$
 $(M.m, N.n) \in E$

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Example

```
public class Delegation {
    public static void main(String args[]) {
        A a = new B();
        a.m();
        Delegation.n();
    }
    public static void n() {
        new C().m();
    }
}

interface I {
    public String str1 = "Printing I string";
    public void m();
}

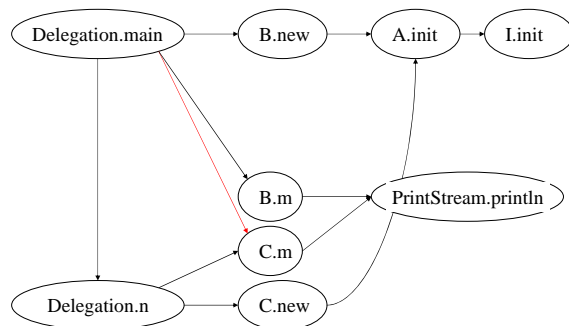
class A implements I {
    public void m() {System.out.println(str1);}
}

class B extends A {
    public B() {super();}
}

class C extends A {
    public void m() {System.out.println("Printing C string");}
}
```

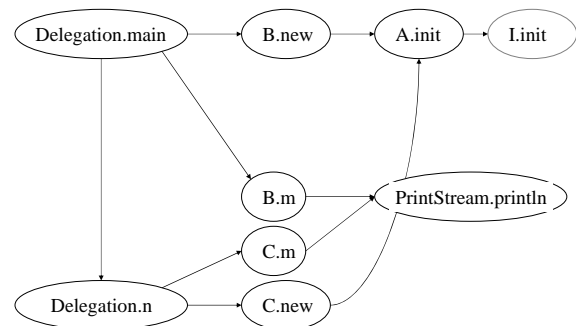
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XTA on Example



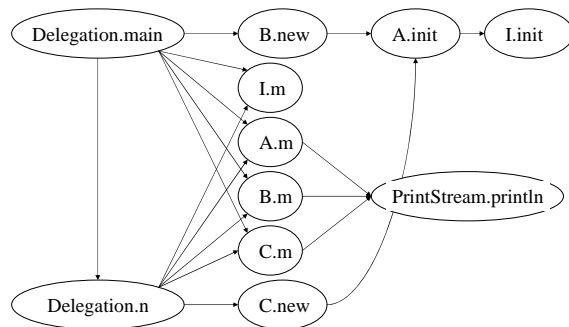
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XTA on Example



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RA vs XTA on Example



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Increasing complexity



- Number of type separating sets S (M number of methods, F number of fields):
 - CHA: 0
 - RTA: 1
 - XTA: $M + F$
- Practical observations on benchmarks:
 - All algorithms RA...XTA scale (1 Mio. Loc)
 - XTA one order of magnitude slower than RTA
 - Correlation to program size rather weak

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Increasing precision



- Practical observations on benchmarks:
 - RTA as baseline: all instantiated (wherever) classes are available in all methods
 - XTA on average:
 - only ca. 10% of all classes are available in methods ☹
 - < 3% fewer reachable methods ☺
 - > 10% fewer call edges
 - > 10% more monomorphic call targets

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Source

- Frank Tip and Jens Palsberg:
Scalable Propagation-Based Call Graph Construction Algorithms.
ACM Conf. on Object-Oriented Programming Systems, Languages and Application – OOPSLA 2000.
- David Grove, Greg DeFouw, Jeffrey Dean, and Craig Chambers:
Call Graph Construction in Object-Oriented Languages.
OOPSLA 1997.

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Conclusion on Call Graphs so far

- Approximations
 - Relatively fast, feasible for large systems
 - Relatively imprecise, conservative
- What is a good enough approximation of certain client analyses
- Answer depends on client analyses (e.g., different answers for software metrics and clustering vs. program optimizations)

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Outline

- Inter-Procedural analysis
- Call graph construction (fast)
- Call graph construction (precise)
- Call graph construction (fast and precise, not today – requires SSA)

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Classic P2A: Introduction

- We try to find all objects that each reference variable may point to (hold a reference to) during an execution of the program.
- Hence, to each reference variable v in a program we associate a set of objects, denoted $Pt(v)$, that contains all the objects that variable v may point to. The set $Pt(v)$ is called the points-to set of variable v .
- Example:


```

A a,b,c;
X x,y;
s1: a = new A( ); // Pt( a ) = {o1}
s2: b = new A( ); // Pt( b ) = {o2}
    b = a; // Pt( b ) = {o1, o2}
    c = b; // Pt( c ) = {o1, o2}
      
```
- Here oi means the object created at allocation site si .
- After a completed analysis, each variable v is associated with a points-to set $Pt(v)$ containing a set of objects that it may refer to

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Name Schema revisited

- The number of objects appearing in a program is in general infinite (countable), hence, we don't have a well-defined set of data values.
- For example, consider the following situation


```

while ( x > y ) {
    A a = new A( );
    ...
}
      
```

 The number of A objects is in cases like this impossible to decide. (Think if x or y depends on some input values).
- From now on, each object **creation point** (`new A()`, `a.clone()`, "hello") represents a unique object (identified by the source code location).
- Again, many run-time objects are mapped to a single abstract object.
- Finitely many abstract objects

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Object Transport as Set Constraints

- Objects can flow between variables due to assignments and calls. Calls will be treated shortly.
- Certain statements generates constraints between points-to sets. We will consider:


```

l = r           ⇒ Pt(r) ⊆ Pt(l)      (Assignment)
site i: l = new A() ⇒ {oi} ⊆ Pt(l)    (Allocation)
      
```
- That is, each assignment can be interpreted as a constraint between the involved points-to sets.
- Each statement in the program will generate constraints, as before equations in DFA, we will have a system of constraints.
- We are looking for the *minimum solution* (minimum size of the points-to sets) that satisfies the resulting system of constraints, i.e., the minimum fixed point of the dataflow equations

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Example

A Simple Program	Generated set constraints
<pre> public A methodX(A param){ A a1 = param; s1 : A a2 = new A(); A a3 = a1; a3 = a2 ; return a3 ; } </pre>	<pre> 1: Pt(param) ⊆ Pt(a1) 2: o1 ∈ Pt(a2) 3: Pt(a1) ⊆ Pt(a3) 4: Pt(a2) ⊆ Pt(a3) </pre>

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Object Transport in terms of P2G edges

- Each constraint can be represented as a relation between nodes in a graph.
- A *Points-to Graph* P2G is a directed graph having variables and objects as nodes and assignments and allocations as edges


```

l = r ⇒ Pt(r) ⊆ Pt(l) ⇒ r → l (Assignment)
site i: l = new A() ⇒ {oi} ⊆ Pt(l) ⇒ oi → l (Allocation)
      
```
- Previous example revisited


```

1: Pt(param) ⊆ Pt(a1)
2: o1 ∈ Pt(a2)
3: Pt(a1) ⊆ Pt(a3)
4: Pt(a2) ⊆ Pt(a3)
      
```
- P2G is our data-flow graph and the objects are our data values to be propagated.
- P2G initialization: $\forall oi \rightarrow l$, add oi to $Pt(l)$.
- P2G propagation: $\forall r \rightarrow l$, add let $Pt(l) = Pt(l) \cup Pt(r)$

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Analysis Approximation: Flow-Insensitivity

- An analysis is *flow-sensitive* if the order of execution of statements is taken in account.
- In the code example below illustrates flow-sensitivity.


```

(1) s1 : f = new A()
(2)     a = f
(3) s2 : f = new A()
          //insensitive: Pt(a)={o1,o2}
          //sensitive: Pt(a)={o1}
(4)     b = f
          //insensitive: Pt(b)={o1,o2}
          //sensitive: Pt(b)={o2}
      
```
- Our approach would have generated the following *set* of constraints


```

o1 ∈ Pt(f), Pt(f) ⊆ Pt(a), o2 ∈ Pt(f), Pt(f) ⊆ Pt(b)
      
```
- Constraints (1) and (3) yield $Pt(f) = \{o1, o2\}$ and consequently that both a and b have $Pt = \{o1, o2\}$.
- Thus, a consequence of using our set constraint approach is *flow-insensitivity*.
- A flow-sensitive analysis requires that each *definition* of a variable has a node and a points-to set. This makes the graph much larger and the analysis more costly.
- Forward reference Static Single Assignment (SSA) representation

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Method Calls: Object Transfer

```
s1 : A a = new A() // o1 -> a
s2 : X x1 = new X() // o2 -> x1
      a.storeX( x1 ) // x1 -> x4 , x4 -> x3 , x3 -> f
      X x2 = a.loadX() // f -> x2

classA {
  X f ;
  private void setX (X x3) {f = x3;}
  private X getX() {return f;}
  public void storeX (X x4) { this.setX(x4);}
  public X loadX() {return this.getX();}
}
```

- Involved object transport
 - Argument passing, i.e., assigning arguments to parameters (e.g. $x1 \rightarrow x4$).
 - A call $a.m()$ involves an implicit assignment $a \rightarrow this_m$.
 - The return assignment involves an implicit step too $f \rightarrow x2$.
- The above approach contains many implicit steps that are hard to perform automatically.
- The procedural method representation gives a more explicit view of the object transport.

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Representation of Methods

OO Definition	Procedural Definition
<pre>classA { public R mName(P1 p1,P2 p2){ ... return Rexpr; } }</pre>	<pre>mName(A this, P1 p1, P2 p2, R ret) { ... ret = Rexpr ; }</pre>
OO Invocation $l = a.mName(x,y);$	Procedural Invocation $mName(a,x,y,l);$

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Advantage of Procedural Representation

- Given a call site $l=r0.m(r1,...,rn)$
 - Represented as $m(r0,r1,...,rn,l)$
 - Targeted at method `public R mName(P1 p1,P2 p2) in classA`
 - Represented as $m(A\ this, P1\ p1,..., Pn\ pn, R\ ret)$
- We add the following P2G edges
- $r0 \rightarrow this, r1 \rightarrow p1, \dots, rn \rightarrow pn, ret \rightarrow l$
- Thus, each resolved call site results in a well-defined set of inter-procedural P2G edges.

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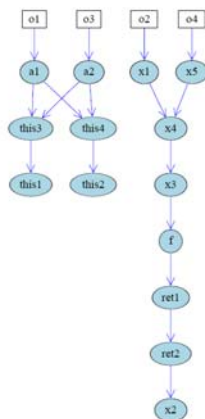
Previous Example Revisited / Extended

```
class Main {
  static procedure main (Main this , String[] args) {
    s1 : A a1 = new A( ) ; // o1 --> a1
    s2 : X x1 = new X( ) ; // o2 --> x1
          storeX ( a1 , x1 ) ; // a1 --> this3 , x1 --> x4
          X x2 ;
          loadX ( a1 , x2 ) ; // a1 --> this4 , ret2 --> x2
    s3 : A a2 = new A( ) ; // o3 --> a2
    s4 : X x5 = new X( ) ; // o4 --> x5
          storeX ( a2 , x5 ) ; // a2 --> this3 , x5 --> x4
          loadX ( a2 , x2 ) ; // a2 --> this4 , ret2 --> x2
  }
}

classA {
  X f ;
  procedure setX(A this1, X x3 ) { f = x3 } // x3 --> f
  procedure getX(A this2, Xret1 ) { ret1 = f } // f --> ret1
  procedure storeX(A this3, X x4 ) { setX(this3,x4) }
    // this3 --> this1, x4 --> x3
  procedure loadX(A this4 , Xret2 ) {getX( this4, ret2 ) }
    // this4 --> this2 , ret1 --> ret2
}
```

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P2G Generated



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Resolving Call Targets

- The procedural method representation makes it quite easy to generate a set of Call Graph edges once the target method has been identified. The problem is to find target methods.
- Static calls and constructor calls are easy, they always have a well-defined target method.
- Virtual calls are much harder; to accurately decide the target of a call site during program analysis is in general impossible.
- Any points-to analysis involves some kind of conservative approximation where we take into account all possible targets.
- The trick is to narrow down the number of possible call targets.

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Resolving Polymorphic Calls

Two approaches to resolve a call site $a.m()$

- Static Dispatch: Given an *externally derived* conservative call graph (discussed before) we can approximate the actual targets of any call site in a program. By using such a call graph we can associate each call site $a.m()$ with a set of pre-computed target methods $T_1.m(), \dots, T_n.m()$.
- Dynamic Dispatch: By using the currently available points-to set $Pt(a)$ itself, we can, for each object in the set, find the corresponding dynamic class and, hence, the target method definition of any call site $a.m()$.

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Static Dispatch

- Given a conservative call graph we can construct a function $staticDispatch(a.m())$ that provides us with a set of possible target methods for any given call site $a.m()$.
- We can then proceed as follows:


```
for each call site  $l = r0.m(r1, \dots, rn)$  do
    let targets = staticDispatch( $r0.m(\dots)$ )
    for each method  $m(A \ this, p1 \ p1, \dots, pn \ pn, R \ ret) \in targets$  do
        add P2G edges  $r0 \rightarrow this, r1 \rightarrow p1, \dots, rn \rightarrow pn, ret \rightarrow l$ 
```
- Advantage: We can immediately resolve all call sites and add corresponding P2G edges.
- Disadvantage: The precision of the externally derived call graph influences the points-to-analysis.
- We refer to P2Gs where no more edges are to be added as *complete*. Complete P2Gs are much easier to handle as will be discussed shortly.

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Dynamic Dispatch

- Given the points-to set $Pt(a)$ of a variable a we can resolve the targets of a call site $a.m()$ using a function $dynamicDispatch(A, m)$ that returns the method executed when we invoke the call $m()$ with signature m on an object O_a of type A .
- We can then proceed as follows:


```
for each call site  $l = r0.m(r1, \dots, rn)$  (or  $m(r0, r1, \dots, rn, l)$ ) do
    for each object  $O_a \in Pt(r0)$  do
        1. Let  $m = signatureOf(m())$ 
        2. Let  $A = typeOf(O_a)$ 
        3. Let  $m(A \ this, p1 \ p1, \dots, pn \ pn, R \ ret) = dynamicDispatch(A, m)$ 
        4. Add P2G edges  $r0 \rightarrow this, r1 \rightarrow p1, \dots, rn \rightarrow pn, ret \rightarrow l$ 
```
- Advantage: We avoid using an externally defined call graph.
- Disadvantage: The P2G is not complete since we initially don't know all members of $Pt(a)$.
- Hence, the P2G will change (additional edges will be added) during analysis.

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Propagating a Complete P2G

- In this approach we use work list to store variable nodes that need to be propagated.


```
1. For each variable  $v$  let  $Pt(v) = \emptyset$  //  $O(\#v)$ 
2. For each allocation edge  $oi \rightarrow v$  do //  $O(\#o)$ 
    (a) let  $Pt(v) = Pt(v) \cup \{oi\}$ 
    (b) add  $v$  to worklist
3. Repeat until worklist empty //  $O(\#v^2 \#o)$ 
    (a) Remove first node  $p$  from worklist
    (b) For each edge  $p \rightarrow q$  do //  $O(\#v)$ 
        i. Let  $Pt(q) = Pt(q) \cup Pt(p)$ 
        ii. If  $Pt(q)$  has changed, add  $q$  to worklist
```
- Time complexity: Let $\#v$ be the number of variable nodes and $\#o$ the number of (abstract) objects.
- A node is added to the work list each time it is changed.
- In the worst case this can happen $\#o$ times for each node, thus, we have $O(\#v^2 \#o)$ number of work list iterations.
- Each such iterations may update every other variable node (hence $O(\#v)$ within the loop). Thus, an upper limit is $O(\#v^2 \#o)$.

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Optimizing the Analysis

- The high time complexity $O(\#v^2 \#o)$ encourages optimizations. Optimizations can basically be done in two different ways:
- We can reduce the size of P2G by identifying points-to sets that must be equal. This idea will be exploited in
 - Removal of strongly connected components
 - Removal of single dominated subgraphs.
- We can speed up the propagation algorithm by processing the nodes in a more clever ordering:
 - Topological node ordering.
- Other optimizations are possible all three are simple and effective.

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Previous Example Revisited: Results of Points-to Analysis

```
class Main {
    static procedure main(Main this, String[] args) {
        s1: A a1 = new A(); // Pt(a1) = {o1}
        s2: X x1 = new X(); // Pt(x1) = {o2}
        storeX(a1, x1);
        X x2; // Pt(x1) = {o2, o4}
        loadX(a1, x2);
        s3: A a2 = new A(); // Pt(a2) = {o3}
        s4: X x5 = new X(); // Pt(x5) = {o4}
        storeX(a2, x5);
        loadX(a2, x2);
    }
}

class A {
    X f; // Pt(f) = {o2, o4}
    procedure setX(A thi s1, X x3) { f = x3; } // Pt(thi s1) = {o1, o3}, Pt(x3) = {o2, o4}
    procedure getX(A thi s2, X r1) { r1 = f; } // Pt(thi s2) = {o1, o3}, Pt(r1) = {o2, o4}
    procedure storeX(A thi s3, X x4) { setX(thi s3, x4); }
    // Pt(thi s3) = {o1, o3}, Pt(x4) = {o2, o4}
    procedure loadX(A thi s4, X r2) { getX(thi s4, r2); }
    // Pt(thi s4) = {o1, o3}, Pt(r2) = {o2, o4}
}
```

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Limitations of Classic Points-to Analysis

- In the previous example we found that $Pt(A, \mathcal{F}) = \{o2, o4\}$. However, from the program code it is obvious that we have two instances of class *A* (*o1* and *o2*) and that $Pt(o1, \mathcal{F}) = \{o2\}$ whereas $Pt(o3, \mathcal{F}) = \{o4\}$. Hence by having a common points-to set for field variables in different objects the different object states are merged.
- Consider two *List* objects created at different locations in the program. We use the first list to store *String* objects and the other to store *Integer*. Using ordinary points to analysis we would find that both these list store both strings and objects.
- Conclusion: Classic points-to analysis merges the states in objects created at different locations and, as a result, can't distinguish their individual states and content.
- Context-sensitive approaches would let each object has its own set of fields. This would however correspond to object/method inlining and increase the number of P2G nodes exponentially and reduce the analysis speed accordingly.
- Flow-sensitivity would increase precision as well, at the price of adding new nodes for every definition of a variable. Once again, increased precision at the price of performance loss.
- The trade-off between precision and performance is a part of everyday life in data-flow analysis. In theory we know how to increase the precision, unfortunately not without a significant performance loss.

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Outline

- Inter-Procedural analysis
- Call graph construction (fast)
- Call graph construction (precise)
- Call graph construction (fast and precise)

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