

# Nominal Anti-Unification with Atom-Variables

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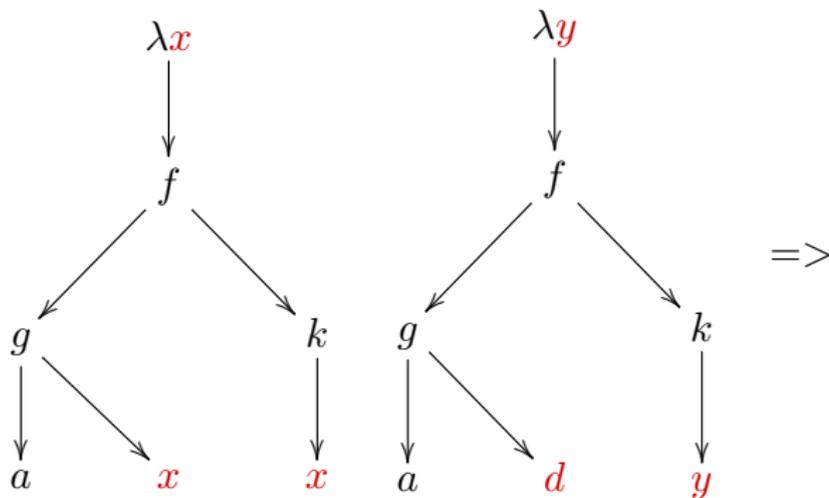
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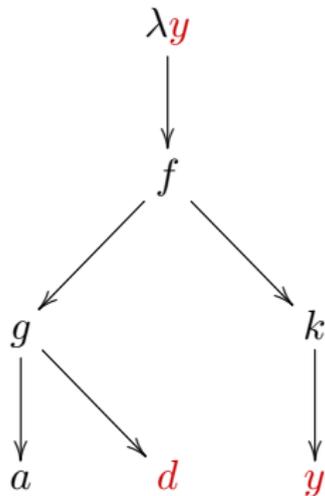
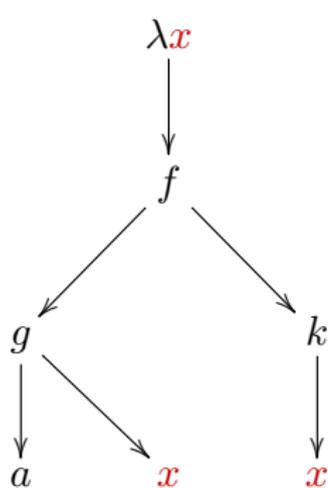
# Examples for Term Generalization of lambda-expressions

Generalize the two terms:



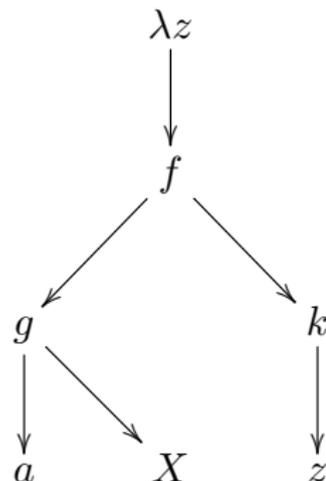
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Generalize the two terms:



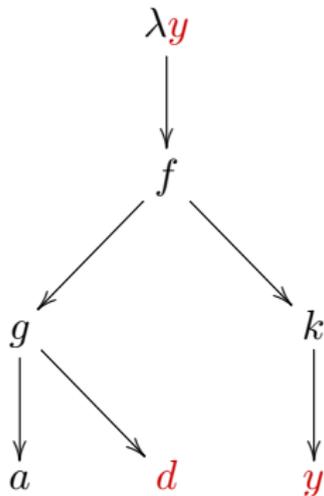
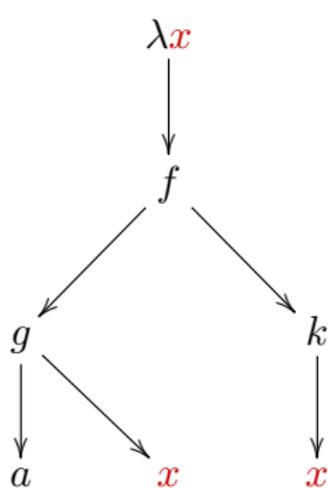
$\Rightarrow$

Generalization:



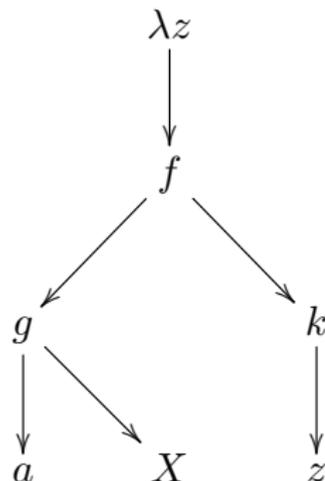
# Examples for Term Generalization of lambda-expressions

Generalize the two terms:



$\Rightarrow$

Generalization:



**Task of Anti-Unification:**

Given a set of terms.

Compute a representation

of all least general generalizations (lgg-s)

# Nominal Term Generalization

Our contribution:

- Anti-Unification,  
in an extended lambda-calculus  $NL_a$  modulo  $\sim_\alpha$   
with grammar:

$$S ::= a \mid f(S_1, \dots, S_n) \mid \lambda a.S$$

- Our general language for the input and algorithm is  $NL_A$ :

$$\begin{aligned} s & ::= W \mid \pi \cdot X \mid f(s_1, \dots, s_n) \mid \lambda W.s \\ W & ::= \pi \cdot A \\ \pi & ::= \emptyset \mid (W_1 W_2) \circ \pi \end{aligned}$$

$A$ : atom-variables standing for atoms

$\pi$ : permutations (of atoms)

$W$ : suspensions

## Some Related Work

Plotkin, A note on inductive generalization, MI, 1971

Urban, Pitts, Gabbay Nominal unification, CSL 2003

Baumgartner, Kutsia, Levy, Villaret. Nominal anti-unification, RTA 2015

S, Sabel, Kutz: Nominal unification with Atom-Variables, JSC 2019

Baumgartner et.al.: Nominal anti-unification algorithm.

- Baumgartner et. al.'s algorithm: results in a **single generalization**. **Striking Problem in Baumgartner et.al.:** There are no *least* general generalizations!
- Since there are infinitely properly increasing chains of more and more general generalizations.
- The problem:  
It is always possible to refine constraints by adding components of the form  $a\#X$ , where  $a$  is an atom that does not occur in the problem.

### **Our approach:**

- Exploit atom-variables to avoid infinite increasing chains;
- And provide a more general formulation

A **term-in context** is a pair  $(t, C)$ , where

$t$  is a  $NL_A$ - term:

$C$  is a  $NL_A$  freshness context, i.e. a set of constraints.  
(of the form:  $A\#s$ .)

ATOMANTIUNIFICATION: State:  $(L, M, \Delta, \theta)$

$L$  Term-pairs to generalize with a  
default generalization-variable

$M$  unsolved term-pairs for later merge

$\Delta$  The freshness context (i.e. constraints)

$\theta$  Substitution; i.e. refinement of the generalization vars.

# ATOMANTIUNIFICATION: Rules of the algorithm

(Dec): Decomposition

$$\frac{\{X:f(s_1,\dots,s_n) \triangleq f(t_1,\dots,t_n)\} \cup \Gamma, M, \nabla, \theta}{\Gamma \cup \{X_1:s_1 \triangleq t_1, \dots, X_n:s_n \triangleq t_n\}, M, \nabla, \theta \cup \{X \mapsto f(X_1, \dots, X_n)\}}$$

where  $X_i$  are fresh variables

(Abs): Abstraction

$$\frac{\{X:\lambda W_1.s \triangleq \lambda W_2.t\} \cup \Gamma, M, \nabla, \theta}{\Gamma \cup \{Y:(W_1 B) \cdot s \triangleq (W_2 B) \cdot t\}, M, \nabla \cup \{B \# \lambda W_1.s, B \# \lambda W_2.t\}, \theta \cup \{X \mapsto \lambda B.Y\}}$$

where  $Y$  is a fresh variable, and  $B$  is a fresh atom-variable

(SusA): SuspensionA

$$\frac{\{X:W_1 \triangleq W_2\} \cup \Gamma, M, \nabla, \theta \quad \nabla \vDash W_1 = W_2}{\Gamma, M, \nabla, \theta \cup \{X \mapsto W_1\}}$$

(SusYY): SuspensionYY

$$\frac{\{X:\pi_1 \cdot Y \triangleq \pi_2 \cdot Y\} \cup \Gamma, M, \nabla, \theta \quad \nabla \vDash \pi_1 = \pi_2}{\Gamma, M, \nabla, \theta \cup \{X \mapsto \pi_1 \cdot Y\}}$$

(Mer): Merging

$$\frac{\Gamma, \{Z_1:s_1 \triangleq t_1, Z_2:s_2 \triangleq t_2\} \cup M, \nabla, \theta \quad \text{EQVM}(\{(s_1, t_1) \preceq (s_2, t_2)\}, \nabla) = \pi}{\Gamma, M \cup \{Z_1:s_1 \triangleq t_1\}, \nabla, \theta \cup \{Z_2 \mapsto \pi \cdot Z_1\}}$$

where  $(Z_1, Z_2)$  is  $(X, Y)$  or  $(A, B)$

(Solve)

$$\frac{\{X:s \triangleq t\} \cup \Gamma, M, \nabla, \theta}{\Gamma, M \cup \{X:s \triangleq t\}, \nabla, \theta} \quad \text{If } \text{Head}(s) \neq \text{Head}(t) \text{ and if } s \text{ and } t \text{ are not both suspensions of atom-variables.}$$

(SolveYY)

$$\frac{\{X:\pi_1 \cdot Y \triangleq \pi_2 \cdot Y\} \cup \Gamma, M, \nabla, \theta \quad \nabla \not\equiv \pi_1 = \pi_2}{\Gamma, M \cup \{X:\pi_1 \cdot Y \triangleq \pi_2 \cdot Y\}, \nabla, \theta}$$

(SolveAB)

$$\frac{\{X:W_1 \triangleq W_2\} \cup \Gamma, M, \nabla, \theta \quad \nabla \not\equiv W_1 = W_2}{\Gamma, M \cup \{A:W_1 \triangleq W_2\}, \nabla, \theta \cup \{X \mapsto A\}}$$

$A$  is a fresh atom-variable.

## Properties of ATOMANTIUNIFICATION

- It computes a single generalization
- The algorithm requires simple exponential time to compute a solution.  
The number of rule applications is polynomial,  
and the solution requires polynomial space.

## ATOMANTIUNIFICATION problems

- requires exponential time:  
checking equivalence of permutations,  
and merging
- It is only weakly complete:  
It is complete if one ignores the freshness constraints.

# An example as a hurdle to completeness of ATOMANTIUNIFICATION

- Input:  $(\emptyset, \{X : (f(A), A, B) \triangleq (c, A, B)\})$
- ATOMANTIUNIFICATION computes the generalization  $(\emptyset, (Y, A, B))$ , however, it is not an lgg.
- Adding  $A\#Y$  or  $B\#Y$  violates the generalization property:  $(f(a), a, a)$  not covered.
- $(\{B\#\lambda A.Y\}, (Y, A, B))$  is the lgg:  
 $(\{B\#\lambda A.Y\}, (Y, A, B))$  allows the instance  $(f(a), a, a)$ , since  $a\#\lambda a.f(a)$  holds.

## Pros and Cons of using and refining freshness constraints

- (+) For a fixed input problem: The tree of properly refining freshness constraint  $C$  is finite.
- (-) However, we did not succeed in constructing a (terminating) algorithm that can construct maximal properly decreasing chains of constraints.

We propose an extension of the expressive power of freshness constraints:

## EQR-Constraints

Permit more general freshness constraints based on **equivalence** relations of atom-variables induced by semantic instantiations.

### Example for EQR-freshness constraints:

The constraint  $A \# \lambda B.A$

has an equivalent EQR-freshness context:

$$(\{A = B\} \implies \text{True})$$

$$\wedge$$

$$(\{A \neq B\} \implies \text{False})$$

- EQR-freshness contexts are strictly more general than freshness contexts (in  $NL_A$ )
- EQR-freshness contexts may be large. (exponentially larger than (usual) freshness contexts.)
- There is an algorithm for refining EQR-freshness contexts using the same set of atom-variables
- The refinement steps are sufficiently fine-grained and terminate.

# The Anti-Unification-Algorithm for EQR-Constraints

ATOMANTIUNIFLGG:

- ① Run ATOMANTIUNIFICATION.
- ② Then: Iteratively refine the EQR-freshness constraints and check.

**Theorem** Algorithm ATOMANTIUNIFLGG terminates, and is complete and it computes a single lgg of given terms.

The complexity is exponential

Slightly changing their approach and algorithm:

**Result:** a unitary nominal anti-unification algorithm with usual freshness constraints.

**idea:** Apply the distinct-names-assumptions  
(OR: do not take atom names too literally)

## Conclusion and Further Work

- We constructed a nominal anti-unification algorithm that is complete and outputs a single least general generalisation. (for atom-variables, and with EQR-constraints)
  
- We constructed a unitary nominal anti-unification algorithm improving the algorithm and result of Baumgartner, Kutsia, Levy, Villaret, RTA 2015. (Only atoms, no atom-variables)

- Is `ATOMANTIUNIFICATION` with refinement of freshness constraints finitary or unary?
- Extending the applicability of the methods to other higher-order languages. E.g. Haskell with recursive `let` constructs