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Lecture 8: Implementation Techniques         Dutline         1. Dowling-Gallier algorithm         2. Full sets         3. Search method for stable models         4. Implementing approximations         5. Branch&bound algorithm	<ul> <li>1</li> <li>T-79.5102 / Autumn 2007 Implementation Techniques</li> <li>Data structures</li> <li>The iteration sequence of the operator T<sub>P</sub> can be implemented more efficiently using certain precomputed data structures.</li> <li>The following data structures are used for a positive program P</li> <li>1. An array occurs[a] of sets of rules indexed by atoms a ∈ Hb( and precomputed as {r   r = h ← B ∈ P and a ∈ B} for each</li> <li>2. An array count[r] of integers indexed by rules r ∈ P. The ini value for a rule r = h ← B is the number  B  of body atoms.</li> <li>3. A set of atoms M ⊆ Hb(P) initialized as T<sub>P</sub> ↑ 1 = T<sub>P</sub>(0).</li> <li>The execution of the function LeastModel(M) given on the nex slide will gradually extend M to LM(P).</li> </ul>
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<ul> <li>T-79.5102 / Autumn 2007 Implementation Techniques</li> <li><b>1. DOWLING-GALLIER ALGORITHM</b></li> <li>W. Dowling and J. Gallier [1984] presented their method originally for testing the satisfiability of sets of Horn clauses.</li> <li>Due to close interconnection of Horn clauses with rules and constraints, we present the method for positive programs directly.</li> <li>The iteration sequence T<sub>P</sub> ↑ 0, T<sub>P</sub> ↑ 1, provides a basic method for computing LM(P) = Ifp(T<sub>P</sub>) for a positive program P.</li> <li>The least fixpoint, for which T<sub>P</sub> ↑ n = T<sub>P</sub> ↑ n − 1 holds, is reached in the method is the satisfication of the satisficat</li></ul>	2 T-79.5102 / Autum 2007 Implementation Techniques Algorithm for Computing LM(P) function LeastModel( $M$ : atom set): atom set; var $Q$ : atom set; $a$ : atom; $r$ : rule; Q := M; while $Q \neq 0$ do a := pick(Q); $Q := Q \setminus \{a\}$ ; for $r$ in occurs[ $a$ ] do count[r] := count[r] - 1; if $count[r] := 0$ and $Head(r) \notin M$ then do $Q := Q \cup \{Head(r)\}$ ; $M := M \cup \{Head(r)\}$ ;

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<b>Example.</b> Consider a positive program <i>P</i> having the following rules:						
$r_0:c.$ $r_1:d.$ $r_2:e \leftarrow$	$-c, d.$ $r_3: f \leftarrow c.$ $r_4$	$: g \leftarrow g, f.  r_5 : h \leftarrow d, f.$				
The algorithm computes $\operatorname{LM}(P) = \{c, d, e, f, h\}$ as follows:						
1. $M := \{c, d\}$	11. $a := d$	21. $r := r_4$				
2. $Q := \{c, d\}$	12. $Q := \{f\}$	22. $count[r_4] := 1$				
3. $a := c$	13. $r := r_2$	23. $r := r_5$				
4. $Q := \{d\}$	14. $count[r_2] := 0$	24. $count[r_5] := 0$				
5 $r := r_2$	15. $Q := \{f, e\}$	25. $Q := \{e, h\}$				
6 count $[r_2] := 1$	16. $M := \{c, d, e, f\}$	26. $M := \{c, d, e, f, h\}$				
7. $r := r_3$	17. $r := r_5$	27. $a := e$				
8 count $[r_3] := 0$	18. $count[r_5] := 1$	28. $Q := \{h\}$				
9. $Q := \{d, f\}$	19. $a := f$	29. $a := h$				
10. $M := \{c, d, f\}$	20. $Q := \{e\}$	30. $Q := \emptyset$				

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- > Assuming constant costs for set operations and arithmetics, the algorithm runs in time linear with respect to ||P||.
- $\blacktriangleright$  The implementation of Q determines the way how computation proceeds: e.g., breadth-first (LIFO) or depth-first (FIFO).

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- $\blacktriangleright$  The set of atoms  $a \in \operatorname{Hb}(P)$  that appear in the negative body literals of a normal program P is denoted by NBA(P).
- $\blacktriangleright$  The members of NBA(P) affect the reduct  $P^M$  play a major role when the stable models M of a normal program P are determined.
- $\blacktriangleright$  The stable models  $M \in SM(P)$  can be characterized in terms of sets of *negative default literals*  $\sim a$  based on atoms  $a \in NBA(P)$ .
- $\blacktriangleright$  The least model associated with a normal logic program P and a set of negative default literals F is  $LM(P_F)$  where

$$P_F = \{a \leftarrow B \mid a \leftarrow B, \sim C \in P \text{ and } \sim C \subseteq F\}.$$

**Definition.** A set F of negative default literals is P-full if and only if for all  $a \in NBA(P)$ , the literal  $\sim a \in F \iff a \notin LM(P_F)$ .

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# 3. SEARCH METHOD FOR STABLE MODELS

- The goal is to compute—as efficiently as possible—one or several stable models for a normal program P given as input.
- ➤ The characterization of stable models based on full sets suggests that the search space essentially consists of subsets of NBA(P).
- ➤ Following the general *branch&bound* search strategy, we gradually build a set L of default literals that constrains stable models being computed and try to prune the search space.
  - 1. Assumptions about models are made one by one.
  - 2. At each point of the search space, stable models that satisfy all the assumptions introduced so far (the set L) are approximated.
  - 3. If a conflict is found, the search backtracks, and the search for other models takes place similarly, if a model is found.

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## **Approximation Criteria**

- ➤ Stable models being computed for a normal program *P* are specified in terms of a set of default literals *L* over Hb(*P*):
  - 1. If  $a \in L$ , then  $a \in M$  for stable models M being computed.
  - 2. If  $\sim a \in L$ , then  $a \notin M$  for stable models M being computed  $(\sim a \in F \text{ holds for the respective full sets } F)$ .
- Such a relationship between a stable model  $M \in SM(P)$  and a set of default literals can be understood as a form of *compatibility*.

**Example.** Consider the normal logic program P =

 $\{a \leftarrow \sim b. \ b \leftarrow \sim a. \ c \leftarrow \sim d. \ d \leftarrow \sim c. \ e \leftarrow \sim f. \ f \leftarrow \sim e. \}.$ 

Now, for instance, the set of default literals  $L = \{a, \sim c\}$  is compatible with stable models  $M_1 = \{a, d, e\}$  and  $M_2 = \{a, d, f\}$  in SM(P).

### Lower and Upper Bounds

- ➤ A lower bound  $LB(P,L) \supseteq L$  is a set of literals which is compatible with any stable model  $M \in SM(P)$  compatible with L.
- ➤ An upper bound  $UB(P,L) \subseteq Hb(P)$  is a set of atoms that contains every  $M \in SM(P)$  compatible with L.
- ➤ An approximation Expand(P,L) is the *least* set of literals L' which contains L and is *closed* in the following senses:
  - (i) If a default literal  $l \in LB(P,L')$ , then  $l \in L'$ .
  - (ii) If an atom  $a \notin UB(P,L')$ , then  $\sim a \in L'$ .
- ➤ The approximation Expand(P,L) can be obtained by computing lower and upper bounds iteratively and applying (i) and (ii).

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## 4. IMPLEMENTING APPROXIMATIONS

- For the sake of efficiency, it is important that the bounds LB(P,L)and UB(P,L) can be computed in linear time.
- The resulting approximation at any point of the search space L should be at least as accurate as WFM(P), i.e.,  $WFM(P) \subseteq L$ .
- > Assumptions embodied in L should be taken fully into account.

**Definition.** For a normal logic program P and a set of default literals L over Hb(P), the set of *active rules* of P given L is

 $\operatorname{ActR}(P,L) = \{a \leftarrow B, \sim C \in P \mid L \cap (\sim B \cup C) = \emptyset\}.$ 

**Remark.** The bodies of rules in ActR(P,L) are not falsified by L!

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Lower Bound		Upper Bound
<b>Definition.</b> The lower bound $LB(P,L)$ is the least set of literals $L'$ which contains $L$ and is closed under the following principles:		<b>Definition.</b> The upper bound $UB(P,L) = LM(ActR(P,L)^{\emptyset})$ where the reduct $ActR(P,L)^{\emptyset}$ is $ActR(P,L)$ with all negative literals removed.
P1: If $a \leftarrow B$ , $\sim C \in \operatorname{ActR}(P,L')$ and $B \cup \sim C \subseteq L'$ , then $a \in L'$ .		<b>Example.</b> Consider the following normal program $P_2$ :
P2: If $b \neq a$ for every rule $a \leftarrow B, \sim C$ in $\operatorname{ActR}(P,L')$ , then $\sim b \in L'$ .		$a \leftarrow \sim b.  b \leftarrow \sim a.  c \leftarrow \sim a.  d \leftarrow \sim c.  e \leftarrow \sim d.$
<ul> <li>P3: If a ∈ L' is the head of exactly one rule a ← B, ~C in ActR(P,L'), then the body B ∪ ~C ⊆ L'.</li> <li>P4: If ~a ∈ L', a rule a ← l<sub>1</sub>,, l<sub>n</sub> ∈ ActR(P,L'), and {l<sub>1</sub>,, l<sub>i-i</sub>, l<sub>i+1</sub>,, l<sub>n</sub>} ⊆ L', then the complement l<sub>i</sub> ∈ L'.</li> <li>P5: If for some atom a ∈ Hb(P) both a ∈ L' and ~a ∈ L', then all</li> </ul>		Verify the following upper bounds for $P_2$ : 1. UB $(P_2, \emptyset) = \{a, b, c, d, e\}$ . $\implies$ Expand $(P_2, \emptyset) = \emptyset$ because also LB $(P_2, \emptyset) = \emptyset$ holds. 2. UB $(P_2, \{a\}) =$ LM $(\{a \leftarrow \sim b. \ d \leftarrow \sim c. \ e \leftarrow \sim e. \}^{\emptyset}) = \{a, d, e\}$ .
literals over $\operatorname{Hb}(P)$ belong to $L'$ .		3. UB( $P_2, \{\sim a\}$ ) = { $a, b, c, d, e$ }.
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Example		Further Examples
Let us compute $L_1 = \operatorname{LB}(P_1, \{\sim b\})$ and $L_2 = \operatorname{LB}(P_1, \{b\})$ for $P_1$ :		<b>Example</b> Let us compute approximations $Expand(\mathbf{P}_{a} \mid a)$ and

 $\operatorname{ActR}(P_1, L_2)$ 

 $r_2, r_3, r_4$ 

 $r_2, r_3$ 

 $r_2$ 

**Example.** Let us compute approximations  $Expand(P_2, \{a\})$  and Expand( $P_2$ , { $\sim a$ }) for the preceding program  $P_2$ :

 $a \leftarrow \sim b$ .  $b \leftarrow \sim a$ .  $c \leftarrow \sim a$ .  $d \leftarrow \sim c$ .  $e \leftarrow \sim d$ .

1.  $LB(P_2, \{a\}) = \{a, \sim b, \sim c, d, \sim e\} = Expand(P_2, \{a\}).$ 

2.  $LB(P_2, \{\sim a\}) = \{\sim a, b, c, \sim d, e\} = Expand(P_2, \{\sim a\}).$ 

**Example.** Let us then analyze a normal program  $P_3$  having two rules:

 $a \leftarrow \sim b$ .  $b \leftarrow b$ .

Now  $LB(P_3, \emptyset) = \emptyset$  but  $UB(P_3, \emptyset) = \{a\}$  so that  $\sim b \in Expand(P_3, \emptyset)$ .  $\implies$  Expand $(P_3, \emptyset) = \{a, \sim b\}$  because LB $(P_3, \{\sim b\}) = \{a, \sim b\}$ .

 $r_1: a \leftarrow \sim b$ .  $r_2: c \leftarrow a$ .  $r_3: b \leftarrow a, \sim c, \sim d$ .  $r_4: d \leftarrow c, \sim e$ .

 $L_2$ 

h

С

all

 $a, \sim c, \sim d$ 

Ρ3

Ρ1

Ρ5

 $\operatorname{ActR}(P_1, L_1)$ 

 $r_1, r_2, r_3, r_4$ 

 $r_1, r_2, r_3, r_4$ 

 $r_1, r_2, r_4$ 

 $r_1, r_2, r_4$ 

 $r_1, r_2, r_4$ 

 $L_1$ 

а

С

d

 $\sim e$ 

 $\sim b$ 

P1

Ρ1

Ρ2

Ρ1

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Implementing Bounds	
<ul> <li>The lower and upper bounds can be implemented as linear time algorithms that resemble the Dowling-Gallier presented above.</li> <li>The accuracy of Expand is at least as good as that of the well-founded model. In fact, we have Expand(P,0) = WFM(P).</li> <li>However, assumptions about stable models to compute, a set of default literals L ⊈ WFM(P), make Expand(P,L) more accurate.</li> <li>Example. For the normal program P<sub>4</sub> consisting of <ul> <li>a ← ~b.</li> <li>b ← ~a.</li> <li>a ← b.</li> </ul> </li> <li>we obtain Expand(P<sub>4</sub>,0) = 0 and Expand(P<sub>4</sub>, {b}) = {a,b,~a,~b}.</li> <li>But the conflict is not detected for P<sub>5</sub> = P<sub>4</sub> ∪ {b.}: WFM(P) = {a,b}.</li> </ul>	function Smodels( var A: literal set; I: A := Expand(P,L) if Conflict(P,A) t if Covered(P,A) t l := Choose(P,A) if Smodels(P,A $\cup$ return $\top$ ; else return Smo Remarks. Recall th The sets of literals represented in space
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5. BRANCH&BOUND ALGORITHM	Exa

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- ➤ The purpose of the function Smodels(P,L) is to check whether P has a stable model M which is compatible with L.
- > The underlying algorithm is based on a number of primitives:

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- 1. The function Expand(P,L) returns a tightened approximation of stable models that are compatible with L.
- 2. The function Conflict(P,L) checks whether the approximation L obtained so far is contradictory  $(\{a, \sim a\} \subseteq L \text{ for some } a)$ .
- 3. The function Covered(P,L) checks whether the approximation L obtained so far covers all atoms of NBA(P), i.e., for each  $a \in NBA(P)$ , either  $a \in L$  or  $\sim a \in L$ .
- The function Choose(P,L) implements the search heuristics, i.e., it picks the next *literal* for branching.



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