ASP Solvers:

The Smodels Language

▲□▶ ▲□▶ ▲ 三▶ ▲ 三▶ 三三 - のへぐ

ASP-solvers

▲□▶▲□▶▲≡▶▲≡▶ ≡ めぬる

- There are lots of ASP solvers available: *smodels, dlv, cmodels, assat, nomore, smodels_{cc}, clasp, platypus, ...*
- We'll briefly look at the original smodels implementation

Smodels

▲□▶ ▲□▶ ▲□▶ ▲□▶ ■ ●の00

Origin Helsinki University of Technology

People I. Niemelä, P. Simons, and T. Syrjänen

Web http://www.tcs.hut.fi/Software/smodels/

Input format lparse(gringo)

Extensions • choice rules

- cardinality/weight constraints
- optimization via minimize/maximize statements

- Let Π be a program and $atoms(\Pi)$ the atoms in Π .
- An assignment is a partial mapping $A : atoms(\Pi) \rightarrow \{T, F\}.$
 - \square Think of A as a three-valued model.

▲□▶ ▲□▶ ▲□▶ ▲□▶ ■ ●の00

- Let Π be a program and $atoms(\Pi)$ the atoms in Π .
- An assignment is a partial mapping $A : atoms(\Pi) \rightarrow \{ \mathbf{T}, \mathbf{F} \}.$

Think of A as a three-valued model.

• An assignment A can be denoted as a pair (A^T, A^F) , where

•
$$A^{\boldsymbol{T}}_{-} = \{ v \in atoms(\Pi) \mid A(v) = \boldsymbol{T} \}$$

•
$$A^{\boldsymbol{F}} = \{ v \in atoms(\Pi) \mid A(v) = \boldsymbol{F} \}$$

A D N A 目 N A E N A E N A B N A C N

- Let Π be a program and $atoms(\Pi)$ the atoms in Π .
- An assignment is a partial mapping $A : atoms(\Pi) \rightarrow \{T, F\}.$

Think of A as a three-valued model.

• An assignment A can be denoted as a pair (A^T, A^F) , where

•
$$A_{\underline{r}}^{T} = \{ v \in atoms(\Pi) \mid A(v) = T \}$$

•
$$A^{\boldsymbol{F}} = \{v \in atoms(\Pi) \mid A(v) = \boldsymbol{F}\}$$

We also denote an assignment A by a set of signed objects:
 {Tv | v ∈ A^T} ∪ {Fv | v ∈ A^F}

A D N A 目 N A E N A E N A B N A C N

- Let Π be a program and $atoms(\Pi)$ the atoms in Π .
- An assignment is a partial mapping $A : atoms(\Pi) \rightarrow \{T, F\}.$

Think of A as a three-valued model.

• An assignment A can be denoted as a pair (A^T, A^F) , where

•
$$A^{\mathsf{T}} = \{v \in atoms(\Pi) \mid A(v) = \mathsf{T}\}$$

•
$$A^{F} = \{v \in atoms(\Pi) \mid A(v) = F\}$$

- We also denote an assignment A by a set of signed objects:
 {Tv | v ∈ A^T} ∪ {Fv | v ∈ A^F}
- Example: Given $atoms(\Pi) = \{a, b, c\}$, assignment $A = \{a \mapsto T, c \mapsto F\}$

can be represented as

$$(\{a\},\{c\})$$
 or $\{Ta,Fc\}$.

(Simplified) Smodels Algorithm

A D N A 目 N A E N A E N A B N A C N

Let Π be a program and A a (partial) assignment.

```
smodels(\Pi, A):

A \leftarrow expand(\Pi, A)

if A^T \cap A^F \neq \emptyset then return

if A^T \cup A^F = atoms(\Pi) then exit with A^T

x \leftarrow select(atoms(\Pi) \setminus (A^T \cup A^F))

smodels(\Pi, A \cup \{Tx\})

smodels(\Pi, A \cup \{Fx\})
```

Call: smodels(Π , \emptyset)

smodels algorithm

- Backtracking search, building a binary search tree
- Choices on unassigned atoms
- The search space is pruned by
 - making one choice at a time by appeal to a heuristics (*select*)
 - the set of remaining choices is reduced and conflicts are detected (*expand*)

▲ロ ▶ ▲周 ▶ ▲ 国 ▶ ▲ 国 ▶ ● の Q @

• expand(Π , A): repeat $A \leftarrow atleast(\Pi, A)$ $A' \leftarrow A$ $A \leftarrow A \cup \{F_X \mid x \in atom(\Pi) \setminus atmost(\Pi, A)\}$ until A = A'

▲□▶ ▲□▶ ▲ 三▶ ▲ 三▶ 三三 - のへぐ

• expand(
$$\Pi$$
, A):
repeat
$$A \leftarrow atleast(\Pi, A)$$
$$A' \leftarrow A$$
$$A \leftarrow A \cup \{Fx \mid x \in atom(\Pi) \setminus atmost(\Pi, A)\}$$
until $A = A'$

- atleast amounts to Fitting's operation
 - These are atoms that **must** be true or false, given A.

$$\begin{array}{l} \mathsf{P} \; \mathsf{expand}(\Pi, A):\\ \mathsf{repeat}\\ & A \leftarrow \mathsf{atleast}(\Pi, A)\\ & A' \leftarrow A\\ & A \leftarrow A \cup \{ \mathbf{F}x \mid x \in \mathsf{atom}(\Pi) \setminus \mathsf{atmost}(\Pi, A) \}\\ \mathsf{until} \; A = A' \end{array}$$

- atleast amounts to Fitting's operation
 - These are atoms that must be true or false, given A.
- atmost computes those atoms derivable in the underlying positive program of Π.
 - Its complement contains those atoms that cannot be derived, given *A*.
 - This is called the greatest unfounded set

- ロ ト - 4 回 ト - 4 □ - 4

expand(
$$\Pi$$
, A):
repeat
 $A \leftarrow atleast(\Pi, A)$
 $A' \leftarrow A$
 $A \leftarrow A \cup \{Fx \mid x \in atom(\Pi) \setminus atmost(\Pi, A)\}$
until $A = A'$

- atleast amounts to Fitting's operation
 - These are atoms that must be true or false, given A.
- atmost computes those atoms derivable in the underlying positive program of Π.
 - Its complement contains those atoms that cannot be derived, given *A*.
 - This is called the greatest unfounded set
- expand computes the *well-founded model* of Π .

Heuristics: Lookahead

▲□▶ ▲□▶ ▲ 三▶ ▲ 三▶ 三 のへぐ

- Strengthen propagation by failed literal detection
- Given a program Π , an atom x, and an assignment A
 - if expand($\Pi, A \cup \{Tx\}$) yields a conflict, then add Fx to A
 - if expand($\Pi, A \cup \{F_X\}$) yields a conflict, then add T_X to A
 - if both yield a conflict, "backtrack"

Heuristics: Lookahead

- Strengthen propagation by failed literal detection
- Given a program Π , an atom x, and an assignment A
 - if expand($\Pi, A \cup \{Tx\}$) yields a conflict, then add Fx to A
 - if expand($\Pi, A \cup \{Fx\}$) yields a conflict, then add Tx to A
 - if both yield a conflict, "backtrack"
- Lookahead is also used for selecting the next unassigned atom x

Heuristics: Lookahead

- Strengthen propagation by failed literal detection
- Given a program Π , an atom x, and an assignment A
 - if expand($\Pi, A \cup \{Tx\}$) yields a conflict, then add Fx to A
 - if expand($\Pi, A \cup \{Fx\}$) yields a conflict, then add Tx to A
 - if both yield a conflict, "backtrack"
- Lookahead is also used for selecting the next unassigned atom x
- That is, given that
 - x^+ is the number of atoms that are assigned through expand($\Pi, A \cup \{Tx\}$) and
 - x[−] is the number of atoms that are assigned through expand(Π, A ∪ {Fx}),

select an atom x with a maximal $min(x^+, x^-)$