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On the Relation of Constraint Answer Set Programming Languages and Algorithms

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Abstract
Recently a logic programming language AC was proposed by Mellarkod et al. (2008) to integrate answer set programming (ASP) and constraint logic programming. Similarly, Gebser et al. (2009) proposed a CLINGCON language integrating ASP and finite domain constraints. These languages allow new efficient inference algorithms that combine traditional ASP procedures and other methods in constraint programming. In this paper we show that a transition system introduced by Nieuwenhuis et al. (2006) to model SAT solvers can be extended to model the "hybrid" ACSOLVER algorithm by Mellarkod et al. developed for simple AC programs and the CLINGCON algorithm by Gebser et al. for clingcon programs. We define weakly-simple programs and show how the introduced transition systems generalize the ACSOLVER and CLINGCON algorithms to such programs. Finally, we state the precise relation between AC and CLINGCON languages and the ACSOLVER and CLINGCON algorithms.

Introduction
Mellarkod et al. (2008) introduced a knowledge representation language AC extending the syntax and semantics of answer set programming with constraint processing features. The origins of their work go back to (Baselice, Bonatti, and Gelfond 2005). In a similar vein, Gebser et al. (2009) proposed a CLINGCON language integrating ASP and finite domain constraints. The AC and CLINGCON languages allow not only new modeling features but also novel computational methods that combine traditional ASP algorithms with constraint (logic) programming (CLP/CSP) algorithms. This combined approach opens new horizons for declarative programming applications. For instance, it allows us to reason about problems with variables whose values range over very large domains such as dynamic systems in real time. Mellarkod et al. presented a “hybrid” ACSOLVER system for finding answer sets of AC programs that combines both ASP and CLP computational tools. The key feature of this system is that it processes a “regular” part of a given program using the ASP algorithm SMODELS (Niemelä and Simons 2000) and a “defined” part using CLP tools. Similarly, system CLINGCON (Gebser, Ostrowski, and Schaub 2009) takes advantage of answer set solver CLASP (Gebser et al. 2007) and constraint solver GECODE (http://www.gecode.org). It is intuitively clear that the AC and CLINGCON languages are related as well as the systems ACSOLVER and CLINGCON. This paper puts these relationships in precise mathematical terms.

We show that transition systems (graphs) introduced by Nieuwenhuis et al. (2006) to model and analyze SAT solvers can be adapted to describe constraint answer set solvers ACSOLVER and CLINGCON. By introducing such new transition systems we provide an alternative description of ACSOLVER and CLINGCON and also an alternative proof of their correctness. This abstract view on the systems allows us to state the relation between them in precise terms by studying the underlying graph representations. The ACSOLVER algorithm was proved to be correct for a class of “simple” programs. We define a more general class of weakly-simple programs and demonstrate how newly introduced transition systems immediately capture a class of algorithms for such programs and demonstrate their correctness.

This work clarifies and extends state of the art developments in the area of constraint answer set programming and we believe will promote further progress in the area.

We start by reviewing AC logic programs and a notion of an answer set for such programs. We introduce a new class of weakly-simple programs. We then review a transition system introduced by Lierler (2008) to model “simple” programs. We extend this transition system to model the ACSOLVER algorithm and show how the newly defined graph can characterize the computation behind the system ACSOLVER. In the subsequent section we introduce the CLINGCON language and formally state its relation to the AC language. At last we define a graph suitable for modeling the system CLINGCON and state a formal result on the relation between the ACSOLVER and CLINGCON algorithms.

A preliminary report on some of the results of this paper has been presented at Workshop on Answer Set Programming and Other Computing Paradigms (Lierler and Zhang 2011).

Review: AC Logic Programs
A sort (type) is a non-empty countable collection of strings over some fixed alphabet. A signature Σ is a collection of sorts, properly typed predicate symbols, constants, and variables. Sorts of Σ are divided into regular and constraint
sorts. All variables in Σ are of a constraint sort. A term of Σ is either a constant or a variable. An atom is of the form \( p(t_1, \ldots, t_n) \) where \( p \) is an \( n \)-ary predicate symbol, and \( t_1, \ldots, t_n \) are terms of the proper sorts. A constraint sort is often a large numerical set with primitive constraint relations. The partitioning of sorts induces a partition of predicates of the AC language:

- **Regular predicates** denote relations among constants of regular sorts;
- **Constraint predicates** denote primitive constraint relations on constraint sorts;
- **Defined predicates** denote relations between constants that belong to regular sort and objects that belong to constraint sorts; such predicates can be defined in terms of constraint, regular, and defined predicates;
- **Mixed predicates** denote relations between constants that belong to regular sort and objects that belong to constraint sorts. Mixed predicates are not defined by the rules of a program and are similar to abducible relations of abductive logic programming (Kakas, Kowalski, and Toni 1992).

An atom formed by a regular predicate is called regular. Similarly for constraint, defined, and mixed atoms. We say that an atom is a **non-mixed atom** if it is regular, constraint, or defined.

A **regular program** is a finite set of rules of the form

\[
a_0 \leftarrow a_1, \ldots, a_l, \text{not} a_{l+1}, \ldots, \text{not} a_m, \\
\text{not not} a_{m+1}, \ldots, \text{not not} a_n,
\]

where \( a_0 \) is \( \perp \) or a ground (non-constraint) atom, and each \( a_i \) \((1 \leq i \leq n)\) is a ground (non-constraint) atom. If \( a_0 = \perp \), we often omit \( \perp \) from the notation. This is a special case of programs with nested expressions (Lifschitz, Tang, and Turner 1999). We assume that the reader is familiar with the definition of an answer set of a logic program and refer to the paper by Lifschitz et al. (1999) for details. A **choice rule** construct \{ \( a \) \} (Niemelä and Simons 2000) of the LParse1 language can be seen as an abbreviation for a rule \( a \leftarrow \text{not not} a \) (Ferraris and Lifschitz 2005). We adopt this abbreviation in the rest of the paper.

An ((AC) logic program) is a finite set of rules of the form \( (1) \) where

- \( a_0 \) is \( \perp \) or a regular or defined atom,
- each \( a_i \), \( 1 \leq i \leq l \), is an arbitrary atom if \( a_0 \) is \( \perp \) or a regular atom,
- each \( a_i \), \( 1 \leq i \leq l \), is a non-mixed atom if \( a_0 \) is a defined atom,
- each \( a_i \), \( l + 1 \leq i \leq m \), is a non-mixed atom,
- each \( a_i \), \( m + 1 \leq i \leq n \), is a regular atom,
- \( n = m \), if \( a_0 \) is a defined atom.

Rule \( (1) \) is called a **defined rule** if \( a_0 \) is a defined atom. It is easy to see that defined rules of a program neither contain mixed atoms in its body nor contain doubly negated atoms \( (not not a) \). We assume that any mixed atom occurring in AC program is of the restricted form \( n(\vec{r}, V) \), where \( \vec{r} \) is a sequence of regular constants and \( V \) is a variable.

A part of the AC program \( \Pi \) that consists of defined rules is called a **defined** part denoted by \( \Pi_D \). By \( \Pi_R \) we denote a non-defined part of \( \Pi \), i.e., \( \Pi \setminus \Pi_D \). For instance, let signature \( \Sigma \) contain two regular sorts \( \text{step} = \{0\} \), \( \text{action} = \{a\} \) and two constraint sorts \( \text{time} = \{0..200\} \), \( \text{computer} = \{1..2\} \); a mixed predicate \( \text{okComp}(\text{computer}, \text{time}) \), two regular predicates \( \text{okTime}(\text{time}) \), \( \text{okComp}(\text{computer}, \text{time}) \). A sample AC program over \( \Sigma \) follows

\[
\begin{align*}
\text{okComp}(1, T) & \leftarrow T \leq 5, \\
\text{okComp}(2, 106) & \leftarrow \text{okTime}(T) \leftarrow T \leq 10, \\
\text{okComp}(1, T) & \leftarrow \text{okTime}(T) \leftarrow T \geq 100, \\
\text{okComp}(2, T) & \leftarrow \text{okTime}(a(0, 0), T) \leftarrow T \neq 1, \\
\text{okComp}(a(0, 0)) & \leftarrow \{0\}
\end{align*}
\]

The first four rules of the program form its defined part whereas the last three rules form \( \Pi_R \).

Mellarkod et al. (Mellarkod, Gelfond, and Zhang 2008) considered programs of more sophisticated syntax than discussed here. For instance, in (Mellarkod, Gelfond, and Zhang 2008) classical negation may precede atoms in rules. Also signature \( \Sigma \) may contain variables of regular sort. Nevertheless, the AC language discussed here is sufficient to capture the class of programs covered by the ACSOLVER algorithm.

The expression \( a_0 \) is the **head** of a rule \( (1) \). If \( B \) denotes the body of \( (1) \), the right hand side of the arrow, we write \( B^{pos} \) for the elements occurring in the positive part of the body, i.e., \( B^{pos} = \{ a_1, \ldots, a_l \} \) and \( B^{neg} \) for the elements occurring under single negation as failure, i.e., \( B^{neg} = \{ a_{l+1}, \ldots, a_m \} \). We frequently identify the body of \( (1) \) with the conjunction of its elements (in which \( not \) is replaced with the classical negation connective \( \neg \)):

\[
a_1 \wedge \cdots \wedge a_l \wedge \neg a_{l+1} \wedge \cdots \wedge \neg a_m \wedge \neg a_{m+1} \wedge \cdots \wedge \neg a_n.
\]

Similarly, we often interpret a rule \( (1) \) as a clause

\[
a_0 \neg \neg a_1 \vee \cdots \vee \neg a_l \vee a_{l+1} \vee \cdots \vee a_m \neg a_{m+1} \vee \cdots \vee \neg a_n
\]

(3)

in the case when \( a_0 = \perp \) in \( (1) \) \( a_0 \) is absent in \( (3) \). Given a program \( \Pi \), we write \( \Pi^{df} \) for the set of clauses \( (3) \) corresponding to all rules in \( \Pi \).

For an AC program \( \Pi \) over signature \( \Sigma \), by the set **ground**(\( \Pi \)) we denote the set of all ground instances of all rules in \( \Pi \). The set **ground***(\( \Pi \)) is obtained from **ground**(\( \Pi \)) by

- dropping the rules where a constraint atom \( a \) occurs in \( B^{pos} \) (\( B^{neg} \)) and \( a \) is false (true, respectively) under the intended interpretation of its symbols,
- dropping all constraint literals from the remaining rules (here we use the term literal to refer to \( a \) and \( not a \)).

It is easy to see that **ground***(\( \Pi \)) is a regular program.

For instance, let **ground**(\( \Pi \)) consist of two rules

\[
\begin{align*}
\text{okTime}(100) & \leftarrow 100 > 100, \\
\text{okComp}(2, 100) & \leftarrow \text{okTime}(101) \leftarrow 101 > 100, \\
\text{okComp}(2, 101)
\end{align*}
\]
then $ground^*(\Pi)$ is
\[
okTime(101) \leftarrow okComp(2, 101).
\]

We say that a sequence of (regular) constants $\vec{r}$ is specified by a mixed predicate $m$ if $\vec{r}$ follows the sorts of the regular arguments of $m$. For instance, for program (2) a sequence $0$ of constants (of type step) is the only sequence specified by mixed predicate $at$. For a set $X$ of atoms, we say that a sequence $\vec{r}$ of regular constants is bound in $X$ by a (constraint) constant $c$ w.r.t. predicate $m$ if there is an atom $m(\vec{r}, c)$ in $X$. A set $M$ of ground mixed atoms is functional over the underlying signature if for every mixed predicate $m$, every sequence of regular constants specified by $m$ is bound in $M$ by a unique constraint constant w.r.t. $m$. For instance, for the signature of program (2) sets \{at(0, 1)\} and \{at(0, 2)\} are functional whereas \{at(0, 1) at(0, 2)\} is not a functional set because $0$ is bound in $M$ by two different constants 1 and 2 w.r.t. $at$.

Definition 1 For an AC program $\Pi$, a set $X$ of atoms is called an answer set of $\Pi$ if there is a functional set $M$ of ground mixed atoms of $\Sigma$ such that $X$ is an answer set of $\Pi$.

For example, sets of atoms
\[
\{at(0, 1), occurs(a, 0)\}
\]
and
\[
\{on, at(0, 0), occurs(a, 0), okComp(1, 0), \ldots, okComp(1, 5), okComp(2, 106), okTime(0), \ldots, okTime(5), okTime(106)\}
\]
are answer sets of (2).

The definition of an answer set for AC programs presented here is different from the original definition in (Mellarkod, Gelfond, and Zhang 2008), but there is a close relation between them.

Proposition 1 For an AC program $\Pi$ over signature $\Sigma$ such that $\Pi$ contains no doubly negated atoms and the set $S$ of all true ground constraint literals over $\Sigma$, $X$ is an answer set of $\Pi$ if and only if $X \cup S$ is an answer set (in the sense of (Mellarkod, Gelfond, and Zhang 2008)) of $\Pi$.

Weakly-Simple AC Programs

The correctness of the ACSOLVER algorithm was shown for simple AC programs. We start this section by reviewing simple programs. We then define a more general class of programs called weakly-simple. In the next section we state correctness results for ACSOLVER-like algorithms for such programs.

We say that an AC program $\Pi$ is safe (Mellarkod, Gelfond, and Zhang 2008) if every variable occurring in a non defined rule in $\Pi$ also occurs in a mixed atom of this rule. An AC program $\Pi$ is super safe if $\Pi$ is safe and

1. if a mixed atom $m(\vec{c}, X)$ occurs in $\Pi$ then a mixed atom $m(\vec{c}, X')$ does not occur in $\Pi$ (where $X$ and $X'$ are distinct variable names),
2. if a mixed atom $m(\vec{c}, X)$ occurs in $\Pi$ then neither a mixed atom $m'(\vec{c}', X)$ such that $\vec{c} \neq \vec{c}'$ nor a mixed atom $m'(\vec{c}, X)$ such that $m \neq m'$ occurs in $\Pi$.

We note that any safe AC program $\Pi$ may be converted to a super safe program $\Pi'$ so that the resulting program has the same answer sets by a simple syntactic transformation. Lierler and Zhang (2011, Section 7) provide such a transformation. We say that an AC program $\Pi$ is simple if it is super safe, and its defined part contains no regular atoms and has a unique answer set.

For any atom $p(\vec{t})$, by $p(\vec{t})^0$ we denote its predicate symbol $p$. For any AC program $\Pi$, the predicate dependency graph of $\Pi$ is the directed graph that

- has all predicates occurring in $\Pi$ as its vertices, and
- for each rule $(1)$ in $\Pi$ has an edge from $a_0^0$ to $a_i^0$ where $1 \leq i \leq l$.

We say that an AC program $\Pi$ is weakly-simple if

- it is super safe,
- all regular atoms occurring in $\Pi_D$ also occur in $\Pi_R$, and
- each strongly connected component of the predicate dependency graph of $\Pi$ is a subset of either regular predicates of $\Pi$ or defined predicates of $\Pi$.

It is easy to see that any simple program is also a weakly-simple program but not the other way around. For example, program (2) is weakly-simple but not simple.

Review: Abstract Smodels

Most state-of-the-art answer set solvers are based on algorithms closely related to the DPLL procedure (Davis, Logemann, and Loveland 1962). Nieuwenhuis et al. described DPLL by means of a transition system that can be viewed as an abstract framework underlying DPLL computation (Nieuwenhuis, Oliveras, and Tinelli 2006). Lierler (2008) proposed a similar framework, SMT$_1$ for specifying an answer set solver SMODELS. Our goal is to design a similar framework for describing an algorithm behind ACSOLVER. As a step in this direction we review the graph SM$_2$ that underlines an algorithm of SMODELS, one of the main building blocks of ACSOLVER. The presentation follows (Lierler 2008).

For a set $\sigma$ of atoms, a record relative to $\sigma$ is an ordered set $M$ of literals over $\sigma$, some possibly annotated by $\Delta$, which marks them as decision literals. A state relative to $\sigma$ is a record relative to $\sigma$ possibly preceding symbol $\perp$. For instance, some states relative to a singleton set \{a\} of atoms are
\[
\emptyset, a, \neg a, a^\Delta, a \neg a, \perp, a \perp, \neg a \perp, a^\Delta \perp, a \neg a \perp.
\]

We say that a state is inconsistent if either $\perp$ or two complementary literals occur in it. For example, states $a \neg a$ and $a \perp$ are inconsistent. Frequently, we consider a state $M$ as a set of literals possibly with the symbol $\perp$, ignoring both the annotations and the order between its elements. If neither a literal $l$ nor its complement occur in $M$, then $l$ is unassigned by $M$. For a set $M$ of literals, by $M^+$ and $M^-$ we denote the set of positive and negative literals in $M$ respectively. For instance, $\{a, \neg b\}^+ = \{a\}$ and $\{a, \neg b\}^- = \{b\}$.

If $C$ is a disjunction (conjunction) of literals then by $\overline{C}$ we understand the conjunction (disjunction) of the complements of the literals occurring in $C$. In some situations, we
will identify disjunctions and conjunctions of literals with the sets of these literals. We assume that the reader is familiar with the definition of *unfounded* for the class of regular programs (Lee 2005).

By *Bodies*(Π, a) we denote the sets of bodies of all rules of a regular program Π with the head a. We recall that a set U of atoms occurring in a regular program Π is *unfounded* (Van Gelder, Ross, and Schlipf 1991; Lee 2005) on a consistent set M of literals with respect to Π if for every a ∈ U and every B ∈ *Bodies*(Π, a), M |= B (where B is identified with the conjunction of its elements), or U ∪ B⁺∞ ≠ ∅.

Each regular program Π determines its *Sm models* graph SMΠ. The set of nodes of SMΠ consists of the states relative to the set of atoms occurring in Π. The edges of the graph SMΠ are specified by the transition rules

*Unit Propagate*:

M ⇒ M l if C ∨ l ∈ Π⁺ and C ⊆ M

*Decide*:

M ⇒ M ⊥ if l is unassigned by M

*Fail*:

M ⇒ ⊥ if M is inconsistent and different from ⊥, M contains no decision literals

*Backtrack*:

P |→ Q ⇒ P 7 if P |→ Q is inconsistent, and Q contains no decision literals

*Unfounded*:

M ⇒ M ¬a if a ∈ U for a set U unfounded on M wrt Π and the transition rules *All Rules Cancelled* and *Backchain True* whose details we omit in this review. A node is *terminal* in a graph if no edge leaves this node.

The graph SMΠ can be used for deciding whether a regular program Π has an answer set by constructing a path from ⊥ to a terminal node. Following proposition serves as a proof of correctness and termination for any procedure that is captured by the graph SMΠ.

**Proposition 2** For any regular program Π,

(a) graph SMΠ is finite and acyclic,

(b) for any terminal state M of SMΠ other than ⊥, M⁺ is an answer set of Π,

(c) state ⊥ is reachable from ⊥ in SMΠ if and only if Π has no answer sets.

**Abstract ACSOLVER**

In order to present the transition system suitable for capturing ACSOLVER we introduce several concepts.

**Query, Extensions, and Consequences:** Given an AC program Π and a set p of predicate symbols, a set X of atoms is a p-input answer set (or an input answer set w.r.t. p) of Π if X is an answer set of Π ∪ Xp where by Xp we denote the set of atoms in X whose predicate symbols are different from the ones occurring in p. Intuitively set p denotes a set of intensional predicates (Ferraris et al. 2009). The concept of p-input answer sets is closely related to “p-stable models” in (Ferraris, Lee, and Lifschitz 2011).

\{a(1), b(1)\} of atoms and let p be a set \{a\} of predicates, then Xp is \{b(1)\}. The set X is a p-input answer set of a program \{a\} ← b(1). On the other hand, it is not an input answer set for the same program with respect to a set \{a, b\}.

For a set S of literals, by SR, SD, and SC we denote the set of regular, defined, and constraint literals occurring in S respectively. By SR,D and SP,C we denote the unions SR ∪ SD and SD ∪ SC respectively. By At(Π) we denote the set of atoms occurring in a program Π.

For an AC program Π, a (complete) query Q is a (complete) consistent set of literals over At(ΠD) ∪ At(ΠR) ∪ At(ΠC).

For a query Q of Π, a complete query E is a satisfying extension of Q w.r.t. Π if Q ⊆ E and there is a (sort respecting) substitution γ of variables in E by ground terms so that the result of this substitution, Eγ, satisfies the conditions

1. if a constraint literal l ∈ Eγ then l is true under the intended interpretation of its symbols, and

2. there is an input answer set A of Π ∪ E w.r.t. defined predicates of Π such that Eγ ⊇ A and Eγ ∩ D = ∅.

We say that literal l is a consequence of Q and Π if for every satisfying extension E of Q w.r.t. Π, l ∈ E. By Cons(Q, Π), we denote the set of all consequences of Q and Π. If there are no satisfying extensions of Q w.r.t. Π we identify Cons(Q, Π) with the singleton {⊥}.

Let Π be (2) and Q be \{okTime(T), T ≠ 1\}. A set

\{on, okTime(T), T ≠ 1\}

forms a satisfying extension of Q w.r.t. Π. Indeed, consider a substitution \{T/100\}. This is the only satisfying extension of Q w.r.t. Π. Consequently, it forms Cons(Q, Π).

On the other hand, there are no satisfying extensions for a query \{¬on, okTime(T)\} so that \{⊥\} corresponds to Cons(Q, Π).

**The graph ACΠ:** For each constraint and defined atom A of signature Σ, select a new symbol Aξ, called the name of A. By Σξ we denote the signature obtained from Σ by adding all names Aξ as additional regular predicate symbols (so that Aξ itself is a regular atom).

For an AC program Π, we select a set of rules consisting of (i) choice rules \{ξ\} for each constraint and defined atom a occurring in Π, and (ii) ΠR whose mixed atoms are dropped, and constraint and defined atoms are replaced by their names. Note that Πξ is a regular program.

For instance, let Π be (2) then Πξ consists of the rules

\{T ≠ 1ξ\} \{okTime(T)ξ\}

← occurs(a, 0), T ≠ 1ξ, not okTime(T)ξ

occurs(a, 0) \{on\}

(5)

For a set M of atoms over Σξ, by Mξ we denote a set of atoms over Σ by replacing each name Aξ occurring in M with a corresponding atom A. For instance, \{T ≠ 1ξ, okTime(T)ξ\} is \{T ≠ 1, okTime(T)\}.

Let Π be an AC logic program. The nodes of the graph ACΠ are the states relative to the set of atoms occurring in Πξ.

For a state M of ACΠ, by query(M) we denote the largest subset of Mξ over At(ΠD) ∪ At(ΠR) ∪ At(ΠC). Let Π be (2)
and $M$ be a state $\text{occurs}(a,0) \ -on^\Delta \ \text{okTime}(T)^\xi^\Delta$ then $\text{query}(M)$ is $\{\text{on}, \ \text{okTime}(T)\}$.

The edges of the graph $\text{AC}_\Pi$ are described by the transition rules of $\text{SM}_\Pi^c$ and the additional transition rule

$$\text{Query Propagate:}$$

$$M \rightarrow M \bot^l \text{ if } l \in \text{Cons}(\Pi, \text{query}(M)),$$

where we abuse notation and identify $\bot^l$ with $\bot$ itself.

The graph $\text{AC}_\Pi$ can be used for deciding whether a weakly-simple $\text{AC}$ program $\Pi$ has an answer set by constructing a path from $\emptyset$ to a terminal node:

**Proposition 3** For any weakly-simple $\text{AC}$ program $\Pi$,

(a) graph $\text{AC}_\Pi$ is finite and acyclic,

(b) for any terminal state $M$ of $\text{AC}_\Pi$ other than $\bot$, $(M^\xi^\Delta)^+ R$ is a set of all regular atoms in some answer set of $\Pi$,

(c) state $\bot$ is reachable from $\emptyset$ in $\text{AC}_\Pi$ if and only if $\Pi$ has no answer sets.

Proposition 3 shows that algorithms that find a path in the graph $\text{AC}_\Pi$ from $\emptyset$ to a terminal node can be regarded as $\text{AC}$ solvers for weakly-simple programs.

Let $\Pi$ be an $\text{AC}$ program (2). Here is a path in $\text{AC}_\Pi$ with every edge annotated by the name of a transition rule that justifies the presence of this edge in the graph:

$$\emptyset \xrightarrow{\text{Unit Propagate}} \text{occurs}(a,0) \xrightarrow{\text{Decide}} \text{occurs}(a,0) \xrightarrow{-on^\Delta \ \text{Decide}} \text{occurs}(a,0) \xrightarrow{-on^\Delta \ (\text{okTime}(T)^\xi^\Delta) \ \text{Query Propagate}} \text{occurs}(a,0) \xrightarrow{-on^\Delta \ (\text{okTime}(T)^\xi^\Delta) \ \text{Backtrack}} \text{occurs}(a,0) \xrightarrow{-on^\Delta \ (\text{okTime}(T)^\xi^\Delta) \ \text{Decide}} \text{occurs}(a,0) \xrightarrow{-on^\Delta \ -\text{okTime}(T)^\xi^\Delta \ -T \neq 1^\xi}$$

Since the last state in the path is terminal, Proposition 3 asserts that $\text{occurs}(a,0)$ is a set of all regular atoms in some answer set of $\Pi$. Indeed, recall answer set (4).

The **ACsolver algorithm**: We can view a path in the graph $\text{AC}_\Pi$ as a description of a process of search for a set of regular atoms in some answer set of $\Pi$ by applying the graph’s transition rules. Therefore, we can characterize an algorithm of a solver that utilizes the transition rules of $\text{AC}_\Pi$ by describing a strategy for choosing a path in this graph. A strategy can be based, in particular, on assigning priorities to transition rules of $\text{AC}_\Pi$, so that a solver never follows a transition due to a rule in a state if a rule with higher priority is applicable. A strategy may also include restrictions on rule’s applications.

We use this approach to describe the ACsolver algorithm (Mellarkod, Gelfond, and Zhang 2008, Fig.1). The ACsolver selects edges according to the priorities on the transition rules of the graph $\text{AC}_\Pi$ as follows:

**Backtrack, Fail >>**

**Unit Propagate, All Rules Cancelled, Backchain True >>**

**Unfounded >>**

**Query Propagate >>**

**Decide.**

Note that ACsolver also only follows a transition due to the rule Query Propagate if there are no satisfying extensions of $\text{query}(M)$ w.r.t. $\Pi_D$, i.e., $\text{Cons}(\Pi, \text{query}(M)) = \{\bot\}$.

Mellarkod et al. (2008) demonstrated the correctness of the ACsolver algorithm for the class of safe canonical programs by analyzing the properties of its pseudocode. Proposition 3 provides an alternative proof of correctness for this algorithm for a more general class of weakly-simple programs that relies on the transition system $\text{AC}_\Pi$. Furthermore, Proposition 3 encapsulates the proof of correctness for a class of algorithms that can be described using $\text{AC}_\Pi$. For instance, it immediately follows that the ACsolver algorithm modified to follow an arbitrary transition due to the rule Query Propagate is still correct.

### The CLINGCON Language

Consider a subset of the AC language, denoted $\text{AC}^-$, so that any AC program without defined atoms is an $\text{AC}^-$ program. It is easy to see that for any $\text{AC}^-$ program, its defined part is empty. The language of the constraint answer set solver CLINGCON defined in (Gebser, Ostrowski, and Schaub 2009) can be seen as a syntactic variant of the $\text{AC}^-$ language.

We now review the clingcon programs and show how they map into $\text{AC}^-$ programs. For a signature $\Sigma$, a clingcon variable is an expression of the form $p(\bar{r})$, where $p$ is a mixed predicate and $\bar{r}$ is a sequence of regular constants. For any clingcon variable $p(\bar{r})$, by $p(\bar{r})^s$ we denote its predicate symbol $p$ and by $p(\bar{r})^s$ we denote its sequence of regular constants $\bar{r}$.

We say that an atom is a clingcon atom over $\Sigma$ if it has the following form

$$v_1 \circ \cdots \circ v_k \circ c_1 \circ \cdots \circ c_m \circ v_{k+1} \circ \cdots \circ v_l \circ c_{m+1} \circ \cdots \circ c_n, \ (6)$$

where $v_i$ is a clingcon variable; $c_i$ is a constraint constant; $\circ$ is a primitive constraint operation; and $\circ$ is a primitive constraint relation.

A clingcon program is a finite set of rules of the form (1) where (i) $a_0$ is $\bot$ or a regular atom, (ii) each $a_i$, $1 \leq i \leq m$ is a regular atom or clingcon atom, and (iii) each $a_i$, $m+1 \leq i \leq n$ is a regular atom.

Any clingcon program $\Pi$ can be rewritten in $\text{AC}^-$ using a function $\nu$ that maps the set of clingcon variables occurring in $\Pi$ to the set of distinct variables over $\Sigma$. For a clingcon variable $v$, $v^\nu$ denotes a variable assigned to $v$ by $\nu$.

For each occurrence of clingcon atom (6) in some rule $r$ of $\Pi$ (i) add a set of mixed atoms $v_i^\nu (v_i^\nu, v_i^\nu)$ for $1 \leq i \leq l$ to the body of $r$, and (ii) replace (6) in $r$ by a constraint atom $v_1^\nu \circ \cdots \circ v_k^\nu \circ c_1 \circ \cdots \circ c_m \circ v_{k+1}^\nu \circ \cdots \circ v_l^\nu \circ c_{m+1} \circ \cdots \circ c_n$.

We denote resulting $\text{AC}^-$ program by $\text{ac}(\Pi)$.

For instance, let clingcon program $\Pi$ over $\Sigma_1$ consist of a single rule

$$\leftarrow \text{occurs}(a,0), \ at(0) \neq 1.$$

Given $\nu$ that maps $at(0)$ to $T$, $\text{ac}(\Pi)$ has the form

$$\leftarrow \text{occurs}(a,0), \ at(0, T), \ T \neq 1.$$

**Proposition 4** For a clingcon program $\Pi$ over signature $\Sigma$, a set $X$ is a constraint answer set of $\Pi$ according to the definition in (Gebser, Ostrowski, and Schaub 2009) iff there is a functional set $M$ of ground mixed atoms of $\Sigma$ such that $X \cup M$ is an answer set of $\text{ac}(\Pi)$.
Note that ac(Π) is a weakly-simple program (in fact, it is a simple program). It follows that a class of algorithms captured by the graph AC_1 is applicable to clingcon programs after minor syntactic transformations. Nevertheless the graph AC_1 is not suitable for describing the CLINGCON system. In the next section we present another graph suitable for this purpose.

The Basic CLINGCON Algorithm

The CLINGCON system is based on tight coupling of the answer set solver CLASP and the constraint solver GECODE. Recall that CLASP starts its computation by building a propositional formula called completion (Clark 1978) of a given program so that its propagation relies not only on the program but also on the completion. Furthermore, it implements such backtracking search techniques as backjumping, learning, forgetting, and restarts. Lierler and Truszczynski (2011) introduced the transition system SML(ASP)_{F,Π} and demonstrated how it captures the CLASP algorithm. It turns out that SML(ASP)_{F,Π} augmented with the transition rule Query Propagate is appropriate for describing CLINGCON. The graph SML(ASP)_{F,Π} extends a simpler graph SM(ASP)_{F,Π}. These extensions are essential for capturing such advanced features of CLASP and CLINGCON as conflict-driven backjumping and learning. To simplify the presentation we review the graph SM(ASP)_{F,Π} and show that augmenting it with the rule Query Propagate captures basic CLINGCON algorithm implementing a simple backtracking strategy in place of conflict-driven backjumping and learning. This abstract view on CLINGCON allows us to compare it to ACSOLVER in formal terms.

Abstract basic CLINGCON: We write Head(Π) for the set of nonempty heads of rules in a program Π. For a clause C = ¬a_1 ∨ ... ∨ ¬a_l ∨ a_{l+1} ∨ ... ∨ a_m we write C^r to denote the rule

\[ \leftarrow a_1, \ldots, a_l, \text{not } a_{l+1}, \ldots, \text{not } a_m. \]

For a set F of clauses, we define F^r = \{ C^r \mid C \in F \}. For a set A of atoms, by Π(A) we denote a program Π extended with the rules \{ a \} for each atom a ∈ A.

The transition graph SM(ASP)_{F,Π} for a set F of clauses and a regular program Π is defined as follows. The set of nodes of SM(ASP)_{F,Π} consists of the states relative to At(F ∪ Π). There are five transition rules that characterize the edges of SM(ASP)_{F,Π}. The transition rules Unit Propagate, Decide, Fail, Backtrack of the graph SM_{F^r,Π} and the transition rule

Unfounded':

\[ M \Rightarrow M \bar{\alpha} \text{ if } \left\{ \begin{array}{l} a \in U \text{ for a set } U \text{ unfounded on } M \\ \text{w.r.t. } Π(At(F ∪ Π) \setminus \text{Head}(Π)) \end{array} \right\} \]

Lierler and Truszczynski (2011) demonstrated how SM(ASP)_{ED-Comp(Π^c),Π} models basic CLASP (without conflict driven backjumping and learning) where ED-Comp(Π^c) denotes classified completion with the use of auxiliary atoms.

We now define the graph CON_{F,Π} for AC programs that extends SM(ASP)_{F,Π} in a similar way as AC_1 extends SM_1.

For an AC logic program Π and a set F of clauses, the nodes of CON_{F,Π} are the states relative to the set At(F ∪ Π^c). The edges of CON_{F,Π} are described by the transition rules of SM(ASP)_{F,Π^c} and the transition rule Query Propagate of AC_Π.

For an AC program Π, a set F of clauses is Π-safe if

1. \[ F \models \neg \alpha, \text{ for every } a \in At(Π^c) \setminus \text{Head}(Π^c), \] and
2. for every answer set X of Π^c there is a model F of Π such that X = M^+ ∩ Head(Π^c).

In fact, a set F of clauses is Π-safe if it is Π^c-safe according to the “saftens” definition given in (Lierler and Truszczynski 2011).

Proposition 5 For any weakly-simple AC program Π and a Π-safe set F of clauses,

(a) graph CON_{F,Π} is finite and acyclic,

(b) for any terminal state M of CON_{F,Π} other than ⊥, \( (M^c)^+ \cap At(Π) \) is a set of all regular atoms in some answer set of Π,

(c) state ⊥ is reachable from ⊥ in CON_{F,Π} if and only if Π has no answer sets.

The algorithm behind basic CLINGCON is modeled by means of the graph AC_{ED-Comp(Π^c),Π} with the following priorities

Backtrack, Fail >> Unit Propagate >> Unfounded >> Query Propagate >> Decide.

Proposition 3 demonstrates that the CLINGCON algorithm is applicable to a broader class of weakly-simple AC programs.

Following concept helps us to formulate the relation between AC_Π and CON_{F,Π} precisely. An edge M \Rightarrow M' in the graph AC_Π (CON_{F,Π}) is singular if:

- the only transition rule justifying this edge is Unfounded, and
- some edge M \Rightarrow M'' can be justified by a transition rule other than Unfounded or Decide.

It is easy to see that due to priorities of ACSOLVER and CLINGCON, singular edges are inessential. We define AC^-Π (CON^-_{F,Π}) as the graph obtained by removing all singular edges from AC_Π (CON_{F,Π}).

Proposition 6 Let Comp(Π^c) be completion classified in a straightforward way by applying distributivity. For every AC program Π, the graphs AC^-Π and CON^-_{Comp(Π^c),Π} are equal.

It follows that the graph CON^-_{Comp(Π^c),Π} also provides an abstract model of ACSOLVER. Hence the difference between ACSOLVER and basic CLINGCON algorithms can be stated in terms of difference in II-safe formulas Comp(Π^c) and ED-Comp(Π^c) that they are applied to.

Conclusions

In this paper, we designed transition systems AC_Π and CON_{F,Π} for describing algorithms for computing (subsets of) answer sets of AC programs. We used these graphs to specify the ACSOLVER and the basic CLINGCON algorithms. We demonstrated a formal relation between the AC and
CLINGCON languages and the algorithms behind ACSOLVER and CLINGCON. Compared with traditional pseudo-code description of algorithms, transition systems use a more uniform (i.e., graph based) language and offer more modular proofs. The graphs $\mathcal{AC}_\Pi$ and $\mathcal{CON}_F$ offer a convenient tool to describe, compare, analyze, and prove correctness for a class of algorithms. In fact we formally show the relation between the subgraphs of $\mathcal{AC}_\Pi$ and $\mathcal{CON}_F$. Furthermore, the transition systems for ACSOLVER and CLINGCON result in new algorithms for solving a larger class of $\mathcal{AC}$ programs — weakly-simple programs introduced in this paper. Neither the ACSOLVER nor CLINGCON procedures, respectively, can deal with such programs. In the future we will consider ways to use current ASP/CLP technologies to design a solver for weakly-simple programs.

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