

一个带破圈启发方法的回答集编程系统*

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An Answer Set Programming System with Cycle Breaking Heuristic

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Abstract: Answer set programming (ASP) is a logic programming paradigm under answer set semantics, which can be utilized in the field of non-monotonic reasoning and declarative problem solving, etc. This paper proposes and implements a cycle breaking heuristic and a bottom-restricted look-ahead procedure for ASP, and the resulting system is called LPS. The experimental results show that, relative to other state-of-the-art ASP systems, LPS could efficiently solve logic programs in phase transition hard-job-regions, and these programs are generally considered difficult to compute. In addition, by applying the so-called dynamic variable filtering (DVF) technique, LPS could greatly reduce the search tree size during the computation.

Key words: answer set programming; heuristic; look-ahead; logic program; phase transition

摘要: 回答集编程(answer set programming, ASP)是一种回答集语义下的逻辑编程范例,可应用于非单调推理,叙述式问题求解等领域.本文为 ASP 提出并实现了一种破圈启发方法与一种基部限制式前向搜索过程,所得到的系统称为 LPS.实验结果显示,相对于其他经典的 ASP 系统, LPS 能够有效地解决处于相变难区域中的逻辑程序,通常这些程序被认为是计算困难的.除此以外,通过使用被称为动态变元过滤(dynamic variable filtering, DVF)的技术, LPS 可以在计算过程中极大地缩小搜索树的尺寸.

关键词: 回答集编程;启发方法;前向搜索;逻辑程序;相变

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1 Introduction

Answer set programming (ASP) is a logic programming paradigm under answer set semantics (also called stable model semantics)^[1,2]. In ASP, problems are encoded as logic programs, the corresponding answer sets of the logic programs give solutions to the original problems. Many applications like non-monotonic reasoning, reasoning

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about actions, declarative problem solving^[3], can be expressed by logic programs, and then solved by an ASP system.

Unfortunately, checking whether a logic program has an answer set is NP-complete^[4]. It follows that computing answer sets for logic programs is a computational hard task. In the last decade, much work has been devoted to the implementation of ASP, most of these systems like Smodels, DLV^[5,6] are based on backtrack search algorithms, which essentially construct a binary search tree and have to handle exponential search space. For such algorithms, heuristics are crucial, good heuristics can significantly improve system performance.

In this paper we propose so-called *cycle breaking* heuristic for ASP, the idea is motivated by Ref.[7], where new upper bounds for computing answer sets are obtained through analyzing the number of *cycles* of a logic program. In particular, Ref.[7] suggests that a good heuristic should break the most number of *even* cycles. As we shall see later, our heuristic matches the purpose quite good. Another important feature of ASP systems is the use of look-ahead procedures^[5,6], the basic idea is to discover inconsistency before selecting a new branching node, and then backtrack as soon as possible, avoiding falling into a deep dead end. It has been shown that look-ahead can greatly speed up the computation, however, it often consumes most of the running time of a system, many restricted look-ahead procedures are then proposed to reduce the running cost. In this paper we present a look-ahead procedure for ASP, which restricts the atoms being looked on *bottoms*. Intuitively, bottoms are the “root” of the dependency graph of a logic program, the truth values of the atoms outside bottoms depend on those inside, therefore restricting the looking over bottoms is considered to be a good idea.

The rest of the paper is organized as follows. Section 2 provides some basic concepts of ASP. Section 3 presents the basic algorithm for computing an answer set. Sections 4 and 5 describe the cycle breaking heuristic and the *bottom-restricted* look-ahead procedure. In Section 6, the experimental results are presented and some related work is briefly discussed in Section 7, we draw conclusions and show future directions in Section 8.

2 Preliminaries

A *logic program* is a finite set of rules of the form: $a \leftarrow b_1, \dots, b_m, \text{not } c_1, \dots, \text{not } c_n$, where a, b_i 's, c_j 's are atoms. We call a the *head* of the rule and the other atoms the *body* of the rule. For convenience, let $\text{Head}(r) = a$, $\text{Pos}(r) = \{b_1, \dots, b_m\}$ and $\text{Neg}(r) = \{c_1, \dots, c_n\}$. A *literal* is an atom or an atom preceded by *not*. The former is called a positive literal and the latter is called a negative literal. If $\text{Neg}(r) = \emptyset$ then r is a *definite* rule. A *definite* logic program P is a set of definite rules, by $\text{Cn}(P)$ we refer to its *unique minimal closure*. By $\text{At}(P)$ we denote the set of atoms occurring in a logic program P . The *Gelfond-Lifschitz reduct* P^A of P with respect to a set $A \subseteq \text{At}(P)$ is obtained by deleting each rule $r \in P$ such that $\text{Neg}(r) \cap A \neq \emptyset$, and removing all negative literals in the remaining rules. Note that a Gelfond-Lifschitz reduct must be a definite program. A set of atoms $A \subseteq \text{At}(P)$ is said to be an *answer set*^[1] of P iff $A = \text{Cn}(P^A)$. The well-founded model^[9] of a logic program P is a unique ordered pair $I(P) = \langle I(P)^+, I(P)^- \rangle$ in which $I(P)^+$ contains atoms which must be *true* and $I(P)^-$ contains atoms which must be *false*. The truth values of atoms which are neither in $I(P)^+$ nor in $I(P)^-$ are *uncertain*. Well-founded model can be characterized by operator γ_P , which is defined as $\gamma_P(A) = \text{Cn}(P^A)$. Note that $A_1 \subseteq A_2$ implies $\gamma_P(A_2) \subseteq \gamma_P(A_1)$, i.e. γ_P is *anti-monotone*. It follows that $\gamma_P^2(A) = \gamma_P(\gamma_P(A))$ is *monotone*, the well-founded model of a program P can be represented as $\langle \text{lfp}(\gamma_P^2), \text{gfp}(\gamma_P^2) \rangle - \text{At}(P)$ in which $\text{lfp}(\gamma_P^2)$ is the least fixed-point of γ_P^2 and $\text{gfp}(\gamma_P^2)$ the greatest fixed-point of γ_P^2 ^[10]. Consider a logic program $P_1 = \{c \leftarrow \text{not } b, b \leftarrow \text{not } c, a \leftarrow\}$, the answer sets of it are $\{a, c\}$ and $\{a, b\}$ while its well-founded model is $\langle \{a\}, \emptyset \rangle$. A *simplification*^[11] of a logic program P under its well-founded model $I(P)$, denoted by $P \setminus I(P)$, is obtained by deleting each rule $r \in P$ such that either $\text{Head}(r) \in I(P)^+$ or $\text{Pos}(r) \cap I(P)^- \neq \emptyset$ or $\text{Neg}(r) \cap I(P)^+ \neq \emptyset$, and remove all positive literals a where $a \in I(P)^+$ and all negative literals *not* a in which $a \in I(P)^-$ from the bodies of the

remaining rules.

A *dependency graph* of a logic program P , denoted by $G(P)$, is a directed graph $\langle V, E \rangle$ with labeled edges. Let $V = At(P)$, there is a *positive* (*negative*, respectively) edge from node p to node q if there exists a rule $r \in P$, $Head(r) = q$ and $p \in Pos(r)$ ($p \in Neg(r)$, respectively). A *strongly connected component* of a directed graph $\langle V, E \rangle$ is a set of nodes $V' \subseteq V$ such that for any $u, v \in V'$, there is a path from u to v , and V' is not a proper subset of any such sets. A *bottom*^[7] is a strongly connected component S of a directed graph, and there is no other strongly connected component S' such that there is a path from S' to S . Note that a strongly connected component can be computed in linear time^[12]. An *odd* (*even*, respectively) *cycle* of a dependency graph is a simple cycle containing an odd (non-zero even, respectively) number of *negative* edges. Both odd and even cycles are called *negative cycles*, simple cycles containing no negative edges are called *positive cycles*.

Odd cycles and even cycles are of great interest in investigating logic programs. It has been shown that a logic program which has no odd cycles^[13] (called *call-consistent* program) at least has one answer set, and a logic program that has no even cycles^[14] has at most one answer set. Furthermore, in the latter case, if the well-founded model of the logic program is also its answer set, then it is the only one, otherwise, it has no answer sets^[15]. Informally speaking, odd cycles eliminate answer sets while even cycles generate answer sets. Based on these results, an algorithm for computing answer sets concerning cycles is proposed in Ref.[7], in next section we shall describe a slightly modified version.

3 Basic Algorithm for Computing Answer Sets

A *signed atom* a^* is an atom a with sign $*$ in $\{+, -\}$. Define $\neg\neg\neg a^- = a^+$ and $\neg a^+ = a^-$. Let σ be a *finite consistent set* of signed atoms, by consistency we mean there is no atom a such that both a^+ , a^- are in σ . Define $c^+ = \{b^+ | b^+ \in \sigma\}$, similarly, $c^- = \{b^- | b^- \in \sigma\}$. For a set of signed atoms \emptyset , let $|\emptyset| = \{b | b^* \in \emptyset\}$.

Definition 1^[11]. Let P be a logic program, and σ a finite consistent set of signed atoms with $|\emptyset| \subseteq At(P)$, let P_σ be the program obtained from P by deleting all rules $r \in P$ which $Pos(r) \cap |\sigma^-| \neq \emptyset$ or $Neg(r) \cap |\sigma^+| \neq \emptyset$, and removing positive literal a from the bodies of the remaining rules if $a \in |\sigma^+|$ or negative literal $not\ a$ if $a \in |\sigma^-|$.

For instance, let P be P_1 mentioned in previous section, $\sigma = \{b^+, c^-\}$, we have $P_\sigma = \{b \leftarrow, a \leftarrow\}$. Intuitively, $|\sigma^+|$ ($|\sigma^-|$, respectively) stands for atoms assumed to be true (false, respectively). Furthermore, consider the dependency graph $G(P_\sigma)$, it is a subgraph of $G(P)$ and there are no edges going out from atoms in $|\sigma|$, in some sense they break the outgoing edges, we call these atoms *breaking nodes*.

Proposition 2^[11]. Let P be a logic program, and σ a finite consistent set of signed atoms with $|\sigma| \subseteq At(P)$. Suppose S is an answer set of P , $|\sigma^+| \subseteq S$ and $|\sigma^-| \cap S = \emptyset$, then S is an answer of P_σ .

Proposition 2 implies that each answer set of P is an answer set of either $P_{\{a^+\}}$ or $P_{\{a^-\}}$ where $a \in At(P)$, however, the inverse generally does not hold. The proposition below shows an interesting property about bottoms.

Proposition 3^[7]. Given a logic program P and its well-founded model $I(P)$, if $Q = P \setminus I(P)$ is not empty, then every bottom of Q must have a pair of nodes a and b , such that there is a negative edge from a to b .

Proposition 3 shows the existence of a negative cycle in each bottom of the well-founded simplification of a program, which plays an important role in the proof of theorem 4. Having introduced the above definitions and propositions, we present our basic algorithm $Ans(P)$ in Fig.1, where $ComputeAns(P_\sigma)$ performs binary search, $Bottoms(Q)$ returns the set of all bottoms of logic program Q and $I(P_\sigma)$ returns the well-founded model of P_σ . The soundness of the algorithm is guaranteed by the theorem below:

Theorem 4^[7]. A logic program P has an answer set if and only if $Ans(P)$ returns True.

Proof: (Sketched) \Rightarrow . By induction on k , the number of odd and even cycles in $G(P)$. For $k=0$, it is well-known

P has a unique answer set $I(P)^+$. For $k > 0$, trivial if $I(P)^+$ is the answer set of P . Suppose $I(P)^+$ is not an answer set for P , pick a breaking node a from $\cup Bottoms(Q)$ on a negative cycle (such a node always exists). Note that an answer set of P is either an answer set of $P_{\{a^+\}}$ or $P_{\{a^-\}}$ and both of them have less negative cycles than k , by induction assumption, $ComputeAns(P_{\{a^+\}})$ ($ComputeAns(P_{\{a^-\}}$, respectively) returns true if $P_{\{a^+\}}$ ($P_{\{a^-\}}$, respectively) has an answer set which is also an answer set of P_\emptyset (i.e. P). Therefore if P has an answer set then $Ans(P)$ returns true. \Leftarrow . Trivial. \square

It has been shown in Ref.[7] that if a program P has at most k even cycles, then P has at most 2^k answer sets, and can be computed in $2^{2k}O(n^k)$ time, where n is the size of the given program. Moreover, restricting the number of odd cycles cannot reduce the complexity. These facts imply that $ComputeAns(P_\sigma)$ should break the most number of even cycles when selecting a breaking node, such that the two resulting logic programs have less even cycles and then easier to compute. In $ComputeAns(P_\sigma)$, we pick breaking nodes from bottoms which at least break one negative cycle, unfortunately, whether it breaks an even cycle is an NP-complete problem^[7], we shall propose a heuristic for this problem in next section.

Procedure $Ans(P)$

Input: A logic program P ;

Output: True if P has an answer set otherwise False.

1. **return** $ComputeAns(P_\emptyset)$;

Procedure $ComputeAns(P_\sigma)$

Input: A logic program P_σ ;

Output: True if P_σ has an answer set that is also an answer set of P_\emptyset otherwise False.

1. **if** $I(P_\sigma)^+$ is an answer set of P_σ **then**

2. **if** $I(P_\sigma)^+$ is an answer of P_\emptyset **then return** True **else return** False

3. **end if**

4. $Q := P_\sigma \setminus I(P_\sigma)$;

5. Pick $a \in \cup Bottoms(Q)$ on a negative cycle;

6. **if** $ComputeAns(P_{\sigma \cup \{a^+\}})$ **then return** True **else return** $ComputeAns(P_{\sigma \cup \{a^-\}})$

Fig.1 Basic algorithm for computing an answer set

4 Cycle Breaking Heuristic

From the above sections, we know that choosing a node that breaks many even cycles can significantly speed up the computation. However, it is very difficult to pick such a node since even deciding whether it breaks one even cycle is NP-complete. A compromise is to prefer breaking negative cycles, no matter which kind of cycles (odd or even) they are. In addition, observe that in a directed graph, a node that has more degrees (occurrences) would probably break more cycles, similar idea has been adopted by the famous MOMS heuristic^[16,17], which prefers picking atoms in shorter open clauses. Based on the above motivations, we propose our heuristic as follows.

Define the *length* of a rule r to be the number of *unassigned* literals in its body with respect to well-founded semantics, denoted by $L(r)$. Let r be a rule containing some unassigned atoms in program P (i.e. open rules), for each unassigned atom $a \in At(P)$ we define $w_1(a) = \sum_{head(r)=a} \alpha^{-L(r)}$, $w_2(a) = \sum_{a \in Pos(r)} \alpha^{-L(r)}$, $w_3(a) = \sum_{a \in Neg(r)} \alpha^{-L(r)}$.

These functions calculate the weights contributed by different occurrences of atom a , more precisely, they calculate the weights of a when it is a head, positive literal and negative literal respectively. The value of α is empirically set to 5, means that every 5 occurrences of an atom in a length $i+1$ rule are counted as 1 occurrence in rule which has length i , this value is obtained by a large number of experimenting, and adopted by many researchers^[16-18]. The evaluation function $w(a)$ is defined as: $w(a) = w_1(a) + w_2(a) + \beta w_3(a)$ where β is an empirically

good value used to emphasize negative edges since we prefer breaking negative cycles, this value could be modified when dealing different logic programs. In this paper we use a fixed value $\beta=1.3$. LPS selects a breaking node a such that $w(a)$ is the greatest, moreover, if $w_2(a)>w_3(a)$ then it first branches a , otherwise it first branches $not a$. Alternating branching order is employed by most ASP systems and has been proved quite useful in some instances.

This function is applied on *bottoms*, since picking a node from a bottom at least breaks one cycle (since bottoms are strongly connected components), furthermore, since there exists at least one negative cycle in each bottom of the well-founded reduction of a program, so the heuristic is expected to efficiently break negative cycles and therefore greatly reduce the depth of the search tree.

For example, let $P_2=\{a\leftarrow not b, b\leftarrow not a, c\leftarrow a, c\leftarrow b\}$, its dependency graph is shown in Fig.2, it is easy to see it has one bottom $\{a,b\}$. In the binary search if we first pick c then the truth values of a,b are still not determined, however, if we first choose an atom from the bottom $\{a,b\}$, then an answer set is immediately derived. Note that in the latter case, the only even cycle of the program is broke when we choose a or b according to our heuristic.

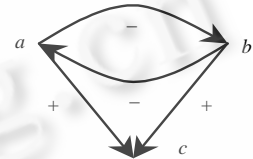


Fig.2 Dependency graph of P_2

5 Bottom-Restricted Look-Aheads

The essential of look-ahead^[5,6] is to discover a dead end as early as possible, thus avoid large scale backtracking. Simply speaking, before choosing a new node in the binary search, look-ahead procedure assumes *true* and then *false* for an *unassigned* atom and performs consistency checking, if both checkings are inconsistent, the search backtracks since there is a contradiction, else if exactly one of the checking is consistent then it picks the atom with the consistent value, otherwise the procedure looks at next unassigned atom. As mentioned before, look-ahead procedure is computational expensive, since for each unassigned atom the consistency checking is performed twice. In LPS look-ahead is applied on bottoms instead of all unassigned atoms, thus called *bottom-restricted* look-ahead. Intuitively, bottoms are the “root” of the dependency graph of a program, the truth values of the atoms outside bottoms depend on those inside. A large program may have small bottoms, therefore looking on bottoms is believed strong enough to discover many contradictions in advance.

We implemented two versions of bottom-restricted look-ahead in LPS: one adopts so-called *static variable filtering* (SVF) while another one adopts *dynamic variable filtering* (DVF). These two techniques are originally proposed in SAT^[19], where they appear to have different performances. Roughly speaking, look-ahead with SVF at most looks every atom once and then return to the binary search, however, if an atom is picked during the looking, then the current partial model is changed and new information about the contradiction may be inferred, so look-ahead with DVF looks unassigned atoms repeatedly until no contradictions could be discovered.

Bottom-restricted look-ahead with DVF is shown in Fig.3 and the SVF version could be obtained by removing the **Do-Until** loop. The function $Conflict(P, \langle |\sigma^+|, |\sigma^-| \rangle)$ in Fig.2 is from Ref.[5], it returns true when P has no answer set S such that $|\sigma^+| \subseteq S$ and $|\sigma^-| \subseteq At(P) \setminus S$, the mechanism behind it is to expand the current partial model $\langle |\sigma^+|, |\sigma^-| \rangle$ by some sophisticated inference rules and see whether $|\sigma^+| \cap |\sigma^-| \neq \emptyset$, if so then the current partial model is consistent, details about this function could be found in Ref.[5]. In Fig.4, we present the improved basic algorithm for LPS, which integrates the bottom-restricted look-ahead and the cycle breaking heuristic.

To see how look-ahead procedure speeds up the computing, consider a program $P_3=\{a_1\leftarrow not b_1, b_1\leftarrow not a_1, \dots, a_n\leftarrow not b_n, b_n\leftarrow not a_n, c\leftarrow not c\}$. P_3 has no answer sets and $n+1$ bottoms (Fig.5), without look-ahead procedure, in the worst case the algorithm may has to go through 2^n partial models before it is sure that there is no

answer sets, however, if we enable look-ahead procedure, the algorithm can return false at one step since conflict checking is performed on each bottom thus the rule $c \leftarrow \text{not } c$ is discovered to be unsatisfiable for any partial model. In this extreme example, the look-ahead procedure saves exponential time.

```

Procedure Look-ahead( $P_\sigma$ )
Input: A logic program  $P_\sigma$ ;
Output: A logic program  $P_\sigma$ , if no dead end is discovered otherwise False.
1. do
2.    $\sigma := \sigma$ ;  $Q := P_\sigma \setminus I(P_\sigma)$ ;
3.   for each  $a \in \cup \text{Bottoms}(Q)$ 
4.     if  $\text{Conflict}(P_\sigma, \langle \sigma^+ \cup \{a\}, \sigma^- \rangle)$  &&
5.        $\text{Conflict}(P_\sigma, \langle \sigma^+, \sigma^- \cup \{a\} \rangle)$  then
6.         return False;
7.     else if  $\text{Conflict}(P_\sigma, \langle \sigma^+ \cup \{a\}, \sigma^- \rangle)$  then
8.        $\sigma = \sigma \cup \{a^-\}$ ;
9.     else if  $\text{Conflict}(P_\sigma, \langle \sigma^+, \sigma^- \cup \{a\} \rangle)$  then
10.       $\sigma = \sigma \cup \{a^+\}$ ;
13.    end if
14.  end for
15. until  $\sigma = \sigma$ 
16. return  $P_\sigma$ 
    
```

Fig.3 Bottom-Restricted look-ahead with DVF

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Procedure ComputeAns( $P_\sigma$ )
Input: A logic program  $P_\sigma$ ;
Output: True if  $P_\sigma$  has an answer set that is also an answer set of  $P_\emptyset$  otherwise False.
1. if Look-ahead( $P_\sigma$ )=False then return False
2.  $P_\sigma := \text{Look-ahead}(P_\sigma)$ ;
3. if  $I(P_\sigma)^+$  is an answer set of  $P_\sigma$  then
4.   if  $I(P_\sigma)^+$  is an answer of  $P_\emptyset$  then return True else return False
5. end if
6.  $Q := P_\sigma \setminus I(P_\sigma)$ ;
7. Pick  $a \in \cup \text{Bottoms}(Q)$  such that  $w(a)$  is the greatest
8. If  $w_2(a) > w_3(a)$  then  $b := a^+$  else  $b := a^-$ 
9. if  $\text{ComputeAns}(P_{\sigma \cup \{b\}})$  then return True else return  $\text{ComputeAns}(P_{\sigma \cup \{b\}})$ 
    
```

Fig.4 Basic algorithm integrated with cycle breaking heuristic and bottom-restricted look-ahead

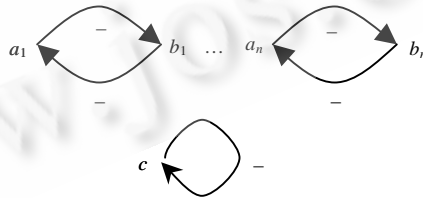


Fig.5 Dependency graph of P_3

6 Experimental Results

In this section we compare LPS with ASP systems Smodels, DLV and Nomore++^[20]. The platform is a Celeron4 1.7GHz PC with 256M memory, running Fedora 2.4.22. LPS is compiled by g++ version 3.3.2 with optimization parameter $-O3$. In all tables, times are in second, tree means the search tree size, i.e. the number of the breaking nodes. We first experiment randomly generated benchmarks^[21]. Let $k-LP(N,L)$ be the classes of programs that have N atoms and L rules with fixed length k , the authors of Ref.[21] discovered so-called hard-job-regions:

classes with $L/N=4\sim 6$, i.e. these programs are generally difficult to solve. The following experiments are done with 3- $LP(200,L)$ classes, which are considered to be moderate, i.e. they would not be too easy to distinguish the systems and too hard to compute.

The first benchmark consists of 282 programs from 11 classes where $L/N=3\sim 8$ with a step of 0.5, following the “easy-hard-easy” pattern. Results are described in Table 1, LPS (SVF) outperforms other solvers in the regions of $L/N=4.5\sim 7$ which fully cover the hard-job-regions, in easy-job-regions, Smodels is the best one. Note that LPS (DVF) has the smallest search tree size during the experiment (DLV provides no information about the search tree) and the running times are quite close to LPS (SVF).

Table 1 Experimental results on “easy-hard-easy” pattern

L/N	#Num	Smodels		LPS (SVF)		LPS (DVF)		Nomore++		DLV	
		Time	Tree	Time	Tree	Time	Tree	Time	Tree	Time	Tree
3	25	2.51	706	19.71	6181	15.58	386	26.41	387	68.85	–
3.5	24	29.54	7007	100.13	21415	88.97	1552	144.11	1725	748.37	–
4	24	115.27	22086	191.18	33332	174.74	2495	344.13	3282	1599.78	–
4.5	23	348.81	46266	237.89	31411	257.73	2476	548.36	4080	2100.10	–
5	27	257.70	35223	207.74	22820	218.33	1808	551.25	3424	1832.41	–
5.5	26	234.63	24454	187.55	16902	241.01	1333	586.01	2951	1420.56	–
6	26	144.04	13275	118.27	8878	144.07	747	415.85	1766	630.84	–
6.5	24	74.95	5825	66.42	4301	92.33	371	273.92	980	268.54	–
7	27	35.78	2548	35.74	1947	49.70	182	140.81	439	107.20	–
7.5	28	16.11	1088	23.98	1116	34.43	110	92.10	254	55.40	–
8	28	8.5	540	16.93	677	22.58	72	56.51	145	33.73	–

From the above results, it is natural to conjecture that LPS may work more efficient in hard-job-regions, so we select three hard classes with $L/N=4.5, 5$ and 5.5 , for each class we experiment 1000 programs, the top 10 hard programs (according to Smodels) are shown in Tables 2~4 respectively. As we expected, LPS completely outperforms other systems on these top hard programs, where LPS (SVF) solved 24 out of 30 programs with the minimal times, and LPS (DVF) outperforms LPS (SVF) on the remaining 6 programs, and still, LPS (DVF) possesses the smallest search tree sizes.

We analyzed some hard programs and found that they contain complicated bottoms, it follows that the cycles occurring in these components are quite complex. They seem to be the reason why these instances are so difficult to solve. In some sense, the experimental results support our conjecture that LPS performs better than other systems in hard-job-regions. Furthermore, LPS (DVF) can efficiently prune the search space, thus has the minimal search tree sizes, this is mainly because LPS (DVF) chooses most breaking nodes during look-ahead instead of binary search.

Table 2 Top 10 out of 1000 hard programs, $k=3, N=200, L/N=4.5$

Instance No.	Smodels		LPS (SVF)		LPS (DVF)		Nomore++		DLV	
	Time	Tree	Time	Tree	Time	Tree	Time	Tree	Time	Tree
095	1122.85	165489	212.30	37953	314.33	3042	937.07	7246	1625.67	–
303	798.74	162475	384.66	49971	523.74	4190	951.59	7226	>3600	–
198	778.73	129200	167.66	22098	187.73	1783	350.82	2700	1938.20	–
703	635.92	113574	129.63	17838	136.12	1430	270.96	2230	627.89	–
536	598.45	53235	394.81	47904	398.52	3683	966.46	7747	>3600	–
400	561.42	71972	392.93	59896	499.34	4507	1348.54	10671	2262.15	–
540	516.70	51230	148.94	18566	122.32	1470	280.82	2121	588.04	–
695	493.15	58198	208.11	26933	192.13	1741	289.39	2178	2359.9	–
700	426.55	64440	340.29	50340	382.31	3712	914.39	6741	2390.54	–
592	367.32	62212	175.79	21790	206.08	1790	376.80	2890	1544.44	–

Table 3 Top 10 out of 1000 hard programs, $k=3, N=200, L/N=5$

Instance No.	Smodels		LPS (SVF)		LPS (DVF)		Nomore++		DLV	
	Time	Tree	Time	Tree	Time	Tree	Time	Tree	Time	Tree
174	779.34	93711	156.29	17549	195.98	1449	801.88	4250	2107.70	–
263	774.74	101593	286.10	34440	430.23	3063	648.61	3877	2748.91	–
021	691.43	103662	164.72	19392	209.54	1599	626.59	4250	1563.71	–
574	653.86	96238	139.36	16483	128.51	1192	388.87	2461	993.45	–
953	619.80	56523	207.76	27351	205.26	2250	694.94	3976	720.61	–
444	604.15	50463	126.99	14833	160.70	1616	562.09	3771	1525.78	–
529	603.35	78678	277.26	34310	445.43	2331	693.48	4306	1559.89	–
028	587.69	70425	376.70	41145	448.02	3265	891.20	5478	1748.90	–
196	553.17	72574	100.78	12306	119.61	1066	578.18	3659	1328.63	–
076	535.12	56781	266.77	33033	354.12	2766	584.74	3619	1557.45	–

Table 4 Top 10 out 1000 hard programs, $k=3, N=200, L/N=5.5$

Instance No.	Smodels		LPS (SVF)		LPS (DVF)		Nomore++		DLV	
	Time	Tree	Time	Tree	Time	Tree	Time	Tree	Time	Tree
255	817.47	63135	351.13	40027	500.73	3060	970.41	4858	>3600	–
214	683.31	73163	143.70	13261	188.16	1135	618.79	2857	1051.08	–
396	649.78	64759	70.84	7171	86.96	871	802.62	4133	1826.96	–
567	624.98	58082	236.67	24210	311.78	1987	1275.24	7038	2091.13	–
534	546.12	61502	202.66	17859	74.31	670	214.27	1106	235.80	–
233	439.80	39295	108.60	8480	152.77	794	558.04	2857	1393.03	–
409	408.06	41100	236.06	24718	329.77	2117	875.90	4941	1939.24	–
929	400.39	50450	280.71	27233	342.84	2078	617.15	3428	1107.26	–
855	379.38	34575	189.74	16482	147.48	1343	530.13	2990	651.68	–
182	348.43	20305	124.17	13879	195.94	1190	681.01	3406	566.30	–

Table 5 presents experimental results on some real-world benchmarks. Roughly speaking, the performance of LPS is close to other systems, though it is not the most efficient one in terms of running time. This is reasonable, since ASP is NP-Complete, the cycle breaking heuristic and the bottom-restricted look-ahead would not be efficient for all benchmarks. Moreover, Smodels is highly optimized for real-world applications, it is not surprising that it is the best one in Table 5. A good result is that, LPS (DVF) still appear more efficient than other systems in terms of tree sizes.

Table 5 Experimental results on bounded model checking

Instance	#Atom	#Rule	Smodels		LPS (SVF)		LPS (DVF)		Nomore++		DLV	
			Time	Tree	Time	Tree	Time	Tree	Time	Tree	Time	Tree
DP-8	691	1112	0.16	9	0.47	26	3.82	9	0.97	12	4.42	–
DP-10	1103	1790	2.12	296	12.77	339	16.73	89	14.31	116	28.43	–
DP-12	1611	2628	347.97	103557	429.36	82456	903.77	9931	565.47	32073	677.33	–
Elavator-1	1103	1596	0.34	17	2.41	55	5.30	19	3.32	15	3.54	–
Elavator-2	3195	4465	4.58	38	26.39	72	37.12	22	21.11	35	43.75	–
Elavator-3	7824	10660	149.14	130	221.23	166	241.18	59	198.43	107	565.9	–
Elavator-4	6437	8957	1164.34	1384	1613.22	1928	1825.12	279	907.11	796	>3600	–
Hartstone-50	854	1138	2.17	142	4.45	173	5.03	77	3.92	103	5.63	–
Hartstone-75	1254	1663	10.94	192	15.77	203	23.54	99	16.68	188	34.55	–
Hartstone-100	1654	2188	26.25	242	35.34	277	38.91	66	21.11	139	58.13	–

7 Related Work

In this section we briefly discuss features of other ASP system. Smodels uses several sophisticated inference rules for characterizing answer sets, during the binary search, it chooses an atom which maximizes the current partial model and restricts look-ahead by removing some unassigned atoms if they have been derived during consistency checking. DLV is designed and optimized for disjunctive logic programs^[6,22], this may explain why DLV appears not so good during the above experiments since all programs are not disjunctive. Several heuristics are

incorporated into DLV^[18], one of these heuristics looks close to ours, however, the most important difference is that the heuristics in DLV do not concern the heads at all, while in our heuristic the heads do contribute to the evaluation function. For restricting the look-ahead, DLV defines so-called look-ahead equivalent literals, i.e. if two literals are look-ahead equivalent, then it is sufficient to only look one of them thus avoiding unnecessary looking. Nomore++^[20] is an operator-based system, it adopts several semantics operators to compute answer sets and treats heads and bodies as equitably computational objects. It follows that its look-ahead not only perform on atoms but also on bodies, Ref.[20] shows that this hybrid look-ahead may save exponentially many choices on some instances. The heuristics in Nomore++ are operator-based and runtime configurable, by combing different semantics operators with choice operator, Nomore++ could enable several heuristics during the computation. It is worth to mention some SAT-based ASP systems like^[23], which compute answer sets by SAT solvers. Generally speaking, SAT-based ASP systems are more efficient than the systems shown above, however, SAT-based ASP systems do not concern any ASP heuristics or implementation techniques at all, thus beyond the scope of this paper.

8 Conclusions and Future Directions

We present an ASP system called LPS, which employs a bottom-restricted look-ahead procedure together with an effective cycle breaking heuristic. The experimental results show that in phase transition hard-job-regions, LPS generally performs better than other ASP systems, i.e. LPS is good at solving random hard programs to some extent. Further, LPS with DVF (dynamic variable filtering) appears to be an efficient ASP system in terms of search tree size, this means DVF is theoretically a good approach for reducing search space.

It is worth to point out that random hard phenomena get more attention in recent years, they are believed to have strong connections with cryptography^[24] (hard random propositional formulas or logic programs could be considered as one-way functions). Since LPS is good at solving random hard instances, it could be applied to the area of cryptography in addition to real-world applications. Our future work will focus on improving LPS, some advanced techniques like conflict recording^[25], non-chronological backtracking^[26] will be added, and various heuristics for real-world applications will also be studied in LPS.

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