

Learning and Exploiting Progress States in Greedy Best-First Search

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Abstract

Previous work introduced the concept of *progress states*. After expanding a progress state, a greedy best-first search (GBFS) will only expand states with lower heuristic values. Current methods can identify progress states only for a single task and only *after* a solution for the task has been found. We introduce a novel approach that learns a description logic formula characterizing all progress states in a classical planning domain. Using the learned formulas in a GBFS to break ties in favor of progress states often significantly reduces the search effort.

1 Introduction

Theoretical properties of optimal state-space search algorithms like A* or IDA* have been extensively studied and are comparatively well understood [Martelli, 1977; Pearl, 1984; Dechter and Pearl, 1985; Korf *et al.*, 2001; Helmert and Röger, 2008; Holte, 2010]. A corresponding theory for suboptimal search algorithms like greedy best-first search (GBFS) [Doran and Michie, 1966] has received growing attention only in the last couple of years [Wilt and Ruml, 2014; Wilt and Ruml, 2015; Wilt and Ruml, 2016; Heusner *et al.*, 2017; Heusner *et al.*, 2018].

The main insight of Heusner *et al.* [2017] is that every run of a GBFS can be partitioned into different episodes defined by so-called *high-water mark benches*, and the *state-space topology* can be partitioned in the same way. All states s on a bench share the same *high-water mark* value, which is the largest heuristic value that needs to be considered to reach a goal state from s . *Progress states* are states that must be expanded to reach the next high-water mark bench. Exploiting knowledge of high-water mark benches or progress states during search gives rise to many applications. The only known algorithm that computes high-water mark benches does so *a posteriori*, i.e., it computes the benches of a problem, after a plan was found [Heusner *et al.*, 2018]. At this point, the high-water mark information is futile.

Inspired by Ståhlberg *et al.* [2021], who successfully learn to characterize unsolvable states using description logic [Baader *et al.*, 2003], we present an approach that learns to

characterize progress states. First, we verify for the GRIPPER and MICONIC planning domains that the set of progress states for the perfect delete-relaxed heuristic h^+ [Hoffmann and Nebel, 2001; Imai and Fukunaga, 2015] can be compactly represented with a description logic formula. Then, we present a method to learn such formulas automatically for any path-independent heuristic.

Our pipeline works as follows: for a given domain and heuristic, we fully expand the reachable state spaces of several small tasks and annotate all states with their heuristic value. Based on the heuristic values, we determine for each state whether it is a progress state. Next, we compute a set of description logic features and evaluate each of them on a subset of states. Then, we adapt a decision tree [Breiman *et al.*, 1984] learning algorithm to learn simple formulas over the description logic features in disjunctive normal form (DNF) which predict whether a state is a progress state. Finally, we show a first use case for the learned progress state classifier: we use our formulas to break ties in a greedy best-first search.

We evaluate our method using the h^+ and h^{FF} heuristics [Hoffmann and Nebel, 2001] and show that our approach successfully learns useful formulas to identify progress states. We observe a trade-off between the quality of the formulas and the time to evaluate them. Most importantly, we show that exploiting progress states is beneficial: it significantly reduces the number of expansions required to find a plan.

2 Background

We first introduce the used planning formalism and provide a short introduction to description logic. Then, we formally define progress states and explain how decision trees work.

2.1 Classical Planning

Throughout this paper, we work with STRIPS tasks [Fikes and Nilsson, 1971] defined in the *Planning Domain Description Language* (PDDL) [McDermott *et al.*, 1998]. A task Π is a tuple $\langle \mathcal{O}, \mathcal{P}, \mathcal{A}, s_I, \gamma \rangle$, where \mathcal{O} is a set of objects and \mathcal{P} is a set of first-order predicates. A predicate $p \in \mathcal{P}$ with arity k grounded to $p(o_1, o_2, \dots, o_k)$ with $o_i \in \mathcal{O}$ is called a *fact*. Let F be the set of all facts. Then any $s \subseteq F$ is called a state and the set of all states \mathcal{S} is called a *state space*. $s_I \in \mathcal{S}$ is the *initial state* and $\gamma \subseteq F$ is the *goal condition*. All states $s \supseteq \gamma$ are *goal states*. \mathcal{A} is a set of *action schemas* that can be grounded using \mathcal{O} . We call grounded action schemas *actions*.

An action a is a tuple $\langle pre, add, del \rangle$ with $pre, add, del \subseteq F$ and is associated with a cost $cost(a) \in \mathbb{R}_0^+$. Action a is applicable in state s if $pre \subseteq s$. Applying a in s , written as $s[[a]]$, leads to the successor state $(s \setminus del) \cup add$. An action sequence $\pi = \langle a_1, a_2, \dots, a_n \rangle$ is applicable in state s if every action a_i is applicable in the state $s[[a_1]][[a_2]] \dots [[a_{i-1}]]$. The cost of an action sequence is the summed-up cost of its actions. A state s' is *reachable* from s if there is an applicable action sequence starting in s and ending in s' . The *reachable state space* $\mathcal{S}_R \subseteq \mathcal{S}$ is the set of all states reachable from s_I . An applicable action sequence starting in state s and ending in a goal state is called an *s-plan*. The objective in classical planning is to find an s_I -plan, i.e., a *plan* for the given task.

2.2 Description Logic

Description logic (DL) is a family of knowledge representation formalisms [Baader *et al.*, 2003] which uses the notions of *concepts*, classes of objects that share some property, and *roles*, relations between these objects. Interpreting concepts and roles for a planning state yields a *denotation*, i.e., a set of objects $O \subseteq \mathcal{O}$ for a concept, and a set of object pairs $\{\langle o_1, o_2 \rangle, \langle o_3, o_4 \rangle, \dots\} \subseteq \mathcal{O} \times \mathcal{O}$ for a role.

Concepts and roles are recursively defined and interpreted for a state $s \in \mathcal{S}$. At its base are the *universal concept* \top and the *bottom concept* \perp with semantics

$$\top(s) = \mathcal{O} \quad \text{and} \quad \perp(s) = \emptyset,$$

as well as *primitive* concepts and roles. A primitive concept $C_{p,i}$ for a k -ary predicate $p \in \mathcal{P}$ and its i -th argument is interpreted in s as

$$C_{p,i}(s) = \{o_i \mid \exists o_1, \dots, o_k \text{ s.t. } p(o_1, \dots, o_k) \in s\}.$$

Accordingly, a primitive role $R_{p,i,j}$ for a k -ary predicate $p \in \mathcal{P}$ and its i -th and j -th arguments is interpreted as

$$R_{p,i,j}(s) = \{\langle o_i, o_j \rangle \mid \exists o_1, \dots, o_k \text{ s.t. } p(o_1, \dots, o_k) \in s\}$$

in s . Let X and X' be two concepts resp. two roles. They can be combined to form new concepts and roles via grammar rules. Examples are negation, union, and intersection which are interpreted in a state s as

$$\begin{aligned} (\neg X)(s) &= \mathcal{O} \setminus X(s) \quad \text{resp.} \quad (\neg X)(s) = \mathcal{O} \times \mathcal{O} \setminus X(s) \\ (X \sqcup X')(s) &= X(s) \cup X'(s), \text{ and} \\ (X \sqcap X')(s) &= X(s) \cap X'(s). \end{aligned}$$

We use the same grammar as Francès *et al.* [2021a]. For details, we refer to their extended paper [Francès *et al.*, 2021b].

We use two functions to convert denotations to integers. For a concept or role X , $|X(s)|$ is the *size* of the set $X(s)$. The concept distance cd between concepts C_1 and C_2 over role R is the smallest $n \in \mathbb{N}_0$ with $x_0 \in C_1(s)$, $x_n \in C_2(s)$, and all $(x_i, x_{i+1}) \in R(s)$.

The complexity $\mathcal{K}(X)$ of the universal concept, the bottom concept, any primitive concept, and any primitive role is 1. The complexity of composed concepts and roles is defined as

$$\begin{aligned} \mathcal{K}(\neg X) &:= 1 + \mathcal{K}(X) \\ \mathcal{K}(X \sqcup X') &= \mathcal{K}(X \sqcap X') := 1 + \mathcal{K}(X) + \mathcal{K}(X') \\ \mathcal{K}(|X|) &:= 1 + \mathcal{K}(X) \\ \mathcal{K}(cd(C_1, R, C_2)) &:= 1 + \mathcal{K}(C_1) + \mathcal{K}(R) + \mathcal{K}(C_2). \end{aligned}$$

2.3 Progress States

Let Π be a planning problem with a state s . A heuristic $h : \mathcal{S} \rightarrow \mathbb{R}_0^+ \cup \{\infty\}$ estimates the cost of an optimal s -plan. Let P be the set of all acyclic s -plans. The *high-water mark* of s is

$$hwm(s) = \min_{\pi \in P} \max_{a_i \in \pi} h(s[[a_1]] \dots [[a_i]]).$$

Heusner *et al.* [2018] define a state s as progress state iff its high-water mark is larger than the high-water mark of at least one of its successor states. Counter-intuitively, this definition excludes goal states for goal-aware heuristics. We therefore adapt the definition: a state s is a *progress state* iff $\gamma \subseteq s$ or $h(s) > \min_{s' \in succ(s)} hwm(s')$, where $succ(s)$ is the set of all successor states of s and γ is the goal condition.

2.4 Decision Trees

A binary decision tree is a machine learning model with a binary tree structure [Breiman *et al.*, 1984]. Let \mathcal{C} be a set of classes and let F be a list of features. A decision tree assigns a class $c \in \mathcal{C}$ to a vector $v \in \mathbb{R}^F$. Each internal tree node i is associated with a feature $f(i) \in \{1, \dots, F\}$ and a threshold $\tau(i) \in \mathbb{R}$. Each leaf node l is associated with a class $class(l) \in \mathcal{C}$. To assign a class to an input vector v , the decision tree is traversed from the root node to a leaf node. At every internal node i , if $v[f(i)] \leq \tau(i)$, then the traversal continues at the first child node, otherwise it continues at the second one. Once a leaf node l is reached, the input is labeled as $class(l)$.

Decision trees are greedily constructed given some training data $\langle D, L \rangle$ with the feature matrix $D \in \mathbb{R}^{M \times F}$ and the label vector $L \in \mathcal{C}^M$ where M is the number of training samples. Each node n is associated with a non-exclusive submatrix $D_n \in \mathbb{R}^{M' \times F}$ and $L_n \in \mathbb{R}^{M'}$. The root node is associated with the whole training data D and L and is initially a leaf node. A leaf node l is associated with the most frequent class in L_l . During training, the algorithm chooses a leaf node l and searches through combinations of features f' and thresholds τ' . f' and τ' are used to group the data points $\langle D_l[i], L_l[i] \rangle$ for $i \in \{1, \dots, M'\}$ into two sets using the test $D_l[i][f'] \leq \tau'$. The quality of the groups is evaluated using a metric (e.g., Gini impurity [Breiman *et al.*, 1984]). The leaf is associated with the combination of the best split ($f(l) = f'$ and $\tau(l) = \tau'$) and two child leaves are added to it, one per split data set. This transforms l into an internal node.

The algorithm continues until all leaves contain only labels from the same class or a maximum tree depth is reached.

3 Handcrafted Formulas

An important first step is to verify that our design space, DNF formulas over description logic features, is expressive enough to compactly characterize progress states. Therefore, we manually create description logic formulas for two standard planning domains, GRIPPER and MICONIC. In contrast to the experiment section, we only consider h^+ here, because it is easier to reason about than h^{FF} . The presented formulas have been verified empirically on all training and validation instances (see below). We slightly simplify the description logic notation here by using constant objects directly and

by defining $p_s(\cdot) = C_{p,0}(s)$ and $p_s(\cdot, o) = \{o' \mid \langle o', o \rangle \in R_{p,0,1}(s)\}$ for predicate p , state s , concept C and role R .

3.1 GRIPPER

An instance in the GRIPPER domain features a single robot with two grippers. The robot can move between two rooms *roomA* and *roomB*, and the goal is to move all balls from *roomA* to *roomB*. A GRIPPER state s is a progress state iff

$$s \models (|at_s(\cdot, roomA)| = 0) \vee \\ ((|at_robby_s(roomA)| > 0) \wedge (|free_s(\cdot)| > 0)) \vee \\ ((|at_robby_s(roomB)| > 0) \wedge (|carry_s(\cdot)| > 0)).$$

The formula describes three types of progress states: (i) there are no balls in *roomA*, (ii) the robot is in *roomA* and has a free gripper or (iii) the robot is in *roomB* and carries at least one ball. In case (i), we are either in a goal state or moving to *roomB* makes progress; in case (ii), picking up a ball makes progress (unless case (i) holds); and in case (iii), dropping a ball makes progress.

3.2 MICONIC

In MICONIC, there is an elevator that operates in a fully connected graph, bringing passengers from an origin floor to a destination floor. Let f_s be the floor for which $lift_at_s(f_s)$ holds in state s . A MICONIC state s is a progress state iff

$$s \models (|origin_s(\cdot, f_s) \cap (\overline{boarded_s(\cdot)} \cup \overline{served_s(\cdot)})| > 0) \vee \\ (|destin_s(\cdot, f_s) \cap (\overline{served_s(\cdot)} \cup \overline{boarded_s(\cdot)})| > 0) \vee \\ (|destin_s(\cdot, f_s) \cap (\overline{boarded_s(\cdot)} \cup \overline{served_s(\cdot)})| = 0).$$

The formula describes three types of progress states: (i) the first line describes states where the elevator is at a floor that is the origin of at least one passenger that is not boarded and has not been served yet; (ii) the second line describes states where the elevator is at a floor that is the destination of at least one passenger that is boarded and has not been served yet; and (iii) the last line describes states where the elevator is at a floor that is not the destination of any passenger that is still waiting (i.e., not boarded and not served). In case (i), progress is made by letting a passenger board; in case (ii), progress is made by letting a passenger leave the elevator; and in case (iii), a goal state has been reached or progress is made by moving to another floor.

4 Learning to Characterize Progress States

In this section we describe our pipeline that labels states as progress states, calculates description logic features for them, and learns a DNF formula that identifies progress states.

4.1 Generating Labeled States

To label the states of a planning instance w.r.t. a heuristic h as progress or non-progress states, we first expand the reachable state space. This includes states reachable only on paths through goal states. For each state s , we track its predecessors $pred(s)$, successors $succ(s)$, and heuristic estimate $h(s)$.

In a second iteration, we compute the high-water marks. Initially, we set $hwm(g) = h(g)$ for all goal states g and regard $hwm(s)$ as undefined for all other states s . We keep

an open list of states ordered by high-water mark values that initially contains all goal states. Upon retrieving a state s from the open list, we insert all its predecessors $p \in pred(s)$ with undefined high-water mark into the open list with a high-water mark value of $hwm(p) = \max(h(p), hwm(s))$. The algorithm guarantees that a state is only inserted once into the open list, namely after its successor with lowest high-water mark value has been retrieved from the open list. When this process terminates, only unsolvable states have an undefined high-water mark, which we treat as ∞ from here on.

In a third iteration over all states, we label all states s that are goal states or that have at least one successor state with a lower high-water mark value (then $hwm(s) > \min_{s' \in succ(s)} hwm(s')$) as progress states. All remaining states (which includes those with $hwm(s) = \infty$) are non-progress states. Note that it is possible to combine the second and third iteration into a single iteration, but considering two separate iterations simplifies the presentation.

4.2 Generating Description Logic Features

The next step of our pipeline takes as input a set of states labeled as “progress state” or “non-progress state”, and a set of grammar rules that defines which features can be generated. The feature generation is an iterative process. In iteration i , only features of complexity i are generated. A new feature is pruned if it has the same denotation as a previously generated feature for all states. In the first iteration, all primitive concepts and roles, as well as the universal and bottom concepts are generated. In every succeeding iteration, the existing features are combined using the provided rules. The feature generation stops once all features up to a given complexity are generated or a time limit is reached.

As the decision tree learns formulas over Boolean features rather than sets of objects, we convert the denotation of each state feature to a Boolean value by first converting it to an integer i (unneeded for concept distance features) and then testing if i is greater than 0. Our feature language would be more expressive if we included comparisons between integer values derived from a feature or to other constants than 0. However, we observed that it is sufficient in most domains to check if a feature generates an empty set, and this restriction speeds up the feature generation significantly. After all features are generated, we iterate over the set of states and store for each state its label and the Boolean evaluation of every feature.

4.3 Learning DNF Formulas

We use the set of labeled states annotated with Boolean features to learn a DNF formula that characterizes progress states. There is no guarantee that our generated features perfectly separate progress from non-progress states. Our grammar might not be expressive enough or we might need a feature with a complexity that is infeasible to generate. Thus, we need a learning algorithm which allows imperfect separations. We decided to train decision trees.

Since our training data is already in an appropriate format (every state is represented by an equally-sized Boolean vector), we can apply a standard algorithm to obtain a decision tree. We do not limit the depth of the generated tree,

Domain	Max. Complexity					# Features	
	1.	2.	3.	4.	5.	Min	Max
BARMAN	7	6	6	6	6	1218	2455
BLOCKSWORLD	10	10	10	10	10	15332	16611
CHILDSNACK	8	8	7	7	7	444	532
DRIVERLOG	9	8	8	7	7	892	1313
FLOORTILE	8	8	7	7	7	1644	3441
GRIPPER	12	12	9	9	9	422	1656
MICONIC	8	8	7	7	7	332	494
VISITALL	12	12	11	11	11	1692	2118

Table 1: Fives times for each domain, we generate description logic features using the progress state labels for h^{FF} . For the i -th repetition, we use states from i state spaces. Left: for each feature set the maximum complexity over its features. Right: the minimum and maximum number of features in the feature sets of a domain.

so if it possible to separate the training data with the generated features, then the tree will separate them. After training the decision tree, we convert it to a DNF: we collect all paths in the tree which classify a state as progress state (e.g., $\langle \text{feature1} > 0 = \top, \text{feature2} > 0 = \perp \rangle$). Each path becomes a clause in the DNF and each decision in the path becomes a literal. The path in the previous example implies a clause with the literals “feature1 > 0” and “feature2 = 0”.

The automatically generated DNFs can be long and may contain redundancy, which makes them slow to evaluate and hard to interpret. We simplify the DNFs with SymPy [Meurer *et al.*, 2017]. If several features separate the data equally well, we break ties first in favor of already used features and second in favor of lower complexity feature. This leads to a significantly lower number of distinct features in a formula which speeds up its evaluation and increases the probability that SymPy finds a simpler, logically equivalent formula.

5 Experiments

We explained above how we can use description logic to generate state features that generalize across all instances of a planning domain and how we can learn DNF formulas to predict whether a state is a progress or non-progress state. Now, we evaluate our pipeline on common planning domains and show that it learns useful formulas. To verify that the formulas learned important information, we use it as tie-breaker in a greedy best-first search (GBFS).

5.1 Setup

We evaluate our approach on the BARMAN, BLOCKSWORLD, CHILDSNACK, DRIVERLOG, FLOORTILE, GRIPPER, MICONIC, and VISITALL domains for the FF heuristic (h^{FF}) and the perfect delete-relaxed heuristic (h^+) [Hoffmann and Nebel, 2001], using the formalization based on operator counting by Imai and Fukunaga [2015] for the latter. For each domain, we define a parameter space (e.g., a range for the number of balls in Gripper) and use it to generate small training and validation instances with PDDL task generators [Seipp *et al.*, 2022]. As our test sets, i.e., for the GBFS runs with and without learned formulas, we use the union of Autoscale 21.11 tasks for optimal and satisficing planning [Tor-

ralba *et al.*, 2021]. The test sets are disjoint from and generally more challenging than our training and validation tasks. For each training and validation instance and each heuristic, we generate the labeled state space with a memory limit of 3.5 GiB and a time limit of 5 hours.

To generate the description logic features, we use the same concepts and roles as Francès *et al.* [2021a]. We use the *size* and the *concept distance* functions to convert the description logic features to integers. We observed that generating the features for a single state space can lead to features that are specific to that state space and to formulas that do not generalize to unseen tasks. Thus, we generate and evaluate 5 feature sets per domain, where the i -th feature set uses states from the first i training instances from that domain. We also observed that the feature generation often exceeds the available memory. As a solution, we select from each training instance a subset of up to 10 000 progress and up to 10 000 non-progress states. We impose no limit on the complexity of the generated features but instead limit the feature generation procedure to 24 hours and 3.5 GiB of memory.

The last step of our pipeline produces the DNF formulas. For each domain, we train a decision tree using each of the 5 generated feature sets. In some domains, one class (progress or non-progress states) heavily outnumbers the other class. Thus, we weight all samples such that both classes have the same impact on the final formula. Furthermore, since some training instances have state spaces of significantly different sizes, we additionally weight all samples to enforce that each instance has the same impact on the final formula. We evaluate all resulting formulas on the validation instances of their domain. Finally, for each domain we select the formula with the highest *F1* score on the validation instances. In case of a tie, the formula with the better *F1* score on the training instances and then the one trained on fewer state spaces is preferred. As the decision tree is a greedy learning algorithm, training is usually quick. We limit training of the DNF formulas to 10 minutes and their validation to 2 hours, both with a memory limit of 3.5 GiB.

Learning to identify progress states is an important contribution on its own. There are many ways in which a search can be improved if it knows which states are progress states or not. We will show that even imperfect formulas contain sufficient information to improve the search. We execute GBFS with h^{FF} and h^+ as baselines and then compare it to GBFS with the same heuristic that uses the progress state information to break ties if two states have the same heuristic value. For GRIPPER and MICONIC, we described perfect, hand-crafted formulas above. To assess the quality of the learned formulas, we also compare them to their perfect counterparts.

We implemented the state space labeling and the final search in Fast Downward [Helmert, 2006]. To generate and evaluate the description logic features, we use the DLPlan library [Drexler *et al.*, 2022], and we adapt scikit-learn [Pedregosa *et al.*, 2011] to train the decision trees. To run our experiments, we use Downward Lab [Seipp *et al.*, 2017]. All steps are executed on a single core of an Intel Xeon Silver CPU. Our benchmarks, code, experiment data, and supplemental results are available online [Ferber *et al.*, 2022].

Domain	Mean Validation F1					First DNF			Best DNF		
	1.	2.	3.	4.	5.	$\mathcal{K}(F)$	Clauses	Literals	$\mathcal{K}(F)$	Clauses	Literals
BARMAN	79	80	81	79	77	7	316	4059	6	1459	25189
BLOCKSWORLD	83	–	–	–	–	10	72	1009	10	72	1009
CHILDSNACK	70	65	73	81	72	8	28	240	7	116	1243
DRIVERLOG	81	81	89	85	85	9	95	829	8	1087	17270
FLOORTILE	75	85	85	85	88	8	10	91	7	65	926
GRIPPER	96	96	96	98	98	11	3	3	9	2	4
MICONIC	98	98	97	98	99	8	1	1	7	23	274
VISITALL	70	70	73	73	73	12	2413	42496	11	3744	74060

Table 2: Left: the mean F1 score (in %) on the validation data for each formula trained for h^{FF} . The i -th formula is trained on i state spaces. Middle: for the DNF trained on the first state space, the maximum complexity in its features, its number of clauses, and its number of literals. Right: for the DNF with the highest F1 score, the maximum complexity in its features, its number of clauses, and its number of literals.

5.2 Results

We start our analysis by inspecting the output of the feature generation step. Table 1 shows for each domain the maximum complexity of features generated when using states from only one or up to five state spaces and the progress state labels for h^{FF} . The results for h^+ qualitatively look the same, so we only report them in the online supplement. We clearly see that using more state spaces causes the feature set to have lower complexity features. This is because more states cause more computational effort per feature. But more importantly, a more diverse training set helps to differentiate similar concepts and roles. As a result, fewer features are pruned and more features are created per complexity level. Thus, there is a trade-off between identifying the meaning of a feature more precisely and generating more complex features. Table 1 also shows that the number of generated features varies significantly between different domains. For example, we generate more than 16 000 features in BLOCKSWORLD but fewer than 500 in MICONIC.

To compare different choices for the complexity trade-off, Table 2 shows the validation performance for DNF formulas trained on up to five state spaces. We see that training on a single state space already produces formulas with a good F1 score. Using more state spaces to generate the formula improves the performance only moderately. For BLOCKSWORLD, where a huge amount of features are generated (see Table 1), using more than one state space exhausts the memory.

Table 2 also shows that all DNFs (almost) always contain at least one of the most complex features available. This suggests that features of high complexity contain essential information. Comparing the number of clauses and literals between the first and best DNFs, it is apparent that the DNF trained on only a single state space requires significantly fewer clauses and literals — there are only two domains where formulas with more than 100 clauses are learned, and only one of them has more than 1000 clauses. On the other hand, the best DNF has more than 100 clauses in four domains, and three of those contain even more than 1000 clauses. A large DNF takes more time to evaluate and hence slows down search. Identifying the perfect trade-off between accuracy and evaluation speed remains future work.

Domain	h^+	h_{LOpt}^+	h_{L*}^+
GRIPPER (17)	137.68	57.03	57.03
MICONIC (14)	82.17	51.07	53.15

Table 3: Geometric mean of the expansions on the commonly solved instances for GRIPPER and MICONIC using the perfect delete-relaxed heuristic without formulas (h^+), with perfect formulas (h_{LOpt}^+), and with learned formulas (h_{L*}^+) in GBFS. The values in brackets show the number of commonly solved instances.

Below, we focus on the DNF formulas with the best F1 score, but results for other formulas can be found in the supplement. To indicate that h^{FF} is enhanced by the best formula we write h_{L*}^{FF} and for h^+ we write h_{L*}^+ .

To show that it is indeed beneficial to characterize progress states, we use them to break ties in GBFS. Table 3 compares a standard GBFS with h^+ against a GBFS with h^+ which is enhanced by perfect progress state formulas. Using the perfect progress state formulas significantly reduces the required number of expansions, which indicates that it pays off to expand progress states as early as possible. We also see that the learned formula for GRIPPER is as good as the perfect formula. For MICONIC, the learned formula is slightly worse than the perfect one but still significantly better than performing a search without progress state information.

Finally, we evaluate the learned formulas for both h^+ and h^{FF} on all domains. Figure 1 compares the expansions required by GBFS with h^{FF} (left) and h^+ (right) to GBFS with the same heuristic and tie-breaking with the progress state information. Both plots show that exploiting the learned DNFs requires significantly fewer expansions in most planning instances. Plain h^{FF} requires fewer expansions than our enhanced h_{L*}^{FF} in only 43 tasks, whereas the enhanced h_{L*}^{FF} requires fewer expansions in 215 tasks. The plain h^+ heuristic requires fewer expansions in 7 tasks, whereas our enhanced h_{L*}^+ requires fewer expansions in 51 tasks. This shows that our learned formulas successfully identify progress states and that exploiting progress states is a useful enhancement for a heuristic that guides GBFS.

Evaluating the progress state formulas in every state can be computationally expensive. Table 4 shows the geometric mean over the runtime on the commonly solved instances

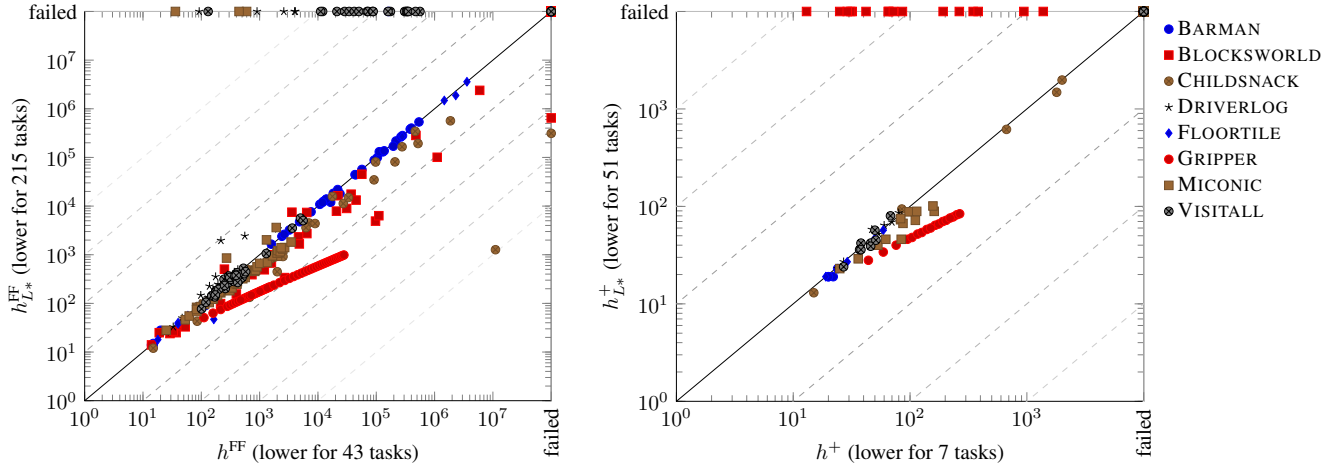


Figure 1: Scatter plot of the required expansions for the commonly solved instances of (Left) h^{FF} and h^{FF} enhanced with our progress state formulas (h_{L*}^{FF}), and (Right) h^+ and h^+ enhanced with our progress state formulas (h_{L*}^+).

Domain	h^{FF}	h_{L*}^{FF}	h^+	h_{L*}^+
BARMAN	3.4	19.1	161	154
BLOCKSWORLD	0.3	0.5	–	–
CHILDSNACK	3.5	6.8	88	87
DRIVERLOG	0.1	3.2	244	252
FLOORTILE	0.4	0.9	11	11
GRIPPER	1.1	0.6	301	266
MICONIC	0.2	1.5	103	95
VISITALL	0.0	10.7	305	303

Table 4: Geometric mean over the runtime on the commonly solved instances for the h^{FF} and h^+ heuristics, respectively.

for our proof of concept implementation. For an expensive heuristic like h^+ , using the formulas does not only reduce the number of expansions, but also the runtime. However, in combination with a fast-to-evaluate heuristic like h^{FF} , the time to evaluate a formula offsets the time saved due to the reduced number of expansions. Finding a sweet spot that optimizes the trade-off between formula complexity and accuracy will be the focus of future work.

6 Conclusions

Heusner *et al.* [2017] showed that knowing which states in a GBFS are progress states bears great potential for improved algorithms. So far, however, this knowledge could only be computed a posteriori, making the information futile. We presented the first algorithm which learns a generalized model to identify progress states. The validation experiments verified that the formulas learned by our approach generalize across instances and that they are able to separate progress states from non-progress states with high accuracy. We furthermore showed that exploiting the learned information significantly decreases search effort, even with a simple baseline method that breaks ties in favor of progress states.

We believe this is only the first step in this direction of research. We observed that the selection of training instances

and states from these instances has a huge impact on the quality of the learned formula, and so far we have not optimized the selection. It is furthermore unclear which feature complexity and which rules are necessary for a given domain. Even more importantly, our approach ignores the fact that a less accurate but faster-to-evaluate formula can be preferable in a search. Adapting the learning pipeline to produce formulas for specific use cases will be challenging future work.

Further ideas for future work include identifying not only progress states but also *crater states* and more sophisticated methods to exploit the learned knowledge, including methods that clear the open list when a progress state is encountered and search algorithms that actively avoid craters and search towards progress states.

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