

Learning Generalised Policies for Numeric Planning

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Abstract

We extend Action Schema Networks (ASNs) to learn generalised policies for numeric planning, which features quantitative numeric state variables, preconditions and effects. We propose a neural network architecture that can reason about the numeric variables both directly and in context of other variables. We also develop a dynamic exploration algorithm for more efficient training, by better balancing the exploration versus learning tradeoff to account for the greater computational demand of numeric teacher planners. Experimentally, we find that the learned generalised policies are capable of outperforming traditional numeric planners on some domains, and the dynamic exploration algorithm to be on average much faster at learning effective generalised policies than the original ASNs training algorithm.

Introduction

Generalised planning is broadly concerned with the representation, synthesis, and learning of plans, policies, heuristics, and other forms of control knowledge applicable to *many* problem instances (Srivastava, Immerman, and Zilberstein 2011; Hu and Giacomo 2011; Celorrio, Aguas, and Jonsson 2019). Interest in generalised planning has steadily increased in recent years, fueled in part by advances in machine learning, and by the development of new formalisms to represent and reason about generalised planning tasks and their solutions (Toyer et al. 2018; Francès et al. 2019; Garg, Bajpai, and Mausam 2019; Bonet and Geffner 2020; Aguas, Jiménez, and Jonsson 2020; Toyer et al. 2020; Shen, Trevizan, and Thiébaux 2020; Karia and Srivastava 2021; Aguas, Jiménez, and Jonsson 2021; Ståhlberg, Bonet, and Geffner 2022; Lin et al. 2022).

An important limit of existing work on generalised planning is that it only allows for primitive forms of quantitative information to be modelled,¹ even though such information is core to many real world problems – for example, modelling a delivery robot requires modelling how much weight it can hold, and modelling flights require reasoning about the the product of distance travelled and fuel consumption per unit of distance. This is despite the existence of the vibrant

field of *Numeric planning*, which extends classical planning formalisms to allow modelling numeric fluents, conditions and effects (Fox and Long 2003), and typically handles them using new heuristic search, optimisation, or satisfiability modulo theory based techniques (Hoffmann 2003; Coles et al. 2013; Scala et al. 2016a,b, 2020; Kuroiwa et al. 2022; Leofante 2023).

In this paper, we extend a state of the art generalised planning approach, namely *Action Schema Networks* (ASNs) (Toyer et al. 2018, 2020) to handle numeric planning problems described in PDDL2.1 (Fox and Long 2003). ASNs is a recent deep learning architecture capable of representing policies for generalised planning, and designed to learn from smaller planning tasks and then apply that knowledge to tackle larger, more complex challenges within the same domain. The architecture exploits the relational structure of planning problems and domains, and its connectivity reflects the precondition-effect relationships captured in the domain’s action schemas. This scheme makes it possible to share weights between policy networks instantiated for different problems in a domain, and learn a single set of parameters which can be transferred to problems of arbitrary size in that domain. ASNs are trained using an imitation learning algorithm, which iteratively explores the state space of the training problems using a teacher planner. Experiments with classical and probabilistic domains have shown that ASNs can outperform conventional planners when the domain has simple tricks that are key to solving larger problems but which can be learned from small problems.

To extend ASNs to numeric planning, We first propose a network module that enables ASNs to directly reason about numeric fluents. Then, we illustrate why such reasoning is not always sufficient for learning effective generalised policies. We argue that reasoning about the interaction between individual fluents is crucial, and allow ASNs to perform this interaction reasoning through numeric comparisons in the problem. To cope with the increased length of numeric plans and run-time of numeric planning teachers in comparison with their classical planning counterparts, we also propose a new training algorithm which offers greater control over the exploration versus learning balance.

Finally, we evaluate our proposed techniques on a representative set of benchmarks from the latest International Planning Competition, featuring both simple and linear nu-

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¹We are only aware of a single exception (Lin et al. 2022).

meric planning domains. Our results show that our extensions to ASNNets allow it to learn generalised policies capable of solving problem instances significantly more complex than those seen in training, and outperform non-learning planners in several domains. We also find that the greater control offered by our new training algorithm allows ASNNets to be trained much more quickly without compromising the effectiveness of the learned generalised policies.

Numeric Planning

As in PDDL2.1 (Fox and Long 2003), a numeric planning problem, denoted as $P = \langle D, I \rangle$, consists of a *domain* D and an *instance* I . The domain includes predicates \mathcal{P} , functions \mathcal{F} , and action schemas \mathcal{A} ; the instance I comprises objects O and additional elements. Each predicate $p \in \mathcal{P}$ and function $f \in \mathcal{F}$ applies to object arguments o_1, \dots, o_n from O to form ground proposition $p(o_1, \dots, o_n)$ and fluents $f(o_1, \dots, o_n)$ respectively.² Through grounding, the sets \mathcal{P}, \mathcal{F} , and O define the set of all possible propositions P and fluents F , which encode a state space S where a state is an assignment of boolean values to each proposition and real values to each fluent.

A *comparison schema* has the form $\xi \triangleright \gamma$, where $\triangleright \in \{<, =, >, \geq\}$, ξ is an arithmetic expression over \mathcal{F} , and γ is a real constant. Once grounded, ξ becomes an arithmetic expression over F , and the ground *comparison* is a mapping from S to a truth value. Each action schema $\alpha \in \mathcal{A}$ has a precondition $\text{pre}(\alpha)$ that is a conjunction of comparison schemas and predicates. The effect of α , $\text{eff}(\alpha)$, is a schema to assign boolean values to propositions and/or increase/decrease/assign the value of arithmetic expressions over \mathcal{F} . Given objects O , action schemas in \mathcal{A} ground to a set of actions A . The problem P is *linear* if all arithmetic expressions are linear, and *simple* if furthermore the numeric action effects only involve increasing or decreasing fluents by a constant.

The instance I , in full, is a tuple $I = \langle O, s_0, G, M \rangle$. The initial state s_0 is any state in S , the goal G is a conjunction of comparisons and propositions, and the plan metric M is an optional arithmetic expression over F . An action is applicable in a state when its precondition is satisfied, and its application yields a new state according to its effect. Propositions and fluents not included in the effect remain unchanged. A plan is a sequence of actions. It is an *executable* plan if when iteratively applied in s_0 , each subsequent state satisfies the next action precondition, and it is a *goal achieving* plan if the final state satisfies G . A *valid* plan is an executable goal achieving plan. The cost of a valid plan is the value of M at the final state if M is provided, or otherwise the number of actions in the plan. For this paper, a *generalised policy* for a domain D is a mapping from instances to executable (but not necessarily goal achieving) plans. The *effectiveness* of a generalised policy over a finite set of instances measures the proportion of instances the policy maps to goal-achieving plans. The more effective a generalised policy is, the larger this number is.

²The value of n , or *arity*, is dependent on the particular predicate or function.

Numeric Action Schema Networks

Action Schema Networks are a state of the art approach for learning generalised policies for classical planning problems (Toyer et al. 2020). Core to its effectiveness is its ability to generalise from a small set of training problems to much larger and unseen problems in the same domain, thereby amortising training time. This original approach is unable to perform numeric reasoning effectively due to a lack of architectural components dedicated to numeric reasoning. In this section we propose *Numeric Action Schema Networks* (ν -ASNNets) for learning generalised policies for numeric planning.

A ν -ASNNet is a neural network with weights θ that takes in input vectors describing the current state s and outputs a probability distribution $\pi^\theta(a|s)$ over all applicable actions a . For each problem instance, a ν -ASNNet is constructed with the weights θ shared between all instances in the same domain. That is, the network architecture is *instance-dependent*, but the weights are *instance-agnostic* through a *weight-sharing* mechanism that we describe later.

For each instance, the ν -ASNNet architecture includes layers of network modules that alternate between encoding action and state information, as shown in Figure 1. Each *action layer* contains an *action module* for each action $a \in A$. The last layer of a ν -ASNNet is always an action layer whose outputs determine π^θ , and we inherit the assumption from ASNNets that the first layer is also always an action layer. Each *state layer* contains one *state module* for each piece of state information, namely propositions, fluents, and comparisons. Each network fixes a hidden dimension d_h , and networks modules propagate forward a hidden representation vector in \mathbb{R}^{d_h} to connected modules in the next layer. Network modules are connected sparsely to modules in adjacent layers through a notion of relatedness.

Definition 1 (relatedness) An action a is related to a proposition p , fluent f , or comparison c at position k , denoted by $R(a, p/f/c, k)$, if $p/f/c$ is a ground instance of the k th unique predicate/function/comparison schema appearing in the action schema of a , respectively.

Example 1 Consider the following action schema for a robotic arm picking up an object of a given weight, subject to a limit on the total load carried by the arm:

```
pickup( $r, o$ ):
  prec : clear( $o$ ), weight( $o$ ) + load( $r$ )  $\leq$  limit( $r$ )
  eff :  $\neg$ clear( $o$ ), holding( $r, o$ ), load( $r$ ) += weight( $o$ )
```

The action $\text{pickup}(r_1, o_1)$ for a particular robot arm r_1 and object o_1 is related at position 1 to the proposition $\text{clear}(o_1)$, at position 2 to the proposition $\text{holding}(r_1, o_1)$, at position 1 to the fluent $\text{weight}(o_1)$, at position 2 to the fluent $\text{load}(r_1)$, at position 3 to the fluent $\text{limit}(r_1)$, and at position 1 to the comparison $\text{weight}(o_1) + \text{load}(r_1) \leq \text{limit}(r_1)$.

The notion of relatedness extends naturally to the various network modules introduced below.

Action modules. The action module for $a \in A$ in the l th action layer takes an input vector u_a^l and produces a hidden representation

$$\phi_a^l = \sigma(W_a^l \cdot u_a^l + b_a^l)$$

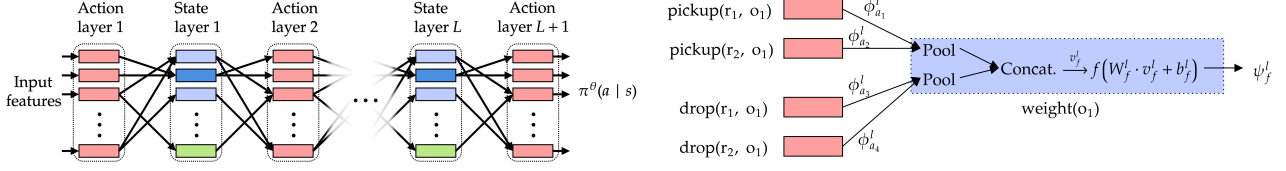


Figure 1: (Left) Overview of an ν -ASNet with L state layers and $L + 1$ action layers, with colours in state layer indicating the different types of modules. (Right) example fluent module for the fluent $\text{weight}(o_1)$ in a problem instance with two robot arms r_1 and r_2 from a domain where $\text{weight}(o)$ occurs in two action schemas $\text{pickup}(r, o)$ and $\text{drop}(r, o)$.

where $W_a^l \in \mathbb{R}^{d_h \times d_a^l}$ and $b_a^l \in \mathbb{R}^{d_h}$ are the learnt weight matrix and bias vector for the action module, σ is a non-linearity, d_h is the fixed hidden representation size, and d_a^l is the dimension of u_a^l . The input vector is constructed by concatenating the hidden representation of all related state modules in the previous layer, ignoring relatedness position

$$u_a^l = \left[\psi_1^{l-1T} \dots \psi_M^{l-1T} \right]^T$$

where ψ_j^{l-1} is the hidden representation produced by a related state module in the preceding state layer. Since all ψ_j^{l-1} have dimension d_h , u_a^l has dimension $M \cdot d_h$.

The related state modules of an action $a \in A$ can be determined by enumerating all the predicates, functions, and comparison schemas in its action schema grounding them using the same objects used to ground a . If we impose an ordering on these constructs (e.g. using position), the structure and dimension of u_a^l across all actions with the same schema is fixed even across different problem instances of the domain. Such structure is the key to weight-sharing. For ν -ASNets of different instances of a domain, all actions modules with action schema α at layer l share the weight matrix W_α^l and bias vector b_α^l . This allows us to apply the same set of weights to any instance in a domain, as all actions in these problem instances are grounded from the same set of action schemas.

Fluent modules. In each state layer, there is one fluent module for each fluent in the problem. Fluent modules allow the network to reason directly about the quantitative components of the state space. Like action modules, a fluent module for fluent $f \in F$ in the l th state layer computes a hidden representation

$$\psi_f^l = \sigma(W_f^l \cdot v_f^l + b_f^l)$$

where $W_f^l \in \mathbb{R}^{d_h \times d_f^l}$ and $b_f^l \in \mathbb{R}^{d_h}$ are the learned weight matrix and bias vector, σ is the same nonlinearity as before, v_f^l is the input feature vector, and d_f^l is the dimension of v_f^l .

Like action modules, weight-sharing requires that the input vectors v_f^l have a similar structure for fluents derived from the same function. Unlike action modules, the number of actions related to a fluent is not instance-agnostic, so simple concatenation of hidden representations of related action modules from the preceding layer is not sufficient. We treat this similarly to how proposition modules are constructed

in the original ASNets. From all the actions related to f at various positions, we extract their action schemas and enumerate all unique pairs $\{(\alpha_1, k_1), \dots, (\alpha_S, k_S)\}$ of action schemas and position pairs. These pairs are only dependent on the function of f . We then construct the input feature by

$$v_f^l = \begin{bmatrix} \text{pool}(\{\phi_\alpha^l \mid \text{op}(a) = \alpha_1 \wedge R(a, f, k_1)\}) \\ \vdots \\ \text{pool}(\{\phi_\alpha^l \mid \text{op}(a) = \alpha_S \wedge R(a, f, k_S)\}) \end{bmatrix}$$

where $\text{op}(a)$ is the action schema of a and pool is a pooling function to combine multiple \mathbb{R}^{d_h} vectors into one \mathbb{R}^{d_h} vector. Like the original ASNets we use the element-wise max function for pool . The structure of the resulting input feature is only dependent on domain information, namely the action schemas and functions, and hence enables weight-sharing – all fluent modules across different ν -ASNets at the same layer l with the same function share the same weight matrix and bias vector.

Proposition modules. Like the original ASNets, in each state layer we include a proposition module for each proposition in the problem. Proposition modules are almost identical to fluent modules, with the same computation for hidden representation, construction of input feature, and weight-sharing property.

Comparison modules. By including fluent values in the network and fluent values in the network input, ν -ASNets are able to learn generalised policies that reason directly on the value of each fluent. Such reasoning is unfortunately not always sufficient. Consider a domain D whose sole numeric component involves robotic arms with load limits lifting up items of varying weight, and suppose two problem instances I and I' differ only in that all the load limit and item weights in I' are double their counterparts in I . There is no practical difference between the problem $P = \langle D, I \rangle$ and $P' = \langle D, I' \rangle$, and an ideal generalised policy should produce the same plan for both problems. Fluent modules are unable to recognise this “symmetry” between P and P' – weights learned by training on P would not apply directly to P' .

More generally, the value of fluents are often only meaningful in the context of other fluent values, and it is valuable to allow learned generalised policies to reason about the interaction between fluents. In particular, fluents interact in comparisons in action preconditions, which we capture

through *comparison modules*. Let $\text{comp}(a)$ denote the comparisons in action a and $C = \bigcup_{a \in A} \text{comp}(a)$, in each state layer there is one comparison module for each comparison. A comparison module for the comparison $c \in C$ in the l th state layer computes a hidden representation

$$\psi_c^l = \sigma(W_c^l \cdot v_c^l + b_c^l)$$

where $W_c^l \in \mathbb{R}^{d_h \times d_c^l}$ and $b_c^l \in \mathbb{R}^{d_h}$ are the learned weight matrix and bias vector, σ is the same nonlinearity as before, v_c^l is the input feature vector, and d_c^l is the dimension of v_c^l .

Similarly to fluent modules, we employ a pooling mechanism to enable weight-sharing between comparisons that share the same schema. From all actions related to c at various positions, we extract their action schemas and enumerate all unique pairs $\{(\alpha_1, k_1), \dots, (\alpha_S, k_S)\}$ of action schema and position pairs. Again, these pairs only depend on the comparison schema of c , and we construct the input feature using them by

$$v_c^l = \begin{bmatrix} \text{pool}(\{\phi_a^l \mid \text{op}(a) = \alpha_1 \wedge R(a, c, k_1)\}) \\ \vdots \\ \text{pool}(\{\phi_a^l \mid \text{op}(a) = \alpha_S \wedge R(a, c, k_S)\}) \end{bmatrix}$$

where pool is the same pooling function as before. The resulting input feature v_c^l is again only dependent on domain information, and hence enables weight-sharing where all comparison modules across different ν -ASNs at the same layer l with the same comparison schema share the same weight matrix and bias vector.

Input. The first and last action layers take minor exceptions to the above as they are the input and output layers of the network. For the first layer, there is no preceding state layer and the input vector u_a^1 is a vector encoding state and heuristic information relevant to the current action. Specifically,

$$u_a^1 = [v_p^T \ v_f^T \ v_c^T \ g_p^T \ g_f^T \ m \ c_a \ c_{\mathcal{L}}^T]^T$$

where v_p, v_f, v_c are the values of the related propositions, fluents, and comparisons of the action in the current state respectively, g_p and g_f indicate if the related propositions and fluents appear in the goal or not, m is a boolean value indicating if the action a is applicable in the current state, c_a is the number of times a has been applied so far, and $c_{\mathcal{L}}$ is a boolean vector encoding landmark information. For v_c , we treat the value of a comparison as the boolean value indicating if it is satisfied.

The lack of goal input g_c for comparisons is a consequence of the lack of overlap between comparisons in action preconditions and comparisons in the goal. We also cannot include goal comparisons directly in the input as there is no notion of relatedness between them and actions. In domains where all instances of interest have goals with the same structure, one can define a “reach” action whose precondition is the original goal and effect is a proposition “goal-reached” which also replaces the goal. This special action would allow ν -ASNs to reason about goal comparisons.

The inclusion of c_a and $c_{\mathcal{L}}$ is to compensate for the *receptive field problem* discussed in the original paper (Toyer

et al. 2020). Essentially, longest chain of related action and state modules the network can reason about is limited by its fixed and finite depth. The inclusion of heuristic information can effectively address this problem. The action count c_a helps the network apply different actions at the same state depending on what actions have been applied before and avoid cycling between adjacent states, and is found to be empirically beneficial (Toyer et al. 2020). The numeric landmark encoding $c_{\mathcal{L}}$ in ν -ASNs is an extension of the LM-cut landmarks from the original ASNs, and is derived from hybrid landmarks extracted from an AND/OR graph structure (Scala et al. 2017). Each such landmark ℓ has a target t_{ℓ} along with a set of actions A_{ℓ} and contributions for each action $\{\lambda_{\ell}^a \mid a \in A_{\ell}\}$, and represents the inequality

$$\sum_{a \in A_{\ell}} \lambda_{\ell}^a y_a \geq t_{\ell}$$

where y_a is the number of times action a is applied from the current state. Given a set of hybrid landmarks, the resulting landmark encoding $c_{\mathcal{L}}$ is a vector in $\{0, 1\}^3$, where $c_{\mathcal{L}}^{(1)} = 1$ if the action a appears as the only action in A_{ℓ} for any landmark, $c_{\mathcal{L}}^{(2)} = 1$ if the action a appears in any A_{ℓ} with other actions, and $c_{\mathcal{L}}^{(3)} = 1$ if a does not appear in any A_{ℓ} .

We have also experimented with other encodings of hybrid landmarks, specifically encodings that take into account the contribution and target of landmarks. We additionally experimented with removing all numeric components of the numeric problem and encoding the LM-cut landmarks of the resulting classical planning problem (Helmert and Domshlak 2009), which were used in the original ASNs. We did not find experimental success for either.

Output. The last layer of the network is the output layer. The output ϕ_a^{L+1} of each action module in the last layer is only a single real number, and the resulting output of the network is the masked softmax of all individual outputs,

$$\pi^{\theta}(a \mid s) = \frac{m_a \exp(\phi_a^{L+1})}{\sum_{a' \in A} m_{a'} \exp(\phi_{a'}^{L+1})}$$

where m_a is a boolean mask of whether the action a is applicable in the current state s , and $\pi^{\theta}(a \mid s)$ is the probability of selecting action a in state s .

Given weights θ for a domain, the resulting generalised policy for a domain produces a plan for each instance by repeatedly selecting and applying actions using π^{θ} , starting at the initial state, and terminating upon reaching a goal state, a state with no applicable action, or a fixed length limit. Like ASNs, we use π^{θ} during training by sampling from it, and during evaluation by greedily selecting the action with maximum probability and breaking ties deterministically.

Miscellaneous. We have introduced comparison modules and fluent modules together, along with their implications for network input and action modules. It is important to note that they can and are designed to be applied separately. We term the network with only comparison modules *C-ASNs* and the network with only fluent modules *F-ASNs*. This specialisation can be empirically beneficial, we hypothesise

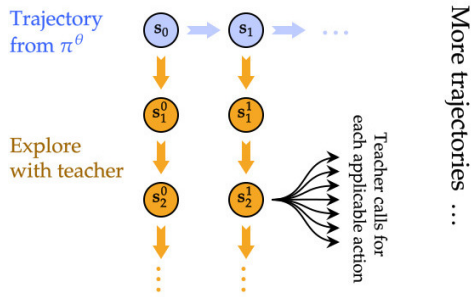


Figure 2: The exploration algorithm used by the original ASNs, where all states from the sampled trajectories are explored with the teacher planner. For all states shown, the teacher planner is called on all applicable actions to mark the actions leading to the lowest-cost plans.

that it allows the network to focus on a particular form of reasoning, reduces computation burden, and potentially reduces overfitting. To disambiguate, we will use B-ASNs to refer to the network with both modules, and ν -ASNs to refer to the collection of architectural variations.

It is worth noting that ν -ASNs make no requirement on the form of comparisons or numeric effects appearing in the problem, and can be applied to the full numeric fragment of PDDL 2.1 (Fox and Long 2003). Similarly, we introduced action preconditions as conjunctions of comparisons and propositions only for simplicity, and they can in fact be conjunctions of any quantifier free formula involving conjunction, disjunction, and negation of propositions and comparisons. Likewise, the goals do not have to be conjunctions of comparisons and propositions either. However, in order for ν -ASNs to generalise, the goals of the problem instances it is applied to should have a consistent structure such that experiences from reaching one instance’s goal is useful to reaching the goal of other instances.

Additionally, we also include skip connections (Toyer et al. 2020). That is, we append the output of each action/proposition/fluent module to the input of the same module in the next layer but omit these in equations for simplicity. Skip connections make it easier for the network to propagate information across layers.

Dynamic Exploration

Like the original ASNs, for each domain, we train ν -ASNs on a small number of training problem instances under the supervision and guidance of a teacher planner. We use the state-of-the-art ENHSP numeric planner as our teacher planner (Scala et al. 2020). The original algorithm trains over a number of epochs. Each epoch involves an *exploration phase* and a *learning phase*. The exploration algorithm used by the original ASNs first uses the current network weights θ to sample a number of trajectories from π^θ , then explores all the states from these trajectories by calling the teacher planner on them, and adding all the resulting states to a multiset state memory S_{mem} , as illustrated in Figure 2. For each state added to S_{mem} , the teacher plan-

ner is called on the resulting state of applying each applicable action, and the actions leading to the lowest-cost plans are marked. The learning phase then updates the weights through mini-batch gradient descent to choose the marked actions using the cross-entropy-based loss

$$\mathcal{L}_\theta = -\frac{1}{|\mathcal{B}|} \sum_{s \in \mathcal{B}} \sum_{a \in A} \left[(1 - y_{s,a}) \cdot \log(1 - \pi^\theta(a | s)) + y_{s,a} \log \pi^\theta(a | s) \right] + \frac{1}{2} \lambda \|\theta\|^2$$

where \mathcal{B} is a mini-batch of states from S_{mem} , $y_{s,a}$ is a binary variable indicating if action a is marked at state s , and the last term is an ℓ_2 regulariser.

In our preliminary experiments we find that the original exploration algorithm is inadequate for effectively learning generalised policies for numeric planning. We observe that for numeric planning, the plan lengths of even simple training problems tend to be longer than that of classical planning, and ENHSP to be much slower than teachers used by ASNs for classical planning. In each epoch, the number of states added to S_{mem} in the original exploration algorithm is quadratic in the plan lengths of the training problems, resulting in a large number of states added for numeric planning. This has a number of downstream consequences in the original exploration algorithm:

1. For each state s added to S_{mem} , the teacher is called for each applicable action in s . An increased number of states added to S_{mem} therefore leads to a significant increase in calls to the teacher planner, especially in domains where states tend to have many applicable actions.
2. The learning phase uses states in S_{mem} for training. We observe that the size of S_{mem} is sometimes a few orders of magnitude larger than the number of states used in the learning phase. In this case many states added to S_{mem} are rarely used, wasting memory and exploration effort.
3. The alternating exploration and learning phases mean that states added to S_{mem} early are explored using historical network weights and hence less useful for learning than more recently added states.

We propose a *dynamic exploration* algorithm to address these problems by removing stale states and dynamically adjusting the amount of exploration performed based on the time spent on the learning phase, as shown in Algorithm 1. In each exploration phase, dynamic exploration first uses the current network weights to generate T_{gen} trajectories for each training problem, i.e. calling the network iteratively starting at the initial state and sampling the action to apply from π^θ (line 8). Each such trajectory terminates upon reaching a goal, a fixed length limit, or a state with no applicable actions. The states in these trajectories are added to an initially empty S_{traj} multiset (line 3). We then repeatedly randomly remove (or *explore*) states from S_{traj} and call the teacher planner for each removed state until a termination condition is met (lines 12 to 13). All states in the resulting teacher plan are added to S_{mem} (line 14). This can be understood as asking the teacher planner to guide the network back onto a valid trajectory. If S_{traj} ever becomes empty,

Algorithm 1: Dynamic exploration. We group new states in S_{expl} and add to S_{mem} with epoch number

Data: A set of training problems P_{train} ; current ν -ASNets weights θ ; epoch number n

```

1 Procedure genTraj()
2   for  $\zeta \in P_{\text{train}}$  do
3      $S_{\text{traj}}.\text{extend}(\text{runPolicy}(s_0(\zeta), \pi_\theta))$ 
4 Procedure explore( $n, S_{\text{mem}}$ )
5    $S_{\text{traj}} \leftarrow \emptyset$ 
6    $S_{\text{expl}} \leftarrow \emptyset$ 
7   for  $i = 1, \dots, T_{\text{gen}}$  do
8     genTraj()
9   while not terminate() do
10    if  $|S_{\text{traj}}| = 0$  then
11      genTraj()
12     $s \leftarrow S_{\text{traj}}.\text{popRandom}()$ 
13     $S_{\text{expl}}.\text{extend}(\text{teacherPlan}(s))$ 
14   $S_{\text{mem}}.\text{extend}((S_{\text{expl}}, n))$ 
15  while  $|S_{\text{mem}}| > N_{\text{mem}}$  do
16     $S_{\text{mem}}.\text{popOldestEpoch}()$ 

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it is refilled by generating one trajectory from each training problem from the initial state.

The termination condition (line 9) is based on the average time t_{learn} spent on recent learning phases and an hyperparameter r to control the ratio between spent on exploration and learning. We terminate exploration when either $r \cdot t_{\text{learn}}$ time has elapsed in the current exploration phase or the number of states explored reaches an upper bound e_{max} , but never before at least e_{min} states have been explored. For the first exploration phase where t_{learn} is undefined, we terminate once at least one state from each problem and at least e_{min} states overall have been explored.

To avoid a size explosion of S_{mem} and ensure its states are recent, we group states in S_{mem} by the epoch they are added in. Whenever the number of states in S_{mem} exceeds a limit N_{mem} , we repeatedly remove the oldest group till the size of S_{mem} falls back under the limit (line 16).

To address point 1, we adopt an alternative action marking method. Whenever a state s is added to S_{mem} , instead of checking if each applicable action can lead to a lowest-cost plan, this option only marks the action selected by the teacher planner, saving the need to call the teacher planner for each applicable action.

Experimental Evaluation

For evaluation, we implemented ν -ASNets and dynamic exploration based on the original implementation of ASNets. Code is available at (Wang and Thiébaux 2024).

Experimental Setup

Teacher planner and benchmark domains. We use the state-of-the-art numeric planner ENHSP-20 (Scala et al. 2020) as the teacher planner, and use benchmarks from the

International Planning Competition 2023 Numeric Track³. These domains only include simple or linear numeric planning problems, and do not fully demonstrate the applicability of ν -ASNets to the entire numeric fragment of PDDL 2.1. For each domain, we use the 3 to 6 smallest instances for training. ENHSP has a wide set of configurations based on heuristic, search algorithm, and the use of methods such as redundant constraints and helpful actions. From these configurations, we select as the teacher ENHSP configuration one that produces short plans quickly (within one second) for the training instances. We do not experiment with domains where such a teacher configuration could not be found. The resulting benchmark domains and teacher configuration used for each domain are shown in Table 1. We classify domains by the proportion of reasoning that is numeric versus propositional into heavily numeric and hybrid domains.

For the domain Counters, we do not use the IPC instances, but instead use a set of evaluation instances with 2 to 60 counters where all the counters start with value 0. We only use one of these evaluation instances for training, and include for training another two instances similar to it but with different initial states. This set up allows us to better understand how the network would generalise across dimensions (number of counters) unvaried during training.

Baselines and ν -ASNet variations. We use ENHSP as the baseline for comparison using all configurations that are used as teacher for at least one domain, and report its results for the best and teacher configurations of each domain. We also compare within the ν -ASNets variations, namely the baseline network without either fluent or comparison modules (N-ASNets), F-ASNets, C-ASNets, and B-ASNets. We also compare the original ASNets training algorithm and the dynamic exploration algorithm for training F-ASNets and C-ASNets, and use superscripted O or D on the ν -ASNets variation to denote them. For fairness, we enable alternative action marking method with the original algorithm.

Hyperparameters. For each domain, we train the network and evaluate it once on each problem instance. Unless otherwise specified, the hyperparameters for ν -ASNets are fixed across domains and architectural variations. These hyperparameters are based on the original ASNets hyperparameters with minor tuning. We use three action layers and two state layers, with a hidden representation size (d_h) of 15 and an ELU as the non-linearity σ (Clevert, Unterthiner, and Hochreiter 2016). When using dynamic exploration, in each exploration phase we collect $T_{\text{gen}} = 2$ trajectories initially, terminate exploration with parameters $r = 1$, $e_{\text{min}} = 10$, and $e_{\text{max}} = 1000$, and impose an $N_{\text{mem}} = 15000$ limit on the size of S_{mem} . When using the original algorithm, we collect two trajectories per problem and explore all states within the collected trajectories. After exploring, the learning phase performs weight optimisation using the Adam optimiser ($\beta_1 = 0.9$, $\beta_2 = 0.99$, and $\epsilon = 10^{-7}$). Mini-batch gradient descent is performed with a learning rate of 0.0003, batch size of 50, and 60 batches per epoch. We addition-

³<https://ipc2023-numeric.github.io/>

Domain	Classification	Teacher	Numeric ASNet						ENHSP		
			B^D	F^D	C^D	F^O	C^O	N^O	best	teacher	best
Block Grouping (20, 4)	HN, simple	hadd-gbfs	15 (8.0)	11 (8.0)	17 (8.0)	10 (15.5)	15 (9.8)	2 (24.0)	17	20	20
Counters (59, 1)	HN, simple	hrmax-astar	9 (0.2)	7 (0.1)	14 (6.4)	10 (0.1)	17 (1.3)	1 (14.8)	17	8	39
Delivery (20, 4)	hybrid, simple	hadd-astar	5 (8.0)	5 (5.3)	20 (1.9)	9 (9.9)	18 (6.7)	17 (3.3)	20	8	16
Drone (20, 4)	HN, linear	hadd-astar	9 (8.0)	4 (8.0)	3 (8.0)	7 (24.0)	3 (24.0)	0 (24.0)	9	11	19
FO-Counters (20, 3)	HN, linear	hrmax-astar	4 (3.6)	5 (1.6)	3 (8.0)	6 (7.3)	3 (10.2)	2 (15.4)	6	4	5
MPrime (20, 4)	hybrid, simple	hmrp-ha-gbfs	16 (3.1)	19 (1.8)	12 (7.9)	18 (4.4)	16 (24.0)	6 (24.0)	19	16	18
Rover (20, 4)	hybrid, simple	hmrp-ha-gbfs	7 (8.0)	4 (8.0)	4 (8.0)	5 (24.0)	4 (24.0)	4 (24.0)	7	7	7
TPP (20, 3)	hybrid, linear	hadd-gbfs	0 (8.0)	0 (8.0)	19 (8.0)	0 (24.0)	20 (24.0)	16 (22.7)	20	4	4
Zenotravel (20, 6)	hybrid, linear	hadd-gbfs	0 (8.0)	0 (8.0)	17 (0.6)	0 (24.0)	16 (0.8)	16 (0.6)	17	20	20

Table 1: Number of instances solved (coverage) by each system, with the ν -ASNet training time in hours shown in parenthesis. The number of instances for evaluation and the number of evaluation instances seen during training are shown in parenthesis after the domain. We also show the classification of the domain (see text) by heavily numeric (HN) versus hybrid (hybrid) and simple versus linear. We additionally show the teacher configuration used for each domain in the format {heuristic}-{search.algorithm}, with the optional “ha” indicating the use of helpful actions.

ally apply an ℓ_2 regulariser with a coefficient of 0.005 and a dropout probability of 0.1. We stop training when all collected trajectories reach the goal for 20 consecutive epochs.

Computational limits. When training, we apply a time limit of 8 hours for dynamic exploration and 24 hours for the original training algorithm. During evaluation, we apply a 1800 seconds time limit per problem for ENHSP and ν -ASNet. Training of ν -ASNet and evaluation of ENHSP is performed on a virtual machine with 32GB of memory and a single dedicated core clocked at 4.5 GHz. Evaluation of ν -ASNet is performed on the same virtual machine with only 8GB of memory.

Results

Table 1 shows the coverage achieved by the ν -ASNet variations and ENHSP by domain. The learned generalised policies are able to achieve coverages competitive with ENHSP, and outperform it on several domains, namely Delivery, FO-Counters, MPrime and TPP. Interestingly, except for FO-Counters, the other three domains all involve some forms of graph traversal and logistics. On Block Grouping, Rover, and Zenotravel, the ν -ASNet achieve coverages commensurate with ENHSP. On the remaining two domains, Counters and Drone, ν -ASNet are able to generalise from the small training problems to bigger problems and perform similarly with or outperform its teacher, but not the best ENHSP configuration.

To better understand how well ν -ASNet are able to generalise, we examine the particular problem instances to see if it is only generalising to problems with similar size to those seen during training. On Block Grouping, the largest training instance has 10 blocks on a 15 by 15 grid, whereas the largest solved instance has 25 blocks on a 100 by 100 grid. The training instances for Counters all have 4 counters, while ν -ASNet variations are generally able to solve evaluation instances with up to 15 counters. This result on Counters shows that ν -ASNet are able to generalise across factors (number of counters in this case) kept constant during training. The largest Delivery training instance has 10

items to deliver, while the largest solved instance has 42. Similar scales of generalisation are achieved on other domains, and demonstrate the strong generalisation capabilities of ν -ASNet.

Table 1 also shows the training time of various ν -ASNet variations. By comparing the training times of C^D -ASNet with C^O -ASNet and F^D -ASNet with F^O -ASNet, our results show that dynamic exploration is able to achieve much lower training times than the original exploration algorithm. This is not a consequence of the lower training time limit we apply for dynamic exploration, as the trend continues even when neither training methods reach their respective time limits, see e.g. in Delivery or FO-Counters. Furthermore, we do not observe any notable reduction in coverage for dynamic exploration when compared to training with the original algorithm. This suggests that dynamic exploration consistently enables learning generalised policies faster without compromising the effectiveness of the learned generalised policies. The only notable exception on Counters with C-ASNet is likely due to a high variance in training time that we found during multiple training runs.

Figure 3 shows the plan cost and evaluation runtime produced by the learned generalised policies and ENHSP with the teacher configuration. When both produce valid plans, they produce plans with similar costs. On Zenotravel ν -ASNet tend to produce better plans, while on Block Grouping ENHSP tends to produce better plans. For runtime, when both produce plans quickly (less than 10 seconds), ENHSP tends to be quicker than the generalised policies. This is likely due to the higher constant overhead required by ν -ASNet to construct the network and load the weights. On more complex problem instances, ν -ASNet tend to produce plans faster than the ENHSP teacher configuration. The large number of points on the top line in the runtime plot demonstrates that the learned generalised policies are able to solve many instances the teacher cannot solve.

Why do we need F-ASNet or C-ASNet? Results in Table 1 show that the specialisation in reasoning offered by F-ASNet or C-ASNet often allow one of them to perform

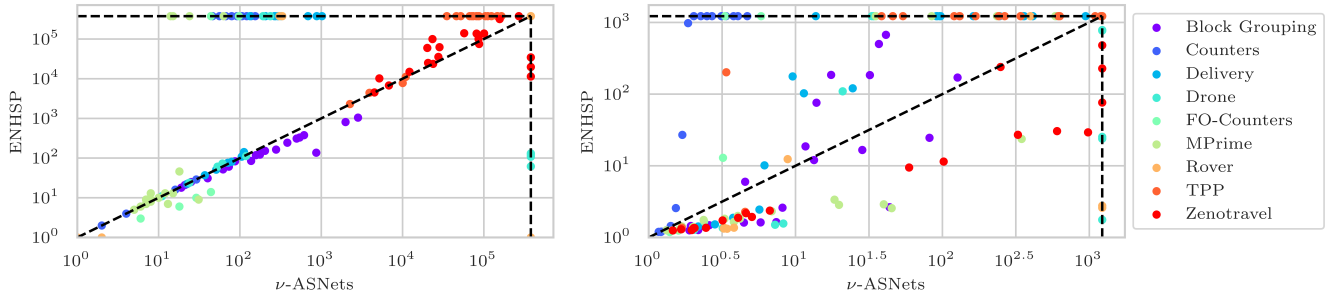


Figure 3: Plan cost (left) and runtime in seconds (right) for each problem instance of the best ν -ASNets variation versus the teacher ENHSP configuration. Points in the bottom-right triangle favour ENHSP and on the top-left triangle favour ν -ASNets. Problems unsolved by a system have value set to the maximum of the axis. A constant of 1 is added to ensure all points lie within view.

better than if they are both included. For example, comparison modules alone on Delivery or fluent modules alone in MPrime achieve notably higher coverages than when the other is included.

Why do fluent modules result in coverages of 0 in TPP and Zenotravel? In both domains, whenever fluent modules are included in the network, the learned generalised policies fail to solve any problem. In these domains, except for fluents used to help the plan metric, all the other fluents are only meaningful in context of each other. We suspect that when fluent modules are included, the network attempts to learn to reason directly on the fluent values, but receive conflicting information on how to do so on the different training instances. This results in training never being able to converge, and ultimately the coverage of 0.

Which ν -ASNets variation is the best? The best ν -ASNets variation by coverage depends on the nature of the domain. Generally, C-ASNets tend to perform well on all benchmark domains, while F-ASNets and B-ASNets perform well on particular domains such as MPrime and Rover respectively.

How can N-ASNets perform well on some domains? N-ASNets is not equipped with network modules for numeric reasoning, but it still has the capability for classical planning reasoning. Unsurprisingly, this allows it to still be effective on hybrid domains where there is a sizeable classical planning component. However, on heavily-numeric domains its unsuitability for numeric reasoning is clear from the poor coverage it achieves.

How do ν -ASNets compare with the IPC 23 competition planners? On domains where we use the same problem sets as IPC 23 (i.e. all but counters), ν -ASNets achieve better coverages than the reported coverage⁴ of the IPC 23 competition planners except on Drone and Rover.

How beneficial are the numeric landmarks? Table 2 shows that the impact of numeric landmark varies notably by domain. For Block Grouping, Counters, and FO-Counters,

Domain	C^D	$C^D(\text{NL})$	F^D	$F^D(\text{NL})$
Block Grouping	17	1	11	2
Counters	14	2	7	1
Delivery	20	20	5	4
Drone	3	2	4	5
FO-Counters	3	2	5	1
MPrime	12	14	19	16
Rover	4	4	4	4
TPP	19	17	0	0
Zenotravel	17	17	0	0

Table 2: Number of instances solved (coverage) depending on the use of numeric landmarks. NL (no landmark) in parenthesis indicates that the landmark was not used.

numeric landmarks are essential for ν -ASNets to perform well. This is expected, as on these domains the goal includes a heavy amount of numeric information, which the network would be uninformed about without numeric landmarks. On other domains, not providing the network with numeric landmarks has little impact on coverage.

Related Work

Existing work on generalised planning is severely limited when it comes to dealing with numeric information. Popular approaches based on Qualitative Numeric Planning (QNP) (Srivastava, Immerman, and Zilberstein 2011; Bonet and Geffner 2020), can represent a fixed number of positive numeric variables that can only be increased or decreased by a positive non-deterministic amount in action effects, and boolean combinations of comparisons of these variables with 0 in action preconditions, initial states and goals. Other approaches allow for incrementing or decrementing a finite set of positive counters by a constant in a deterministic fashion (Srivastava, Immerman, and Zilberstein 2010; Srivastava et al. 2015). The more recent Generalised Integer Numeric Planning (GLINP) (Lin et al. 2022) supports non-simple numeric effects, but is limited to integer variables. In contrast our work support the full numeric fragment (level 2) of PDDL2.1 (Fox and Long 2003), including nonlinear effects

⁴<https://ipc2023-numeric.github.io/results/presentation.pdf>

(Scala et al. 2016a) and numeric variables whose number grows with the number of objects. On the other hand, the above works provide guarantees on the effectiveness of generalised policies, whereas our learning approach cannot.

Independently and concurrently to our work, Tariq, Valenzano and Soutchanski (2023) experimented with handling numeric planning problems with the original ASNet policy representation. They reduced the set of values each fluent takes to a finite range, which they then manually discretised into consecutive intervals, each represented by a new predicate. Numeric conditions in the action schemas are then compiled into a disjunction over these predicates. As Tariq et. al observe, this approximation of the original numeric problem creates a large number of related propositions for each action, which leads to impractically large networks and compromises the sparseness of the ASNets policy representation. The empirical evaluation conducted by Tariq et al. only used four unseen test instances per domain. These are only marginally larger than the training instances and solved within less than a second by both ENHSP and ASNets. In contrast, we have proposed an architecture that treats fluents and comparisons as first-class citizens and explicitly reasons about them. Its performance is competitive with ENHSP over the latest numeric planning competition instances.

Conclusion and Future Work

We have introduced ν -ASNets and its variations, neural network architectures for learning generalised policies for numeric planning based on ASNets. The network is able to reason directly about numeric values through fluent modules, and about numeric contexts through comparison modules. We also introduced dynamic exploration, which trains ν -ASNets much faster than the original algorithm used by ASNets, without harming the effectiveness of the learned generalised policies.

Our work leaves significant room for future research. A common trait in plans for numeric planning problems is the repetition of actions. We believe that a network architecture capable of not just predicting the action to apply, but also the number of times to apply it, can be highly effective. Additionally, while our work has focused on numeric planning, the method we use to construct comparison modules can in principle be applied to other components of actions modelled in PDDL, to include constructs such as action effects and universal/existential quantifiers. This leads to a network that can potentially learn generalised policies for a much more expressive class of problems than numeric planning.

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