Action Schema Networks – IPC Version

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Abstract

Action Schema Networks (ASNets) are neural network architectures for generalized reactive policies that leverage the relational nature of planning problems represented in (P)PDDL. An ASNet policy can be learned by imitating the actions selected by a traditional non-learning planner. Even when the training set consists of only a few small problems, ASNets are able to solve significantly larger instances within the same domain. In this paper, we describe the ASNet planner submitted to the Learning Track of the International Planning Competition (IPC23). It closely follows the original implementation (Toyer et al. 2020), but focuses on solving deterministic planning problems and employs newer PDDL parsers and eagerexecution-based machine learning frameworks.

1 ASNets

In this section we briefly describe the core ideas of ASNets. For a complete and comprehensive explanation, we refer the reader to (Toyer et al. 2020, 2018). Action Schema Networks (ASNets) attempt to connect the world of automated planning to deep learning by encoding the relational structure of factored planning problems into neural networks, where neurons represent actions and propositions. Connections between each layer are derived from a graph representation based on the action schema of the given domain. Because all problems from the same domain share the same action schema, weights of the policy network trained on a small set of problems can be shared with networks instantiated for larger problems from the domain. Therefore, training with simpler problems makes solving harder problems with policy networks possible using the learned parameters.

ASNets are policy networks that take in the current state s and return a policy $\pi^{\theta}(a|s)$. The current state is encoded as a feature vector and fed into an alternating sequence of action and proposition layers. Each action layer consists of action modules representing actions, and each proposition layer is formed by proposition modules corresponding to propositions. The connections between modules are determined by the *relatedness* of actions and propositions.

Definition 1 (*Relatedness*) A proposition *p* is related to an action *a* at position *k* if the lifted proposition (predicate) is

the *k*-th unique predicate in the precondition or effect of the lifted action (action schema).

A proposition module is connected to an action module in the next layer, if and only if the underlying action and proposition are related. Similarly, action modules are connected to related proposition modules in the next layer. Because the relatedness only depends on predicates and action schemas, propositions and actions grounded from the same action schema share the same number of connections.

Each action module at layer l with underlying action a takes in a vector $u_a^l \in \mathbb{R}^{d_a^l}$ and outputs another vector $\phi_a^l \in \mathbb{R}^{d_h}$, where d_h is the dimension of the hidden vector. The outputs of related proposition modules $p_1, ... p_M$ in the previous layer are vectors $\psi_{p_1}^{l-1}, ..., \psi_{p_M}^{l-1}$. These vectors, together with the output vector of the same action module from the previous layer ϕ_a^{l-1} are concatenated to form the input vector u_a^l , which is then fed into an affine transform $f(W_a^l \cdot u_a^l + b_a^l) = \phi_a^l$. All actions a from the same action schema i in this layer share the same weights and bias, namely $W_a^l = W_i^l, b_a^l = b_i^l$.

Similarly, a proposition module at layer l with underlying proposition p takes in a vector $v_p^l \in \mathbb{R}^{d_p^l}$ and outputs another vector $\psi_p^l \in \mathbb{R}^{d_h}$. Instead of concatenation, the outputs of related action modules $\phi_{a_1}^{l-1}, ..., \phi_{a_N}^{l-1}$ from the previous layer are grouped by their position of relatedness and then passed through a max pooling function. Then the outputs of the pooling functions are concatenated with the output vector of the same proposition module from the previous layer ψ_p^{l-1} . Finally, the resulting vector v_p^l is fed into an affine transform $f(W_p^l \cdot v_p^l + b_p^l) = \psi_a^l$. The weights W_p^l and bias b_p^l are shared by all propositions grounded from the same predicate in this layer.

Because weights are shared between actions grounded from the same action schema (same for propositions of the same predicate), the learned weights are generic transformations that can be applied to the instantiated networks for of problems of any size.

The input features to the first action layer are encodings of the current states. For each action module with related propositions $p_1, ..., p_M$, the input vector u_a^1 consists of four parts: a vector $v \in \{0, 1\}^M$ where $v_i = 1$ iff the corresponding proposition p_i is true in current state; a vector

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 $g \in \{0,1\}^M$ where $g_i = 1$ iff the proposition appears unnegated in the goal; a scalar m = 1 if the action is applicable in current state; and finally, an extension part including features derived from other heuristics. The first half of the extension part is a one-hot label vector $c \in \{0,1\}^3$ derived from LM-cut landmarks (Helmert and Domshlak 2009). The vector has $c_1 = 1$ iff the action appears as the only action in at least one landmark; $c_2 = 1$ iff the action appears in a landmark containing two or more actions; and $c_3 = 1$ otherwise. The second half of the extension is an integer recording the number of times the action has appeared in the plan previously.

The last layer is also an action layer but with different outputs. The output dimension of each action module is one. Specifically, each action module outputs a single scalar value $\phi_{a_1}^{L+1}, ..., \phi_{a_N}^{L+1}$. The final policy is then the log-scaled probability distribution:

$$\pi^{\theta}(a_i|s) = \frac{m_i \exp(\phi_{a_i}^{L+1})}{\sum_{j=1}^N m_j \exp(\phi_{a_j}^{L+1})}.$$

The model is then trained through imitation learning. Namely, the model will try to imitate the action selection policy of a teacher planner. At each training step, the trainer samples a set of states (so-called "rollouts") and tries to find a plan from the states to the goal by choosing actions according to the learnt policy π^{θ} . Then the trainer compares these plans to solutions found by the teacher planner and learns from their differences. Originally ASNets used SSiPP (Trevizan and Veloso 2014)) for stochastic problems and fastdownward (Helmert 2006) for deterministic problems as teachers. The teacher outputs a Q-value $Q^{teach}(s, a)$, which represents the cost of reaching the goal from state s when taking action a and acting optimally afterwards. Therefore, the best action will have the lowest Q-value. Using the Qvalues, the trainer calculates the cross-entropy loss between the learnt policy π^{θ} and the teacher's policy. By minimising the loss, the model learns to make similar decisions as the teacher planner.

2 Changes to ASNets

In this section we describe our changes and improvements on the original implementation of ASNets.

2.1 ML Framework

The learning module of the original ASNets is based on TensorFlow 1.x (TF1) (Abadi 2016), which is a deprecated ML framework with very different runtime behaviour compared to the latest frameworks. We updated the implementation by replacing the learning framework with TensorFlow 2.x (TF2) (Singh et al. 2020), which has faster and simpler API and more supported functions. Some major differences between TF1 and TF2 includes:

• TF1 uses lazy execution and requires manual graph construction representing the model architecture (Abadi 2016), and then manually compiles the graph by passing the input and output tensors to a *Session*. TF2 executes eagerly (Singh et al. 2020). The construction of the graph

can be done in the background by calling higher-level API modules.

- TF1 variables are based on implicit global namespaces. Model variables are stored in the graph even if no Python variable is pointing to it. TF2 variables are managed the same way as Python variables, making it faster and more memory-efficient.
- TF2 provides a well-defined high-level API called Keras (Gulli and Pal 2017) to easily build and train a neural network while remaining highly customisable.

We re-implemented the model and model trainer using the Keras API. The behaviours of the action and proposition modules are defined in separate classes. Connections between the modules are defined in a network extended from the Keras Layer class, which provides a well-defined API for training and evaluation. The shared weights are managed through a weight manager where weights are defined as TensorFlow variables with learnable values. Then we used a Keras Loss Class to implement the loss function described in (Toyer et al. 2020) to use the class interface in training directly. Finally, we rewrote the trainer with the new models and components while keeping the data generation and training mechanism unchanged.

2.2 Training

Originally ASNets were designed for solving Stochastic Shortest Path problems (SSP) (Bertsekas and Tsitsiklis 1996). Because we only consider deterministic domains in this competition, we simplified and optimised the training for the deterministic setting.

First, for the teacher planner, we followed the configuration of the deterministic version of the original ASNets, which used fast-downward (Helmert 2006) in solving deterministic problems. In other words, we use fast-downward to compute the teacher's policy $Q^{teach}(s, a)$ for state *s* and action *a* during the training process. And for the solver algorithm used by fast-downward, we chose the LAMA-first algorithm (Richter and Westphal 2010), whereas the original version of ASNets used A-star with LM-cut (Helmert and Domshlak 2011). In stochastic problems, the policy is generally represented by a probability distribution function, while in deterministic settings, we simply take the best action. Therefore, the Q-value computed by the teacher becomes a one-hot label indicating which action is the optimal selection and rejecting all other actions.

Because ASNets can be trained on small problems and then reuse the same learnt weights for solving larger problems, it is unnecessary to train the model on large problems directly. However, in this competition, we don't know the domains and sizes of problems in advance. Hence we cannot set the boundary between "easy" and "hard" problems¹. Because generating training data for the model requires solving many sub-problems (namely, finding plans from some intermediate states to the goal), if a training problem is too hard

¹The competition provides problems in increasing order of difficulty

to solve, the learning script might waste too much time solving these sub-problems instead of training the model. Therefore, the program has to decide by itself whether to use the given training problem or not.

We set the timeout threshold of the fast-downward solver to 1 minute. In other words, if the teacher fails to find a plan for the given state within one minute, it will skip the problem and train the model with data from other problems. Besides, we train the model in an incremental manner: we first train the model with only the two easiest problems. The learnt policy is used as a basic planner. Then we append one more problem to the training set and train a new model. We repeat this process until all problems are in the training set or the teacher fails to return plans for all new problems. We record the problems that failed to be solved by the planner and then try to train the model again with fewer rollouts per step and a longer time limit (2 minutes) for the teacher planner. The incremental training set strategy guarantees we can have a benchmark model before the trainer runs out of memory or fails to learn from harder problems.

We used a stepped decreasing learning rate starting from $1e^{-3}$. The learning rate drops to $1e^{-4}$ after 500 steps and further drops to $1e^{-5}$ after 1000 steps. The L1 regularisation parameter is set to 0, and the L2 regularisation parameter is set to $2e^{-4}$. The model has 3 action layers and 2 proposition layers with hidden vector size 16 and dropout rate 0.1. We save the trained model after every epoch if the evaluation result is better than the previous one.

2.3 Evaluation

The original ASNets select an action to execute in a state s by identifying the action with the highest probability in their last layer, i.e., $\arg \max_{a \in A(s)} \pi^{\theta}(a|s)$, breaking ties arbitrarily. While effective in practice, an ASNet does not guarantee that the generated plans will be loop-free. For example, in a Blocks World problem, the arg max strategy cannot guarantee that picking up block A from the table, placing A back on the table, and repeating the cycle indefinitely will never happen. The solution proposed in (Toyer et al. 2020) is to sample the action to be executed in s according to $\pi^{\theta}(\cdot|s)$, guaranteeing that loops will be eventually broken; however, this method performs worse than the arg max strategy in practice because the probability of sampling an unfavorable action is higher than the probability of the arg max strategy resulting in a loop. To address these issues, we use a new execution method based on the arg max strategy that ensures loop-free execution while preserving its performance in the absence of loops. To do so, the planner keeps track of the set E(s) of actions previously executed in each visited state s. An action for state s is then chosen according to $\arg \max_{a \in A(s) \setminus E(s)} \pi^{\theta}(a|s)$, i.e., we select the highest probability action that has not yet been executed. If there are no applicable actions remaining, i.e., E(s) equals A(s), the planner returns a failure to find a plan. This strategy behaves as the arg max strategy until a loop is detected, in which case it systematically executes different actions from highest to lowest probability according to $\pi^{\theta}(\cdot|s)$.

3 Issues and Solutions

The outcomes of the competition significantly deviated from our expectations and our own experiment results. In the original paper, the model's performance surpasses LAMA-first on *blocksworld* after sufficient training. Although we acknowledge that the model might not perform as well as the original paper due to limited training time and resources, and a different training set, we did not anticipate the planner failing in multiple domains. Therefore, we thoroughly checked our implementation and training/evaluation pipeline. After careful inspection and validation, we identified three major issues that led to the unexpected poor performance:

- 1. There was a bug in the implementation of the loop-free execution strategy mentioned in Section 2.3. The original approach utilized a dictionary that stored states and the number of times they appeared in the plan as string-integer pairs. The string representation of states was used as the dictionary keys. The loop-free execution strategy ensured that actions were not repeated in the same state, which prevented loops in the plan. However, some states were incorrectly mapped to the same string while converting states to string representations. As a result, the planner treated different states as if they were the same and employed an unnecessary loop-breaking strategy. Consequently, the plan generated was not as good as it could be and in some cases, no valid solution was found.
- 2. One of the pre-processing steps was to quickly find a superset of all grounded actions applicable to a given problem. This reachability analysis method did not work correctly when the problem contained negative preconditions. As a result, it failed to return all reachable actions, which caused an error when the fast-downward planner returned these missing actions.
- 3. As we mentioned in Section 1, the encoding of actions contained an extension part based on LM-cut landmarks. These values were computed using the SSIPP. For this reason, SSIPP was instantiated for each problem. However, the SSIPP planner had an upper bound on the number of atoms in the problem. This upper bound was set to 16384 by default, which was much lower than the upper bound of atoms in the problems provided. Therefore, when the input problem became large, the planner failed to parse the problem and quit without solving the problem.

The competition's planner performs poorly due to a combination of these issues. However, it's important to note that these problems are not connected to the methodology or core concepts underlying ASNets. To obtain an unbiased evaluation of ASNets we need to address these issues first. This was achieved by the following changes:

 We fixed the string representation of states. The states are now represented as a string of tuples of propositions. This ensures that no different states are mapped to the same key. The trade-off is that this method may result in a long key string for complex problems, leading to higher memory usage.

Domain	#Train	Agile Scores	Quality Scores	Coverage
blocksworld	30	22.25	27.66	37
childsnack	12	0	0	0
childsnack-s	8	2.67	2.11	3
ferry	40	32.32	41.66	48
floortile	20	15.0	19.13	24
miconic	52	35.76	44.27	61
rovers	16	4.5	4.33	6
satellite	32	29.66	38.19	45
sokoban	13	1.53	1.04	2
spanner	31	9.97	10.0	10
transport	50	20.93	28.06	35
overall(ex. childsnack-s)	297	171.92	214.34	268

Table 1: The column "#Train" indicates the number of problems that the planner is trained on before reaching the time limit. Agile scores and quality scores are computed using the same scripts from the competition. Coverage is the number of problems solved among the 90 testing problems per domain. "childsnack-s" is the planner trained with the same settings as others but ignores the first 12 training problems.

- 2. We fixed the bug in the reachability algorithm. Due to the time limit in the competition, we optimised for a faster approximation of a set of applicable actions that returns more false positives in domains with negative preconditions.
- 3. We increased the upper bound of atoms to the maximum acceptable range of 65535.

After having resolved the issues, we re-evaluated the planner's performance under the same time and resource constraints as in the competition, namely training with a single-core CPU, 32 GiB memory, and 24 hours, and testing with a single-core CPU, 8 GiB memory, and 0.5 hours.

In order to determine the agile and quality scores, we utilized the scripts and data provided by the competition. The experiment outcomes are presented in Table 1.

4 Ablation Analysis

In this section, we analyse the competition results after reevaluation of the model's performance. First of all, it's important to note that the planners used by participants in the learning track have different "learning components." Huzar learned a Graph Neural Network that helps the planner shrink the set of available actions and then solve the problem using LAMA; GOFAI learned to find partial grounding of problems and then solve the problem with a portfolio; Vanir learns a hierarchical policy, which is a rooted tree with each node representing a decomposition of parent problems and sketch rules (Drexler, Seipp, and Geffner 2023). Muninn, the most similar competition entry to our planner, learned a neural network policy by optimising two value functions. It is expected that ASNets is no match for planner portfolios or variants of LAMA. After analysing the logs of learning and planning for each domain, we offer the following observations:

Improved Testing Scores After fixing the pipeline, the results significantly improved in almost all domains. The total coverage increased from 40 to 268 and the quality score rose from 29 to 214. The only domain for which results did not

improve is *childsnack*, where we still cannot solve any problem, due to insufficient training time.

Insufficient Training Time The planner faces a significant issue with limited time and resources, which hinders its ability to exploit and fully learn from the training set. Unfortunately, due to the 24-hour training time limit, the trainer times out on all domains before completing the full training set. As shown in Table 1, the trainer can only utilize a maximum of 52 problems in the miconic domain, and in the worst-case scenario, it can only use 12 problems in the childsnack domain. As expected, the planner's performance on *childsnack* is the worst. One observation is that the first 12 problems of childsnack have at most 2 "children". However, even the easiest problem in the test set (easy/p01.pddl) has four children sitting on two tables. The planner failed to learn enough information to generalise to such cases. To demonstrate that the planner can learn useful domain knowledge given enough training examples in the domain, we trained the planner on *childsnack* with the same settings, but we started training from the 13th problem in the training set. As the planner now has more information regarding the domain, we immediately saw an improvement in the quality of the resulting planner, as shown in Table 1. Additionally, we note that the original implementation of ASNets (Toyer et al. 2020) is optimized for multi-core processing. Unfortunately, since we only had a single-core CPU in this competition, the learner was operating in a sub-optimal situation, adversely impacting the training process and making it notably slower.

Overfitting training strategy As described in section 2.2, the training strategy used by the planner iteratively adds new problems to the training set. This approach ensures that the planner retains knowledge from simpler problems while also learning new knowledge from more complex ones. To check this, we validated the domain knowledge by solving training set problems using the learnt domain knowledge. And it turns out the planner can always solve all the trained problems, regardless of its performance in the testing set. This observation potentially indicates that the model is overfit-

ted. Besides, the training time per epoch increases as the total number of problems increases. As a result, when we retrain the model with a new problem and all previously seen problems, the trainer spent a significant amount of time overfitting to seen problems. To illustrate this point, we conducted an experiment in which we trained the planner on *blocksworld* using the same settings but extended the training time to 3 days. However, we could only increase the number of problems the planner was trained on from 30 to 34. This result demonstrated that overfitting was more significant than we had anticipated. Therefore, developing a better training strategy for the planner to work on most of the trainable problems is necessary.

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