

# Fast Downward Stone Soup 2023

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Fast Downward Stone Soup (Helmert, Röger, and Karpas 2011) is a sequential portfolio planner, built on top of the Fast Downward planning system (Helmert 2006, 2009). It participated in three previous International Planning Competitions (IPC): 2011 (Helmert et al. 2011), 2014 (Röger, Pommerening, and Seipp 2014), and 2018 (Seipp and Röger 2018). In the last IPC, Fast Downward Stone Soup was the winner of the satisficing and the cost-bounded tracks.

In this planner abstract, we present the Fast Downward Stone Soup portfolio submitted to the sequential optimal and satisficing tracks of IPC 2023. After two IPCs (2014 and 2018) without participating in the optimal track, this is the first time after 12 years that an optimal version of Fast Downward Stone Soup enters the competition. The procedure used for building the portfolios is the same as in 2011, 2014, and 2018. Therefore, we only briefly explain this procedure here and refer the reader to the original paper for more details (Helmert, Röger, and Karpas 2011). We highlight also the configurations used for the optimal track portfolio, and the new additions for the satisficing track.

## Building the Portfolios

The Stone Soup algorithm requires the following information as input:

- A set of *planning algorithms*  $\mathcal{A}$ . We use a different set of Fast Downward configurations depending on the track, which we describe below.
- A set of *training instances*  $\mathcal{I}$ , for which portfolio performance is optimized. We use a set of 7330 instances, described below.
- Complete *evaluation results* that include, for each algorithm  $A \in \mathcal{A}$  and training instance  $I \in \mathcal{I}$ ,
  - the *runtime*  $t(A, I)$  of the given algorithm on the given training instance on our evaluation machines, in seconds (we did not consider anytime planners), and
  - the *plan cost*  $c(A, I)$  of the plan that was found.

We use time and memory limits of 10 minutes and 8 GiB to generate this data. If algorithm  $A$  fails to solve instance  $I$  within these bounds, we set  $t(A, I) = c(A, I) = \infty$ .

The procedure computes a portfolio as a mapping  $P : \mathcal{A} \rightarrow \mathbb{N}_0$  which assigns a time limit (possibly 0 if the algorithm is not used) to each component algorithm. It is a

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build-portfolio(algorithms, results, granularity, timeout):  
  portfolio := { $A \mapsto 0 \mid A \in \text{algorithms}$ }  
  repeat [timeout/granularity] times:  
    candidates := successors(portfolio, granularity)  
    portfolio :=  $\arg \max_{C \in \text{candidates}}$  score( $C$ , results)  
  portfolio := reduce(portfolio, results)  
  return portfolio
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Figure 1: Stone Soup algorithm for building a portfolio.

simple hill-climbing search in the space of portfolios, shown in Figure 1.

In addition to the algorithms and the evaluation results, the algorithm takes two parameters, *granularity* and *timeout*, both measured in seconds. The timeout is an upper bound on the total time for the generated portfolio, which is the sum of all component time limits. The granularity specifies the step size with which we add time slices to the current portfolio.

The search starts from a portfolio that assigns a time limit of 0 seconds to all algorithms. In each hill-climbing step, it generates all possible *successors* of the current portfolio. There is one successor per algorithm  $A$ , where the only difference between the current portfolio and the successor is that the time limit of  $A$  is increased by the given granularity.

We evaluate the quality of a portfolio  $P$  by computing its *portfolio score*  $s(P)$ . The portfolio score is the sum of *instance scores*  $s(P, I)$  over all instances  $I \in \mathcal{I}$ . The function  $s(P, I)$  is similar to the scoring function used for the International Planning Competitions since 2008. The only difference is that we use the best solution quality among our algorithms as reference quality (instead of taking solutions from other planners into account): if no algorithm in a portfolio  $P$  solves an instance  $I$  within its allotted runtime, we set  $s(P, I) = 0$ . Otherwise,  $s(P, I) = c_I^*/c_I^P$ , where  $c_I^*$  is the lowest solution cost for  $I$  of any input algorithm  $A \in \mathcal{A}$  and  $c_I^P$  denotes the best solution cost among all algorithms  $A \in \mathcal{A}$  that solve the instance within their allotted runtime  $P(A)$ .

In each hill-climbing step the search chooses the successor with the highest portfolio score. Ties are broken in favor of successors that increase the timeout of the component algorithm that occurs earliest in some arbitrary total order.

The hill-climbing phase ends when all successors would

exceed the given time bound. A post-processing step reduces the time assigned to each algorithm by the portfolio. It considers the algorithms in the same arbitrary order used for breaking ties in the hill-climbing phase and sets their time limit to the lowest value that would still lead to the same portfolio score.

## Training Benchmark Set

As benchmarks, we used all tasks and domains from previous IPCs, from Delfi (Katz et al. 2018), and from the 22.03 Autoscale collection Torralba, Seipp, and Sievers (2021), leading to a set of 92 domains with 7330 tasks. We used Downward Lab (Seipp et al. 2017) to run all planners on all benchmarks on Intel Xeon Silver 4114 2.2 GHz processors, imposing a memory limit of 8 GiB and a time limit of 30 minutes for optimal planners and 5 minutes for satisficing and agile planners. For each run, we stored its outcome (plan found, out of memory, out of time, task not supported by planner, error), the execution time, the maximum resident memory, and if the run found a plan, the plan length and plan cost. This data set is online available<sup>1</sup>. As training data for our optimal (satisficing) portfolio, we selected from each domain the 30 tasks which are solved by the fewest optimal (satisficing) planners, which results in 2377 remaining tasks.

## Planning Algorithms

### Satisficing Track

For the satisficing track, we collect our input planning algorithms from several sources. First, we use the component algorithms of the previous Fast Downward Stone Soup portfolios that participated in the sequential satisficing track of other IPCs (Helmert et al. 2011; Röger, Pommerening, and Seipp 2014; Seipp and Röger 2018). In total, this gives us 87 different configurations.

Second, we add different combinations of  $h^{cea}$  (Helmert and Geffner 2008),  $h^{CG}$  (Helmert 2006),  $h^{FF}$  (Hoffmann and Nebel 2001), and  $h^{LM}$  (Richter, Westphal, and Helmert 2011) with greedy best-first search (GBFS) but using different open-lists:

- For each one heuristic  $h \in \{h^{cea}, h^{CG}, h^{FF}\}$ , we use one configuration of (eager) GBFS with the  $\epsilon$ -greedy open-list (Röger and Helmert 2010) ordered by  $h$ . We also add the same configurations but using lazy GBFS (Richter and Helmert 2009).
- We use configurations using lazy GBFS alternating between three open-lists:  $[h^{FF}, h, \text{PAR}(h^{FF}, h)]$  where  $h \in \{h^{cea}, h^{CG}, h^{LM}\}$  and  $\text{PAR}(h^{FF}, h)$  is a Pareto open-list using  $h^{FF}$  and  $h$  (Röger and Helmert 2010).
- We add two other configurations using lazy GBFS: one alternating between  $[\epsilon\text{-greedy}(h^{FF}), h^{LM}]$  open-lists, and one alternating between  $[h^{FF}, \epsilon\text{-greedy}(h^{LM})]$ . For each of these two configurations, we also include configurations with additional open-lists only containing states generated by preferred operators.

<sup>1</sup>URL to be published

All configurations using  $h^{LM}$  have one version using reasonable orders and one without them.

This makes for a total of 18 new configurations. Overall, this leaves us with 105 planner configurations as input for the hill-climbing procedure.

### Optimal Track

For the optimal track, we distinguish whether planning algorithms support conditional effects or not. We use  $A^*$  with the following heuristics without support for conditional effects:

- the blind heuristic
- BJOLP (Domshlak et al. 2011)
- Cartesian abstractions (Seipp and Helmert 2018):
  - for subtasks induced by goals and fact landmarks (one of four different limits: 10s, 60s, 300s generation time or 1 million state-changing transitions in all abstractions)
  - for subtasks only induced by goals (1 million transitions)
  - for subtasks only induced by fact landmarks (1 million transitions)
- $h^2$  (Haslum and Geffner 2000)
- $h^{\max}$  (Bonet and Geffner 2001)
- pattern databases (PDBs) (Culberson and Schaeffer 1998; Edelkamp 2001), all combined in the canonical PDB heuristic (Haslum et al. 2007), computed with the following pattern collections:
  - CEGAR with a maximum size of 1 million states in individual PDBs and 10 million states in all PDBs, computation time limit of 10s/60s/300s with enabling stagnation after 2s/12s/20s, enabling blacklisting after 75% of the computation time limit or on stagnation, computing wildcard plans (Rovner, Sievers, and Helmert 2019)
  - hill climbing (Haslum et al. 2007) (thus leading to iPDB) with a computation time limit for the main loop of the algorithm of 10s/60s/300s
  - interesting patterns of size 1/2/3 (Pommerening, Röger, and Helmert 2013)
- LM-cut (Helmert and Domshlak 2009)
- merge-and-shrink heuristics (Helmert et al. 2014; Sievers and Helmert 2021) with bisimulation-based shrinking (Nissim, Hoffmann, and Helmert 2011) and a size limit of 50000 states per abstraction, exact label reduction (Sievers, Wehrle, and Helmert 2014), one of two merge strategies: SCC-DFP or SCC-sbMIASM (Sievers, Wehrle, and Helmert 2016), and a computation time limit for the main loop of the algorithm of 10s/60s/300s
- operator-counting heuristics (Pommerening et al. 2014) with different types of constraints:
  - post-hoc optimization constraints over interesting patterns (size 1/2/3) (Pommerening, Röger, and Helmert 2013)
  - delete relaxation constraints, leading to  $h^+$  (exact IP model and LP relaxation) (Imai and Fukunaga 2015)

- state equation constraints (Pommerening et al. 2014)
- state equation constraints and LM-cut constraints (Pommerening et al. 2014)
- state equation constraints, LM-cut constraints, and delete relaxation constraints (LP relaxation) (Pommerening et al. 2014)
- diverse potential heuristics and potential heuristics optimized for the initial state or all states (Seipp, Pommerening, and Helmert 2015)

This makes for a total of 37 planning algorithms. All of them prune operators during successor generation using atom-centric stubborn sets (Röger et al. 2020). However, this feature is turned off after the first 1000 expansions if less than 20% of operators are pruned.

For tasks with conditional effects, we cannot use stubborn sets pruning and we only use those of the above algorithms that support conditional effects: blind,  $h^2$ ,  $h^{\max}$ , all merge-and-shrink variants. This is a total of 11 planning algorithms.

### Resulting Portfolios

The resulting satisficing portfolio uses 41 of the 144 possible algorithms, running them for 8–135 seconds. On the training set, the portfolio achieves an overall score of 1999.93, which is much better than the best component algorithm with a score of 1650.40. If we had an oracle to select the best algorithm (getting allotted the full 1800 seconds) for each instance, we could reach a total score of 2073.

### Executing Sequential Portfolios

In the previous sections, we assumed that a portfolio simply assigns a runtime to each algorithm, leaving their sequential order unspecified. With the simplifying assumption that all planner runs use the full assigned time and do not communicate information, the order is indeed irrelevant. In reality the situation is more complex.

First, the Fast Downward planner uses a preprocessing phase that we need to run once before we start the portfolio, so we do not have the full 1800 seconds available.<sup>2</sup> Therefore, we treat per-algorithm time limits defined by the portfolio as relative, rather than absolute values: whenever we start an algorithm, we compute the total allotted time of this and all following algorithms and scale it to the actually remaining computation time. We then assign the respective scaled time to the run. As a result, the last algorithm is allowed to use all of the remaining time.

Second, in the satisficing setting we would like to use the cost of a plan found by one algorithm to prune the search of subsequent planner runs (in the optimal track we stop after finding the first plan). We therefore use the best solution found so far for pruning based on  $g$  values: only paths in the state space that are cheaper than the best solution found so far are pursued.

Third, planner runs often terminate early, e.g., because they run out of memory or find a plan. Since we would like

<sup>2</sup>The preprocessing phase consists of converting the input PDDL task (Fox and Long 2003) into a SAS<sup>+</sup> task (Bäckström and Nebel 1995) with the Fast Downward translator component.

to use the remaining time to continue the search for a plan or improve the solution quality, we sort the algorithms by their coverage scores in decreasing order, hence beginning with algorithms likely to succeed quickly.

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For a portfolio planner, not those who *combined* the components deserve the main credit but those who *contributed* them. We therefore wish to thank all Fast Downward contributors and the people who came up with the algorithms we use in our portfolio.

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