

Ninjas' Revenge: the secret genetic technique

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Abstract

Knapsack problems are centered around finding an optimal distribution of values within a certain number of containers or bags, so to speak. They are classified as NP-hard and have been the subject of interest for many studies, where different kinds of solutions have been proposed. In this report we detail our own solution for a particular instance of a knapsack problem, where we make use of genetic algorithms to find the optimal assignment of multiple items. We compared this approach against more traditional exhaustive search methods, and our experiments showed that the genetic algorithm was capable of obtaining optimal results in reasonable times as long as the number of items was not exceedingly large.

Keywords: Knapsack problem, packing problems, genetic algorithms

1. Problem statement and goals

One of the most popular problems studied in the area of constrained combinatorial optimization is the knapsack problem. Given a finite set of objects with associated weights and values, the objective is to maximize the value of a collection formed by these items without exceeding a predefined weight limit.

The knapsack problem has diverse practical applications which makes it particularly interesting. For example, it has been applied to production and inventory management [Ziegler \(1982\)](#), financial models [Mathur et al. \(1983\)](#) and queueing operations in computer systems [Gerla and Kleinrock \(1977\)](#) or even manufacturing [Bitran and Tirupati \(1989\)](#).

There are several variations to the knapsack problem. We could consider there are multiple copies of each item or take into consideration their volume in addition to their weight. However, we will focus on the simplest case: the one dimensional 0-1 knapsack problem, where the only constraint is the weight and the number of copies of each item is limited to 1.

The premise for this particular problem is that a group of 18 ninjas must sneak into a palace and carefully choose which items they should put in their bag, the goal here being to maximize the value of the bag's contents. Although it might seem silly on the surface, this is actually an NP-hard problem. Thus, no known algorithm achieves optimal solutions in

polynomial time for all cases of this problem. Still, a sufficiently good solution can be found quickly by resorting to heuristic methods. The one that we will explore is an effective and commonly used metaheuristic known as “genetic algorithm”.

Genetic algorithms (GAs) belong to the family of techniques known as evolutionary algorithms. They draw inspiration from natural selection and genetic processes to provide solutions to complex optimization problems and model evolutionary systems.

The main process can be described as an evolutionary cycle. They first initialise a population of chromosomes, which represent candidate solutions. Some of these individuals will be drawn from the population through a selection mechanism and compared according to their fitness. In the reproduction phase, the genetic material of the best individuals will be combined to form offspring. Also, some values of the offspring’s chromosomes will be mutated. Finally, a replacement strategy is set to generate the next generation population. This process is repeated for several generations or until some convergence criterion is met.

We can describe formally the 0-1 knapsack problem as follows: given a set of n items and a *knapsack* with:

$$\begin{aligned} p_j &= \text{profit of item } j, \\ w_j &= \text{weight of item } j \\ c &= \text{capacity of the knapsack,} \end{aligned}$$

we will select a subset of items so as to obtain:

$$\begin{aligned} &\text{maximize } z = \sum_{j=1}^n p_j x_j \\ &\text{subject to } \sum_{j=1}^n w_j x_j < c, \end{aligned}$$

where we have that:

$$x_j = \begin{cases} 1 & \text{if item } j \text{ is selected;} \\ 0 & \text{otherwise} \end{cases}, \quad j \in N = \{1, \dots, n\}$$

A solution to the knapsack problem can be represented as a sequence of binary values $[x_1, x_2, \dots, x_n]$. We will apply GAs to the eighteen different cases, each of them specified by a file containing a value for n and c in the first line and values for p_j and w_j in subsequent lines for each object x_j .

The goal of this project is to find the optimal solution to each of the eighteen knapsack problems proposed by means of applying a GA approach. To this end, we will have to choose and implement suitable mechanisms in each of the phases of the evolutionary cycle. The description of these methods, as well as the reasons for their election, are detailed in Section 3.

2. Previous work

In the last few decades, there has been an impressive amount of research into knapsack problems as well as bin-packing problems, the latter being an special case where the number

of containers is not fixed but rather the goal is to minimize them instead. This interest stems mainly from their simple premise and structure, which are deceptively complex and allow for many different approaches, some more efficient than others naturally.

Martello et al. (1997) [Martello et al. \(2000\)](#) explore different kinds of exact algorithms that could be used to solve knapsack problems, mainly focused on the combination of branch-and-bound algorithms and dynamic programming. Moreover, the application of genetic algorithms to these types of problems has also proved to be particularly promising. Falkenauer (1998) [Falkenauer \(1998\)](#) describes a Grouping Genetic Algorithm to solve bin-packing problems with great success, where its performance was superior to previous approaches and could be extended to other types of grouping problems. Mohamadi (2010) [Mohamadi \(2010\)](#) on the other hand proposes an improved representation scheme and genetic algorithm to solve a set of 100 sample problems, and the results attest to its efficiency.

All in all, previous research shows that there is merit in using genetic algorithms to solve packing problems, and their performance may even surpass other tried and tested methods based on more traditional techniques.

3. Implementation details

There are several aspects to consider when applying genetic algorithms to solve a problem, each one with its own methods. In this section, we define the different methods used in our implementation for each one of these aspects.

3.1. Population initialization

In order to initialize the population, we used a random initialization approach that has been traditionally used in GAs for its simplicity and efficiency. However, we also introduced a parameter for the initialization range to give the algorithm better chances of finding the optimal solution, in case we already know where this solution should be. This initialization range, in conjunction with the possibility of sorting the input items by their value/weight ratio, can highly improve the algorithm's performance.

3.2. Fitness function and selection

To be able to assess how well an individual (i.e. possible solution) fits our objective we must first define the fitness function. For our particular problem, we defined the following fitness function:

$$F(x) = \begin{cases} -\infty & \text{if } x \bullet w > c \\ x \bullet v & \text{otherwise} \end{cases}$$

where x is the bitstring representing the individual, w is the list of item weights, v is the list of item values and c is the knapsack capacity.

Once we have the fitness function we are able to proceed to the selection step. We decided to implement the tournament and elitism selection methods for our algorithm. Particularly, we use a fixed value of $k = 2$ for both tournament and elitism. We decided to use tournament

selection instead of other methods such as standard roulette wheel or rank selection because of its simplicity and because seems to outperform the other approaches [Razali et al. \(2011\)](#). On the other hand, elitism was implemented because of its usage in the tournament selection method.

3.3. Crossover and mutation

Several crossover techniques can be applied in GAs such as one-point, 2-point, multi-point (with more than two cut points) and uniform crossovers. In this work, we have implemented the one-point, 2-point and uniform crossovers. The reason behind this choice is the simplicity of these techniques and their good performance, especially the 2-point crossover [Hasançebi and Erbatur \(2000\)](#); [DeJong and Spears \(1990\)](#); [Adeli and Cheng \(1993\)](#).

For the mutation step, we implemented a simple bit-flip where a random number of genes between 1 and the size of the chromosome are flipped. In addition, we set the mutation probability to 0.5 since lower values made the algorithm get stuck in local minimums more frequently.

3.4. Replacement strategy

The final step of the GA is the replacement strategy used to obtain the new generation. In our case, we used elitism to replace the worst individuals of the current population. With this strategy, the old generation might be entirely replaced. In addition to the elitism, we also decided to remove repeated individuals before applying the replacement. This decision was made to ensure that the diversity of the population is maintained. Additionally, we defined an extra case that does not use elitism. This is when there is no good solution (i.e. the fitness value is $-\infty$) neither in the parents nor the offspring. In this case, we decided to obtain the new generation by selecting, randomly, half of the individuals from μ and the other half from λ .

4. Results and Discussion

The main objective of this project was to implement a genetic algorithm that was able to achieve optimal, or near-optimal, solutions for the knapsack problem in a reasonable amount of time when compared with the exhaustive search method. To achieve this, we tried different methodologies such as sorting the items by their value/weight ratio or by using different crossover methods.

First, we started by comparing the three crossover methods implemented (i.e. one-point, two-point and uniform). In [Table 1](#), we can see the results obtained with each method in three different problem sizes. We can see how the worst method, in terms of finding the solution, is the one-point crossover since it is only able to find it for small sizes of the problem. The uniform crossover seems to outperform the two-point crossover method both in execution time and the number of times the best solution is found for small and medium-size problems. However, the two-point method is the only one able to find the best solution for the three files. Although these results are not concluding since we have not tested enough sizes, they might indicate that for small and medium-sized problems the usage of uniform

crossover should be preferred over the other two options, while the two-point method should be used when we have a greater problem size.

Sol	#items	One-point		Two-point		Uniform	
		Time (s)	Best found	Time (s)	Best found	Time (s)	Best found
6	50	56.206	1/3	76.890	1/3	65,544	2/3
9	100	131.852	0/3	109.723	1/3	43.720	3/3
14	300	114.531	0/3	270.457	1/3	121.396	0/3

Table 1: Comparison between implemented crossover methods using the sort option, 10000 max generations, 600 stall generations, a population size of $\max(\#items^2, 2000)$ and tournament selection.

Once we tested the different crossover methods, we proceeded to compare the algorithm without sorting the items by value/weight rate and with the sorting. To be able to compare them we fixed the set of parameters to avoid biasing the results as much as possible.

Looking at Table 2, it is possible to observe the following:

1. It was possible to get the optimal value for $\sim 89\%$ of the files tried.
 - (a) For sol_17 and sol_18 we obtained sub-optimal solutions close to the optimal solution but, due to the restriction of stall generations, the algorithm was not able to produce better results.
2. There is a strong correlation between the number of items and the time it takes to find the optimal solution. That is, a Pearson correlation of 0.9873 for the runs without sort and 0.9979 for the runs with sort. However, certain files do not follow this trend. For instance, if we compare the results from sol_12 and sol_3 we see how the problem with more items (i.e. sol_12) needs more or less the same time as the solution with fewer items (i.e. sol_3) and even finds the optimal with more frequency. This is a behaviour that we would usually not expect.
3. Sorting the items using their value/weight ratio helped to improve the time required to find the optimal solution significantly. In some cases, the improvement achieved is ~ 56 times faster than the not sorted result. The mean speedup is around ~ 6 times.

It is interesting to find that sorting the items by their value/weight ratio yielded excellent results. We think the reason for this is that sorting the items ends up placing the best ratios in the lower section of the bitstring (i.e. the bits belonging to the right half). If we combine this with a relatively small population range, it is likely to get the first generation with a high fitness value and it is even possible to find the optimal solution among the initial generation. In the reproduction step, having this configuration helps next generations to improve their fitness score faster. This can be observed in the time required to find the solution in all the files tried. There might be some cases where the best solution is found by combining low value/weight ratio items or that there is not much difference in ratio between items. This could cause the sorting to have a lower impact on the final performance but it is still higher than the no sorting option.

Sol	#items	Best value	Best value found	Time without sort (s)	Best found	Time with sort (s)	Best found
1	4	19	19	0.0378	4/4	0.0375	4/4
2	19	12248	12248	2.445	4/4	0.430	4/4
3	30	99798	99798	34.501	0/3	24.437	2/3
4	40	99924	99924	4.382	3/3	0.674	3/3
5	45	23974	23974	117.021	1/3	2.102	3/3
6	50	142156	142156	247.373	0/3	76.890	1/3
7	50	5345	5345	73.624	3/3	18.885	3/3
8	60	99837	99837	12.803	3/3	34.425	3/3
9	100	99837	99837	94.960	0/3	109.723	1/3
10	100	1333930	1333930	321.234	0/2	23.198	2/2
11	100	10892	10892	322.240	1/2	128.105	1/2
12	200	100236	100236	49.784	2/2	24.183	2/2
13	200	1103604	1103604	632.185	0/2	512.575	1/2
14	300	1688692	1688692	417.575	0/2	196.617	1/2
15	400	3967180	3967180	216.064	0/2	258.675	1/2
16	500	54939	54939	614.917	0/2	191.009	1/2
17	1000	109899	109894	699.159	0/2	643.658	0/2
18	10000	1099893	1099132	4548.846	0/1	7,208.676	0/1

Table 2: Comparison between the algorithm with the sort option and without it using 10000 max generations, 600 stall generations, a population size of $\max(\#items^2, 2000)$, tournament selection and two-point crossover.

Regarding the evolution of the fitness value throughout the different generations, most of the plots obtained (can be found within the project folder) share the same trend. The fitness tends to flatten at the beginning after a few epochs/generations when a high value is reached. The flattening happens because we keep the best value found in subsequent iterations. However, we think that because of the mutation process that happens in each generation, the GA can break out and find even higher fitness values. This can be observed in the graphics by the layered shape obtained. As an example, Figure 1 shows the fitness progression for the last file `ninja_18_10000`, where 10000 items were used.

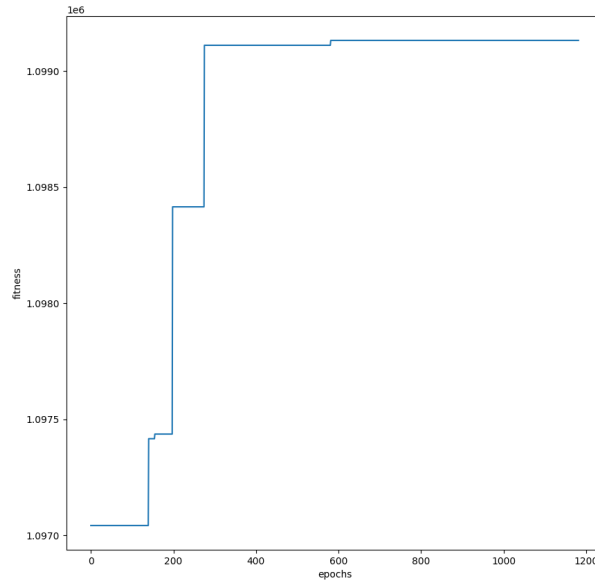


Figure 1: Evolution of the fitness value for the `ninja_18_10000` file using the sort option, 10000 max generations, 600 stall generations, a population size of 2000, tournament selection and two-point crossover.

5. Strengths and weaknesses

After seeing the results obtained from our GA implementation we have detected some strengths and weaknesses in it.

Starting with its strengths, we have seen how the algorithm can find the optimal solution in a reasonable amount of time in most cases. This is a very good outcome since we would usually expect to obtain sub-optimal results. Also, we noticed that the algorithm can find multiple optimal solutions for some of the problems. Being able to find several optimal solutions to one problem opens the possibility to introducing new factors when deciding

which one best suits us. This can be considered a strength compared to the standard exhaustive search algorithm since the latter is only able to find one possible solution. Another strong point of our solution is its simplicity, since it does not require complex functions or procedures to achieve good results. We consider that this makes the algorithm easy to understand and thus adaptable to other problems that can be represented as bitstrings.

Following with its weaknesses, we have seen how the algorithm struggles when the size of the problem increases dramatically (e.g. 1000 or 10000 items). In these cases, we observed how the exhaustive search outperformed our algorithm in terms of solution quality and time complexity since it could find the optimal solution in less time. Also, for some problems, we observed how successive executions performed considerably different when looking at the execution time and the solution obtained. From this, we infer that the GA is highly dependent on the initialization of the initial population.

6. Conclusions and future work

The genetic algorithm implemented has proven to be a good heuristic for generating approximate solutions of the unidimensional 0-1 knapsack problem. The results achieved were optimal or near-optimal in the eighteen scenarios studied despite the imposition of stopping criteria.

We also noted how the performance of the system improves drastically when the data is preprocessed suitably. Having a good understanding of the problem representation and the internal mechanisms in a GA is essential for this preliminary step.

Another aspect that determines the performance of the system is of course the strategies implemented in each phase of the evolutionary cycle. We relied on several methods to choose from at each step which makes these systems very flexible and adaptable. This in turn, along with a proper selection of parameters, allowed us to tailor the algorithm to the problem.

Nonetheless, there are ways to further optimize our solution. As mentioned earlier, GAs are highly sensitive to the population seeding technique. The random initialization of the population can generate individuals with poor fitness, which increases the time needed to achieve optimality. Instead, we could resort to better seeding techniques such as Sorted Population [Yugay et al. \(2008\)](#), which creates a large initial population, sorts it in ascending order according to their fitness and chooses a percentage of individuals that have above average fitness. This and other initialization techniques like Nearest Neighbour or Gene Bank were described and compared by [Hassanat et al. \(2018\)](#).

As we have seen in Section 5, one of the main weaknesses of our algorithm is that its execution time increases dramatically with the problem size. One way to mitigate this issue would be to parallelize the reproduction process. In this way, we could randomly split the population into multiple batches and then perform the reproduction step simultaneously for each batch. Given that, we only would need to wait for the slowest batch to finish and then join all the resulting offspring together to perform the replacement step.

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