Genetic Algorithms

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October 6, 2013

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1 Introduction

Genetic algorithms are a biologically-inspired approach to heuristic search which mimic natural selection. Unlike many other evolutionary strategies and evolutionary programming, they are not designed to solve a specific problem, but are designed to solve the problem of optimisation which is made difficult by substantial complexity and uncertainty[2].

The complexity of the task should make it such that discovering an optimum solution is a long, maybe even impossible, task. At the same time the uncertainty needs to be reduced so that the knowledge of *available* options can be increased.

The initial design for a genetic algorithm was a method for moving from one population of chromosomes to another using a form of natural selection. This algorithm also included methods for crossover, mutation and inversion. This idea of having a large population was the distinguishing feature from any past attempts which had only considered the parent and one offspring, where the offspring was simply a mutation of the parent[3].

1.1 Evolutionary Algorithms

As their name suggests, an evolutionary algorithm applies elements from the biological theory of evolution to the problem of optimisation. These elements include:

- Reproduction
- Mutation
- Recombination
- Selection

Typically, a population of candidate solutions are generated to which a fitness function can be applied. The population is then subject to some form of evolution, and this process is repeated until a halting criteria is met.

Genetic algorithms are a type of evolutionary algorithm with a focus on the genetic evolution of solutions. Candidate solutions for genetic algorithms, known as *chromosomes* are encoded as a series of *genes*. These genes are a representation of the choices which need to be optimised for the solution and can be as simple as a single bit or as complex as a real number, depending on the problem.

There are many other forms of both evolutionary and genetic algorithms which this report with mention in later sections.

2 Basic Genetic Algorithm Principals

The basic principals of genetic algorithms are to represent candidate solutions as a population of chromosomes, from this population the fittest members can be picked out and used in the next generation and to create new member of the population through reproduction and/or mutation.

This cycle repeats with the aim to produce better performing individuals in each generation until the optimum solution is either reached or gotten close enough to that any future improvement is unnecessary or unwanted due to other constraints such as processing time. The latter of these allows a genetic algorithm to come up with a "good" solution in a reasonable amount of time.

Reproduction and mutation are an important part of genetic programming, and too of evolutionary programming. Without these parts the algorithm would quickly reach a local optimum for the initial population and would not improve past this.

2.1 Chromosome Representation

One of the key parts parts in implementing a genetic algorithm is the representation of chromosomes. This is very dependent of the problem the genetic algorithm needs to optimise and can have a knock on affect on the efficiency and accuracy of the algorithm.

Sometimes a simple solution is enough to represent the problem, binary strings are a commonly suggested approach. However sometimes more complex representations are required, potentially any data structure can be used as a chromosome but lists and trees are the common choices as they are easy to perform crossover¹ and mutation on.

As a very simple example, to maximise y in: y = f(x), one could represent the value of x as a binary string, an example of which is shown in figure 1.

00101010

Figure 1: Chromosome representation as a binary string

2.2 Fitness Function

To fulfil the step of natural selection; the process of choosing the "best" members of a population, there needs to be a way of evaluating each chromosome, such that they can be compared to one another.

The function for doing so is known as a fitness function, which typically returns either a single number or a list of numbers, depending on the problem. Each chromosome can then be ranked in order of fitness and the top members of a population can then be selected.

As the value returned from a fitness function, with be specific to the domain it has be used within, it is necessary to rescale the fitness value to ensure uniformity when genetic algorithms are applied over several different domains at the same time.

¹A term used instead of reproduction in genetic algorithms.

2.3 Selection

Selection simulates the "survival of the fittest" nature of biology. However, it is beneficial not to keep some lesser performing members of the population to avoid getting trapped in local optimum. Most selection algorithms will introduce an element of randomness into the selection to deal with this.

This report will discuss the effects of different selection algorithms in Section 3.

- 2.4 Crossover
- 2.5 Mutation
- 2.6 Termination

3 The Effects of Selection Algorithms on GAs

The selection of the "best" individuals is an important step in genetic algorithms and as such several selection algorithms have been put forward.

The simplest is to just pick the top individuals of a generation to carry on their genes in the next generation, either through directly copying them, crossover or mutation.

This has its flaws and doesn't mirror the biological process of evolution. The main flaw of this approach is that it can be beneficial to include weaker members of a population to avoid getting stuck in local optimum. By including other, seemingly worse, members of a population, the effects of crossover and mutation can overcome this problem.

Another, commonly used, technique is to use a tournament-based selection algorithm, which involves selecting the best member from a pool made up of random members of the population. The size of the pool has a big effect on the selection of individuals, and is generally inversely proportional to the number of weaker members selected.

Tournament selection is one of the closest to biological natural selection, and has the benefits that it is efficient, especially on parallel architectures.

Another popular selection algorithm is roulette-wheel selection, where the chance a particular individual has of being selected is proportional to its fitness. Those with a higher fitness have a higher probability of being selected. The equation used to calculate the probability is shown in equation 1, where i is the individual in question, f_x is the fitness of individual x and N is the number of individuals in the population.

$$p_i = \frac{f_i}{\sum_{j=1}^N f_j} \tag{1}$$

There have been studies into the effects of different selection algorithms on genetic algorithms. Goldberg and Deb[1] compared four methods of selection; proportional reproduction, ranking selection, tournament selection and the GENITOR algorithm.

They analysed each selection scheme in terms of:

- 1. growth ratio; the expected ratio for the members of the best class to the umber of members of the population.
- 2. Takeover time; the approximate number of generations it takes to converge, this gives an idea of how long it would take before mutation, rather than crossover, becomes the primary method for exploring the search space.
- 3. Time complexity; the standard big O measure for time complexity.

It was found that linear ranking and binary tournament (a tournament of only two individuals) selection had similar growth rates, but that tournament selection with larger pool sizes had higher growth rates, though a non-linear ranking function can also achieve this.

Takeover times all converged in around $O(\log n)$ generations. Most interestingly the time complexities had a large range, from $O(n^2)$ to O(n), with

tournament selection being an O(n). This is particularly interesting as tournament selection is a very easy algorithm to make parallel, this fits in with genetic algorithms nicely as they improve drastically from parallel implementation.

These results should be taken with some scepticism as they are formulated from the pure mathematics of the selection techniques and not from running experiments and the authors even mention that it is only designed as a simple method to better understand the *expected* behaviour and suggest looking at better methods, citing the inherently noisy nature of genetic algorithms.

3.1 Genetic Drift

Rogers and Prügel-Bennett[4] defined a method of analysing selection schemes using genetic drift, a term borrowed from biology, which describes the change in frequency of an allele (gene variation) through random sampling of the population.

Genetic drift is a phenomenon observed in genetic algorithms due to the nature of selection and, unlike analysis methods like convergence time, leads to a exact analytical solution. Though previous attempts to calculate genetic drift were often approximations or difficult to generalise to other cases, according to the authors.

In genetic algorithms, genetic drift is a measure of the change in fitness variation within a population and is worked out using probabilistic analysis of the selection scheme. The inherent problem with this is that genetic algorithms are a form of subsymbolic learning, which is very difficult to apply probabilistic methods to, due to the randomness used within them.

The authors attempt to deal this by using results from experiments averaged over 100,000 runs. Again though, a lot will depend on the nature of the problem, the representation of chromosomes and even the way in which the selection scheme is implemented. Like with the work of Goldberg and Deb, this only provides an indication of the performance of a selection scheme, rather than a full view. It does, however, seem to give some useful insights into different selection schemes and may help rule out less useful selection schemes.

4 Code Example

A simple python implementation is shown in figure 1.

```
import random, math, time
from graph import Graph
MUTATION = 0.001
CROSSOVER = 0.6
class Chromosome:
  @ class method
  def create(cls, graph):
     genes = [i for i in xrange(graph.num_nodes())]
     random.shuffle(genes)
     return Chromosome(genes, graph)
  def __init__(self , genes , graph):
     while not self.valid(genes):
       genes = self.make_valid(genes)
     self.genes = genes
     self.graph = graph
  def valid(self, genes):
     for i in xrange(len(genes)):
       if i not in genes:
          return False
     return True
   \begin{array}{l} def \ make\_valid(self \,, \ genes) \colon \\ missing = \left[ \, i \ for \ i \ in \ xrange(len(genes)) \ if \ i \ not \ in \ genes \right] \\ double = \left[ \, i \ for \ i \ in \ xrange(len(genes)) \ if \ genes.count(i) > 1 \right] \\ \end{array} 
     random.shuffle(missing)
     random.shuffle(double)
     for i in double:
       genes [genes.index(i)] = missing.pop()
     return genes
  def = str_{-}(self):
     return "{} - {}".format(self.genes, Chromosome.fitness(self))
  def __repr__(self):
     return self._str_()
  @classmethod
  def fitness(cls, c):
     return sum ([graph.distance(n1, n2)
                    for (n1, n2) in zip(c.genes, c.genes[1:])])
class BasicGA(object):
  def __init__(self , graph , population):
     self.graph = graph
     self.pop_size = len(population)
     self.population = population
     {\tt self.mutants\_rate} \ = \ {\tt int} \, (\, {\tt math.ceil} \, (\, {\tt self.pop\_size} \ * \ {\tt MUTATION}) \, )
     self.crossover_rate = int(math.ceil(self.pop_size * CROSSOVER))
  def run(self):
     """Runs the GA once."""
     best = self.select()
```

```
offspring = self.crossover(best)
    mutants = self.mutate(best)
    cur_best = self.population[0]
    self.population = best + offspring + mutants
    return cur_best
  def select(self):
    """ Selects the best individuals from a population."""
    retained_pop = self.pop_size - self.mutants_rate - self.
    crossover_rate
    return sorted (self.population, key=Chromosome.fitness)[0:
    retained_pop]
  def crossover(self, population):
    """ Performs crossover on a certain amount of the population."""
    return [self.perform_crossover(population)
             for i in xrange(self.crossover_rate)]
  {\tt def} perform_crossover(self, population):
    Perform crossover by:
    - Choosing a crossover point randomly.
    - Choosing two parents.
    - Slicing the genes of both parents together at the crossover
    point.
    point = int(random.random() * self.graph.num_nodes())
    p1 = random.choice(population)
    p2 = random.choice(population)
    return Chromosome(p1.genes[:point] + p2.genes[point:], self.
    graph)
  def mutate(self, population):
    """ Performs mutation of a certain number of the population."""
    return [self.perform_mutation(population)
             for i in xrange(self.mutants_rate)]
  {\color{red} \textbf{def} \hspace{0.1cm}} perform\_mutation \hspace{0.1cm} (\hspace{0.1cm} \hspace{0.1cm} self \hspace{0.1cm}, \hspace{0.1cm} population \hspace{0.1cm} ):
   """Randomly swap two genes to mutate the chromosome."""
   target = random.choice(population)
   genes = [i for i in target.genes]
   pos1 = int(random.random() * len(genes))
   pos2 = int(random.random() * len(genes))
   temp = genes[pos1]
   genes[pos1] = genes[pos2]
   genes [pos2] = temp
   return Chromosome (genes, self.graph)
class TournamentGA(BasicGA):
  def __init__(self , graph , num, tournament_size):
    super(TournamentGA, self).__init__(graph, num)
    {\tt self.tournament\_size} \ = \ tournament\_size
  def select(self):
    Selects the best members of the population using a tournament-
    based
    algorithm.
    retained\_pop = self.pop\_size - self.mutants\_rate - self.
    c \, rossover\_rate
    temp = self.population
```

```
best = []
     while len(best) != retained_pop:
       tournament = random.sample(temp, min(self.tournament_size,
     len(temp)))
       b = min(tournament, key=Chromosome.fitness)
       best.append(b)
     return best
if __name__ == "__main__":
  graph = Graph(25, 10, seed="genetic-algorithm")
  # Reset the random seed from graph creation.
  #random.seed()
  pop = [Chromosome.create(graph) for _ in xrange(100)]
  \begin{array}{l} {\rm gas} = \left[ {\rm BasicGA} \left( {\rm graph} \,, \, \, {\rm pop} \right), \,\, {\rm TournamentGA} \left( {\rm graph} \,, \, \, {\rm pop} \,, \,\, 2 \right), \\ {\rm TournamentGA} \left( {\rm graph} \,, \, \, {\rm pop} \,, \,\, 10 \right) \right] \end{array}
  for ga in gas:
    avg = []
     improving = True
     runs = 0
     while improving:
       runs += 1
       best = ga.run()
       avg.append(Chromosome.fitness(best))
       if len(avg) > 100:
          median = avg[len(avg)/2 + 1]
          improving = Chromosome. fitness (best) != median and len (avg)
     print "{} found best: {} in {} runs".format(ga.__class__.
     __name__ ,
                                                                str(best),
```

Listing 1: A Python implementation of a simple Genetic Algorithm

```
import random
class Node:
  def __init__(self , id , graph):
    self.x = int(random.random() * graph.size)
    self.y = int(random.random() * graph.size)
    self.id = id
  def distance(self, other):
    return pow(self.x - other.x, 2) + pow(self.y - other.y, 2)
  def = str_{-}(self):
    return "(\dot{i}, \dot{j})".format(self.x, self.y)
class Graph:
  nodes = None
  def __init__(self, nodes, size, **kwargs):
   if 'seed' in kwargs:
      random.seed(kwargs['seed'])
    self.size = size
    self.nodes = [Node(i, self) for i in xrange(nodes)]
  def distance(self, n1, n2):
    return self.nodes[n1].distance(self.nodes[n2])
  def num_nodes(self):
```

```
return len(self.nodes)

def __str__(self):
    s = ""
    for x in xrange(0, self.size):
        for y in xrange(0, self.size):
        added = False
        for node in self.nodes:
        if node.x == x and node.y == y:
            s += str(node.id)
            added = True
    if not added:
        s += " \n"
    s += "\n"
    return s
```

Listing 2: The Graph Code for listing 1

References

- [1] David E. Goldberg and Kalyanmoy Deb. A comparative analysis of selection schemes used in genetic algorithms. In *Foundations of Genetic Algorithms*, pages 69–93, 1991.
- [2] John H. Holland. Adaptation in natural and artificial systems: an introductory analysis with applications to biology, control, and artificial intelligence. MIT Press, April 1992.
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