

Chapter 7

Genetic Programming

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Abstract Welcome to genetic programming, where the forces of nature are used to automatically evolve computer programs. We give a flavour of where GP has been successfully applied (it is far too wide an area to cover everything) and interesting current and future research but start with a tutorial of how to get started and finish with common pitfalls to avoid.

7.1 Introduction

Getting computers to automatically solve problems is central to artificial intelligence, machine learning and the broad area covered by what Turing called “machine intelligence” [1]. As we shall show, this is what genetic programming is actually doing today.

Genetic programming [2] works by applying the power of evolution by natural selection [3] to artificial populations inside your computer, cf. Figure 7.1. Unlike in nature, you decide who is fit, who survives and who has children. Like nature, children are not identical to their parents but suffer random mutations and can be created by fusing together the genetic characteristics of their parents. Unlike other approaches to evolving expressions, genetic programming works because it has defined a way of representing expressions whereby they can be randomly mutated

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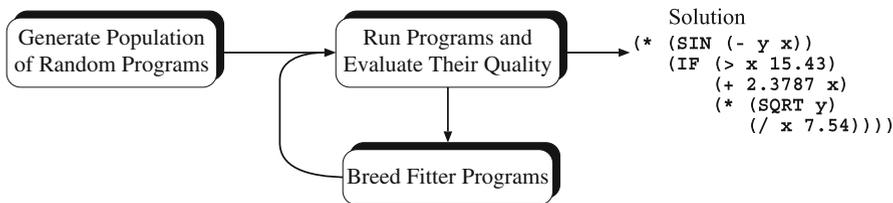


Fig. 7.1 The basic control flow for GP, where survival of the fittest is used to find solutions.

and still be syntactically correct expression which can be evaluated. Like nature, many mutants are not as fit as their parents but, like nature, every so often, a mutant is created which is better. Similarly children produced by sex have genes which are a random combination of parental genes. Again, every once in a while an improved combination is found and the offspring program is selected for prospers and in subsequent generations copies of it spread through the evolving population.

Genetic programming can be thought of as like domesticated animals and plants, where improvements have been made by breeders progressively selecting preferred characteristics. (Darwin studied the records of breeders of domesticated pigeons.) Thus you too must impose a direction on evolution, e.g. to control a robot, design a radio aerial or find a genetic component of breast cancer survival, you must select programs that are better at doing it. For example, given the cause of death and life span of 253 Swedish women cancer patients, you might select a program which correctly predicted more cases of survival for more than 8 years after surgery than one which was less accurate.

With large populations and/or many generations, selecting individual programs becomes too tedious to do by hand. Instead we pass the job to a computerized automatic “fitness function”. On your behalf, it prefers better programmes over the less good. Ultimately it is your fitness function which guides the evolution of your population by selecting who will survive and who will have children. The fitness function is literally a matter of life or death.

There are several fine books on GP ([2, 4] and [5] leap to mind); however, we strongly encourage *doing* GP as well as reading about it. There are many good free (unsupported) GP implementations (e.g. lilGP, ECJ, Beagle¹ and TinyGP) but it is not so hard to write your own.

7.1.1 Overview

The next section describes the main parts of genetic programming, while Section 7.3 describes how you put them together to get a working system. Next Sections 7.4 and 7.5 describe advanced GP techniques. We survey the enormous variety of

¹ Darwin was the naturalist onboard HMS Beagle for 5 years [6].

applications of GP in Section 7.6. This is followed by a collection of troubleshooting suggestions (Section 7.7) and by our conclusions (Section 7.8).

7.2 Representation, Initialization and Operators in Tree-Based GP

7.2.1 Representation

In artificial intelligence it has become accepted wisdom that how information about the application and its solution is stored (i.e. represented internally within the computing system) and manipulated by it, is crucial to successful implementation. Huge effort is spent by very clever people on designing the correct representation.

Genetic programming has ignored this. In GP, the evolved program contains the solution and “representation” refers to the language evolution uses to write the program. The same representation might be used in a program evolved to predict breast cancer survival as one evolved to find insider trading in a stock market.

In GP, programs are usually expressed as *syntax trees* rather than as lines of code. For example Figure 7.2 shows the tree representation of the program $\max(x+x, x+3*y)$. The variables and constants in the program (x , y and 3) are leaves of the tree. In GP they are called *terminals*, while the arithmetic operations ($+$, $*$ and \max) are internal nodes called *functions*. The sets of allowed functions and terminals together form the *primitive set* of a GP system.

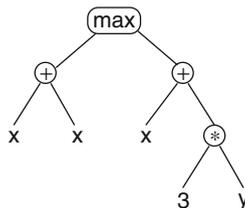


Fig. 7.2 GP syntax tree representing $\max(x+x, x+3*y)$.

It is common to represent expressions in *prefix* notation, e.g. $\max(x+x, x+3*y)$ becomes $(\max(+ x x) (+ x (* 3 y)))$. Usually the number of arguments a function takes is known, e.g. *sin* has one argument but $*$ has two. When the arity is known, the brackets in prefix-notation expressions are not needed. This means trees can be represented as simple linear sequences. Usually this is much more efficient than tree-based representation of programs, which require the storage and management of numerous pointers. In effect, the function’s name gives its arity and from the arities the brackets can be inferred. For example, the

expression $(\max (+ x x) (+ x (* 3 y)))$ can be written unambiguously as $\max + x x + x * 3 y$.

The choice of whether to use such a linear representation or an explicit tree representation is typically guided by convenience, efficiency, the genetic operations being used (some may be more easily or more efficiently implemented in one representation), and other data one may wish to collect during runs. (It is sometimes useful to attach additional information to nodes, which may be easier to implement if they are explicitly represented).

Tree representations are the most common in GP. However, there are other important representations including linear [4, 7–9] and graph [10–12] based programs.

7.2.2 Initializing the Population

As with other evolutionary algorithms, in GP the individuals in the initial population are typically randomly generated. There are a number of different approaches to generating this random initial population, e.g. [13]. However, we will describe two of the simplest methods (the `full` and `grow` methods) and the most widely used combination of the two known as *Ramped half-and-half* [2].

In both the `full` and the `grow` methods, the initial individuals are generated so that they do not exceed a maximum depth you decide. The *depth* of a node is the number of edges that need to be traversed to reach the node starting from the tree’s root node (depth 0). The depth of a tree is the depth of its deepest leaf (e.g. the tree in Figure 7.2 has a depth of 3). The `full` method generates full trees (i.e. all leaves are at the same depth). It does this by choosing at random from the available functions (known as the function set) until the maximum tree depth is reached. Then the tree is finished by adding randomly chosen leaves from the available terminals (known as the terminal set). Figure 7.3 shows a series of snapshots of the construction of a full

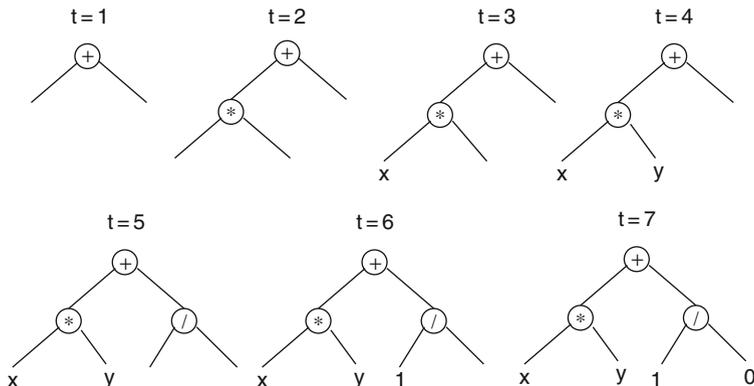


Fig. 7.3 Creation of a full tree having maximum depth 2 using `full` initialization ($t = \text{time}$) [53].

tree of depth 2. The children of the $*$ and $/$ nodes must be leaves or otherwise the tree would be too deep. Thus, at steps $t = 3$, $t = 4$, $t = 6$ and $t = 7$ a terminal must be chosen. (In this example leafs x , y , 1 and 0 were randomly chosen).

Although, the `full` method generates trees where all the leaves are at the same depth, this does not necessarily mean that all initial trees will have an identical number of nodes (often referred to as the *size* of a tree) or the same shape. This happens only if all the functions have the same arity (i.e. have the same number of inputs.) Nonetheless, even when mixed-arity primitive sets are used, the range of program sizes and shapes produced by the `full` method may be limited. The `grow` method creates trees of more varied sizes and shapes. Nodes are selected from the whole primitive set (i.e. functions and terminals) until the depth limit is reached. Once the depth limit is reached only terminals may be chosen (just as in the `full` method). Figure 7.4 illustrates growing a tree with depth limit of 2. In Figure 7.4 ($t = 2$) the first argument of the $+$ root node happens to be a terminal. This prevents that branch from growing any more. The other argument is a function ($-$). It can grow one level before its arguments are forced to be terminals to ensure that the resulting tree does not exceed the depth limit. C++ code for a recursive implementation of both the `full` and the `grow` methods is given below.

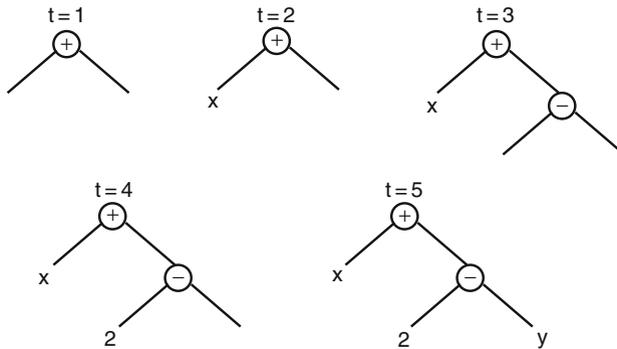


Fig. 7.4 Creation of a five node tree using the `grow` initialization method with a maximum depth of 2 ($t = \text{time}$). A terminal is chosen at $t = 2$, causing the left branch of the root to be closed at that point even though the maximum depth had not been reached [5].

Because neither the `grown` or the `full` method provides a very wide array of sizes or shapes on their own, Koza proposed a combination called *ramped half-and-half* [2]. Half the initial population is constructed using `full` and half is constructed using `grow`. This is done using a range of depth limits (hence the term “ramped”) to help ensure that we generate trees having a variety of sizes and shapes.

While these methods are easy to implement and use, the sizes and shapes of the trees generated are highly sensitive to the number of functions, the number of inputs they have and the number of terminals. This makes it difficult to control the sizes and shapes of the trees. For example, if there are many more terminals than

functions, the `grow` method will almost always generate very short trees regardless of the depth limit. Similarly, if the number of functions is considerably greater than the number of terminals, then the `grow` method will be like the `full` method.

```
//Choose desired depth uniformly at random between min and max depth.
//Choose either full or grow.
SubInit(rnd(max_depth-min_depth)+min_depth, rnd(2), min_depth);

void Individual::SubInit(int depth, BOOL isfull, int min_depth) {
  if (depth <= 0)
    i=rand_terminal(); // terminal required
  else if (isfull || min_depth>0)
    i=rand_function(); // function required
  else { //grow: terminal allowed 50% of the time
    if (rnd(2)) // terminal required
      i=rand_terminal();
    else // node required
      i=rand_function();
  }
  SETNODE(code[ip],i); //store opcode in Individual
  ip++;

  for(int a=0;a<argnum(i);a++) {
    SubInit(depth-1,isfull, min_depth-1, tree);
  }
}
```

C++ code fragment to create a random tree. For efficiency the tree is flattened and stored in array `code` (access is via macro `SETNODE`). `SubInit` recursively calls itself until it reaches a leaf of the tree. (Based upon Andy Singleton's GPquick.)

The initial population need not be entirely random. If something is known about likely properties of the desired solution, trees having these properties can be used to seed the initial population.

7.2.3 Selection

As with other evolutionary algorithms, in GP better individuals are more likely to have more child programs than inferior individuals. Tournament selection is most often used, followed by fitness-proportionate selection [4], but any standard evolutionary algorithm selection mechanism (e.g. stochastic universal sampling) can be used.

In *tournament selection* a number of individuals are chosen at random from the population. These are compared with each other and the best of them is chosen to be the parent. When doing crossover, two parents are needed and, so, two selection tournaments are made. Note that tournament selection only looks at which program is better than another. It does not need to know how much better. This effectively

automatically rescales fitness, so that the selection pressure is constant. Thus, a single extraordinarily good program cannot immediately swamp the next generation with its children. If it did, this would lead to a rapid loss of diversity with potentially disastrous consequences for a run. Conversely, tournament selection amplifies small differences in fitness to prefer the better program even if it is only marginally superior to the other individuals in a tournament.

Tournament selection, due to the random selection of programs to be included in the tournament, is inherently noisy. So, while preferring the best, tournament selection does ensure that even below average programs have some chance of having children. Since tournament selection is easy to implement and provides automatic fitness rescaling, it is commonly used in GP.

7.2.4 Recombination and Mutation

Crossover (recombination) and mutation in GP are very different from crossover and mutation in other evolutionary algorithms. The most commonly used form of crossover is *subtree crossover*. Given two parents, subtree crossover randomly (and independently) selects a *crossover point* (a node) in each parent tree. Then, it creates the offspring by replacing the subtree rooted at the crossover point in a copy of the first parent with a copy of the subtree rooted at the crossover point in the second parent [2], as illustrated in Figure 7.5.

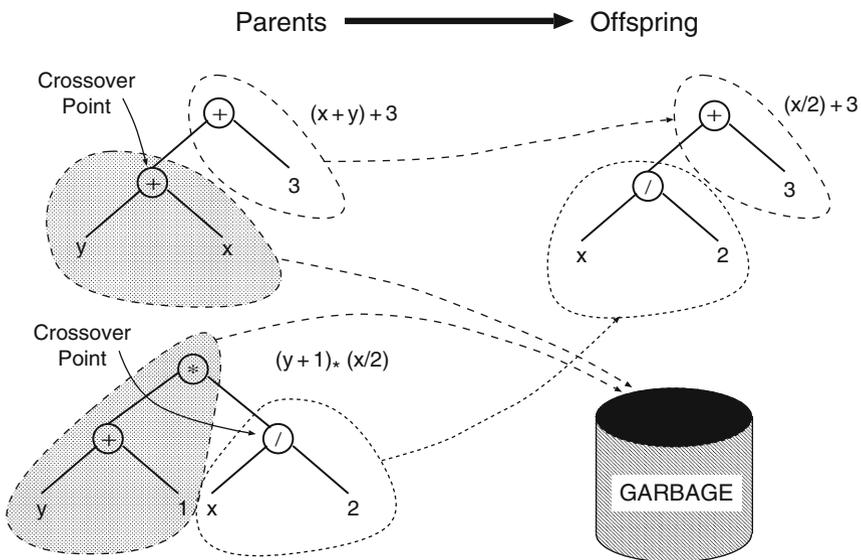


Fig. 7.5 Example of subtree crossover. Note that the trees on the left are actually *copies* of the parents. So, their genetic material can freely be used without altering the original individuals [5].

Typical GP primitive sets lead to trees with an average arity of at least two. This means most of the program will be leaves. So if crossover points were chosen uniformly, crossovers would frequently swap very small subtrees (even just leaves), i.e. exchange only very small amounts of genetic material. Whereas in nature (and many GAs) often both parents contribute more-or-less equally to their offspring's genetic code. To counter this, Koza suggested the widely used approach of choosing functions 90% of the time and leaves 10% of the time [2]. Many other types of crossover and mutation of GP trees are possible (see [5, pp 42–44]).

The most commonly used form of mutation in GP is *subtree mutation*. It randomly selects a mutation point in a tree and substitutes the subtree rooted there with a randomly generated subtree (cf. Figure 7.6 and [15]).

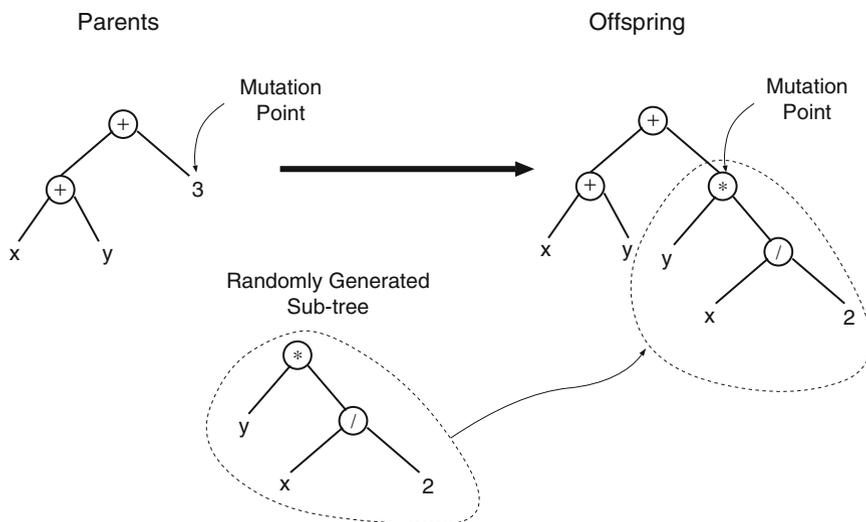


Fig. 7.6 Example of subtree mutation [5].

Another common form of mutation is *point mutation*, which is GP's rough equivalent of the bit-flip mutation used in genetic algorithms [14]. In point mutation, a random node is selected and the primitive stored there is replaced with a different random primitive of the same arity taken from the primitive set. If no other primitives with that arity exist, nothing happens to that node (but other nodes may still be mutated). When subtree mutation is applied, it changes exactly one subtree. On the other hand, every node in the tree has a small probability of being mutated by point mutation. This means point mutation independently changes a random number of nodes.

In GP normally only one genetic operator is used to create each child. Which one is used is chosen at random. Typically, crossover is applied with the highest probability, the *crossover rate* often being 90% or higher. On the contrary, the *mutation rate* is much smaller, typically being in the region of 1%. If the sum of all

the probabilities comes to less than 100% the remaining offspring are created simply by copying better individuals from the current population. (This is known as *reproduction*.)

7.3 Getting Ready to Run Genetic Programming

7.3.1 Step 1: Terminal Set

GP is not typically used to evolve programs in the familiar languages people normally write programs in. Instead simpler programming languages are used. Indeed GP can usually be thought of as evolving executable expressions rather than fully fledged programs. The first two preparatory steps, the definition of the terminal and function sets, specify the language. Together they define the ingredients that are available to GP to create computer programs.

Typically the terminal set contains *the program's inputs* (e.g. x , y , cf. Table 7.1). It may also contain *functions with no arguments*. They might be needed because they return different values each time they are used, such as a function which returns random numbers or returns the distance from a robot to an obstacle or because the function produces *side effects*. Functions with side effects may change some global data structures, draw on the screen, print to a file, control the motors of a robot, etc.

Table 7.1 Examples of primitives in GP function and terminal sets.

Function set		Terminal set	
<i>Kind of primitive</i>	<i>Example</i>	<i>Kind of primitive</i>	<i>Example</i>
Arithmetic	$+$, $*$, $/$	Variables	x , y
Mathematical	\sin , \cos , \exp	Constant values	3, 0.45
Boolean	AND, OR, NOT	0-arity functions	<code>rand</code> , <code>go_left</code>
Conditional	IF-THEN-ELSE		
Looping	FOR, REPEAT		

Often an evolved program will need access to constants. We do not know in advance what their values will be, so GP chooses some randomly. In some implementations the number of constants is limited and it may be that new ones cannot be created during the GP run. Instead their values must be chosen as the population is initialized. Typically this is done by a special terminal that represents an *ephemeral random constant*. Every time it is chosen (either at the start or when a new subtree is created by mutation), a different random value is generated. This is used for that *particular* terminal and remain fixed for the rest of the run.

7.3.2 Step 2: Function Set

The function set typically contains only the arithmetic functions ($+$, $-$, $*$, $/$). However, all sorts of other functions and constructs typically encountered in computer

programs can be used, see Table 7.1. Sometimes specialized functions or terminals, which are designed to solve particular problems are used. For example, if the goal is to evolve art, then the function set might include such actions as `select_from_pallet` and `paint`.

For GP to work effectively, most function sets are required to have an important property known as *closure* [2]. Closure can be broken down into *type consistency* and *evaluation safety*. Finally the primitive set must be able to (i.e. must be *sufficient* to) express solutions to the problem.

7.3.2.1 Type Consistency

Crossover (Section 7.2.4) can mix and join nodes arbitrarily. So it is important that *any* subtree can be used as any argument for every function in the function set. The simplest way to achieve this is to ensure all functions return values of the same type and that each of their arguments also have this type. For example, `+`, `-`, `*`, and `/` can be defined so that they each take two integer arguments and return an integer. The terminals would also be integers. (Automatic-type conversion, e.g., between Booleans and integers, default values and polymorphic functions, can also be used to ensure that crossover always produces syntactically correct and runnable programs.) Sections 7.4 and 7.5 will describe safe ways to extend GP.

7.3.2.2 Evaluation Safety

The purpose of evaluation safety is to ensure evolved programs can run and thereby be assigned fitness even when they run into errors. For example, an evolved expression might divide by 0 or call `MOVE_FORWARD` when facing a wall or precipice. It is common to use *protected* versions of numeric functions that can otherwise throw exceptions, such as division, logarithm, exponential and square root. The protected version of a function first tests for potential problems with its input(s) before executing the corresponding instruction. If a problem is spotted then some default value is returned. Protected division (often notated with `%`) checks to see if its second argument is 0. If so, `%` typically returns the value 1 (regardless of the value of the first argument). (The decision to return the value 1 provides the GP system with a simple way to generate the constant 1, via an expression of the form `(% x 0)`. This combined with a similar mechanism for generating 0 via `(- x x)` ensures that GP can easily construct these two important constants.) Similarly, in a robotic application a `MOVE_AHEAD` instruction can be modified to do nothing if a forward move is illegal or if moving the robot might damage it. (Braitenberg permitted his imaginary robots to make dangerous moves as a way of weeding the poor control program from the better [16].)

An alternative to protected functions is to trap run-time exceptions and strongly reduce the fitness of programs that generate such errors. However, if the likelihood of generating invalid expressions is very high, this can lead to too many individuals

in the population having nearly the same (very poor) fitness. This makes it hard for selection to choose which individuals might make good parents.

7.3.2.3 Sufficiency

By sufficiency we mean it is possible to express a solution to the problem using the elements of the primitive set. For example, $\{\text{AND}, \text{OR}, \text{NOT}, x_1, x_2, \dots, x_N\}$ is a sufficient primitive set for logic problems, since it can produce all Boolean functions of the variables x_1, x_2, \dots, x_N . The primitive set $\{+, -, *, /, x, 0, 1, 2\}$ is unable to represent transcendental functions, such as $\sin(x)$. When a primitive set is insufficient, GP can often develop programs that approximate the desired solution. Which may be good enough for the user's purpose. Adding a few unnecessary primitives in an attempt to ensure sufficiency tends not to slow down GP overmuch.

7.3.3 Step 3: *Fitness Function*

The task of the fitness measure is to choose which parents are to have offspring. That is, which parts of the search space we have just sampled (which is what the current population has done for us) are worth exploring further. The fitness function is our primary (and often sole) mechanism for giving a high-level statement of requirements to GP.

Fitness can be measured in many ways. For example, in terms of the amount of *error* between its output and the desired output; the amount of *time* (fuel, money, etc.) required; the *accuracy* of the program in recognizing patterns or classifying objects; the *payoff* a game-playing program produces.

Fitness evaluation normally requires executing all the programs in the population, typically multiple times. While one can compile the GP programs that make up the population, the overhead of building a compiler is usually substantial, so it is much more common to use an interpreter to evaluate the evolved programs. Interpreting a program tree means executing the nodes in the tree in an order that guarantees that nodes are not executed before the value of their arguments (if any) is known. This is usually done by traversing the tree recursively starting from the root node and postponing the evaluation of each node until the values of its children (arguments) are known. Other orders, such as going from the leaves to the root, are possible. If none of the primitives have side effects, the two orders are equivalent. Figure 7.7 contains C++ code fragments which implements top down recursive tree evaluation using a linear data structure for speed.

In some problems we are interested in the *output* produced by a program. In other problems we are interested in the actions performed by a program composed of functions with side effects. In either case the fitness of a program typically depends on the results produced by its execution on many different inputs or under a variety of different conditions. For example, the program might be tested on all possible combinations of inputs x_1, x_2, \dots, x_N . Alternatively, a robot control program might

```

//Flatted tree is stored in array of nodes.
//evalnode has three components. It trades space for speed.
//node (not shown) stores the same information in one byte. It
//is used to store the population. node is expanded to evalnode
//before fitness testing.
//typedef allows code to be compiled for several applications.

typedef float retval;
typedef retval (*EVALFUNC)(); // evaluation code with global pointer
typedef struct evalnode {
    EVALFUNC ef;
    evalnode* jump;
    retval value;
} evalnode; // node type

evalnode* IP; //global pointer
evalnode* ExprGlobal; //Array holding flatten tree
//some old Sun compilers incorrectly optimise EVAL (workaround via FTP)
#define EVAL ((++IP)->ef)()
#define GETVAL IP->value
#define TRAVERSE() IP=(++IP)->jump

retval D0Eval() { return data[0]; } //data holds inputs D0-D9. Typically it
retval D9Eval() { return data[9]; } //is different for each training case
retval ConstEval() {return GETVAL;}
retval AddEval() {return EVAL + EVAL;}
retval SubEval() {return EVAL - EVAL;}
retval MulEval() {return EVAL * EVAL;}
retval DivEval() {// "Protected" division
    const retval numerator = EVAL;
    const retval denominator = EVAL;
    if (denominator !=0) return numerator/denominator;
    else return 1;
}
retval IflteEval() {//IfLTE(condition1islessthan,condition2,dothis,dohat)
    retval rval=EVAL;
    if (rval<=EVAL) {
        rval=EVAL;
        TRAVERSE(); // Jump the third expression
    } else {
        TRAVERSE(); // Jump the second expression
        rval=EVAL;
    }
    return rval;
}

retval Individual::evalAll() {// eval the whole expression anew
    IP=ExprGlobal-1; // start at beginning of flatten tree
    return EVAL;
}

```

Fig. 7.7 Example fast interpreter (based on Andy Singleton's GPquick). The tree is linearized and functions and terminals within it are replaced by pointers to C++ functions which implement them. On a typical modern computer the GP individual and the interpreter are held in fast cache. Since the tree is flatten into the traditional depth first order, the interpreter runs from top of the tree to the rightmost terminal in one forward pass. This avoids backtracking. Continuous forward motion suits typical cache architectures.

be tested with the robot in a number of starting locations. These different test cases typically contribute to the fitness value of a program incrementally and for this reason are called *fitness cases*.

Despite all this sophistication and the computational work it does, the fitness function ultimately boils down to just one bit of information: does this mutated program beget another child? We do not know the correct answer to this question. So we add noise (e.g. via tournament selection). Fortunately it is not necessary to get the answer right all the time, or even most of the time, just as long as we are right occasionally. We sometimes lose sight of this hard truth. Sometimes it may be better to accept a less accurate calculation of fitness. If by doing so we reduce the time taken to calculate a fitness value. Thus allowing us to take the life or death decision more times.

7.3.4 Step 4: GP Parameters

The most important control parameter is the *population size*. It is impossible to make general recommendations for setting *optimal* parameter values, as these depend too much on the details of the application. However, genetic programming is in practice robust and it is likely that many different parameter values will work. As a consequence, one need not typically spend a long time tuning GP for it to work adequately. Some possible parameter settings are given in the tableau in Table 7.2.

Table 7.2 Typical parameters for example genetic programming run.

Objective:	Record your problem here
Function set:	For example, +, -, % (protected division) and ×; all operating on floats
Terminal set:	For example, x and constants chosen randomly between -5 and $+5$
Fitness:	e.g. sum of absolute errors for a number of fitness cases. The number of fitness cases may be limited by the amount of training data available to evaluate the fitness of the evolved individuals. In other cases, e.g. 22-bit even parity [17], there can be too much training data. Then the fitness function may use a fraction of the training data. This does not necessarily have to be done manually as there are a number of algorithms that dynamically change the test set as the GP runs (see [5, Sect. 10.1]).
Selection:	Tournament size 7
Initial pop:	Ramped half-and-half (Section 7.2.2) depth range of 2–6
Parameters:	As a rule one prefers to have the largest population size that your system can handle gracefully. Normally the population size should be at least 500 and people often use much larger populations. (However, some prefer much smaller populations. Typically these rely on mutation rather than crossover and run many more generations.) Traditionally, 90% of children are created by subtree crossover. However, the use of a 50–50 mixture of crossover and a variety of mutations also appears to work well [5, Chapter 5]. Some implementations do not require arbitrary limits of tree size. Even so, because of bloat (the uncontrolled growth of program sizes during GP runs [5, Sect. 11.3]), it is common to impose either a size or a depth limit or both (see Section 7.7.6).
Termination:	10–50 (The most productive search is usually performed in those early generations.)

7.3.5 Step 5: When to Stop and How to Decide Who is the Solution

The last step, is choosing when to stop the GP and how to decide which of the thousands of programs that it has evolved to use. Typically we stop either when an acceptable solution has been found or a maximum number of generations has been reached. Typically, the single best-so-far individual is used. Although one might wish to study additional individuals, e.g. to look for particularly short or elegant solutions.

7.4 Guiding GP with A Priori Knowledge

As described so far, GP is essentially knowledge free: a powerful search mechanism (evolution) is set free to search the space of all expressions which can be formed from the function and terminal set. Contrast this with traditional methods, such as linear regression, in which a very basic search mechanism is used to search a very restricted set of expressions. Linear regression is often extended to more complex forms (polynomial regression, log regression, etc.), but this still leaves a vast gap between the complete search of GP and the very restricted parameter search of classical regression.

In many applications, the user will know a great deal about the form of acceptable solutions. It can be highly desirable to incorporate this knowledge into the search, since it can save the user time, e.g. by enabling the user to exclude solutions which will not be useful for some reason or to impose a preference ordering on solutions. Including the user's background knowledge can also increase data efficiency and so allow more complex models to be learnt than could possibly be justified solely by the available data. In some cases, such restriction may be essential, because it may not be possible to provide meaningful fitness values for all the solutions a GP system could evolve. Finally, GP search may be more efficient if the user's knowledge can be used to increase the concentration of solutions in the search space. This increase may be non-trivial (in example 3 below it is of many orders of magnitude) but it is nevertheless the least important reason.

In principle, a wide range of mechanisms could be used to restrict the search space; in practice, most available systems use some form of grammar, generally an extension of Context Free Grammars (CFG [18]). The constraints may range over a wide range of complexity, for example

1. Strongly typed systems (Section 7.3.2.1, [19]): in which only type-consistent expressions can be evolved.
2. Extended process models: in many domains, such as ecological modelling [20], there is a known sub-model of processes which are certainly occurring (zooplankton are eating phytoplankton, for example), but there may also be other unknown processes occurring which require adaptation of this process model to fit the data.

3. Dimensional consistency [21]: For example, in physics, equations must be consistent in time (t), length (l) and mass (m) dimensions. For example, integrating Newton’s second equation of motion gives $s = ut + \frac{1}{2}at^2$. s has dimensions of length. u is a velocity and hence has dimension l/t . $\frac{1}{2}$ is a pure number and so has no dimensions. a is an acceleration and hence has dimension l/t^2 . Putting these together, the right-hand side gives $l/t \times t + l t^{-2} \times t^2 = l$. Which is indeed the same as the dimensions of the left-hand side (l). Dimensionally inconsistent formulae may fit the data well but they are nevertheless unacceptable.

7.4.1 Context-Free Grammars in GP

CFG-based GP systems are the most widely used, and the simplest to explain, so we take them as our base case. From the user’s perspective, a CFG-GP system is very similar to a standard GP system. However, instead of just providing a list of function and terminal symbols², the user must provide a grammar specifying the ways in which they may be used. That is the only change really required; the user does not have to do anything special with respect to the GP operators (selection, crossover, mutation); the system takes care of those. For some systems, there may be one further difference. In a grammar-based system, the fitness function can be defined in the same way as for standard GP. However, the grammar defines how to build up more complex expressions from simpler ones. In many cases, it is easier to define how to build up the meanings of the expressions (i.e. how to evaluate them) at the same time—we call this “providing a semantics for the grammar”. If this is done, the fitness function definition may reduce to just a few lines of code, defining how these values contribute to the fitness. We give a brief example in Section 7.4.1.1 below.

The grammar provides an additional way for the user to interact with the evolving population. When the CFG-GP system is first run, it may not produce the results the user desires, e.g. the evolved solution may not fit the data sufficiently well. Or it may not be acceptable to the user for some other reason. However, the CFG-GP runs may help the user see how the problem may be solved. Frequently, it is possible to incorporate this insight into the grammar so as to achieve more useful results in subsequent runs. This ability to interact with the solution space, through grammar definitions, is one of the primary practical benefits of grammar-based GP systems.

Of course, the implementation of CFG-GP is a little more complex than simple tree GP, though this is generally not visible to the user. In GP, the individuals of the evolutionary population are expression trees; in CFG-GP, the individuals are parse trees from the grammar, i.e. CFG-GP individuals are paths through the user-supplied

² A word of caution: GP and grammar terminology were both developed before grammar-based GP systems and use some of the same words. Unfortunately, when they came together in grammar-based GP, some inconsistencies arose. Thus, in a CFG-GP system, a (GP) function symbol is a terminal (in grammar terms), though it is not a member of the GP terminal set. Unfortunately there does not seem to be any reasonable way to resolve this inconsistency.

grammar, starting from the grammar’s start symbol (the root of the parse tree). Eventually the path will reach terminals of the grammar (i.e. symbols which cannot be expanded further). The list of terminals (in the order they were encountered) is the output of the grammar. Typically, this list is an executable program (written in the language specified by the user’s grammar). It is then run in order to find the fitness of the CFG-GP individual.

Initialization, which requires ensuring grammar consistency while guaranteeing to stay within the depth bound, is also a little complex; most systems use a variant of the grow-tree algorithm described in Section 7.2.2. This is combined with a counting mechanism, to ensure that it is always possible to complete a parse tree within the remaining depth. Crossover and mutation are defined in ways that preserve grammar consistency. Mutation replaces a subtree from the grammar with a random subtree. It creates the random subtree in the same way as the initial population is created, except that it starts from the location in the grammar occupied by the subtree it has just removed, rather than at the root. Crossover is essentially like normal GP crossover, except that crossover is only allowed between nodes with the same grammar non-terminal. This ensures, as with mutation, that the offspring is consistent with the grammar.

7.4.1.1 Example of CFG-GP: Strong Typing in GP

We use Strongly Typed GP [19] as a simple example of grammar use. A GP problem requiring two types, arithmetic and Boolean, might use a grammar such as in Table 7.3. Thus the first “arithmetic” rule says that an arithmetic expression may consist of the sum of two arithmetic expressions, or (| means or) the difference of two arithmetic expressions, and so on. The second “interaction” rule says that a Boolean expression may be formed by comparing two arithmetic expressions, with any of the comparison operators <, = or >. Thus the grammar permits arbitrarily complex nesting of arithmetic and Boolean expressions, but guarantees that they are combined in meaningful ways.

In systems which also support semantic specification within the rules, a rule such as $A \rightarrow A * A$ would be expanded to include variables. These variables represent the values generated by the rule (such as $A(A_0) \rightarrow A(A_1) * A(A_2)$). Extra (semantic) rules then give the values of those variables (such as $val(A_0) = val(A_1) * val(A_2)$). Of course, in simple cases like this, where the meaning of ‘*’ is already built into

Table 7.3 An example grammar for Boolean and arithmetic types. The following six rules define how non-terminal symbols A (the start symbol, representing arithmetic expressions) and B (representing Boolean expressions) can be expanded into 15 (grammar) terminals + - * / x 0 if(, ,) < = > & ∨ ¬ true false.

<i>Arithmetic rules</i>	<i>Interaction rules</i>	<i>Boolean rules</i>
$A \rightarrow A + A A - A A * A A / A$	$A \rightarrow \text{if}(B, A, A)$	$B \rightarrow B \& B B \vee B \neg B$
$A \rightarrow x 0$	$B \rightarrow A < A A = A A > A$	$B \rightarrow \text{true} \text{false}$

the language in which the GP system is written, the advantage is limited. In more complex domains or problems where other properties of the expression in addition to its value may be needed, semantic specification can greatly simplify coding the problem.

7.4.2 Variants of Grammar-Based GP

7.4.2.1 More Powerful Grammars

Perhaps the most important issue is that the user's knowledge may not be expressible in context-free form. This has led to a wide range of extended-grammar systems. They fall into two main classes: Context Sensitive Grammars (CSG) [22] and attribute and other semantic grammars [23].

CSG permit more precise syntactic restrictions on the search space; for some problem domains, this greater expressiveness is important for encoding the problem.

Attribute grammars extend the semantic specification we described in Section 7.4.1.1. In some problems, the semantics may allow us to decide early in the evaluation process, that the individual will have low fitness. For example, in a constraint problem, we may know that if a constraint is breached early in evaluating an individual, the violation is only going to get worse as we continue with its evaluation. Thus semantic constraints may be used to short-circuit fitness evaluation. But even more intriguingly, they may be used to avoid creating poor individuals at all. For example, when we come to cross over individuals, the semantic values attached to the nodes in an individual might indicate that a crossover at a particular point would automatically breach a constraint. A system based on semantic grammars can then simply abort the crossover, never creating the potentially poor individual. In general, if it is difficult to express the user's knowledge about the search space in a CFG, consider using either a CSG or a semantic grammar.

7.4.2.2 More Flexible Representations: GE and Tree Adjunct Grammars

The reduction in search space size provided by a CFG representation can be beneficial for search; but it comes at a cost. The CFG reduces not only the number of formulae that can be represented in the search space, but also the links between them. Paradoxically, in some cases this sparser search space might be *more* difficult for evolution to search than the original search space. Two approaches have been introduced to avoid this problem. In the first [24, 25], the CFG representation is linearized: instead of representing the individuals directly as grammar parse trees, a coding scheme represents them as linear strings. This approach has led to one of the most widely used GP systems, known as Grammatical Evolution (GE) [25]. The other [26] uses an alternative representation from natural language study, Tree Adjunct Grammars (TAGs [27]); unlike CFG trees, any rooted subtree of a TAG tree is

syntactically and semantically meaningful, so that there is much more flexibility in transforming one TAG tree to another.

In both cases, the syntactic flexibility provides an additional benefit: it is relatively easy to implement new operators (often analogous to biological processes that occur in DNA evolution) which may simplify search in particular domains. Practically, this means that where search with standard-GP and CFG-GP systems has stagnated, it may be worth investigating GE or TAGs. They may be able to solve problems which are beyond the reach of more classical GP systems.

7.4.2.3 Grammar Learning

A number of more experimental grammar-based systems [28–30] refine the grammar describing the search space as search proceeds. (These systems are usually based on probabilistic grammars: each grammar rule option has a probability attached to it, indicating the probability that it will be used in generating an individual.) This has two consequences. It can make for faster search. But more importantly, it means that the grammar at the end of the search space may give an explicit representation of the space of solutions (rather than the implicit representation given by the best individuals in a final GP population).

This explicit representation may be of value in its own right, especially in applications such as scientific research, where the desired outcome is better understanding of the processes in the domain, rather than simply predictive models. The use of probabilistic grammars means that the understanding may be quite sensitive, going beyond just the content of the grammar rules. In some parts of the grammar, the probabilities may converge close to either 1.0 or 0.0, indicating that that aspect of the grammar is important in defining a solution to the problem; in others, the probabilities may be more widely spread, indicating that that aspect of the grammar is not particularly important to the problem solution.

7.5 Expanding the Search Space in Genetic Programming

In Section 7.4 we described some of the ways in which you can give the evolutionary process a helping hand. For example, by providing domain-specific data types or by constraining programs to conform to an appropriate grammar. But, in the context of a particular problem or a particular set of program representations, we do not necessarily know *how* to give evolution a helping hand. In which case it can be useful to expose more, rather than less, of the system's decisions about data and control architecture to evolution. Doing so will often incur new costs, some from the added complexity of the system and some from the expansion of the space of programs which GP is searching [31]. However in many cases these costs can be justified by improvements to problem-solving power or scalability.

We will discuss some of the ways in which researchers have expanded the purview of the evolutionary process in GP. In Section 7.5.1, we first examine the evolution of data structures and the ways in which they can be accessed and manipulated by evolving programs. We then turn to program and control structure. Section 7.5.2 describes how GP can be used to evolve programs that use subroutines, macros, and more exotic techniques for controlling the flow of execution. The concept of “development” (here development means the evolved programs build *other* structures which then produce the desired behaviors) provides for even more evolutionary flexibility (Section 7.5.3). The last part of this section (Section 7.5.4) describes mechanisms by means of which the evolutionary processes themselves can be allowed to evolve.

7.5.1 *Evolving Data Structures and Their Use*

The earliest and simplest GP applications evolved programs that used single, simple data types. The use of multiple—but still simple—data types has been helped in a variety of ways, for example by the use of strong typing (see Section 7.4 and [19]). An important technique for evolving programs that use more complex data types is *indexed memory*, which was first presented by Teller [32]. An indexed memory is simply an array of variables of some simple type, accessed using integer indices. Teller showed that by including indexed memory `read` and `write` functions in the GP function one can evolve programs that use memory in relatively complex and useful ways. He also showed that the inclusion of indexed memory was useful in expanding the space of programs over which GP can search, e.g. it can be shown to include programs for all Turing computable functions.

Indexed memory can be used by evolving programs to implement a wide variety of more complex data structures, but modern software engineering practice suggests that it is even more useful for human programmers—and hence possibly also for GP—to have access to higher level data structures. Langdon has investigated the extent to which GP can evolve, and subsequently use, more abstract data structures including stacks, queues and lists [33]. He showed that GP can indeed evolve and subsequently solve problems using such data structures and that GP with abstract data types can outperform GP with indexed memory on several problems, including a context-free language recognition problem and the problem of implementing a simple four function calculator.

Alternative program representations provide additional opportunities for the evolution and use of data structures. For example, approaches based on polymorphic functional representations, initially developed by Yu using Haskell [34], have recently been extended by Binard and Felty, using a version of the λ -calculus to which they have added an operation of abstraction on types [35]. They showed how their system could evolve and use abstractions for Boolean and list data types which were not explicitly present in their initial environments.

To some extent, the ways in which data types and program syntax are interrelated determine the ways in which GP can discover and use complex data structures. Strongly typed GP and polymorphic GP provide two approaches but they do not exhaust the possibilities. For example, in the Push programming language all communication between instructions is accomplished via typed global data stacks. It is not specified by placing the instructions next to each other, as in most programming languages. This decouples an evolving program's type structures from its control structures and thereby permits greater flexibility (for good or ill) in the expression of programs that manipulate multiple data types [36, 37].

7.5.2 Evolving Program and Control Structure

Most interesting programs that are written by humans involve the use of control structures not available in the simplest GP systems. These include mechanisms that support iteration, recursion, and the definition and use of reusable code modules. As with data structures, GP researchers have developed a range of techniques for evolving programs that evolve and use these powerful control abstractions.

Limited forms of iteration are relatively easy to handle through the use of primitive functions that simply repeat the execution of a subexpression some specified number of times. This was demonstrated in Koza's first book using `do_until` structures [2]. A variety of more sophisticated techniques, such as the "restricted iteration creation" operations of Koza and Andre [38], have been developed to help GP systems incorporate iteration into evolving programs. Both iteration and recursion present challenges with respect to nontermination; this is generally handled either by imposing execution limits or by using primitives that are naturally self-limiting, such as the `foldr` function in Haskell [39]. While the search space of recursive programs appears to be rugged, and several early attempts to evolve recursive programs produced negative results (e.g. [40]), more recent research has been increasingly successful (e.g. [37, 39, 41–44]).

Modular structures can also be incorporated into evolving programs in several ways. A common approach, pioneered by Koza [45], is to simultaneously evolve a "main program" (sometimes called a "result producing branch") and one or more "automatically defined function" (ADF) branches that can be called by the main program and possibly by each other (usually with restrictions to prevent nonterminating recursion). This approach has been shown to provide dramatic advantages in certain problem areas with exploitable regularities. In the original ADF framework the number of ADFs and the numbers of arguments that they take are specified manually, but the subsequent development of "architecture altering operations" brought these decisions, as well, under evolutionary control [46].

A variety of other approaches to the evolution and use of modules have also been developed. For example, in "evolutionary module acquisition" the code for modules is not evolved in separate branches but rather is extracted from the main programs of relatively successful individuals in the population [47–49]. "Automatically defined

macros” allow GP to evolve not only function modules but also control structure modules that execute code conditionally or repeatedly [50]. And several researchers have shown how GP can be used to evolve object-oriented programs in which functionality is modularized through the use of classes and objects [51, 52].

More radical forms of control structure evolution have also been explored. For example, the inclusion of combinators (higher order functions studied in the theory of functional programming languages) in the function set can allow GP to explore a large space of control architectures while imposing minimal constraints on program syntax [37, 53]. Perhaps the greatest flexibility—and therefore potentially the most intractable search space—is provided by the Push programming language, in which programs can contain arbitrary code-manipulation instructions and thereby transform their own code in arbitrary ways during execution [36, 37]. All of these innovations have been demonstrated to be useful in certain circumstances, but further study is required to determine exactly when.

7.5.3 *Evolving Development*

In nature an organism’s genes do not interact directly with its environment; rather, they direct the construction of proteins which form the organism’s body. It is that body—the phenotype—that interacts with the organism’s environment.

Several GP techniques have been inspired by the biological distinction between genotype and phenotype and by the process, called ontogeny, by which the genotype leads to the phenotype (e.g. [54–57]). In the most common approach, *developmental GP*, the programs produced by GP are structure-building programs, and it is the structures that are built by these programs, rather than the programs themselves, that are tested for fitness in the problem environment.

Typically one begins the developmental process with an “embryo” that consists of a minimal structure of the appropriate kind. The functions in the GP function set then, when executed, augment this embryo. For example, if the desired structure is a neural network then the embryo might consist of a single input node connected to a single output node and the functions in the GP function set might add additional nodes and connections [58]. Or if the desired structure is an electrical circuit then the embryo might consist of a voltage source connected to a load resistance and the functions in the GP function set might add components and wires [46].

The developmental approach has been successful in a wide range of application areas, ranging from the evolution of control systems [59] to the evolution of quantum circuits [60]. Part of its appeal comes from the way that it facilitates the application of GP to the evolution of structures that are not themselves best viewed as computer programs; developmental GP still evolves computer programs, but the programs build (develop) structures that might be quite different in nature from computer programs. Other attractions of developmental GP may derive from ways in which it affects the GP search process. For example, one might expect mutation to have a different range of effects when applied early in a developmental process than when applied to the fully developed phenotype (as is done in standard GP).

Whether this will be the case, and whether a developmental approach will therefore help or hinder, will depend on the specific problem and program representations. In biology, however, evolution often proceeds through adjustments to developmental programs and timing [61], and there is recent evidence that it can also lead to desirable properties such as robustness and self-repair in GP [62].

When the phenotype is not a computer program, developmental GP makes it easier to apply GP by making it easier to choose the function set. Because of the freedom that one has in designing a structure-building function set, developmental GP also allows you to experiment with different genotype-to-phenotype mappings, some of which may be more successful than others.

7.5.4 Evolving Evolutionary Mechanisms

The most radical expansions of the GP search space involve evolutionary control of the evolutionary process itself. There is a long history of research on self-adaptive mechanisms in evolutionary computation (e.g. see [63, 64]). In most areas outside of GP this means that numerical parameters of the evolutionary algorithm—for example mutation rates—are themselves encoded in the evolving genomes and are thereby subject to variation and selection. Similar strategies can also be applied to GP, but because GP involves the evolution of programs it is natural to ask whether a GP process can also usefully evolve its own utility programs—for example its utility program for performing mutation—and other aspects of the overall evolutionary algorithm along with the main problem-solving programs that are its primary targets.

Several approaches to self-adaptation in GP have been explored. These include several “meta-GP” approaches, in which programs implementing genetic operators (like mutation and crossover) co-evolve with problem-solving programs in separate populations [11, 65, 66]. In “autoconstructive evolution” these evolving auxiliary functions are encoded in the problem-solving programs themselves; much as in biology. Code for reproduction (mate selection, mutation, recombination, etc.) can be intermingled with, and can interact with, code for survival (problem-solving performance) in an individual’s genome [36].

The attractions of these techniques, which allow a GP system to evolve itself as it runs, stem from the possibility that the resulting systems will be adapted to their problem environments and therefore more effective than hand-designed systems. As with the other expansions to the GP search space discussed above, however, there are significant associated costs and many open research questions about how and when these costs can be overcome or justified.

7.6 Applications

There are more than 5000 recorded uses of GP. These include an enormous number of applications. It is impossible to list them all. However we shall start with a discussion of the general kinds of problems where GP has proved successful (Section 7.6.1) and the important area of symbolic regression (Section 7.6.2). Next

come sections which review the main application areas of GP: Image and Signal processing (Section 7.6.3) Finance (Section 7.6.4) Industrial Process Control (Section 7.6.5) Medicine and Bioinformatics (Section 7.6.6) Hyper-heuristics (Section 7.6.7) Entertainment and Computer Games (Section 7.6.8) and Art (Section 7.6.9). We conclude with a description of some of the human-competitive results automatically generated by GP (Section 7.6.10).

7.6.1 *Where GP Has Done Well*

If one or more of the following apply, GP may be suitable.

- The interrelationships among the relevant variables are unknown or poorly understood. GP can help discover which variables and operations are important; provide novel solutions to individual problems; unveil unexpected relationships among variables; and, sometimes GP can discover new concepts. These might then be taken and applied as in a conventional way.
- Finding the size and shape of the ultimate solution is a major part of the problem.
- Many training data are available in computer-readable form.
- There are good simulators to test the performance of tentative solutions to a problem, but poor methods to directly obtain good solutions.
In many areas there are tools to evaluate a completed design, (e.g. how far will this bridge bend under the forecast load.) Such tools solve the *direct problem* of working out the behaviour of a solution. However, the knowledge held within them cannot be easily used to solve the *inverse problem* of designing an artefact from its requirements. GP can exploit simulators and analysis tools and “data-mine” them to solve the *inverse problem* automatically.
- Conventional mathematical analysis cannot give analytic solutions.
- An approximate solution is acceptable.
- Small improvements are highly prized. Even in mature applications GP can sometimes discover small delta improvements, which may be very valuable.

Two examples are NASA’s work on satellite radio aerial design [67] and Spector’s evolution of new quantum computing algorithms that out-performed all previous approaches [68, 69]. Both of these domains are complex, do not have analytic solutions, but good simulators existed which were used to define the fitness of evolved solutions. In other words, people did not know how to solve the problems but they could (automatically) recognize a good solution when they saw one. In both cases GP discovered highly successful and unexpected designs. Also the key component of the evolved quantum algorithm was extracted and applied elsewhere [70].

7.6.2 *Curve Fitting, Data Modelling and Symbolic Regression*

There are many very good tools which will fit curves to data; however, typically they require you to specify the type of curve you want to fit, e.g. a straight line, an

exponential, a Gaussian distribution. Where GP can help is where the form of the curve or underlying model is unknown. In fact the main problem can be discovering the form of the solution or which data to use. This is generally known as *symbolic regression*.

By regression we mean finding the coefficients (e.g. slope and y -intercept) of a predefined function such that the function best fits some data. However, until a good fit is found the experimenter has to keep trying different functions by hand until a good model for the data is found. Sometimes, even expert users have strong biases when choosing functions to fit. For example, in many applications there is a tradition of using linear models, even when the data might be better fit by a more complex model. Since GP does not make this assumption, it is well suited to this sort of discovery task.

For instance, GP can evolve *soft sensors* [71]. The idea is to evolve a function which estimates what a real sensor would measure, based on data from other actual sensors in the system, (e.g. where placing an actual sensor would be expensive.) Experimental data (e.g. from industrial plant) typically come in large tables where numerous quantities are reported. Usually we know which variable we want to predict (e.g. the soft sensor value) and which other quantities we can use to make the prediction (e.g. the real sensor values). If this is not known, then experimenters must decide which are going to be their *dependent variables* before applying GP. Sometimes there are hundreds or even thousands of variables. (In bioinformatics the number of variables may approach a million.) It is well known that in these cases the efficiency and effectiveness of any machine learning or program induction method, including GP, can dramatically drop as most of the variables are typically redundant or irrelevant. This forces the system to waste considerable energy on isolating the key features. To avoid this, it is necessary to perform some form of *feature selection*, i.e. we need to decide which *independent variables* to keep and which to leave out. There are many techniques to do this, its even possible that GP itself can be used to do feature selection [72].

There are problems where more than one output (prediction) is required. For example, Table 7.4 contains data collected from a robot. The left-hand side gives four control variables, while the right-hand side contains six dependent variables measured after the robot obeyed the commands. The *Elvis* robot is shown in Figure 7.8 during the acquisition of a data sample. The roles of the independent and dependent variables are swapped when GP is given the task of controlling the arm given data from the robot's eyes.

There are several GP techniques which might be used to deal with applications where multiple outputs are required, e.g. GP individuals made of multiple trees, linear GP with multiple output registers, graph-based GP with multiple output nodes, and a single GP tree with primitives operating on vectors.

After a suitable data set has been assembled, the GP terminal set (cf. Section 7.3.1) must be defined. Since the independent variables will become the evolved code's inputs, they must be included in the terminal set. Typically some constants are also included. Next is the function set (cf. Section 7.3.2). It is often sufficient to give GP the standard four arithmetical operations (+ - × %) and an `if`. The

Arm actuator				Left eye			Right eye		
	x	y	size	x	y	size	x	y	size
-376	-626	1000	-360	44	10	29	-9	12	25
-372	-622	1000	-380	43	7	29	-9	12	29
-377	-627	899	-359	43	9	33	-20	14	26
-385	-635	799	-319	38	16	27	-17	22	30
-393	-643	699	-279	36	24	26	-21	25	20
-401	-651	599	-239	32	32	25	-26	28	18
-409	-659	500	-200	32	35	24	-27	31	19
-417	-667	399	-159	31	41	17	-28	36	13
-425	-675	299	-119	30	45	25	-27	39	8
:	:	:	:	:	:	:	:	:	:

continues for a total of 691 lines

Table 7.4 Samples showing the size and location of Elvis’s finger tip as apparent to his two eyes, given various right-arm actuator set points. GP inverts the mapping and evolves 4 functions which take data collected by both cameras (which show a target) and output instructions to the four arm motors so that his arm moves to the target.

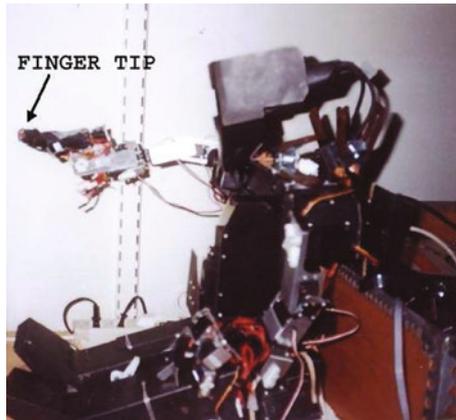


Fig. 7.8 Elvis robot sitting with his right hand outstretched. The apparent position and size of a bright red laser attached to his finger tip is recorded. The data are then used to train a GP to move the robot’s arm to a spot in three dimensions using only his eyes.

terminal and function sets are the raw components from which GP tries to build its solutions.

In virtually all symbolic regression applications the fitness function (cf. Section 7.3.3) must measure how close the outputs produced by each program are to the values of the dependent variables, when the corresponding values of the independent ones are used as inputs for the program. So, symbolic regression fitness functions tend to include summing the errors measured for each record in the data set. Usually either the absolute difference or the square of the error is used.

7.6.3 *Image and Signal Processing*

Ford were among the first to consider using GP for industrial signal processing [66]. They evolved algorithms for pre-processing electronic motor vehicle signals for possible use in engine monitoring and control.

Several applications of GP for image processing have been for military uses. For example, QinetiQ evolved programs to pick out ships using SAR radar from space satellites and to locate ground vehicles from airborne photo reconnaissance. They also used GP to process surveillance data for civilian purposes, such as predicting motorway traffic jams from subsurface traffic speed measurements [67]. Satellite images can also be used for environmental studies and for prospecting for valuable minerals [24].

Zhang has been particularly active at evolving programs with GP to visually classify objects (such as human faces) [76].

To some extent, extracting text from images (OCR) can be done fairly reliably and the accuracy rate on well-formed letters and digits is close to 100%. However, many interesting cases remain [77] such as Arabic [78] and oriental languages, handwriting [79–81] (such as the MNIST examples of handwritten digits from IRS tax returns) and musical scores [82].

The scope for applications of GP to image and signal processing is almost unbounded. A promising area is medical imaging. GP image techniques can also be used with sonar signals [83]. Off-line work on images includes security and verification. For example, [84] have used GP to detect image watermarks which have been tampered with.

7.6.4 *Financial Trading, Time Series Prediction and Economic Modelling*

GP is very widely used in these areas. It is impossible to describe all its applications instead we will just hint at a few. Chen has written more than 60 papers on using GP in finance and economics. He has investigated modelling of agents in stock markets [85], game theory, evolving trading rules for the S&P 500 [86] and forecasting the Hong Kong Hang-Seng index.

The *efficient markets hypothesis* is a tenet of economics. It is founded on the idea that everyone in a market has “perfect information” and acts “rationally”. If the efficient markets hypothesis held, then everyone would see the same value for items in the market and so agree the same price. Without price differentials, there would be no money to be made from the market itself. Whether it is trading potatoes in northern France or dollars for yen, it is clear that traders are not all equal and considerable doubt has been cast on the efficient markets hypothesis. So, people continue to play the stock market. Game theory has been a standard tool used by economists to try to understand markets but is often supplemented by simulations

with both human and computerized agents. GP is increasingly being used as part of these simulations of social systems.

The US Federal Reserve Bank used GP to study intra-day technical trading on the foreign exchange markets to suggest the market is “efficient” and found no evidence of excess returns [87]. This negative result was criticized in [88]. Later work by Neely et al. suggested that data after 1995 are consistent with Lo’s *adaptive markets hypothesis* rather than the *efficient markets hypothesis* [89]. GP and computer tools are being used in a novel data-driven approach to try and resolve issues which were previously a matter of dogma.

From a more pragmatic viewpoint, Kaboudan shows GP can forecast international currency exchange rates [90], stocks and stock returns, house prices and consumption of natural gas. Tsang and his co-workers continue to apply GP to a variety of financial arenas, including betting [91], forecasting stock prices, studying markets, approximating Nash equilibrium in game theory and arbitrage. Dempster and HSBC also use GP in foreign exchange trading [92]. Pillay has used GP in social studies and teaching aids in education, e.g. [93].

7.6.5 Industrial Process Control

Kordon and his coworkers in Dow Chemical have been very active in applying GP to industrial process control. In [94] Kordon describes where industrial GP stands now and how it will progress. Another active collaboration is that of Kovacic and Balic, who used GP in the computer numerical control of industrial milling and cutting machinery [95]. The partnership of Deschaine and Francone is most famous for their use of Discipulus for detecting bomb fragments and unexploded ordinance [96]. Genetic programming has also been used in the food processing industry. For example, Barriere et al. modelled the ripening of camembert [97].

Lewin, Dassau and Grosman applied GP to the control of an integrated circuit fabrication plant [98]. GP has also been used to identify the state of a plant to be controlled (in order to decide which of various alternative control laws to apply). For example, Fleming’s group in Sheffield used multi-objective GP [99] to reduce the cost of running aircraft jet engines.

7.6.6 Medicine, Biology and Bioinformatics

Kell and his colleagues in Aberystwyth have had great success in applying GP widely in bioinformatics [100]. Another very active medical research group is that of Moore and his colleagues at Vanderbilt [101]. Many medical data sets are very wide. Some have many thousands of inputs, but relatively few cases. (For example, a typical GeneChip data set will have tens of thousands of measurements per patient but may cover less than a hundred people [72]). Such wide data sets tend to

be avoided by traditional statistical techniques, where often the first reaction is to try and remove as many attributes as possible. Discarding whole columns of training data is often called “feature selection”. However, as has been repeatedly shown, e.g. by the Aberystwyth and Vanderbilt groups, GP can sometimes be successfully applied directly to very wide data sets.

Computational chemistry is widely used in the drug industry. Some properties of simple molecules can be calculated. However, the interactions between chemicals which might be used as drugs and medicinal targets within the body are beyond exact calculation. Therefore, there is great interest in the pharmaceutical industry in approximate *in silico* models which attempt to predict either favourable or adverse interactions between proto-drugs and biochemical molecules. Since these are computational models, they can be applied very cheaply in advance of the manufacturing of chemicals, to decide which of the myriad of chemicals might be worth for further study. Potentially, such models can make a huge impact both in terms of money and time without being anywhere near 100% correct. Machine learning and GP have both been tried. GP approaches include [102, 103].

7.6.7 GP to Create Searchers and Solvers—Hyper-heuristics

A heuristic can be considered to be a rule-of-thumb or “educated guess” that reduces the search required to find a solution. A meta-heuristic (such as a genetic algorithm) is a non-problem specific heuristic, i.e. a rule-of-thumb which can be tried on a range of problems. A hyper-heuristic is a heuristic to choose other heuristics. The difference between meta-heuristics and hyper-heuristics is that the meta-heuristic operates directly on the problem search space with the goal of finding optimal or near-optimal solutions. Hyper-heuristics operate on the heuristics search space (which consists of the heuristics used to solve the target problem). Their aim is to find good heuristics for a problem, for a certain class of instances of a problem or even for a particular instance of the problem.

GP has been very successfully used as a hyperheuristic. For example, GP has evolved competitive SAT solvers [104], state-of-the-art bin packing algorithms, particle swarm optimizers, evolutionary algorithms and travelling salesman problem solvers [105].

7.6.8 Entertainment and Computer Games

Today, a major usage of computers is interactive games. There has been some work on incorporating artificial intelligence into mainstream commercial games. Naturally the software owners are not keen on explaining exactly how much AI the games contain or giving away sensitive information on how they use AI. However, published work on GP and games includes Othello, Poker, Backgammon [106],

robotics, including robotic football, Corewares, Ms Pac-Man, radio-controlled model car racing, Draughts, and Chess. Funes [107] reports experiments which attracted thousands of people via the Internet who were entertained by evolved Tron players.

7.6.9 *The Arts*

Computers have long been used to create purely aesthetic artefacts. Much of today's computer art tends to ape traditional drawing and painting, producing static pictures on a computer monitor. However, the immediate advantage of the computer screen — movement — can also be exploited. In both cases evolutionary computation can, and has been, exploited. Indeed, with evolution's capacity for unlimited variation, evolutionary computation offers the artist the scope to produce ever changing works. The use of GP in computer art can be traced back at least to the work of Karl Sims and William Latham. Christian Jacob's work provides many examples. Many recent techniques are described in [108].

Evolutionary music has been dominated by Jazz [109], which is not to everyone's taste. Most approaches to evolving music have made at least some use of interactive evolution [110] in which the fitness of programs is provided by users, often via the Internet. The limitation is almost always finding enough people willing to participate [111]. It is surprising given their monetary value that so far little use has been made of GP to generate novel cell phone ring tones.

One of the sorrows of AI is that as soon as it works it stops being AI and becomes computer engineering. For example, the use of computer-generated images has recently become cost-effective and is widely used in Hollywood. One of the standard state-of-the-art techniques is the use of Reynold's swarming "boids" [112] to create animations of large numbers of rapidly moving animals. This was first used in *Cliffhanger* (1993) to animate a cloud of bats. Its use is now commonplace (herds of wildebeest, schooling fish, and even large crowds of people). In 1997 Craig was awarded an Oscar.

7.6.10 *Human Competitive Results: The Humies*

A particularly informative measure of the power of a problem-solving technology is its track record in solving problems that could only be solved previously by means of human intelligence and ingenuity. In order to highlight such achievements by genetic and evolutionary computation an annual competition has been held since 2004 at the Genetic and Evolutionary Computation Conference (GECCO), organized by the Association for Computing Machinery's Special Interest Group on Genetic and Evolutionary Computation (ACM SIGEVO). This competition, known as the "Humies", awards substantial cash prizes to results deemed "human competitive" as assessed by objective criteria such as patents and publications [113].

25 gold, silver and bronze “medals” with cash prizes have been awarded in the Humies competition, totaling \$45,700. Of these, 13 of the medals (4 gold, 7 silver, 2 bronze) have been awarded to teams using GP (as opposed to other genetic and evolutionary computation methods), in application areas including antenna design, quantum circuit design, mechanical engineering, optical system design, game strategy design, computer vision, and pure mathematics. Figure 7.9 shows a gold medal winning result from 2004, an antenna that was designed using GP for NASA’s Space Technology 5 mission [67]. Figure 7.10 shows a silver medal winning result from 2005, a lens system that duplicates the functionality of the patented Nagler lens system but with a novel topology [114].

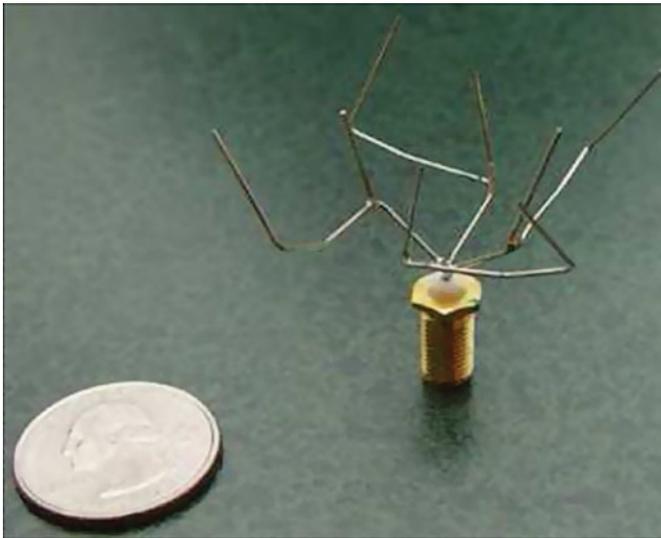


Fig. 7.9 Award winning human-competitive antenna design produced by GP.

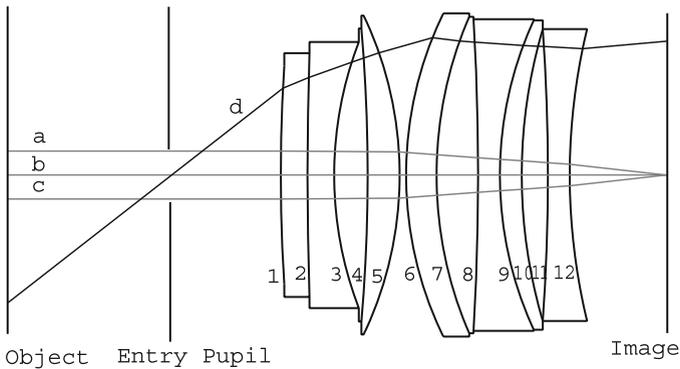


Fig. 7.10 Award winning human-competitive lens designed by GP [114, Figure 12].

7.7 Trouble-Shooting GP

The evolutionary dynamics are often very complex so it often difficult to troubleshoot evolutionary computation system. On the other hand, they can be very resilient and even GP systems with horrendous bugs can evolve valid solutions. Nonetheless, we suggest some general issues to keep in mind. To a large extent the advice in [2, 115] and [33, Chapter 9] also remains sound.

7.7.1 Can You Trust Your Results?

Here we have an interesting divergence between universities and life. Are your results about your GP system? (e.g. is it better than a GP without mutation?) Or about your application? (e.g. have you evolved a better image filter?) [5] describes how to overcome the stochastic nature of evolution and what it means to be better.

In complex applications with powerful techniques, like GP, the danger of learning the training data and so creating a solution which fits it too faithfully is ever present [116]. The classic example is where a neural network was asked to find tanks. It was presented with pictures of fields containing tanks and the same fields without tanks. After prolonged training, it could differentiate between the two. However, the final system failed to find tanks. Eventually the problem was traced to the training images. Since tanks are heavy all the pictures with tanks in them were taken on one day, the tanks were moved, and some time later another set of pictures were taken. The ANN had cheated, it had learnt (using brightness) to distinguish pictures taken early in the day from those taken later and totally ignored the presence or absence of the tanks. While it humorous and is easy to see after the fact, you must ensure the machine learning does not play this joke on you.

7.7.2 Study Your Populations

If you are not getting your desired results, take the time to dig around in the populations and see what is actually being evolved. For example, if you included a particular input or function, is it included in the better individuals? Are they using it in sensible ways? (Sometimes the tree may include an input but it has no or little impact on the program's behaviour, e.g., because it is multiplied by zero.) Is the primitive being multiply used? Similarly, if you are using grammatical evolution, are your evolved individuals using your grammar as you expected? Or is the grammar biasing the system in an undesirable or an unexpected way?

Remember GP is doing genetic search. GP increases the numbers of genes (i.e. terminals and functions) which appear in above average fitness individuals. You might keep a count of the number of times important primitives occur in the population. Mostly gene numbers vary randomly according to how lucky they are. However, look out for primitives that become extinct or (if using mutation) reduce to low

background levels. If this happens, it suggests that the fitness function is driving the current GP population in an unintended direction.

Is the distribution of fitness values of members of the population (particularly that of the better programs) dominated by a few values with large gaps between them? This suggests jumping these gaps may be hard. It also suggests that the next improvement may also be separated from the current best. So finding the next improved solution will require jumping a large gap and so be difficult. Perhaps changing your fitness function to be more continuous will improve performance considerably.

7.7.3 Studying Your Programs

A major advantage of GP is you create *visible* programs. You can see how they work. You can explain how they work to your customers. When presenting GP results, include a slide of the evolved program. The `dot` package (<http://www.graphviz.org/>) is good a starting point for nicely presented graphs. GP trees can be automatically converted to dot (<http://www.cs.ucl.ac.uk/staff/W.Langdon/lisp2dot.html>).

There are methods to automatically simplify expressions (e.g. in Mathematica and Emacs). However, simply things like removing excess significant digits and combining constant terms can make your solution more intelligible. After cleaning up the answer, make sure it still works.

In some cases the details of the trees (e.g. the particular nodes) are less important than the general size and shape. In [117], Daida describes a way to visualize the size and shape of either individual trees or an entire population, cf. Figure 7.11. (A Mathematica implementation is available via <http://library.wolfram.com/infocenter/MathSource/5163/>.)

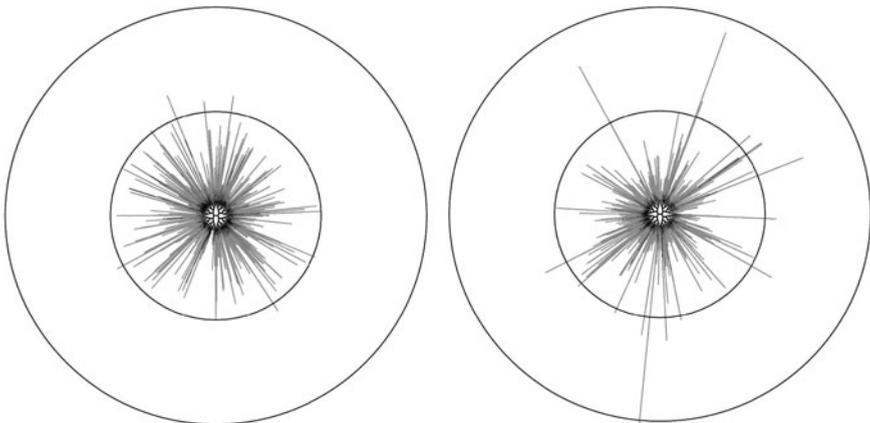


Fig. 7.11 The size and shape of 1,000 individuals in the final generation of runs using a depth limit of 50 (on the left) and a size limit of 600 (on the right). The inner circle is at depth 50, and the outer circle is at depth 100. These plots are from [119].

7.7.4 Encourage Diversity

If GP is to benefit from using a population: the population must be diverse. (Otherwise it might be much more efficient to use a hill climbing or other single point search like simulated annealing.) Sections 7.7.2 and 7.7.3 have described measuring its diversity in terms of its genes (i.e. functions and terminals) and the size and shape of the programs. You can also consider the variation in the programs' behaviour (cf., for Boolean problems, [118]). Studying behaviour, as opposed to genetic makeup, avoids the problem that GP populations often bloat [5, Sect. 11.3]. In bloated populations syntactically (i.e. genetically) evolved programs appear different but the difference is in unused code. However, studying the programs' behaviours will show if the population's phenotypes have converged excessively.

Since mutation and crossover often produce diverse but low fitness individuals you could restrict studies of population diversity to just the subset of the population which is selected to have children.

If you suspect the population has converged excessively you could

- Not use the reproduction operator.
- Add one or more mutation operators.
- Use a weaker selection mechanism, e.g., to reduce the tournament size.
- If you are using the "steady state" approach, i.e. you add new programs to the population immediately, rather than waiting until a whole new population has been created. You could choose who to overwrite at random. (Often people kill the worse member of the population and replace him with the new child. This is fine but tends to increase the convergence of the population.) You might want to protect the best member of the population to ensure he is not deleted.
- Use a generational population model instead of a steady-state model.
- The standard population is panmictic. This means there are no restrictions on which individual mates with and favourable genetic innovations rapidly spread through and may take over the whole population. In contrast a large population may be split into semi-isolated demes [5, Sect. 10.5] which keeps diversity high by slowing the spread of improvements [33, 120].

Demes are often used in conjunction with parallel hardware. The speed at which innovations spread is controlled both by the number of emigrants and how the demes are interconnected. Both all-to-all and toroidal topologies are common. They have very short paths between demes. Arranging demes in a ring gives a longer convergence time. Typically many individuals (say 2%) are transferred between demes for each generation. Over biological timescales, such a high immigration rate is sufficient to prevent a converged population diverging into separate species. However, GP is typically not run for so many generations and, from an engineering standpoint, one has to trade off rapid take up of good solutions versus searching different locations.

- Use fitness sharing or even multi-objective approaches to encourage the formation of many fitness niches.

7.7.5 Approximate Solutions Are Better than No Solution

When GP starts, typically, it starts from nowhere, i.e. well behind the state of the art. For the initial population to evolve, it must contain some programs with an “edge”. With some advantage, even if slight, over the rest of the population. You must design your selection and fitness function to amplify this. A typical fitness function gives a continuous measure of how far a program is from your requirements. The final program will be a descendent of those you select at the beginning so the fitness function must not only prefer better than random but also reward approximate solution to the whole problem which may be refined into a complete solution. Or rather, an approximate solution to the whole problem.

This is true throughout the run. At every generation the fitness function must seek out approximate solutions from which better ones can be evolved. Even standing still may not be enough. Nature tends to the easy thing. In GP this often corresponds to finding new programs which have the same fitness as their parents. Unfortunately within a few generations, they can evolve to become very resistant to change and further progress to your goal (as opposed to theirs) becomes very hard.

Consider (just for illustrative purposes) a problem with just five test cases, four of which are fairly easy and consequently less important, with the fifth being crucial and quite difficult. So the population may contain individuals that can do the four easier tasks, but are unable to make the jump to the fifth. There are several things you could try: (1) weight the hard task more heavily or (2) use a multi-objective approach. However, a more fundamental and probably more successful approach is to redesign how you sort the programs, e.g. (3) divide the task up in some way into sub-tasks, (4) provide more tasks, or (5) change it from being a binary condition (meaning that an individual does or does not succeed on the fifth task) to a continuous condition, so that an individual GP program can partially succeed on the fifth task to a greater or less extent. The idea is to create a smoother gradient for the evolutionary process to follow.

7.7.6 Control Bloat

If you are running out of memory or your execution times seem inordinately long, look the size of your evolved expressions. Often they will be growing over time. It is usually necessary to provide some form of bloat control, cf. [5, Sect. 11.3]. Controlling bloat is also important if one’s goal is to find a comprehensible model, since in practice these must be small. A large model will not only be difficult to understand but also may over-fit the training data [121].

7.7.7 Convince Your Customers

For your work to make an impact it must be presented in a form that can convince others of the validity of its results and conclusions. This might include a pitch within

a corporation seeking continued financial support for a project, the submission of a research paper to a journal or the presentation of a GP-based product to potential customers. [5] contains suggestions on improving written and verbal presentation of artificial evolution experiments. While [122, e.g., Chapter 14] has many suggestions about getting your work accepted (and paid for) by your customers.

The burden of proof is on the users of GP. It is important to use the customer's language. If the fact that GP discovered a particular chemical is important in a reaction or drug design, you should make this stand out during the presentation. A great advantage of GP over many AI techniques is that its results are often simple equations. Ensure these are intelligible to your customer, e.g., by simplification. Also make an effort to present your results using your customer's terminology. Your GP system may produce answers as trees, but if the customers use spreadsheets, consider translating the tree into a spreadsheet formula.

Also, one should try to discover how the customers intend to validate GP's answer. Do not let them invent some totally new data which have nothing to do with the data they supplied for training ("just to see how well it does..."). Avoid customers with contrived data. GP is not omnipotent, it knows nothing about things it has not seen. At the same time you should be scrupulous about your own use of holdout data. GP is a very powerful machine learning technique. With this comes the ever present danger of over-fitting. One should never allow performance on data reserved for validation to be used to choose which answer to present to the customer.

7.8 Conclusions

We have seen how genetic programming works and how to use it. We have hinted at just two of the many exciting research areas. The first widens the application of GP by using formal grammars to capture user knowledge and so guide GP. The second (so far) complementary approach extends GP by doing the opposite! That is, sometimes GP can benefit from having additional freedom, including the freedom to evolve parts of itself. We have skimmed through a few of GP's many applications, including cases where GP has evolved human competitive solutions. Finally we have tried to distill some practical "how to" knowledge into a few pages.

To conclude perhaps the best introduction to genetic programming is to create your own (or borrow someone else's) and evolve things.

References

1. Turing, A.M.: Intelligent machinery. Report for National Physical Laboratory. Reprinted in Ince, D.C. (ed.) (1992) *Mechanical Intelligence: Collected Works of A.M. Turing.* (1948)
2. Koza, J.R.: *Genetic Programming: On the Programming of Computers by Means of Natural Selection.* MIT Press (1992)
3. Darwin, C.: *The Origin of Species*, 1985 edn. John Murray, Penguin classics (1859)

4. Banzhaf, W., Nordin, P., Keller, R.E., Francone, F.D.: *Genetic Programming— An Introduction; On the Automatic Evolution of Computer Programs and its Applications*. Morgan Kaufmann, San Francisco, CA, USA (1998)
5. Poli, R., Langdon, W.B., McPhee, N.F.: A field guide to genetic programming. Published via <http://lulu.com> and freely available at <http://www.gp-field-guide.org.uk>, (2008) (With contributions by Koza, J.R.)
6. Darwin, C.: *Voyage of the Beagle*. Henry Colburn, London, penguin classics, 1989 edition, (1839)
7. Nordin, P., Banzhaf, W., Francone, F.D.: Efficient evolution of machine code for CISC architectures using instruction blocks and homologous crossover. In: Spector, L., Langdon, W.B., O'Reilly, U.-M., Angeline, P.J. (eds.) *Advances in Genetic Programming 3*, chapter 12, pp. 275–299. MIT Press, Cambridge, MA, USA (1999)
8. Foster, J.A.: Review: Discipulus: a commercial genetic programming system. *GP&EM*, **2**(2), 201–203 (2001)
9. Brameier, M., Banzhaf, W.: A comparison of linear genetic programming and neural networks in medical data mining. *IEEE Trans. EC*, **5**(1), 17–26 (2001)
10. Poli, R.: Discovery of symbolic, neuro-symbolic and neural networks with parallel distributed genetic programming. Technical Report CSRP-96-14, University of Birmingham, School of Computer Science, August 1996. Presented at 3rd International Conference on Artificial Neural Networks and Genetic Algorithms, ICANNGA'97
11. Teller, A.: Evolving programmers: The co-evolution of intelligent recombination operators. In: Angeline, P.J., Kinnear, Jr., K.E. (eds.) *Advances in Genetic Programming 2*, chapter 3, pp. 45–68. MIT Press, Cambridge, MA, USA (1996)
12. Miller, J.F., Smith, S.L.: Redundancy and computational efficiency in cartesian genetic programming. *IEEE Trans. EC*, **10**(2), 167–174 (2006)
13. Langdon, W.B.: Size fair and homologous tree genetic programming crossovers. *GP&EM*, **1**(1/2), 95–119 (2000)
14. Goldberg, D.E.: *Genetic Algorithms in Search Optimization and Machine Learning*. Addison-Wesley, (1989)
15. Angeline, P.J.: Subtree crossover: building block engine or macromutation? In: Koza et al., J.R. (eds.) *GP 1997*, pp. 9–17. Morgan Kaufmann, Stanford, 13–16 Jul.
16. Braitenberg, V.: *Vehicles*. MIT Press, (1984)
17. Poli, R., Page, J., Langdon, W.B.: Smooth uniform crossover, sub-machine code GP and demes: A recipe for solving high-order boolean parity problems. In: Banzhaf, W., Daida, J., Eiben, A.E., Garzon, M.H., Honavar, V., Jakiela, M., Smith, R.E. (eds.) *Proceedings of the Genetic and Evolutionary Computation Conference*, volume 2, pp. 1162–1169. Morgan Kaufmann, Orlando, Florida, USA (1999)
18. Chomsky, N.: Three models for the description of language. *IEEE Trans. Inf. Theory*, **2**(3), 113–124 (1956)
19. Montana, D.J.: Strongly typed genetic programming. *Evol. Comput.*, **3**(2), 199–230 (1995)
20. Recknagel et al., F.: Comparative application of artificial neural networks and genetic algorithms for multivariate time series modelling of algal blooms in freshwater lakes. *HydroInformatic*, **4**(2), 125–134 (2002)
21. Ratle, A., Sebag, M.: Genetic programming and domain knowledge: Beyond the limitations of grammar-guided machine discovery. In: Schoenauer, M., Deb, K., Rudolph, G., Yao, X., Lutton, E., JulianMerelo, J., Schwefel, H.-P. (eds.) *Parallel Problem Solving from Nature – PPSN VI 6th International Conference*, volume 1917 of LNCS, pp. 211–220. Springer Verlag, Paris, France (2000)
22. Wong, M.L., Leung, K.S.: Inductive logic programming using genetic algorithms. In: Brahan, J.W., Lasker, G.E. (eds.) *Advances in Artificial Intelligence - Theory and Application II*, pp. 119–124. I.I.A.S., Ontario, Canada (1994)
23. Hussain, T.S., Browse, R.A.: Attribute grammars for genetic representations of neural networks and syntactic constraints of genetic programming. In *Workshop on Evolutionary Computation. Held at the 12 Canadian Conference on Artificial Intelligence*, Vancouver, Canada (1998)

24. Paterson, N.R., Livesey, M.: Distinguishing genotype and phenotype in genetic programming. In: John, R. Koza, (ed.) *Late Breaking Papers at the Genetic Programming 1996 Conference Stanford University July 28–31, 1996*, pp. 141–150. Stanford Bookstore, Stanford University, CA, USA 28–31 (1996)
25. Ryan et al., C.: Grammatical evolution: evolving programs for an arbitrary language. In: Banzhaf et al., W. (eds.) *EuroGP, LNCS 1391*, pp. 83–95. Paris, 14–15 Apr (1998)
26. Hoai et al., N.X.: Representation and structural difficulty in genetic programming. *IEEE Trans. EC*, **10**(2), 157–166 (2006)
27. Joshi, A.K., Levy, L.S., Takahashi, M.: Tree adjunct grammars. *J. Comput. Syst. Sci.* **10**, 136–163 (1975)
28. Ratle, A., Sebag, M.: Avoiding the bloat with probabilistic grammar-guided genetic programming. In: Collet et al., P. (eds), *EA 2001, LNCS 2310*, pp. 255–266
29. Shan, Y., McKay, R.I., Baxter, R., Abbass, H., Essam, D., Hoai, N.X.: Grammar model-based program evolution. In: *Proceedings of the 2004 IEEE Congress on Evolutionary Computation*, pp. 478–485. IEEE Press, Portland, Oregon (2004)
30. Bosman, P.A.N., de Jong, E.D.: Grammar transformations in an EDA for genetic programming. In: Poli, R., Cagnoni, S., Keijzer, M., Costa, E., Pereira, F., Raidl, G., Upton, S.C., Goldberg, D., Lipson, H., de Jong, E., Koza, J., Suzuki, H., Sawai, H., Parmee, I., Pelikan, M., Sastry, K., Thierens, D., Stolzmann, W., Lanzi, P.L., Wilson, S.W., O’Neill, M., Ryan, C., Yu, T., Miller, J.F., Garibay, I., Hollifield, G., Wu, A.S., Riopka, T., Meysenburg, M.M., Wright, A.W., Richter, N., Moore, J.H., Ritchie, M.D., Davis, L., Roy, R., Jakiela, M., (eds.) *GECCO 2004 Workshop Proceedings*, Seattle, Washington, USA 26–30 (2004)
31. Langdon, W.B., Poli, R.: Mapping non-conventional extensions of genetic programming. *Natural Comput.*, **7**, 21–43 (2008)
32. Teller, A.: The evolution of mental models. In: Kinnear, Jr., K.E. (ed.) *Advances in GP*, ch 9, pp. 199–219. MIT Press (1994)
33. Langdon, W.B.: *Genetic Programming and Data Structures: Genetic Programming + Data Structures = Automatic Programming!*, volume 1 of *Genetic Programming*. Kluwer, Boston, (1998)
34. Yu, T., Clack, C.: PolyGP: A polymorphic genetic programming system in haskell. In: Koza, J.R., Banzhaf, W., Chellapilla, K., Deb, K., Dorigo, M., Fogel, D.B., Garzon, M.H., Goldberg, D.E., Iba, H., Riolo, R. (eds.) *Genetic Programming 1998: Proceedings of the Third Annual Conference*, pp. 416–421. University of Wisconsin, Madison, Wisconsin, USA, Morgan Kaufmann. (1998)
35. Binard, F., Felty, A.: Genetic programming with polymorphic types and higher order functions. In: Keijzer, M., Antoniol, G., Congdon, C.B., Deb, K., Doerr, B., Hansen, N., Holmes, J.H., Hornby, G.S., Howard, D., Kennedy, J., Kumar, S., Lobo, F.G., Miller, J.F., Moore, J., Neumann, F., Pelikan, M., Pollack, J., Sastry, K., Stanley, K., Stoica, A., Talbi, El-G., Wegener, I. (eds.) *GECCO '08: Proceedings of the 10th annual conference on Genetic and evolutionary computation*, pp. 1187–1194. ACM, Atlanta, GA, USA (2008)
36. Spector, L., Robinson, A.: Genetic programming and autoconstructive evolution with the push programming language. *GP&EM*, **3**(1), 7–40 (2002)
37. Spector, L., Klein, J., Keijzer, M.: The push3 execution stack and the evolution of control. In: Beyer, H.-G., O’Reilly, U.-M., Arnold, D.V., Banzhaf, W., Blum, C., Bonabeau, E.W., Cantu-Paz, E., Dasgupta, D., Deb, K., Foster, J.A., de Jong, E.D., Lipson, H., Llorca, X., Mancoridis, S., Pelikan, M., Raidl, G.R., Soule, T., Tyrrell, A.M., Watson, J.-P., Zitzler, E. (eds.) *GECCO 2005: Proceedings of the 2005 conference on Genetic and evolutionary computation*, volume 2, pp. 1689–1696. ACM Press, Washington DC, USA (2005)
38. Koza, J.R., Andre, D.: Evolution of iteration in genetic programming. In: Fogel, L.J., Angeline, P.J., Baeck, T. (eds.) *Evolutionary Programming V: Proceedings of the Fifth Annual Conference on Evolutionary Programming*, San Diego, MIT Press, (1996)
39. Yu, T.: Hierarchical processing for evolving recursive and modular programs using higher order functions and lambda abstractions. *GP&EM*, **2**(4), 345–380 (2001)

40. Whigham, P.A., McKay, R.I.: Genetic approaches to learning recursive relations. In: Yao, X. (ed.) *Progress in Evolutionary Computation*, LNAI 956, pp. 17–27. Springer (1995)
41. Brave, S.: Evolving recursive programs for tree search. In: Angeline, P.J., Kinnear, Jr., K.E. (eds.) *Advances in Genetic Programming 2*, chapter 10, pp. 203–220. MIT Press, Cambridge, MA, USA (1996)
42. Wong, M.L., Leung, K.S.: Evolving recursive functions for the even-parity problem using genetic programming. In: Angeline, P.J., Kinnear, Jr., K.E. (eds.) *Advances in GP 2*, ch 11, pp. 221–240. MIT Press (1996)
43. Yu, T.: A higher-order function approach to evolve recursive programs. In: Yu et al., T. (eds.) *GP 2005*, ch 7, pp. 93–108. Springer, Ann Arbor, 12–14 May (2005)
44. Agapitos, A., Lucas, S.M.: Learning recursive functions with object oriented genetic programming. In: Collet, P., Tomassini, M., Ebner, M., Gustafson, S., Ekárt, A. (eds.) *Proceedings of the 9th European Conference on Genetic Programming*, volume 3905 of *Lecture Notes in Computer Science*, pp. 166–177. Budapest, Hungary, Springer (2006)
45. Koza, J.R.: *Genetic Programming II: Automatic Discovery of Reusable Programs*. MIT Press, Cambridge Massachusetts, (1994)
46. Koza et al., J.R.: *Genetic Programming 3: Darwinian Invention and Problem Solving*. Morgan Kaufman (1999)
47. Angeline, P.J., Pollack, J.: Evolutionary module acquisition. In: Fogel, D., Atmar, W. (eds.) *Proceedings of the Second Annual Conference on Evolutionary Programming*, pp. 154–163. La Jolla, CA, USA (1993)
48. Kinnear, K.E. Jr.: Alternatives in automatic function definition: A comparison of performance. In: Kinnear, K.E. Jr., (ed.) *Advances in Genetic Programming*, chapter 6, pp. 119–141. MIT Press, (1994)
49. Walker, J.A., Miller, J.F.: The automatic acquisition, evolution and reuse of modules in cartesian genetic programming. *IEEE Trans. EC* **12**(4), 397–417 (2008)
50. Spector, L.: Simultaneous evolution of programs and their control structures. In: Angeline, P.J., Kinnear, K.E. Jr.: (eds.) *Advances in Genetic Programming 2*, chapter 7, pp. 137–154. MIT Press, Cambridge, MA, USA (1996)
51. Bruce, W.S.: Automatic generation of object-oriented programs using genetic programming. In: Koza et al., J.R. (eds.) *GP 1996*, pp. 267–272. MIT Press, Stanford, 28–31 Jul 1996
52. Lucas, S.: Exploiting reflection in object oriented genetic programming. In: Keijzer et al., M. (eds.) *EuroGP, LNCS 3003*, pp. 369–378. Springer 5–7 Apr (2004)
53. Briggs, F., O’Neill, M.: Functional genetic programming with combinators. In: Pham et al., T.L. (eds.) *Proceedings of the Third Asian-Pacific workshop on Genetic Programming, ASPGP*, pp. 110–127, Military Technical Academy, Hanoi, Vietnam (2006)
54. Banzhaf, W.: Genotype-phenotype-mapping and neutral variation – A case study in genetic programming. In: Davidor, Y., Schwefel, H.-P., Männer, R. (eds.) *Parallel Problem Solving from Nature III*, volume 866 of *LNCS*, pp. 322–332, Springer-Verlag, Jerusalem (1994)
55. Jacob, C.: Genetic L-system programming. In: Davidor, Y., Schwefel, H.-P., Männer, R. (eds.) *Parallel Problem Solving from Nature III*, volume 866 of *LNCS*, pp. 334–343. Springer-Verlag, Jerusalem (1994)
56. Spector, L., Stoffel, K.: Ontogenetic programming. In: Koza et al., J.R. (eds.) *GP-96*, pp. 394–399. Stanford, MIT Press 28–31 Jul (1996)
57. Harding, S.L., Miller, J.F., Banzhaf, W.: Self-modifying cartesian genetic programming. In: Thierens, D., Beyer, H.-G., Bongard, J., Branke, J., Clark, J.A., Cliff, D., Congdon, C.B., Deb, K., Doerr, B., Kovacs, T., Kumar, S., Miller, J.F., Moore, J., Neumann, F., Pelikan, M., Poli, R., Sastry, K., Stanley, K.O., Stutzle, T., Watson, R.A., Wegener, I. (eds.) *GECCO ’07: Proceedings of the 9th annual conference on Genetic and evolutionary computation*, volume 1, pp. 1021–1028. ACM Press, London (2007)
58. Gruau, F.: Genetic micro programming of neural networks. In: Kinnear, Jr., Kenneth, E. (ed.) *Advances in Genetic Programming*, chapter 24, pp. 495–518. MIT Press, (1994)
59. Koza, J.R., Keane, M.A., Streeter, M.J., Mydlowec, W., Yu, J., Lanza, G.: *Genetic Programming IV: Routine Human-Competitive Machine Intelligence*. Kluwer Academic Publishers, (2003)

60. Gottlieb, G.: *Individual development and evolution: The genesis of novel behavior*. Oxford University Press, New York (1992)
61. Gottlieb, G.: *Individual development and evolution: The genesis of novel behavior*. Oxford University Press, New York (1992)
62. Miller, J.F.: Evolving developmental programs for adaptation, morphogenesis, and self-repair. In: Banzhaf et al., W. (eds), *ECAI, LNAI 2801*, pp. 256–265. Springer 14–17 Sep (2003)
63. Angeline, P.J.: Adaptive and self-adaptive evolutionary computations. In: Palaniswami, M., Attikiouzel, Y. (eds.) *Computational Intelligence: A Dynamic Systems Perspective*, pp. 152–163. IEEE Press, (1995)
64. Meyer-Nieberg, S., Beyer, H.-G.: Self-adaptation in evolutionary algorithms. *Parameter Setting in Evolutionary Algorithm*. Springer (2006)
65. Schmidhuber, J.: Evolutionary principles in self-referential learning. on learning now to learn: the meta-meta-meta...-hook. Diploma thesis, Technische Universität München, Germany, 14 May (1987)
66. Edmonds, B.: Meta-genetic programming: Co-evolving the operators of variation. CPM Report 98-32, Centre for Policy Modelling, Manchester Metropolitan University, UK, AytounSt., Manchester, M1 3GH. UK, (1998)
67. Lohn et al., J.: Evolutionary antenna design for a NASA spacecraft. In: O'Reilly et al., U.-M. eds. *GPTP II*, ch 18, pp. 301–315. Springer, Ann Arbor, 13–15 May 2004
68. Langdon, W.B., Buxton, B.F.: Genetic programming for mining DNA chip data from cancer patients. *Genet. Prog. Evol. Mach.*, **5**(3), 251–257, (2004)
69. Spector, L., Barnum, H., Bernstein, H.J., Swamy, N.: Finding a better-than-classical quantum AND/OR algorithm using genetic programming. In: Angeline, P.J., Michalewicz, Z., Schoenauer, M., Yao, X., Zalzal, A. (eds.) *Proceedings of the Congress on Evolutionary Computation*, volume
70. Spector, L., Bernstein, H.J.: Communication capacities of some quantum gates, discovered in part through genetic programming. In: Shapiro, Jeffrey H., Hirota, O. (eds) *Proceedings of the Sixth International Conference on Quantum Communication, Measurement, and Computing (QCMC)*, pp. 500–503. Rinton Press, (2003)
71. Jordaan, E., Kordon, A., Chiang, L., Smits, G.: Robust inferential sensors based on ensemble of predictors generated by genetic programming. In: Yao, X., Burke, E., Lozano, J.A., Smith, J., Merelo-Guervós, J.J., Bullinaria, J.A., Rowe, J., Tiño Ata Kabán, P., Schwefel, H.-P. (eds.) *Parallel Problem Solving from Nature – PPSN VIII*, volume 3242 of *LNCS*, pp. 522–531. Springer-Verlag Birmingham, UK (2004)
72. Langdon, W.B., Buxton, B.F.: Genetic programming for mining DNA chip data from cancer patients. *GP&EM*, **5**(3), 251–257 (2004)
73. Hampo, R.J., Marko, K.A.: Application of genetic programming to control of vehicle systems. In: *Proceedings of the Intelligent Vehicles '92 Symposium*, pp. 191–195. IEEE, Detroit, Mi, USA (1992)
74. D. H. and S.C. Roberts. Incident detection on highways. In U.-M. O'Reilly et al., eds., *GPTP II*, ch 16, pp 263–282. Springer, Ann Arbor, 13-15 May (2004)
75. Ross et al., B.J.: Hyperspectral image analysis using genetic programming. *Appl. Soft Comput.* **5**(2), 147–156 (2005)
76. Zhang, M., Smart, W., Using gaussian distribution to construct fitness functions in genetic programming for multiclass object classification. *Pattern. Recognit. Lett.*, **27**(11), 1266–1274, *Evolutionary Computer Vision and Image Understanding*. (2006)
77. Cilibrasi, R., Vitanyi, P.M.B.: Clustering by compression. *IEEE Trans. Inf. Theory*, **51**(4), 1523–1545 (2005)
78. Klassen, T.J., Heywood, M.I.: Towards the on-line recognition of arabic characters. In *Proceedings of the 2002 International Joint Conference on Neural Networks IJCNN'02*, pp. 1900–1905. IEEE Press, Hilton Hawaiian Village Hotel, Honolulu, Hawaii, (2002)
79. De Stefano et al., C.: Character preclassification based on genetic programming. *Pattern Recogn. Lett.* **23**(12), 1439–1448 (2002)

80. Teredesai, A., Govindaraju, V.: GP-based secondary classifiers. *Pattern Recogn.* **38**(4), 505–512 (2005)
81. Parkins, A.D., Nandi, A.K.: Genetic programming techniques for hand written digit recognition. *Signal Process.*, **84**(12), 2345–2365 Dec (2004)
82. Quintana et al., M.I.: Morphological algorithm design for binary images using genetic programming. *GP&EM*, **7**(1), 81–102 (2006)
83. Martin, M.C. Evolving visual sonar: Depth from monocular images. *Pattern Recognition Letters*, **27**(11), 1174–1180 Evolutionary Computer Vision and Image Understanding. (2006)
84. Usman et al., I.: Image authenticity and perceptual optimization via genetic algorithm and a dependence neighborhood. *IJAMCS*, **4**(1), 615–620 (2007)
85. Chen, S.-H., Liao, C.-C.: Agent-based computational modeling of the stock price-volume relation. *Information Sciences*, **170**(1), 75–100 (2005)
86. Yu, T., Chen, S.-H.: Using genetic programming with lambda abstraction to find technical trading rules. In: *Computing in Economics and Finance*, University of Amsterdam, (2004)
87. Neely et al., C.J.: Is technical analysis in the foreign exchange market profitable? A GP approach. *J. Finan. Quant. Anal.* **32**(4), 405–426 (1997)
88. Marney, J.P., Miller, D., Fyfe, C., Tarbert, H.F.E.: Risk adjusted returns to technical trading rules: a genetic programming approach. In: *7th International Conference of Society of Computational Economics*, Yale, 28–29 (2001)
89. Neely et al., C.J.: The adaptive markets hypothesis: evidence from the foreign exchange market. Working Paper 2006-046B, Federal Reserve Bank of St. Louis Rev (2007)
90. Kaboudan, M.: Extended daily exchange rates forecasts using wavelet temporal resolutions. *New Math. Nat. Comput.* **1**, 79–107 (2005)
91. Tsang et al., E.P.K.: EDDIE beats the bookies. *Softw. Pract. Exper.*, **28**(10), 1033–1043 (1998)
92. Austin et al., M.P.: Adaptive systems for foreign exchange trading. *Quant. Finan.*, **4**(4), 37–45 (2004)
93. Pillay, N.: Evolving solutions to ASCII graphics programming problems in intelligent programming tutors. In: Akerkar, R. ed. *International Conference on Applied Artificial Intelligence (ICAAI'2003)*, pp. 236–243. Fort Panhala, Kolhapur, India TMRF. 15–16 Dec (2003)
94. Kordon, A.: Evolutionary computation in the chemical industry. In: Yu et al., T. (eds.) *Evolutionary Computation in Practice*, ch 11, pp. 245–262. Springer (2008)
95. Kovacic, M., Balic, J.: Evolutionary programming of a CNC cutting machine. *Int. J. Advan. Manufact. Technol.*, **22**(1–2), 118–124 (2003)
96. Deschaine, L.: Using information fusion, machine learning, and global optimisation to increase the accuracy of finding and understanding items interest in the subsurface. *GeoDrill. Int.* **122**, 30–32 May (2006)
97. Barriere, O., Lutton, E., Baudrit, C., Sicard, M., Pinaud, B., Perrot, N.: Modeling human expertise on a cheese ripening industrial process using GP. In: Rudolph, G. (ed.) *PPSN X*, volume 5199 of *LNCS*, Springer, Dortmund (2008)
98. Dassau et al., E.: Modeling and temperature control of rapid thermal processing. *Comput. Chem. Eng.*, **30**(4), 686–697 (2006)
99. Rodriguez-Vazquez et al., K.: Identifying the structure of nonlinear dynamic systems using multiobjective genetic programming. *IEEE Trans. Syst., Man Cyberne., Part A*, **34**(4), 531–545 (2004)
100. Kell, D.: Defence against the flood. *Bioinformatics World*, pp. 16–18. (2002)
101. Moore et al., J.H.: Symbolic discriminant analysis of microarray data in automimmune disease. *Genet. Epidemiol.*, **23**, 57–69 (2002)
102. Barrett, S.J., Langdon, W.B.: Advances in the application of machine learning techniques in drug discovery, design and development. In: Tiwari, A., Knowles, J., Avineri, E., Dahal, K., Roy, R. (eds.) *Applications of Soft Computing: Recent Trends*, Advances in Soft Computing, pp. 99–110. On the World Wide Web, Springer, (2006)
103. Bains et al., W.: Evolutionary computational methods to predict oral bioavailability QSPRs. *Curr. Opin. Drug Discov. Develop.*, **5**(1), 44–51 (2002)

104. Fukunaga, A.: Automated discovery of composite SAT variable selection heuristics. In: AAAI, pp. 641–648. (2002)
105. Keller, R.E., Poli, R.: Cost-benefit investigation of a genetic-programming hyperheuristic. In *Evolution Artificielle, 8th International Conference*, Lecture Notes in Computer Science, Tours, Springer, France (2007)
106. Azaria, Y., Sipper, M.: GP-gammon: genetically programming backgammon players. *Gene. Program. Evolvable Mach.*, **6**(3), 283–300 (2005)
107. Funes, P., Sklar, E., Juille, H., Pollack, J.: Animal-animat coevolution: using the animal population as fitness function. In: Pfeifer et al., R. (eds.) SAB, pp. 525–533. Zurich, Aug 17–21 (1998). MIT Press
108. Machado, P., Romero, J. (eds.): *The Art of Artificial Evolution*. Springer (2008)
109. Spector, L., Alpern, A.: Criticism, culture, and the automatic generation of artworks. In: AAAI-94, pp. 3–8. Seattle, Washington, USA AAAI Press/MIT Press (1994)
110. Takagi, H.: Interactive evolutionary computation: Fusion of the capabilities of EC optimization and human evaluation. *Proc. IEEE*, **89**(9), 1275–1296 (2001)
111. Langdon, W.B.: Global distributed evolution of L-systems fractals. In: Keijzer et al., M. (eds.) EuroGP’2004, LNCS 3003, pp. 349–358. Coimbra, Portugal, Springer 5–7 Apr (2004)
112. Reynolds, C.W.: Flocks, herds, and schools: a distributed behavioral model. *SIGGRAPH Comput. Graph.*, **21**(4), 25–34 (1987)
113. Koza et al., J.R.: Automatic creation of human-competitive programs and controllers by means of genetic programming. *GP&EM*, **1**(1/2), 121–164 (2000)
114. Koza, J.R., Al-Sakran, S.H., Jones, L.W.: Automated re-invention of six patented optical lens systems using genetic programming. In: Beyer, H.-G., O’Reilly, U.-M., Arnold, D.V., Banzhaf, W., Blum, C., Bonabeau, E.W., Cantu-Paz, E., Dasgupta, D., Deb, K., Foster, J.A., de Jong, E.D., Lipson, H., Llorca, X., Mancoridis, S., Pelikan, M., Raidl, G.R., Soule, T., Tyrrell, A.M., Watson, J.-P., Zitzler, E. (eds.) *GECCO 2005: Proceedings of the 2005 conference on Genetic and evolutionary computation*, volume 2, pp. 1953–1960. ACM Press, Washington DC, USA (2005)
115. Kinnear, K.E., Jr.: A perspective on the work in this book. In: Kinnear, Jr., K.E. (ed.) *Advances in Genetic Programming*, chapter 1, pp. 3–19. MIT Press, (1994)
116. Corney, D.P.A.: Intelligent analysis of small data sets for food design. Ph.D. thesis, University College, London (2002)
117. Daida, J.M., Hilss, A.M., Ward, D.J., Stephen, L.: Long. Visualizing tree structures in genetic programming. *Genet. Prog. Evol. Mach.*, **6**(1), 79–110 (2005)
118. McPhee, N.F., Ohs, B., Hutchison, T., Semantic building blocks in genetic programming. In: O’Neill, M., Vanneschi, L., Gustafson, S., Alcazar, A.I.E., De Falco, I., Cioppa, A.D., Tarantino, E. (eds.) *Proceedings of the 11th European Conference on Genetic Programming, EuroGP 2008*, volume 4971 of *Lecture Notes in Computer Science*, pp. 134–145. Springer, Naples (2008)
119. Crane, E.F., McPhee, N.F.: The effects of size and depth limits on tree based genetic programming. In: Yu et al., T. (eds.) *Genetic Programming Theory and Practice III*, ch 15, pp. 223–240. Springer, Ann Arbor, 12–14 May (2005)
120. Spector, L., Klein, J.: Trivial geography in genetic programming. In: Yu et al., T. (eds.) *GTP III*, ch 8, pp. 109–123. Springer, Ann Arbor, 12–14 May (2005)
121. Gelly, S., Teytaud, O., Bredeche, N., Schoenauer, M.: Universal consistency and bloat in GP. *Issue on New Methods in Machine Learning. Theory Appl. Rev. d’Intelligence Artif.*, **20**(6), 805–827 (2006)
122. Yu, T., Davis, D., Baydar, C., Roy, R. (eds.) *Evolutionary Computation in Practice*, volume 88 of *Studies in Computational Intelligence*. Springer, (2008)