

# **The ECJ Owner's Manual**

A User Manual for the ECJ Evolutionary Computation Library

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**Where to Obtain ECJ**

<http://cs.gmu.edu/~eclab/projects/ecj/>

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# Chapter 1

## Introduction

The purpose of this manual is to describe practically every feature of ECJ, an evolutionary computation toolkit. It's not a good choice to learn the system. It's very terse, boring, and long, and not organized as a tutorial but rather as an encyclopedia. Instead, I refer you to ECJ's four tutorials and various other documentation that comes with the system. But when you need to know about some particular gizmo that ECJ has available, this manual is where to look.

### 1.1 About ECJ

ECJ is an evolutionary computation framework written in Java. The system was designed for large, heavy-weight experimental needs and provides tools which provide many popular EC algorithms and conventions of EC algorithms, but with a particular emphasis towards genetic programming. ECJ is free open-source with a modified academic license which requires acknowledgment of use of the system in the body of significant published work.

ECJ is now well over ten years old and is a mature, stable framework which has (fortunately) exhibited relatively few serious bugs over the years. Its design has readily accommodated many later additions, including multiobjective optimization algorithms, island models, master/slave evaluation facilities, coevolution, steady-state and evolution strategies methods, parsimony pressure techniques, and various new individual representations (for example, rule-sets). The system is widely used in the genetic programming community and is reasonably popular in the EC community at large. I myself have used it in over thirty or forty publications.

A toolkit such as this is not for everyone. ECJ was designed for big projects and to provide many facilities, and this comes with a relatively steep learning curve. We provide tutorials and many example applications, but this only partly mitigates ECJ's imposing nature. Further, while ECJ is *extremely* "hackable", the initial development overhead for starting a new project is relatively large. As a result, while I feel ECJ is an excellent tool for many projects, other tools might be more apropos for quick-and-dirty experimental work.

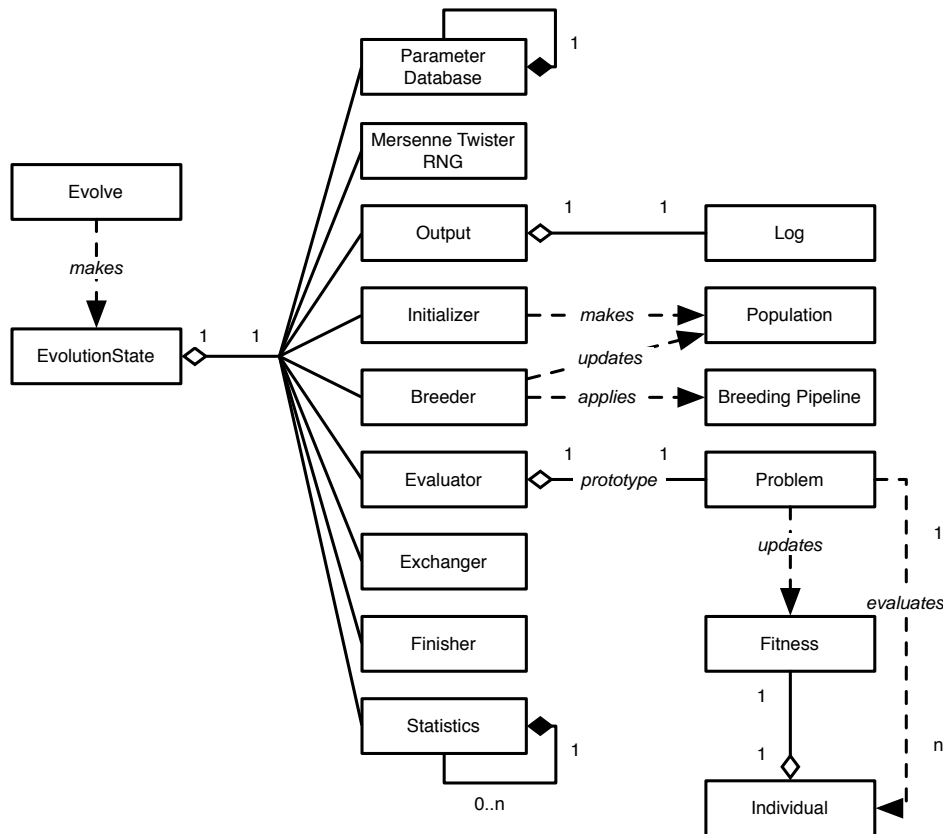
**Why ECJ was Made** ECJ's primary inspiration comes from *lil-gp* [16], from which it owes much. Homage to *lil-gp* may be found in ECJ's command-line facility, how it prints out messages, and how it stores statistics.

Work on ECJ commenced in Fall 1998 after experiences with *lil-gp* in evolving simulated soccer robot teams [5]. This project involved heavily modifying *lil-gp* to perform parallel evaluations, a simple coevolutionary procedure, multiple threading, and strong typing. Such modifications made it clear that *lil-gp* could not be further extended without considerable effort, and that it would be worthwhile developing an "industrial-grade" evolutionary computation framework in which GP was one of a number of orthogonal features. I intended ECJ to provide at least ten years of useful life, and I believe it has performed well so far.



Figure 1.1 Top-Level Loop of ECJ's SimpleEvolutionState class, used for basic generational EC algorithms. Various sub-operations are shown occurring before or after the primary operations. The full population is revised each iteration.





tionaryEvaluator, among others. There are likewise custom breeders and initializers for different functions. The Exchanger provides an opportunity for other hooks, notably internal and external island models. For example, post-breeding exchange might allow external immigrants to enter the population, while emigrants might leave the population during post-evaluation exchange. These singleton operators comprise most of the high-level “verbs” in the ECJ system, as shown in Figure 1.2.

**Parameterized Construction** ECJ is unusually heavily parameterized: practically every feature of the system is determined at runtime from a parameter. Parameters define the classes of objects, the specific subobjects they hold, and all of their initial runtime values. ECJ does this through a bootstrap class called *Evolve*, which loads a *ParameterDatabase* from runtime parameter files at startup. Using this database, *Evolve* constructs the top-level *EvolutionState* and tells it to “setup” itself. *EvolutionState* in turn calls subsidiary classes (such as *Evaluator*) and tells them to “setup” themselves from the database. This procedure continues down the chain until the entire system is constructed.

**State Objects** In addition to “verbs”, *EvolutionState* also holds “nouns” — the state objects representing the things being evolved. Specifically, *EvolutionState* holds exactly one *Population*, which contains some  $N$  (typically 1) *Subpopulations*. Multiple *Subpopulations* permit experiments in coevolution, internal island models, etc. Each *Subpopulation* holds some number of *Individuals* and the *Species* to which the *Individuals* belong. *Species* is a flyweight object for *Individual*: it provides a central repository for things common to many *Individuals* so they don’t have to each contain them in their own instances.

While running, numerous state objects must be created, destroyed, and recreated. As ECJ only learns the specific classes of these objects from the user-defined parameter file at runtime, it cannot simply construct them using Java’s new operator. Instead such objects are created by constructing a *prototype* object at startup time, and then using this object to stamp out copies of itself as often as necessary. For example, *Species* contains a prototypical *Individual*. When new *Individuals* must be created for a given *Subpopulation*, they are copied from the *Subpopulation*’s *Species* and then customized. This allows different *Subpopulations* to use different *Individual* representations.

In keeping with its philosophy of orthogonality, ECJ defines *Fitnesses* separate from *Individuals* (representations), and provides both single-objective and multi-objective *Fitness* subclasses. For historical reasons, *Subpopulation* stores the *Fitness* prototype, rather than keeping it in *Species*. Different *Subpopulations* may likewise use different *Fitness* forms.

**Breeding** A *Species* holds a prototypical *breeding pipeline* which is cloned by the *Breeder* and used per-thread to breed individuals and form the next-generation population. Breeding pipelines are tree structures where a node in the tree filters incoming *Individuals* from its child nodes and hands them to its parents. The leaf nodes in the tree are *SelectionMethods* which simply choose *Individuals* from the old subpopulation and hand them off. There exist *SelectionMethods* which perform tournament selection, fitness proportional selection, truncation selection, etc. Nonleaf nodes in the tree are *BreedingPipelines*, many of which copy and modify their received *Individuals* before handing them to their parent nodes. Some *BreedingPipelines* are representation-independent: for example, *MultiBreedingPipeline* asks for *Individuals* from one of its children at random according to some probability distribution. But most *BreedingPipelines* act to mutate or cross over *Individuals* in a representation-dependent way. For example, the *GP CrossoverPipeline* asks for one *Individual* of each of its two children, which must be genetic programming *Individuals*, performs subtree crossover on those *Individuals*, then hands them to its parent.

A tree-structured breeding pipeline allows for a rich assortment of experimenter-defined selection and breeding processes. Further, ECJ’s pipeline is *copy-forward*: *BreedingPipelines* must ensure that they copy *Individuals* before modifying them or handing them forward, if they have not been already copied. This guarantees that new *Individuals* are copies of old ones in the population, and furthermore that multiple pipelines may operate on the same *Subpopulation* in different threads without the need for locking. ECJ may apply multiple threads to parallelize the breeding process without the use of Java synchronization at all.



Figure 1.3 Top-Level data objects used in evolution.



Figure 1.4 A typed genetic programming parse tree.

**Evaluation** The Evaluator performs evaluation of a population by passing one or (for coevolutionary evaluation) several Individuals to a Problem subclass which the Evaluator has cloned off of its prototype. Evaluation may too be done in multithreaded fashion with no locking, using one Problem per thread. Individuals may also undergo repeated evaluation in coevolutionary Evaluators of different sorts.

In most projects using ECJ, the primary task is to construct an appropriate Problem subclass. The task of the Problem is to assess the fitness of the Individual(s) and set its Fitness accordingly. Problem classes also report if the ideal Individual has been discovered.

**Utilities** In addition to its ParameterDatabase, ECJ also uses a checkpointable Output convenience facility which maintains various streams, repairing them after checkpoint. Output also provides for message logging, retaining in memory all messages during the run, so that on checkpoint recovery the messages are printed out again as before. Other utilities include population distribution selectors, searching and sorting tools, etc.

The quality of a random number generator is important for a stochastic optimization system. As such, ECJ's random number generator was the very first class written in the system: it is a Java implementation of the highly respected Mersenne Twister algorithm [10] and is the fastest such implementation available. Since ECJ's release, the ECJ MersenneTwister and MersenneTwisterFast classes have found their way in a number of unrelated public-domain systems, including the popular NetLogo multiagent simulator [21]. MersenneTwisterFast is also shared in ECJ's sister software, the MASON multiagent simulation toolkit [7].

**Representations and Genetic Programming** ECJ allows you to specify any genome representation you like. Standard representation packages in ECJ provide functionality for vectors of all Java data types; arbitrary-length lists; trees; and collections of objects (such as rulesets).

ECJ is perhaps best known for its support of "Koza"-style tree-structured genetic programming representations. ECJ represents these individuals as forests of parse-trees, each tree equivalent to a single Lisp

s-expression. Figure 1.4 shows a parse-tree for a simple robot program, equivalent to the Lisp s-expression (if (and on-wall (tick > 20) (\* (ir 3) 6) 2.3). In C this might look like (onWall && tick > 20) ? ir(3) \* 6 : 2.3. This notionally says “If I’m on the wall and my tick-count is greater than 20, then return the value of my third infrared sensor times six, else return 2.3”. Such parse-trees are typically evaluated by executing their programs in a test environment, and modified via subtree crossover (swapping subtrees among individuals) or various kinds of mutation (replacing a subtree with a randomly-generated one, perhaps).

ECJ allows multiple subtrees for various experimental needs: Automatically Defined Functions (ADFs — a mechanism for evolving subroutine calls [4]), or parallel program execution, or evolving teams of programs. Along with ADFs, ECJ provides built-in support for Automatically Defined Macros (ADM)s [18] and Ephemeral Random Constants (ERCs [3], such as the numbers 20, 3, 6, and 2.3 in Figure 1.4).

Genetic programming trees are constructed out of a “primordial soup” of function templates (such as on-wall or 2.3. Early forms of genetic programming were typeless: though such templates had a predefined arity (number of arguments), any node could be connected to any other. Many genetic programming needs require more constraints than this. For example, the node if might expect a boolean value in its first argument, and integers or floats in the second and third arguments, and return a float when evaluated. Similarly and might take two booleans as arguments and return a boolean, while \* would take ints or floats as arguments and return a float.

Such types are often associated with the kinds of data passed from node to node, but they do not have to be. For example, typing might be used to constrain certain nodes to be evaluated in groups or in a certain order: for example, a function type-block might insist that its first argument be of type *foo* and its second argument be of type *bar* to make certain that a *foo* node be executed before a *bar* node.

ECJ permits a simple static typing mechanism called *set-based* typing, which is suitable for many such tasks. In set-based typing, the return type and argument types of each node are each defined to be sets of type symbols (for example, {*bool*} or {*foo*, *bar*, *baz*}, or {*int*, *float*}. The desired return type for the tree’s root is similarly defined. A child node is permitted to fit into the argument slot of a parent node if the child node’s return type and type of the that argument slot in the parent are *compatible*. We define types to be compatible if their set intersection is nonempty (that is, they share at least one type symbol).

Set-based typing is sufficient for the typing requirements found in many programming languages, including ones with type hierarchies. It allows, among other things, for nodes such as \* to accept *either* integers or floats. However there are considerable restrictions on the power of set-based typing. It’s often useful for the return type of a node to *change* based on the particular nodes which have plugged into it as arguments. For example, \* might be defined as returning a float if at least one of its arguments returns floats, but returning an integer if both of its arguments return integers. if might be similarly defined not to return a particular type, but to simply require that its return type and the second and third argument types must all match. Such “polymorphic” typing is particularly useful in situations such as matrix multiplication, where the operator must place constraints on the width and height of its arguments and the final returned matrix. In this example, it’s also useful to have an infinite number of types (perhaps to represent matrices of varying widths or heights).

ECJ does not support polymorphic typing out of the box simply because it is difficult to implement many if not most common tree modification and generation algorithms using polymorphic typing: instead, set-based typing is offered to handle as many common needs as can be easily done.

**Out of the Box Capabilities** ECJ provides support out-of-the-box for a bunch of algorithm options:

- Generational algorithms:  $(\mu, \lambda)$  and  $(\mu + \lambda)$  Evolution Strategies, the Genetic Algorithm, Genetic Programming variants, and Differential Evolution
- Steady-State evolution
- Parsimony pressure algorithms
- Spatially-embedded evolutionary algorithms
- Random restarts

- Multiobjective optimization, including the NSGA-II and SPEA2 algorithms.
- Cooperative, 1-Population Competitive, and 2-Population Competitive coevolution.
- Multithreaded evaluation and breeding.
- Parallel synchronous and asynchronous Island Models spread over a grid of computers.
- Internal synchronous Island Models internally in a single ECJ process.
- Massive parallel generational fitness evaluation of individuals on remote slave machines.
- Asynchronous Evolution, a version of steady-state evolution with massive parallel fitness evaluation on remote slave machines.
- Opportunistic Evolution, where remote slave machines run their own mini-evolutionary processes for a while before sending individuals back to the master process.
- Internal synchronous Island Models internally in a single ECJ process.
- A large number of selection and breeding operators

ECJ also has a GUI, though in truth I nearly universally use the command-line.

**Idiosyncracies** ECJ was developed near the introduction of Java and so has a lot of historical idiosyncracies.<sup>1</sup> Some of them exist to this day because of conservatism: refactoring is disruptive. If you code with ECJ, you'll definitely have to get used to one or more of the following:

- No generics at all, few iterators or enumerators, no Java features beyond 1.4 (including annotations), and little use of the Java Collections library. This is part historical, and part my own dislike of Java's byzantine generics implementation, but it's mostly efficiency. Generics are very slow when used with basic data types, as they require boxing and unboxing. The Java Collections library is unusually badly written in many places internally: and anyway, for speed we tend to work directly with arrays.
- Hand-rolled socket code. ECJ's parallel facility doesn't rely on other libraries.
- ECJ loads nearly every object from its parameter database. This means that you'll rarely see the new keyword in ECJ, nor any constructors. Instead ECJ's usual "constructor" method is a method called `setup(...)`, which sets up an object from the database.
- A proprietary logging facility. ECJ was developed before the existence of `java.util.logging`. Partly out of conservatism, I am hesitant to rip up all the pervasive logging just to use Sun's implementation (which isn't very good anyway).
- A parameter database derived from Java's old `java.util.Properties` list rather than XML. This is historical of course. But seriously, do I need a justification to avoid XML?
- Mersenne Twister random number generator. `java.lang.Random` is grotesquely bad, and systems which use it should be shunned.
- A Makefile. ECJ was developed before Ant and I've personally never needed it.

---

<sup>1</sup>It used to have a lot more — I've been weeding out ones that I think are unnecessary nowadays!

## 1.3 Unpacking ECJ and Using the Tutorials

ECJ comes as a single tarball, `ecj.tar.gz`, or as a ZIP file, `ecj.zip`. After unpacking this, you're left with one directory called `ecj`.

In the `ecj` directory you'll find several items:

- A top-level README file, which should be self-explanatory in its importance.
- ECJ's LICENSE file, which describes the primary license (AFL 3.0, a BSD-style academic license).
- A CHANGES log, which lists all past changes to all versions (including the latest).
- A Makefile. ECJ does not use Ant, and in fact you can compile ECJ very straightforwardly by simply compiling all the java files in the `ec` directory. But we provide a helpful Makefile which will compile ECJ and do various other useful tasks.
- The docs directory. This contains most of the ECJ documentation.
- The start directory. This contains various scripts for starting up ECJ: though in truth we rarely use them.
- The `ec` directory. This contains ECJ proper. `ec` is the top-level package for ECJ.

### 1.3.1 The `ec` Directory, the CLASSPATH, and jar files

The `ec` directory is ECJ's top-level package. Every subdirectory is a subpackage, and most of them are headed by helpful README files which describe the contents of the directory. Most packages contain not only Java files and class files but also parameter files and occasional data files: ECJ was designed originally for the class files to be compiled and stored right alongside the Java files in these directories, though it can be used with the separate-build-area approach taken by IDEs like Eclipse.

Because `ec` is the top-level package, you can compile ECJ, more or less, by just sticking its parent directory (the `ecj` directory), in your CLASSPATH. You will also need to add certain **jar files** in order to compile ECJ's distributed evaluation and island model facilities, and its GUI. You can get these jar files from the ECJ main website (<http://cs.gmu.edu/~eclab/projects/ecj/>). Note that none of these libraries is required. For example, if the libraries for the distributed evaluator and island model are missing, ECJ will compile but will complain if you try to run those packages with compression turned on (a feature of the packages). The GUI library is optional to ECJ, so if you don't install its libraries, you can still compile ECJ by just deleting the `ec/display` directory.

#### 1.3.1.1 The `ec/display` Directory: ECJ's GUI

This directory contains ECJ's GUI. It's in a state of disrepair and I suggest you do not use it. ECJ is really best as a command line program. In fact, as mentioned above, you can simply delete the directory and ECJ will compile just fine.

#### 1.3.1.2 The `ec/app` Directory: Demo Applications

This directory contains all the demo applications. We have quite a number of demo applications, many sharing the same subdirectories. Read the provided README file for some guidance.

### 1.3.2 The docs Directory

This directory contains all top-level documentation of ECJ except for the various README files scattered throughout the package. The `index.html` file provides the top-level entry point to the documentation.

The documentation includes:

- Introduction to parameters in ECJ
- Class documentation
- ECJ's four tutorials and post-tutorial discussion. The actual tutorial code is located in the `ec/app` directory.
- An (old) overview of ECJ
- An (old) discussion of ECJ's warts
- Some (old) graph diagrams of ECJ's structure
- This manual

#### 1.3.2.1 Tutorials

ECJ has four tutorials which introduce you to the basics of coding on the system. **I strongly suggest you go through them before continuing through the rest of this manual.** They are roughly:

1. A simple GA to solve the MaxOnes problem with a boolean representation.
2. A GA to solve an integer problem, with a custom mutation pipeline.
3. An evolution strategy to solve a floating-point problem, with a custom statistics object and reading and writing populations.
4. A genetic programming problem, plus some elitism.

As should be obvious from the rest of this manual, this **barely scratches the surface** of ECJ. No mention is given of parallelism, differential evolution, coevolution, multiobjective optimization, list and ruleset representations, grammatical encoding, spatial embedding, etc. But it'll get you up to speed.



## Chapter 2

# ec.Evolve and Utility Classes

ECJ is big. Let us begin.

ECJ's entry point is the class `ec.Evolve`. This class is little more than bootstrapping code to set up the ECJ system, construct basic datatypes, and get things going.

To run an ECJ process, you fire up `ec.Evolve` with certain runtime arguments.

```
java ec.Evolve -file myParameterFile.params -p param=value -p param=value (etc.)
```

ECJ sets itself up entirely using a **parameter file**. To this you can add additional **command-line parameters** which override those found in the parameter file. More on the parameter file will be discussed starting in Section 2.1.

ECJ can also restart from a checkpoint file it created in a previous run, like this:

```
java ec.Evolve -checkpoint myCheckpointFile.gz
```

Checkpointing will be discussed in Section 2.3.

The purpose of `ec.Evolve` is to construct an `ec.EvolutionState` instance, or load one from a checkpoint file; then get it running; and finally clean up. The `ec.EvolutionState` class actually performs the evolutionary process. Most of the stuff `ec.EvolutionState` holds is associated with evolutionary algorithms or other stochastic optimization procedures. However there are certain important utility objects or data which are created by `ec.Evolve` prior to creating the `ec.EvolutionState`, and are then stored into `ec.EvolutionState` after it has been constructed. These objects are:

- The **Parameter Database**, which holds all the parameters `ec.EvolutionState` uses to build and run the process.
- The **Output**, which handles logging and writing to files.
- The **Checkpointing Facility** to create checkpoint files as the process continues.
- The **Number of Threads** to use, and the **Random Number Generators**, one per thread.
- A simple declaration of the **Number of Jobs** to run in the process.

The remainder Section 2 discusses each of these items. It's not the most exciting of topics: but it's important in order to understand the rest of the ECJ process.

## 2.1 The Parameter Database

To build and run an experiment in ECJ, you typically write three things:

- (In Java) A **problem** which evaluates individuals and assigns fitness values to them.
- (In Java) Depending on the kind of experiment, various **components** from which individuals can be constructed — for example, for a genetic programming experiment, you'll need to define the kinds of nodes which can be used to make up the individual's tree.
- (In one or more Parameter Files) Various **parameters** which define the kind of algorithm you are using, the nature of the experiment, and the makeup of your populations and processes.

Let's begin with the third item. Parameters are the lifeblood of ECJ: practically everything in the system is defined by them. This makes ECJ highly flexible; but it also adds complexity to the system.

ECJ loads parameter files and stores them into the `ec.util.ParameterDatabase` object, which is available to nearly everything. Parameter files are an extension of the files used by Java's old `java.util.PropertyList` object. Parameter files usually end in `".params"`, and contain parameters one to a line. Parameter files may also contain blank (all whitespace) lines, which are ignored, and also lines which start with `"#"`, which are considered comments and also ignored. An example comment:

```
# This is a comment
```

The parameter lines in a parameter file typically look like this:

```
parameter.name = parameter value
```

A **parameter name** is a string of non-whitespace characters except for `"="`. After this comes some optional whitespace, then an `"="`, then some more optional whitespace.<sup>1</sup> A **parameter value** is a string of characters, including whitespace, except that all whitespace is trimmed from the front and end of the string. Notice the use of a period the parameter name. It's quite a common convention to use periods in various parameter names in ECJ. We'll get to why in a second.

Here are some legal parameter lines:

```
generations = 400
pop.subpop.0.size          =1000
pop.subpop=                ec.Subpopulation
```

Here are some illegal parameter lines:

```
generations
= 1000
pop subpop = ec.Subpopulation
```

### 2.1.1 Inheritance

Parameter files may be set up to **derive from** one or more other parameter files. Let's say you have two parameter files, `a.params` and `b.params`. Both are located in the same directory. You can set up `a.params` to derive from `b.params` by adding the following line as the very first line in the `a.params` file:

```
parent.0 = b.params
```

---

<sup>1</sup>Actually, you can omit the `"="`, but it's considered bad style.

This says, in effect: “include in me all the parameters found in the b.params file, but any parameters I myself declare will override any parameters of the same name in the b.params file.” Note that b.params may itself derive from some other file (say, c.params). In this case, a.params receives parameters from both (and parameters in b.params will likewise override ones of the same name in c.params).

Let’s say that b.params is located inside a subdirectory called foo. Then the line will look like this:

```
parent.0 = foo/b.params
```

Notice the forward slash: ECJ was designed on UNIX systems. Likewise, imagine if b.params was stored in a sibling directory called bar: then we might say:

```
parent.0 = ../bar/b.params
```

Long story short: parameter files are declared using traditional UNIX path syntax.

A parameter file can also derive from *multiple* parent parameter files, by including each at the beginning of the file, with consecutive numbers, like this:

```
parent.0 = b.params
parent.1 = yo/d.params
parent.2 = ../z.params
```

This says in effect: “first look in a.params for the parameter. If you can’t find it there, look in b.params and, ultimately, all the files b.params derives from. If you can’t find it in any of them, look in d.params and all the files it derives from. If you can’t find it in any of them, look in z.params and all the files it derives from. If you’ve still not found the parameter, give up.”

This is essentially a depth-first search through a tree or DAG, with parents overriding their children (the files they derive from) and earlier siblings overriding later siblings. Note that this multiple inheritance scheme is not the same as C++ or Lisp/CLOS, which use a distance measure!

When you fire up ECJ, you point it at a single parameter file, and you can provide additional parameters at the command-line, like this:

```
java ec.Evolve -file parameterFile.params -p command-line-parameter=value \
               -p command-line-parameter=value ...
```

Furthermore, your program itself can submit parameters to the parameter database, though it’s very unusual to do so. When a parameter is requested from the parameter database, here’s how it’s looked up:

1. If the parameter was declared by the program itself, this value is returned.
2. Else if the parameter was provided on the command line, this value is returned.
3. Else the parameter is looked up in the provided parameter file and all derived files using the inheritance ordering described earlier.
4. Else the database signals failure.

## 2.1.2 Kinds of Parameters

ECJ supports the following kinds of parameters:

- **Numbers.** Either long integers or double floating-point values. Examples:

```
generations = 500
tournament.size = 3.25
minimum-fitness = -23.45e15
```

- **Arbitrary Strings** trimmed of whitespace. Example:

```
crossover-type = two-point
```

- **Booleans.** Any value except for "false" (case-insensitive) is considered to be *true*. It's best style to use lower-case "true" and "false". The first two of these examples are *false* and the second two are *true*:

```
print-params = false
die-silently = fAlSe
pop.subpop.0.perform-injections = true
quit-on-run-complete = whatever
```

- **File Path Names.** Paths can be of three types. **Absolute paths**, which (in UNIX) begin with a "/", stipulate a precise location in the file system. **Relative paths**, which do not begin with a "/", are defined relative to the parameter file in which the parameter was located. You've seen relative paths already used for derived parameter files. Finally, **Execution relative paths** are defined relative to the directory in which the ECJ process was launched. Execution relative paths look exactly like relative paths except that they begin with the special character "\$". Examples of all three kinds of paths:

```
stat.file = $out.stat
eval.prob.map-file = ../dungeon.map
temporary-output-file = /tmp/output.txt
```

- **Class Names.** Class names are defined as the full class name of the class, including the package. Example:

```
pop.subpop.0.species = ec.gp.GPSpecies
```

- **Arrays.** ECJ doesn't have direct support for loading arrays, but has a convention you should be made aware of. It's common for arrays to be loaded by first stipulating the number of elements in the array, then stipulating each array element in turn, starting with 0. The parameter used for the number of elements differs from case to case. Note the use of periods prior to each number in the following example:

```
gp.fs.0.size = 6
gp.fs.0.func.0 = ec.app.ant.func.Left
gp.fs.0.func.1 = ec.app.ant.func.Right
gp.fs.0.func.2 = ec.app.ant.func.Move
gp.fs.0.func.3 = ec.app.ant.func.IfFoodAhead
gp.fs.0.func.4 = ec.app.ant.func.Progn2
gp.fs.0.func.5 = ec.app.ant.func.Progn3
```

The particulars vary. Here's another, slightly different, example:

```
exch.num-islands = 8
exch.island.0.id = SurvivorIsland
exch.island.1.id = GilligansIsland
exch.island.2.id = FantasyIsland
exch.island.3.id = TemptationIsland
exch.island.4.id = RhodeIsland
exch.island.5.id = EllisIsland
exch.island.6.id = ConeyIsland
exch.island.7.id = TreasureIsland
```

Anyway, you get the idea.

### 2.1.3 Namespace Hierarchies and Parameter Bases

ECJ has lots of parameters, and by convention organizes them in a namespace hierarchy to maintain some sense of order. The delimiter for paths in this hierarchy is — you guessed it — the period.

The vast majority of parameters are used by one Java object or another to set itself up immediately after it has been instantiated for the first time. ECJ has an important convention which uses the namespace hierarchy to do just this: the **parameter base**. A parameter base is essentially a path (or namespace, what have you) in which an object expects to find all of its parameters. The prefix for this path is typically the parameter name by which the object itself was loaded.

For example, let us consider the process of defining the class to be used for the global population. This class is found in the following parameter:

```
pop = ec.Population
```

ECJ looks for this parameter, expects a class (in this case, `ec.Population`), loads the class, and creates one instance. It then calls a special method (`setup(...)`, we'll discuss it later) on this class so it can set itself up from various parameters. In this case, `ec.Population` needs to know how many subpopulations it will have. This is defined by the following parameter:

```
pop.subpops = 2
```

`ec.Population` didn't know that it was supposed to look in `pop.subpops` for this value. Instead, it only knew that it needed to look in a parameter called `subpops`. The rest (in this case, `pop`) was provided to `ec.Population` as its *parameter base*: the text to be prepended — plus a period — to all parameters that `ec.Population` needed to set itself up. It's not a coincidence that the parameter base also happened to be the very parameter which defined `ec.Population` in the first place. This is by convention.

Armed with the fact that it needs to create an array of two subpopulations, `ec.Population` is ready to load the classes for those two subpopulations. Let's say that for our experiment we want them to be of different classes. Here they are:

```
pop.subpop.0 = ec.Subpopulation
pop.subpop.1 = ec.app.myapp.MySpecialSubpopulation
```

The two classes are loaded and one instance is created of each of them. Then `setup(...)` is called on each of them. Each subpopulation looks for a parameter called `size` to tell it how many individuals will be in that subpopulation. Since each of them is provided with a different parameter base, they can have different sizes:

```
pop.subpop.0.size = 100
pop.subpop.1.size = 512
```

Likewise, each of these subpopulations needs a "species". Presuming that the species are different classes, we might have:

```
pop.subpop.0.species = ec.vector.VectorSpecies
pop.subpop.1.species = ec.gp.GPSpecies
```

These species objects themselves need to be set up, and when they do, their parameter bases will be `pop.subpop.0.species` and `pop.subpop.1.species` respectively. And so on.

Now imagine that we have ten subpopulations, all of the same class (`ec.Subpopulation`), and *all but the first one* has the exact same size. We'd wind up having to write silly stuff like this:

```

pop.subpop.0.size = 1000
pop.subpop.1.size = 500
pop.subpop.2.size = 500
pop.subpop.3.size = 500
pop.subpop.4.size = 500
pop.subpop.5.size = 500
pop.subpop.6.size = 500
pop.subpop.7.size = 500
pop.subpop.8.size = 500
pop.subpop.9.size = 500

```

That's a lot of typing. Though I am saddened to report that ECJ's parameter files *do* require a lot of typing, at least the parameter database facility offers an option to save our fingers somewhat in this case. Specifically, when the `ec.Subpopulation` class sets itself up each time, it actually looks in not one but *two* path locations for the `size` parameter: first it tacks on its current base (as above), and if there's no parameter at that location, then it tries tacking on a **default base** defined for its class. In this case, the default base for `ec.Subpopulation` is the prefix `ec.subpop`. Armed with this we could simply write:

```

ec.subpop.size = 500
pop.subpop.0.size = 1000

```

When ECJ looks for subpopulation 0's size, it'll find it as normal (1000). But when it looks for subpopulation 1 (etc.), it won't find a size parameter in the normal location, so it'll look in the default location, use what it finds there (500). Only if there's no parameter to be found in either location will ECJ signal an error.

It's important to note that if a class is loaded from a default parameter, this doesn't mean that the default parameter will become its parameter base: rather, the original expect location will continue to be the base. For example, imagine if both of our Species objects were the same class, and we had defined them using the default base. That is, instead of

```

pop.subpop.0.species = ec.vector.VectorSpecies
pop.subpop.1.species = ec.vector.VectorSpecies

```

...we simply said

```

ec.subpop.species = ec.vector.VectorSpecies

```

When the species for subpopulation 0 is loaded, its parameter base is not going to be `ec.subpop.species`. Instead, it will still be `pop.subpop.0.species`. Likewise, the parameter base for the species of subpopulation 1 will still be `pop.subpop.1.species`.

Keep in mind that all of this is *just a convention*. You can use periods for whatever you like ultimately. And there exist a few global parameters without any base at all. For example, the number of generations is defined as

```

generations = 200

```

...and the seed for the random number generator the fourth thread is

```

seed.3 = 12303421

```

...even though there is *no object set up* with the `seed` parameter, and hence no object has `seed` as its parameter base. Random number generators are one of the few rare objects in ECJ which are not specified from the parameter file.

## 2.1.4 Loading Parameters

Parameters are looked up in the `ec.util.ParameterDatabase` class, and parameter names are specified using the `ec.Parameter` class. The latter is little more than a cover for Java strings. To create the parameter `pop.subpop.0.size`, we say:

```
Parameter param = new Parameter("pop.subpop.0.size");
```

Of course, usually we don't want to just make a direct parameter, but rather want to construct one from a parameter base and the remainder. Let's say our base (`pop.subpop.0`) is stored in the variable `base`, and we want to look for `size`. We do this as:

```
Parameter param = base.push("size");
```

Here are some common `ec.util.ParameterDatabase` methods:

### **ec.util.ParameterDatabase Methods**

---

```
public boolean exists(Parameter parameter, Parameter default)
```

If either parameter exists in the database, return true. Either parameter may be null.

```
public String getString(Parameter parameter, Parameter default)
```

Look first in *parameter*, then failing that, in *default parameter*, and return the result as a String, else null if not found. Either parameter may be null.

```
public File getFile(Parameter parameter, Parameter default)
```

Look first in *parameter*, then failing that, in *default parameter*, and return the result as a File, else null if not found. Either parameter may be null.

```
public Object getInstanceForParameterEq(Parameter parameter, Parameter default, Class superclass)
```

Look first in *parameter*, then failing that, in *default parameter*, to find a class. The class must have *superclass* as a superclass, or can be the *superclass* itself. Instantiate one instance of the class using the default (no-argument) constructor, and return the instance. Throws an `ec.util.ParamClassLoadException` if no class is found.

```
public Object getInstanceForParameter(Parameter parameter, Parameter default, Class superclass)
```

Look first in *parameter*, then failing that, in *default parameter*, to find a class. The class must have *superclass* as a superclass, but may not be *superclass* itself. Instantiate one instance of the class using the default (no-argument) constructor, and return the instance. Throws an `ec.util.ParamClassLoadException` if no class is found.

```
public int getBoolean(Parameter parameter, Parameter default, double defaultValue)
```

Look first in *parameter*, then failing that, in *default parameter*, and return the result as a boolean, else *defaultValue* if not found or not a boolean. Either parameter may be null.

```
public int getIntWithDefault(Parameter parameter, Parameter default, int defaultValue)
```

Look first in *parameter*, then failing that, in *default parameter*, and return the result as an int, else *defaultValue* if not found or not an int. Either parameter may be null.

```
public int getInt(Parameter parameter, Parameter default, int minValue)
```

Look first in *parameter*, then failing that, in *default parameter*, and return the result as an int, else *minValue*−1 if not found, not an int, or < *minValue*. Either parameter may be null.

```
public int getIntWithMax(Parameter parameter, Parameter default, int minValue, int maxValue)
```

Look first in *parameter*, then failing that, in *default parameter*, and return the result as an int, else *minValue*−1 if not found, not an int, < *minValue*, or > *maxValue*. Either parameter may be null.

```
public long getLongWithDefault(Parameter parameter, Parameter default, long defaultValue)
```

Look first in *parameter*, then failing that, in *default parameter*, and return the result as a long, else *defaultValue* if not found or not a long. Either parameter may be null.

```

public long getLong(Parameter parameter, Parameter default, long minValue)
    Look first in parameter, then failing that, in default parameter, and return the result as a long, else minValue−1 if not
    found, not a long, or < minValue. Either parameter may be null.

public long getLongWithMax(Parameter parameter, Parameter default, long minValue, long maxValue)
    Look first in parameter, then failing that, in default parameter, and return the result as a long, else minValue−1 if not
    found, not a long, < minValue, or > maxValue. Either parameter may be null.

public float getFloatWithDefault(Parameter parameter, Parameter default, float defaultValue)
    Look first in parameter, then failing that, in default parameter, and return the result as a float, else defaultValue if not
    found or not a float. Either parameter may be null.

public float getFloat(Parameter parameter, Parameter default, float minValue)
    Look first in parameter, then failing that, in default parameter, and return the result as a float, else minValue−1 if not
    found, not a float, or < minValue. Either parameter may be null.

public float getFloatWithMax(Parameter parameter, Parameter default, float minValue, float maxValue)
    Look first in parameter, then failing that, in default parameter, and return the result as a float, else minValue−1 if not
    found, not a float, < minValue, or > maxValue. Either parameter may be null.

public double getDoubleWithDefault(Parameter parameter, Parameter default, double defaultValue)
    Look first in parameter, then failing that, in default parameter, and return the result as a double, else defaultValue if
    not found or not a double. Either parameter may be null.

public double getDouble(Parameter parameter, Parameter default, double minValue)
    Look first in parameter, then failing that, in default parameter, and return the result as a double, else minValue−1 if
    not found, not a double, or < minValue. Either parameter may be null.

public double getDoubleWithMax(Parameter parameter, Parameter default, double minValue, double maxValue)
    Look first in parameter, then failing that, in default parameter, and return the result as a double, else minValue−1 if
    not found, not a double, < minValue, or > maxValue. Either parameter may be null.

```

---

## 2.1.5 Debugging

Your ECJ experiment is loading silently and running, but how do you know you didn't make a mistake in your parameters? How do you know ECJ is using the parameters you stated rather than some default values? If you include the following parameter in your collection:

```
print-params = true
```

...then ECJ will print out all the parameters which were used or tested for existence. For example, you might get things like this printed out:

```

!P: pop.subpop.0.file
P: pop.subpop.0.species = ec.gp.GPSpecies
<P: ec.subpop.species
P: pop.subpop.0.species.pipe = ec.breed.MultiBreedingPipeline
<P: gp.species.pipe
!E: pop.subpop.0.species.pipe.prob
P: pop.subpop.0.species.pipe.num-sources = 2
<P: breed.multibreed.num-sources
P: pop.subpop.0.species.pipe.source.0 = ec.gp.koza.CrossoverPipeline
<P: breed.multibreed.source.0
E: pop.subpop.0.species.pipe.source.0.prob = 0.9
<E: gp.koza.xover.prob

```



A **P** means that a parameter was used. An **E** means that a parameter was tested for existence. An **!** means that the parameter did not exist. A **<** means that the parameter existed in the default base as well as the primary base, but the value of the primary base was the one used. In this last case, the primary base is printed out on the line immediately prior.

There are a few other debugging parameters of less value. At the end of a run, ECJ can dump all the parameters in the database; all the parameters accessed (retrieved or tested for existence); all the parameters used (retrieved); all the parameters *not* accessed; and all the parameters *not* used. Pick your poison. Here are the relevant parameters:

```
print-all-params = true
print-accessed-params = true
print-used-params = true
print-unaccessed-params = true
print-unused-params = true
```

Typically you'd only want to set one of these to true. The most useful one is `print-unaccessed-params`, since by examining the results you can see if a parameter you set was used or not: if not, probably because it wasn't typed right. It also tells you about old, disused parameters. In fact, as I was writing this manual and needed `print-unaccessed-params` examples, I ran the Lawnmower problem (in `ec/app/lawnmower`) and got the following:

```
Unaccessed Parameters
===== (Ignore parent.x references)

gp.fs.2.info = ec.gp.GPFuncInfo
gp.koza.grow.min-depth = 5
gp.tc.0.init.max = 6
gp.koza.mutate.build.0 = ec.gp.koza.GrowBuilder
gp.tc.1.init.max = 6
parent.0 = ../../gp/koza/koza.params
gp.koza.grow.max-depth = 5
gp.tc.2.init.max = 6
gp.koza.mutate.ns.0 = ec.gp.koza.KozaNodeSelector
gp.fs.0.info = ec.gp.GPFuncInfo
gp.koza.half.growp = 0.5
gp.tc.0.init.min = 2
gp.koza.mutate.source.0 = ec.select.TournamentSelection
gp.koza.mutate.tries = 1
gp.tc.1.init.min = 2
gp.fs.1.info = ec.gp.GPFuncInfo
gp.tc.2.init.min = 2
gp.koza.mutate.maxdepth = 17
```

Most of these unaccessed parameters are perfectly fine; standard boilerplate stuff for genetic programming that didn't happen to be used by this application. But then there's the first parameter: `gp.fs.2.info = ec.gp.GPFuncInfo`, and two others like it later. I had deleted the `GPFuncInfo` class from the ECJ distribution well over a *year ago*. But apparently I forgot to remove a vestigial parameter which referred to it. Oops!

By the way, note the request to ignore "parent.x references"—this means to ignore the stuff like `parent.0 = ../../gp/koza/koza.params` that gets printed out with everything else.

## 2.2 Output

ECJ has its own idiosyncratic logging and output facility called `ec.util.Output`. This is largely historical: ECJ predates any standard logging facilities available in Java. The facility is in part inspired by a similar facility that existed in the **lil-gp** C-based genetic programming system. The system has generally worked out well so we've not seen fit to replace it.

The primary reason for the central logging and output facility is to survive checkpointing and restarting from checkpoints (see Section 2.3). Except for the occasional debugging statement which we've forgotten to remove, all output in ECJ goes through `ec.util.Output`.

The output facility has four basic features:

- **Logs**, attached to Files or to Writers, which output text of all kinds. Logs can be *restarted*, meaning that they can be reopened when ECJ is restarted from a checkpoint.
- Two dedicated Logs, the **Message Logs**, which write text out to stdout and stderr respectively.
- The ability to print arbitrary text to any Log.
- Short **Announcements** of different kinds. Announcements are different from arbitrary text in that they are not only written out to Logs (usually the stderr message Log) but are also stored in memory. This allows them to be checkpointed and automatically reposted after ECJ has started up again from a checkpoint.

The least important announcements are simple **messages**. One special kind of message is the **system message** generated by ECJ itself. Next in importance are **warnings**. One special kind of warning, the **once-only-warning**, will be written only once to a Log even if it's posted multiple times. Next are various kinds of *errors*, which can cause ECJ to quit. First, a whole bunch of basic **errors** can pile on before ECJ finally decides to quit. Second, **fatal errors** will cause ECJ to quit immediately rather than wait for more errors to accumulate.

### 2.2.1 Creating and Writing to Logs

There are many methods in `ec.util.Output` for creating Logs. Here are the two most common ones:

---

#### `ec.util.Output` Methods

```
public int addLog(File file, boolean appendOnRestart)
```

Add a log on a given file. If ECJ is restarted from a checkpoint, and *appendOnRestart* is true, then the log will be appended to the current file contents. Else they will be replaced. The Log is registered with `ec.util.Output` and its log number is returned.

```
public int addLog(File file, boolean appendOnRestart, boolean gzip)
```

Add a log on a given file. If ECJ is restarted from a checkpoint, and *appendOnRestart* is true, then the log will be appended to the current file contents. Else they will be replaced. If *gzip* is true, then the log will be gzipped. You cannot have both *appendOnRestart* and *gzip* true at the same time. The Log is registered with `ec.util.Output` and its log number is returned.

---

Two logs are always made for you automatically: a log to **stdout** (log number 0); and another log to **stderr** (log number 1). The stderr log prints all announcements, but the stdout log does not.

To write arbitrary text to a log, here are the most common methods:

---

#### `ec.util.Output` Methods

```
public void print(String text, int log number)
```

Prints a string to a log.

```
public void println(String text, int log number)
    Prints a string to a log, plus a newline.
```

---

To post a message or generate a warning or error (all of which ordinarily go to the stderr log, and are also stored in memory):

### **ec.util.Output Methods**

---

```
public void message(String text)
    Posts a message.

public void warning(String text)
    Posts a warning.

public void warning(String text, Parameter parameter, Parameter default)
    Posts a warning, and indicates the parameters which caused the warning. Typically used for cautioning the user
    about the parameters he chose.

public void warnOnce(String text)
    Posts a warning which will not appear a second time.

public void warnOnce(String text, Parameter parameter, Parameter default)
    Posts a warning which will not appear a second time, and indicates the parameters which caused the warning.
    Typically used for cautioning the user about the parameters he chose.

public void error(String text)
    Posts an error message. The contract implied in using this method is that at some point in the near future you will
    call exitIfErrors().

public void error(String text, Parameter parameter, Parameter default)
    Posts an error message, and indicates the parameters which caused the warning. Typically used for cautioning the
    user about the parameters he chose. The contract implied in using this method is that at some point in the near
    future you will call exitIfErrors().

public void exitIfErrors()
    If an error has been posted, exit.

public void fatal(String text)
    Posts an error message and exits immediately.

public void fatal(String text, Parameter parameter, Parameter default)
    Posts an error message, indicates the parameters which caused the warning, and exits immediately. Typically
    used for cautioning the user about the parameters he chose.
```

---

## **2.2.2 The ec.util.Code Class**

ECJ Individuals, Fitnesses, and various other components sometimes need to write themselves to a file in a way which can *both* be read by humans *and* be read back into Java resulting in perfect copies of the original. This means that neither printing text nor writing raw data binary is adequate.

ECJ provides a utility facility to make doing this task a little simpler. The `ec.util.Code` class encodes and decodes basic Java data types (booleans, bytes, shorts, ints, longs, floats, chars, Strings) into Strings which can be emitted as text. They all have the same pattern:

### **ec.util.Code Methods**

---

```
public static String encode(boolean val)
    Encodes val into a String and returns it.
```

```

public static String encode(byte val)
    Encodes val into a String and returns it.

public static String encode(short val)
    Encodes val into a String and returns it.

public static String encode(int val)
    Encodes val into a String and returns it.

public static String encode(long val)
    Encodes val into a String and returns it.

public static String encode(float val)
    Encodes val into a String and returns it.

public static String encode(double val)
    Encodes val into a String and returns it.

public static String encode(char val)
    Encodes val into a String and returns it.

public static String encode(String val)
    Encodes val into a String and returns it. Obviously encoding a String into a String sounds goofy, but go with us
    here.

```

---

These methods encode their data in an idiosyncratic way. Here's a table describing it:

Data Type	Encoding	Example
boolean	T or F	T
byte	<code>bvalueAsDecimalNumber </code>	<code>b59 </code>
short	<code>svalueAsDecimalNumber </code>	<code>s-321 </code>
int	<code>ivalueAsDecimalNumber </code>	<code>i42391 </code>
long	<code>lvalueAsDecimalNumber </code>	<code>l-342341232 </code>
float	<code>fvalueEncodedAsInteger valuePrintedForHumans </code>	<code>f-665866527 -9.1340002E14 </code>
double	<code>dvalueEncodedAsLong valuePrintedForHumans </code>	<code>d4614256656552045848 3.141592653589793 </code>
char	<code>'characterWithEscapes'</code>	<code>'w' or ' ' or '\n' or '\n' or '\u2FD3'</code>
String	<code>"stringWithEscapes"</code>	<code>"Dragon in Chinese is:\n\u2FD3"</code>

These are of course idiosyncratic,<sup>2</sup> but lacking a Java standard for doing the same task, they do an adequate job. You're more than welcome to go your own way.

### 2.2.2.1 Decoding the Hard Way

To decode a sequence of values from a String, you begin by creating an `ec.util.DecodeReturn` object wrapped around the String:

```
DecodeReturn decodeReturn = new DecodeReturn(string);
```

To decode the next item out of the string, you call:

```
Code.decode(decodeReturn);
```

The type of the decoded data is stored here:

---

<sup>2</sup>The eccentricities in this class stem from it being developed well before Java had any standard way to do such things itself — indeed Java still doesn't have a standard way to do most of this. I might improve it in the future, at the very least, by not requiring type symbols (like `b|`) in front of integer types. And including methods named things like `DecodeReturn.getFloat()` which throws exceptions rather than requiring one to look up type information.

```
int type = decodeReturn.type;
```

... and is one of the following `ec.util.DecodeReturn` constants:

```
public static final byte DecodeReturn.T_ERROR = -1;
public static final byte DecodeReturn.T_BOOLEAN = 0;
public static final byte DecodeReturn.T_BYTE = 1;
public static final byte DecodeReturn.T_CHAR = 2;
public static final byte DecodeReturn.T_SHORT = 3;
public static final byte DecodeReturn.T_INT = 4;
public static final byte DecodeReturn.T_LONG = 5;
public static final byte DecodeReturn.T_FLOAT = 6;
public static final byte DecodeReturn.T_DOUBLE = 7;
public static final byte DecodeReturn.T_STRING = 8;
```

If the type is a boolean (false = 0, true = 1), byte, char, short, int, or long, the result is stored here:

```
long result = decodeReturn.l;
```

If the type is a double or float, the result is stored here:

```
double result = decodeReturn.d;
```

If the type is a String, the result is stored here:

```
double result = decodeReturn.s;
```

To decode the next element out of the String, just call `Code.decode(decodeReturn)` again. Continue doing this until you're satisfied or reach a type of `T_ERROR`.

### 2.2.2.2 Easier Decoding

One of the most common decoding tasks is reading a decoded number or boolean from a single line, often preceded with a preamble, such as:

Evaluated: T

... Or ...

Size of Genome: i13|

The `Code` class has some convenience methods for decoding these lines without having to muck about with a `DecodeReturn`:

#### **ec.util.Code Methods**

---

```
public static float readFloatWithPreamble(String preamble, EvolutionState state, LineNumberReader reader)
    Decodes and returns an encoded single floating-point value from the reader, first skipping past an expected preamble. If the preamble does not exist, or the value does not exist, an error is issued.

public static float readDoubleWithPreamble(String preamble, EvolutionState state, LineNumberReader reader)
    Decodes and returns an encoded double floating-point value from the reader, first skipping past an expected preamble. If the preamble does not exist, or the value does not exist, an error is issued.

public static float readBooleanWithPreamble(String preamble, EvolutionState state, LineNumberReader reader)
    Decodes and returns an encoded boolean value from the reader, first skipping past an expected preamble. If the preamble does not exist, or the value does not exist, an error is issued.
```

```

public static byte readByteWithPreamble(String preamble, EvolutionState state, LineNumberReader reader)
    Decodes and returns an encoded byte from the reader, first skipping past an expected preamble. If the preamble
    does not exist, or the value does not exist, an error is issued.

public static short readShortWithPreamble(String preamble, EvolutionState state, LineNumberReader reader)
    Decodes and returns an encoded short from the reader, first skipping past an expected preamble. If the preamble
    does not exist, or the value does not exist, an error is issued.

public static float readIntegerWithPreamble(String preamble, EvolutionState state, LineNumberReader reader)
    Decodes and returns an encoded integer from the reader, first skipping past an expected preamble. If the preamble
    does not exist, or the value does not exist, an error is issued.

public static long readLongWithPreamble(String preamble, EvolutionState state, LineNumberReader reader)
    Decodes and returns an encoded long from the reader, first skipping past an expected preamble. If the preamble
    does not exist, or the value does not exist, an error is issued.

public static char readCharacterWithPreamble(String preamble, EvolutionState state, LineNumberReader reader)
    Decodes and returns an encoded character from the reader, first skipping past an expected preamble. If the
    preamble does not exist, or the value does not exist, an error is issued.

public static char readStringWithPreamble(String preamble, EvolutionState state, LineNumberReader reader)
    Decodes and returns an encoded string from the reader, first skipping past an expected preamble. If the preamble
    does not exist, or the value does not exist, an error is issued.

```

---

## 2.3 Checkpointing

ECJ supports **checkpointing**, meaning the ability to save the state of the stochastic optimization process to a file at any point in time, and later start a new ECJ process resuming in that exact state. Checkpointing is particularly useful when doing long processes on shared servers or other environments where the process may be killed at any time. ECJ's checkpointing procedure largely consists of applying Java's serialization mechanism to the `ec.EvolutionState` object, which in turn serializes the entire object graph of the current system.

Turn on checkpointing like this:

```
checkpoint = true
```

ECJ typically writes out checkpoint files every  $n$  generations (or, in the steady-state evolution situation, every  $n$  generations' worth of evaluations of individuals). To set  $n = 4$ , you'd say:

```
checkpoint-modulo = 4
```

ECJ writes to checkpoint files named `ec.generation.gz`, where *generation* is the current generation number. If you don't like the `ec` prefix for some reason, change it to, say, `curmudgeon` like this:

```
prefix = curmudgeon
```

Whenever a checkpoint is written, this fact is also added as an announcement. Here's the output of a typical run with checkpointing every two generations.

```

| ECJ
| An evolutionary computation system (version 19)
| By Sean Luke
| Contributors: L. Panait, G. Balan, S. Paus, Z. Skolicki, R. Kicinger, E. Popovici,
|               K. Sullivan, J. Harrison, J. Bassett, R. Hubley, A. Desai, A. Chircop,
|               J. Compton, W. Haddon, S. Donnelly, B. Jamil, and J. O'Beirne

```

```
| URL: http://cs.gmu.edu/~eclab/projects/ecj/
| Mail: ecj-help@cs.gmu.edu
|      (better: join ECJ-INTEREST at URL above)
| Date: July 10, 2009
| Current Java: 1.5.0_20 / Java HotSpot(TM) Client VM-1.5.0_20-141
| Required Minimum Java: 1.4
```

```
Threads:  breed/1 eval/1
Seed: -530434079
Job: 0
Setting up
Initializing Generation 0
Subpop 0 best fitness of generation: Fitness: -1542.1932
Generation 1
Subpop 0 best fitness of generation: Fitness: -1499.354
Checkpointing
Wrote out checkpoint file ec.2.gz
Generation 2
Subpop 0 best fitness of generation: Fitness: -1497.0482
Generation 3
Subpop 0 best fitness of generation: Fitness: -1481.9377
Checkpointing
Wrote out checkpoint file ec.4.gz
Generation 4
Subpop 0 best fitness of generation: Fitness: -1426.816
...
```

Imagine that at this point the power failed and we lost the process. We'd like to start again from the checkpoint file ec.4.gz. We can do that by typing:

```
java ec.Evolve -checkpoint ec.4.gz
```

Notice that we don't provide a parameter file or optional command-line parameters. That's because the parameter database has already been built and stored inside the checkpoint file. When ECJ starts up from a checkpoint file, it starts right where it left off, but first spits out all the announcements that had been produced up to that point, with one exception. See if you can catch it:

```
Restoring from Checkpoint ec.4.gz
```

```
| ECJ
| An evolutionary computation system (version 19)
| By Sean Luke
| Contributors: L. Panait, G. Balan, S. Paus, Z. Skolicki, R. Kicinger, E. Popovici,
|               K. Sullivan, J. Harrison, J. Bassett, R. Hubley, A. Desai, A. Chircop,
|               J. Compton, W. Haddon, S. Donnelly, B. Jamil, and J. O'Beirne
| URL: http://cs.gmu.edu/~eclab/projects/ecj/
| Mail: ecj-help@cs.gmu.edu
|      (better: join ECJ-INTEREST at URL above)
| Date: July 10, 2009
| Current Java: 1.5.0_20 / Java HotSpot(TM) Client VM-1.5.0_20-141
| Required Minimum Java: 1.4
```

```
Threads:  breed/1 eval/1
Seed: -530434079
Job: 0
```

```

Setting up
Initializing Generation 0
Subpop 0 best fitness of generation: Fitness: -1542.1932
Generation 1
Subpop 0 best fitness of generation: Fitness: -1499.354
Checkpointing
Wrote out checkpoint file ec.2.gz
Generation 2
Subpop 0 best fitness of generation: Fitness: -1497.0482
Generation 3
Subpop 0 best fitness of generation: Fitness: -1481.9377
Checkpointing
Generation 4
Subpop 0 best fitness of generation: Fitness: -1426.816
Generation 5
Subpop 0 best fitness of generation: Fitness: -1336.0835
Checkpointing
Wrote out checkpoint file ec.6.gz
Generation 6
Subpop 0 best fitness of generation: Fitness: -1302.0063
...

```

## 2.4 Threads and Random Number Generation

In many cases ECJ supports multiple threads at two stages of the evolutionary process: during *breeding* and during *evaluation*. You can specify the number of threads for each of these processes like this:

```

breedthreads = 4
evalthreads = 4

```

Typically, but not always, you'd want to set these numbers to match the number of cores or processors on your computer. And usually these two numbers should be the same. If you don't know the number of cores, you can let ECJ try to figure it out for you by saying:

```

breedthreads = auto
evalthreads = auto

```

ECJ is still capable of producing replicable results even when threading is turned on: you'll get the same results if you use the same number of evaluation and breeding threads and the same random number generator seeds. Which brings us to...

### 2.4.1 Random Numbers

As befitting its name, stochastic optimization is *stochastic*, meaning involving randomness. This means that a random number generator is central to the algorithms in ECJ, and it's crucial to have a fairly good generator. Unfortunately, Java's default random number generator, `java.util.Random`, is notoriously bad. It creates highly nonrandom sequences, so much so that websites have been developed to show off how awful it is.<sup>3</sup> **Never, ever, use `java.util.Random` in your ECJ code.**

ECJ comes with a high quality random number generator ready for you to use: `ec.util.MersenneTwisterFast`. This is a fast implementation of a famous random number generator, the Mersenne Twister.<sup>4</sup> The Mersenne Twister has a very high period and good statistical randomness qualities.

<sup>3</sup>See for example <http://alife.co.uk/nonrandom/>

<sup>4</sup><http://www.math.sci.hiroshima-u.ac.jp/~m-mat/MT/emt.html>



If you're comfortable with `java.util.Random`, you'll be fine. `ec.util.MersenneTwisterFast` has all the methods that `java.util.Random` has, plus one or two more.

In ECJ, Mersenne Twister is seeded with a single 32-bit integer other than zero (actually, it's a long, but only the first 32 bits are used). You specify this seed with the following parameter:

```
seed.0 = -492341
```

Setting the seed this way gives you control over ECJ's results: if you set the seed to the same value, ECJ will produce the exact same results again. But if you like you can also let ECJ set the seed to the current wall clock time in milliseconds, which is almost always different for different runs:

```
seed.0 = time
```

One reason ECJ's Mersenne Twister implementation is fairly fast is that it's not threadsafe. Thus ECJ maintains one random number generator for each thread used by the program. This means that if you have more than one thread, you'll have more than one random number generator, and each one of them will need a seed. Let's say you've settled on two threads. You can set both random number generator seeds like this:

```
evalthreads = 2  
breedthreads = 2  
seed.0 = -492341  
seed.1 = 93123
```

You can also use wall clock time. Specifically, if you instead do the following:

```
evalthreads = 2  
breedthreads = 2  
seed.0 = time  
seed.1 = time
```

...ECJ will guarantee that the two seeds differ. Last, if you set your threads automatically:

```
evalthreads = auto  
breedthreads = auto
```

...then ECJ will automatically set all the seeds using wall clock time, except the ones you specify by hand. After all, you don't know how many seeds you'll get!

The Mersenne Twister random number generators are stored in an array, located in a variable called `random` in the `ec.EvolutionState` object. The size of the array is the maximum of the number of breed and evaluation threads being used. How do you know which random number generator you should use? Many methods in ECJ are passed a **thread number**. This number is the index into the random number generator array for the thread in which this method is being called. For example, to get a random double, you typically see things along these lines:

```
double d = state.random[threadnum].nextDouble();
```

If you're in a single-threaded portion of the program, you can just use generator number 0.

## 2.4.2 Selecting from Distributions

Selecting from distributions is a common task in stochastic optimization. ECJ has a utility class, `ec.util.RandomChoice`, which makes it easy to set up and select from histogram-style (arbitrary) distributions, such as selecting randomly from a Population by Fitness.

The distributions in question come in the form of arrays of floats, doubles, or special objects which can provide their own float or double values. The values in these arrays are expected to form a probability density function (PDF). The objective is to select indexes in this array proportional to their value. To begin,

you call one of the following methods on your array to have RandomChoice convert it into a Cumulative Density Function (CDF) to make selection easier:

### ec.util.RandomChoice Methods

---

`public static void organizeDistribution(float[] probabilities, boolean allowAllZeros)`

If the array is all zeros, then if *allowAllZeros* is false, then an ArithmeticException is thrown, else the array is converted to all ones. Then the array is converted to a CDF. If the array has negative numbers or is of zero length, an Arithmetic Exception is thrown.

`public static void organizeDistribution(float[] probabilities)`

If the array is all zeros, then if *allowAllZeros* is false, then an ArithmeticException is thrown. If not, then the array is converted to a CDF. If the array has negative numbers or is of zero length, an Arithmetic Exception is thrown.

`public static void organizeDistribution(double[] probabilities, boolean allowAllZeros)`

If the array is all zeros, then if *allowAllZeros* is false, then an ArithmeticException is thrown, else the array is converted to all ones. Then the array is converted to a CDF. If the array has negative numbers or is of zero length, an Arithmetic Exception is thrown.

`public static void organizeDistribution(float[] probabilities)`

If the array is all zeros, then if *allowAllZeros* is false, then an ArithmeticException is thrown. If not, then the array is converted to a CDF. If the array has negative numbers or is of zero length, an Arithmetic Exception is thrown.

`public static void organizeDistribution(Object[] objs, RandomChoiceChooser chooser, boolean allowAllZeros)`

The objects in *objs* are passed to *chooser* to provide their floating-point values (and to set them if needed). If the array is all zeros, then if *allowAllZeros* is false, then an ArithmeticException is thrown, else the array is converted to all ones. Then the array is converted to a CDF. If the array has negative numbers or is of zero length, an Arithmetic Exception is thrown.

`public static void organizeDistribution(Object[] objs, RandomChoiceChooser chooser)`

The objects in *objs* are passed to *chooser* to provide their floating-point values (and to set them if needed). If the array is all zeros, then if *allowAllZeros* is false, then an ArithmeticException is thrown. Then the array is converted to a CDF. If the array has negative numbers or is of zero length, an Arithmetic Exception is thrown.

`public static void organizeDistribution(Object[] objs, RandomChoiceChooserD chooser, boolean allowAllZeros)`

The objects in *objs* are passed to *chooser* to provide their double-floating-point values (and to set them if needed). If the array is all zeros, then if *allowAllZeros* is false, then an ArithmeticException is thrown, else the array is converted to all ones. Then the array is converted to a CDF. If the array has negative numbers or is of zero length, an Arithmetic Exception is thrown.

`public static void organizeDistribution(Object[] objs, RandomChoiceChooserD chooser)`

The objects in *objs* are passed to *chooser* to provide their double-floating-point values (and to set them if needed). If the array is all zeros, then if *allowAllZeros* is false, then an ArithmeticException is thrown. Then the array is converted to a CDF. If the array has negative numbers or is of zero length, an Arithmetic Exception is thrown.

---

These methods rely on two special interfaces, **RandomChoiceChooser** (ec.util.RandomChoiceChooser) and **RandomChoiceChooserD** (ec.util.RandomChoiceChooserD). RandomChoiceChooser requires two methods which map between Objects and floating-point values.

```
public float getProbability(Object obj);
public void setProbability(Object obj, float probability);
```

RandomChoiceChooserD is the same, except that it's used for double values:

```
public double getProbability(Object obj);
public void setProbability(Object obj, double probability);
```

Once the array has been modified, you can then select random indexes from it. This is done by first generating a random floating-point number from 0...1, then passing that number into one of the following methods.

---

### ec.util.RandomChoice Methods

---

```
public static int pickFromDistribution(float[ ] probabilities, float probability)
    Selects and returns an index in the given array which contains the given probability.

public static int pickFromDistribution(double[ ] probabilities, double probability)
    Selects and returns an index in the given array which contains the given probability.

public static int pickFromDistribution(Object[ ] objs, RandomChoiceChooser chooser, float probability)
    Selects and returns an index in the given array which contains the given probability. The chooser will provide the
    floating-point values of each element in the array.

public static int pickFromDistribution(Object[ ] objs, RandomChoiceChooserD chooser, double probability)
    Selects and returns an index in the given array which contains the given probability. The chooser will provide the
    floating-point values of each element in the array.
```

---

## 2.5 Jobs

Perhaps you need to run ECJ 50 times and collect statistics from all 50 runs. ECJ's `ec.Evolve` class provides a rudimentary but extensible jobs facility to do this. You specify the number of jobs as follows:

```
jobs = 50
```

Each job will automatically use a different set of random number generator seeds. Additionally, if there is more than one job, ECJ will prepend each statistics file with `job.jobnumber`. For example, if we ran with just a single job (the default) we'd probably create an output statistics file called `out.stat`. But if we ran with multiple jobs, during the fourth job we'd create the output statistics file as `job.3.out.stat` (jobs start with 0).

Jobs are restarted properly from checkpoints: when you resume from a checkpoint, you'll start up right in that job and continue from there. This is accomplished by storing the job parameter, and the runtime arguments, in the `ec.EvolutionState` object. See extended comments in the `ec.Evolve` source code for more information.

But what if you need more job complexity? For example, what if you want to run ECJ with 10 different parameter settings and 50 runs per setting? You'll need to do some coding.

For example, let's say we want to run 50 jobs, and each job changes the generation length. The first job has 20 generations, the second job has 21 generations, etc. Here's the trick. After ECJ creates the `ec.EvolutionState` object, it calls the `startFresh()` method on this object. The default implementation calls `setup(...)` on this object, then starts running the evolutionary loop. Here's your chance. Let's say you're using the common `EvolutionState` `ec.simple.SimpleEvolutionState` subclass as your `EvolutionState`. Override the `startFresh(...)` method in a custom subclass of `SimpleEvolutionState` along these lines:

```

public class ec.app.myexample.MySimpleEvolutionState extends ec.simple.SimpleEvolutionState
{
    public void startFresh()
    {
        // setup() hasn't been called yet, so very few instance variables are valid at this point.
        // Here's what you can access:  parameters, random, output, evalthreads, breedthreads,
        //                                randomSeedOffset, job, runtimeArguments, checkpointPrefix
        // Let's modify the 'generations' parameter based on the job number
        int jobNum = ((Integer)(job[0])).intValue();
        parameters.set(new ec.util.Parameter("generations"), "" + (jobNum + 20));

        // call super.startFresh() here at the end.  It'll call setup() from the parameters
        super.startFresh();
    }
}

```

Now we need to stipulate that this is our EvolutionState, by changing the state parameter:

```

state = ec.app.myexample.MySimpleEvolutionState
jobs = 50

```

In truth though, in all my experiments, I have personally *always* handled different parameter settings in a completely different way: on the command line using a UNIX script. It's much simpler than mucking with Java code. For example, to run ten runs each of five different population sizes, perhaps you could do this (in the tcsh shell language):

```

@ seed = 92341
foreach size (2 4 8 16 32)
    foreach job (1 2 3 4 5 6 7 8 9 10)
        @ seed = ${seed} + 17
        java ec.Evolve -file ant.params \
            -p seed.0=${seed} \
            -p pop.subpop.0.size=${size} \
            -p stat.file=out.${size}.${job}.stat
    end
end

```

## 2.6 The ec.Evolve Top-level

ECJ's Evolve.java class looks complex but that's just because it has various gizmos to do jobs, checkpoint handling, etc. But in fact, the top-level loop for ECJ can be quite small. Here's all you need to do:

1. Try to load an ec.EvolutionState from a checkpoint file.
2. If it loads call run(...) on it.
3. Otherwise...
  - (a) Load a parameter database.
  - (b) Create a new ec.EvolutionState from the parameter database.
  - (c) Call run(...) on it.
4. Call cleanup(...) and exit.

Here's the basic code. As you can see, it's not very complex.

```
public static void main(String[] args)
{
    EvolutionState state = Evolve.possiblyRestoreFromCheckpoint(args);
    if (state!=null) // loaded from checkpoint
        state.run(EvolutionState.C_STARTED_FROM_CHECKPOINT);
    else
    {
        ParameterDatabase parameters = Evolve.loadParameterDatabase(args);
        state = Evolve.initialize(parameters, 0);
        state.run(EvolutionState.C_STARTED_FRESH);
    }
    Evolve.cleanup(state);
    System.exit(0);
}
```

The source code for `ec.Evolve` is very heavily commented with examples and ideas for customization. Check it out!



## Chapter 3

# ec.EvolutionState and the ECJ Evolutionary Process

As discussed in Section 2, the purpose of the `ec.Evolve` class is simply to set up an `ec.EvolutionState` and get it going. `ec.EvolutionState` is the central object in all of ECJ.

An ECJ process has only one `ec.EvolutionState` instance. Practically everything in ECJ, except for `ec.Evolve` itself, is pointed to somehow from `ec.EvolutionState`, so if you checkpoint `ec.EvolutionState`, the entire ECJ process is written to disk. Various subclasses of `ec.EvolutionState` define the stochastic optimization process. And a great many methods are handed the `ec.EvolutionState` instance, and so have essentially global access to the system.

If you peek inside `ec.EvolutionState`, you will find a number of objects, as shown in Figure 3.1:

- Some familiar objects, placed there by `ec.Evolve` after it created the `ec.EvolutionState`: the **Parameter Database**, **Output**, and array of **Random Number Generators**. Additionally, the number of **breeding threads** and **evaluation threads** (and various checkpoint and job stuff, not shown below):

```
public ec.util.ParameterDatabase parameters;  
public ec.util.MersenneTwisterFast[] random;  
public ec.util.Output output;  
public int breedthreads;  
public int evalthreads;
```

- A **Population**, which holds the individuals in the evolutionary process; plus the current **generation** (iteration) of the evolutionary process and the total **number of generations** to run (how or whether these last two variables are used depends on the evolutionary process in question):

```
public ec.Population population;  
public int generation;  
public int numGenerations;
```

- An **Initializer**, whose job is to create the initial `ec.Population` at the beginning of the run, and a **Finalizer**, whose job is to clean up at the very end of the run.

```
public ec.Initializer initializer;  
public ec.Finalizer finalizer;
```

- An **Evaluator**, whose job is assign quality assessments (fitnesses) to each member of the `Population`, and a **Breeder**, whose job is to produce a new `Population` from the previous-generation's `Population` through some collection of selection and modification operators.



Figure 3.1 Top-Level operators and utility facilities in EvolutionState, and their relationship to certain state objects. A repeat of Figure 1.2. Compare to Figure 4.2.

```
public ec.Evaluator evaluator;
public ec.Breeder breeder;
```

- An **Exchanger**, which optionally exports Population members to other ECJ processes, or imports ones to add to the Population. And finally, a **Statistics** object, whose methods are called at many points in the run to output statistics on the current run performance.

```
public ec.Exchanger exchanger;
public ec.Statistics statistics;
```

## 3.1 Common Patterns

Most of ECJ's classes follow certain patterns which you'll see many times, so it's useful to review them here.

### 3.1.1 Setup

Nearly all classes adhere to the `ec.Setup` interface. This interface is `java.io.Serializable` (which is why ECJ can serialize all its objects) and defines a single additional method:



---

## ec.Setup Methods

```
public void setup(EvolutionState state, Parameter base)
```

Constructs the Setup object from the Parameter Database using *base* as the primary parameter base. Nearly all ECJ classes implement this method.

---

ECJ objects are born from ECJ's Parameter Database, which constructs them with the default (no-argument) constructor. Then they have `setup(...)` called on them, and are expected to construct themselves by loading parameters as necessary from the Parameter Database (`state.parameters`), using the provided Parameter base. Thus the `setup(...)` method is, for all intents and purposes, the constructor.

When implementing `setup(...)` always call `super.setup(...)` if a superclass exists.

### 3.1.2 Singletons and Cliques

Singletons (`ec.Singleton`) are Setups which create a single instance and that's it. For example, `ec.EvolutionState` is a Singleton, as are `ec.Initializer`, `ec.Finalizer`, `ec.Evaluator`, `ec.Breeder`, `ec.Exchanger`, and `ec.Statistics`. Singletons are generally meant to be globally accessible.

Though Singleton are single objects, Cliques (`ec.Clique`) are objects for which only a small number (but usually more than 1) are created. Cliques are also generally meant to be globally accessible. Most Cliques have a globally accessible registry of some sort in which all Clique members can be found.

Because they are global, Prototypes and Singletons usually are set up from a single parameter base (the one provided by `setup(...)`).

### 3.1.3 Prototypes

Prototypes (`ec.Prototype`) are by far the most common objects in ECJ. Prototypes are Setups which follow the following design pattern: only one instance is loaded from the Parameter Database and set up; this object is the *prototype*. Then many objects are created by *deep cloning* the prototype. One example of a Prototype is an Individual (`ec.Individual`): a single prototypical Individual is created when ECJ starts up; and further Individuals are deep cloned from this prototype to fill Populations.

Because they can be deep cloned, Prototypes implement the `java.lang.Cloneable` interface, so you must implement the method:

---

## ec.Prototype Methods

```
public Object clone()
```

Deep-clones the object. Implemented by all Prototypes. Must call `super.clone()`, possibly catching a thrown `CloneNotSupportedException`.

---

Unlike Singletons and Cliques, Prototypes also usually have *two* parameter bases: the primary base provided by `setup(...)`, and a default base. As a result, Prototypes must implement a method which can be called to provide this default base:

---

## ec.Prototype Methods

```
public ec.util.Parameter defaultBase()
```

Returns the default base for the Prototype.

---

The standard way to implement this method is to consult a special **defaults class** in the Parameter's Java package. For example, in the `ec.simple` package the defaults class is `ec.simple.SimpleDefaults`. Here's the entirety of this class:

```

public final class SimpleDefaults implements ec.DefaultsForm
{
    public static final String P_SIMPLE = "simple";
    public static final Parameter base() { return new Parameter(P_SIMPLE); }
}

```

The Parameter returned by `base()` here provides **package default base** for the `ec.simple` package. Now consider `ec.simple.SimpleFitness`, a Prototype in this package. This class implements the `defaultBase()` method like this:

```

public static final String P_FITNESS = "fitness";
public Parameter defaultBase()
{
    return SimpleDefaults.base().push(P_FITNESS);
}

```

Thus, as a result the default parameter base for `ec.simple.SimpleFitness` is `simple.fitness`.

### 3.1.4 The Flyweight Pattern

Many Prototypes follow what is commonly known as the **flyweight pattern**. Prototypes are often great in number and Java is a memory hog: so it's helpful for groups of Prototypes to place shared information common to them in a single central location rather than keep copies of their own. For various reasons (particularly because it's hard to do serialization) ECJ doesn't use static variables to store this common information. Instead groups of Prototypes often all point to a special object which contains information common to all of them. For example, instances of `ec.Individual`, in groups, typically share a common `ec.Species` which contains information common to them. At any particular time there may be several such groups of Individuals, each with a *different* Species.

### 3.1.5 Groups

Groups are similar to Prototypes in that a single object is loaded from the Parameter Database and further objects are created by a cloning procedure. Groups are likewise `java.lang.Cloneable`. However, Groups are different in that there is no prototype per se: the object loaded from the Parameter Database isn't held in reserve but is actively used. It must not just clone another object, but actually create a new, fresh, clean object ready to be used. This is done by implementing the method:

---

#### ec.Group Methods

```
public ec.util.Parameter emptyClone()
```

Returns a pristine, new clone of the Group which has been emptied of members.

---

This method is normally implemented by cloning the object, cleaning out the clone, and returning it in the same pristine state that it would be if it had been created directly from the Parameter Database. At present there are only a few ECJ objects which implement Group: namely, `ec.Population`, `ec.Subpopulation`, and certain specialized subclasses of `ec.Subpopulation`.

## 3.2 Populations, Subpopulations, Species, Individuals, and Fitnesses

Populations, Subpopulations, and Individuals are the “nouns” of an evolutionary system, and Fitnesses are the “adjectives”. They're pretty central to the operation of any evolutionary or sample-based stochastic search algorithm.



Figure 3.2 Top-Level data objects used in evolution. A repeat of Figure 1.3.

In ECJ, an **individual** is a candidate solution to a problem. Some  $M$  Individuals are grouped together into a sample of solutions known as a **subpopulation**. Some  $N$  subpopulations are grouped together into the system's **population**. There's only one population per evolutionary process. The most common scenario is for ECJ to have  $M$  individuals grouped into a single subpopulation, which then is the sole member of ECJ's population. However, coevolutionary algorithms (Section 7.1) typically have  $N > 1$  subpopulations: as do a special and little-used internal island model scheme (see Section 6.2).<sup>1</sup>

Usually ECJ's population is an instance of the class `ec.Population` and its subpopulations are instances of the class `ec.Subpopulation`. Both of these are Groups. Let's say that there's a single subpopulation, which must contain 100 individuals. We'll express this as follows:

```
pop = ec.Population
pop.subpops = 1
pop.subpop.0 = ec.Subpopulation
pop.subpop.0.size = 100
```

Obviously further subpopulations would be `pop.subpop.1`, `pop.subpop.2`, etc. The population is found in an instance variable in the `EvolutionState`:

```
public Population population;
```

The Population is little more than an array of Subpopulations. To get Subpopulation 0, with the `EvolutionState` being `state`, you'd say:

```
Subpopulation theSubpop = state.population.subpops[0];
```

Subpopulations themselves contain arrays of individuals. To get Individual 15 of Subpopulation 0, you'd say:

```
Individual theIndividual = state.population.subpops[0].individuals[15];
```

In addition to an array of individuals, each subpopulation contains a **species** which defines the individuals used to fill the subpopulation, as well as their fitness and the means by which they are modified. Subpopulations also contain some basic parameters for creating initial individuals, though the procedure is largely handled by `Species`.<sup>2</sup> We'll get to creation and modification later.

Species have an odd relationship to Individuals and to Subpopulations. First recall the Flyweight pattern in Section 3.1.4. Individuals are related to a common Species using the Flyweight pattern: they use Species to store a lot of common information (how to modify themselves, for example). Ordinarily you'd think that the Subpopulation would be a good place for this storage. However different Subpopulations can share the same Species. This allows you to, for example, have one Species guide an entire evolutionary run that might have twenty Subpopulations in it.<sup>3</sup> The species of Subpopulation 0 may be found here:

```
Species theSpecies = state.population.subpops[0].species;
```

A Species contains three major elements: first, the **prototypical Individual** for Subpopulations which use that Species. Recall that Individuals are Prototypes and new ones are formed by cloning from a prototypical individual held in reserve. This "queen bee" individual, so to speak, is found here:

```
Individual theProto = state.population.subpops[0].species.i_prototype;
```

A Species also contains a **prototypical Fitness** object. In ECJ *fitnesses* are separate from *individuals*.

---

<sup>1</sup>Because these two techniques use the subpopulations in different ways, they cannot be used together (a rare situation in ECJ).

<sup>2</sup>You might be asking: if Species are responsible for making individuals, why are Subpopulations involved at all? A very good question indeed.

<sup>3</sup>Granted, this isn't very common.

Individuals define the candidate solution, and Fitnesses define how well it has performed. Like Individuals, Fitnesses are also Prototypes. The prototypical Fitness for Subpopulation 0 may be found here:

```
Fitness theProtoFitness = state.population.subpops[0].species.f_prototype;
```

The Species class you pick is usually determined by the kind of Individual you pick, that is, by the kind of representation of your solution. You define the class of the Species for Subpopulation 0, and its prototypical Fitness and prototypical Individual, as follows. For example, let's make Individuals which are arrays of integers, and a simple Fitness common to many evolutionary algorithms:

```
pop.subpop.0.species = ec.vector.IntegerVectorSpecies
pop.subpop.0.species.ind = ec.vector.IntegerVectorIndividual
pop.subpop.0.species.fitness = ec.simple.SimpleFitness
```

By way of explanation, IntegerVectorIndividual, along with various other “integer” vector individuals like LongVectorIndividual, ShortVectorIndividual, and ByteVectorIndividual, requires an IntegerVectorSpecies. And ec.simple.SimpleFitness is widely used for problems such as Genetic Algorithms or Evolution Strategies. The prototypical Individual is never assigned a Fitness (it's null). But once assembled in a Subpopulation, each Individual has its very own Fitness. To get the Fitness of individual 15 in Subpopulation 0, you'd say:

```
Fitness theFitness = state.population.subpops[0].individuals[15].fitness;
```

Last, a Species contains a **prototypical Breeding Pipeline** to modify individuals. We'll get to that in Section 3.5.

Since they're Prototypes, Individuals, Fitnesses, and Species all have default bases. We'll talk about the different kinds of Individuals, Fitnesses, and Species later, plus various default bases for them.

### 3.2.1 Making Large Numbers of Subpopulations

Let's say you're doing an evolutionary experiment (perhaps coevolution, see Section 7.1) which involves 100 Subpopulations. It's going to get very tiresome to repeat...

```
pop = ec.Population
pop.subpops = 100
pop.subpop.0 = ec.Subpopulation
pop.subpop.0.size = 100
pop.subpop.0.species = ec.vector.IntegerVectorSpecies
pop.subpop.0.species.ind = ec.vector.IntegerVectorIndividual
pop.subpop.0.species.fitness = ec.simple.SimpleFitness
....
pop.subpop.1 = ec.Subpopulation
pop.subpop.1.size = 100
pop.subpop.1.species = ec.vector.IntegerVectorSpecies
pop.subpop.1.species.ind = ec.vector.IntegerVectorIndividual
pop.subpop.1.species.fitness = ec.simple.SimpleFitness
....
pop.subpop.2 = ec.Subpopulation
pop.subpop.2.size = 100
pop.subpop.2.species = ec.vector.IntegerVectorSpecies
pop.subpop.2.species.ind = ec.vector.IntegerVectorIndividual
pop.subpop.2.species.fitness = ec.simple.SimpleFitness
....
```

... and so on some 100 times. Even with the help of ECJ's default parameters, you'll still be typing an awful

lot. Population has a simple mechanism to make this easier on you: the parameter...

```
pop.default-subpop = 0
```

This says that if you *do not* specify a Subpopulation in parameters, ECJ will assume its parameters are identical for those of Subpopulation 0. Thus you could simply say:

```
pop = ec.Population
pop.subpops = 100
pop.default-subpop = 0
pop.subpop.0 = ec.Subpopulation
pop.subpop.0.size = 100
pop.subpop.0.species = ec.vector.IntegerVectorSpecies
pop.subpop.0.species.ind = ec.vector.IntegerVectorIndividual
pop.subpop.0.species.fitness = ec.simple.SimpleFitness
...
```

... and be done with it. Note that you can always specify a Subpopulation specially. For example, suppose all of your Subpopulations were exactly like Subpopulation 0 *except* for Subpopulation 19. You can say:

```
pop = ec.Population
pop.subpops = 100
pop.default-subpop = 0
pop.subpop.0 = ec.Subpopulation
pop.subpop.0.size = 100
pop.subpop.0.species = ec.vector.IntegerVectorSpecies
pop.subpop.0.species.ind = ec.vector.IntegerVectorIndividual
pop.subpop.0.species.fitness = ec.simple.SimpleFitness
... pop.subpop.19 = ec.Subpopulation
pop.subpop.19.size = 25
pop.subpop.19.species = ec.vector.FloatVectorSpecies
pop.subpop.19.species.ind = ec.vector.DoubleVectorIndividual
pop.subpop.19.species.fitness = ec.simple.SimpleFitness
...
```

Note that even though Subpopulation 19 shared the same fitness type as the others, we still had to specify it. It's an all-or-nothing proposition: either you say *nothing* about that particular Subpopulation, or you say *everything*.

### 3.2.2 How Species Make Individuals

Species have two ways to create new individuals: from scratch, or reading from a stream. To generate an individual from scratch, you can call (in `ec.Species`):

#### **ec.Species Methods**

---

```
public Individual newIndividual(EvolutionState state, int thread)
    Returns a brand new, randomized Individual.
```

---

The default implementation of this method simply clones an Individual from the prototype and returns it. Subclasses of Species override this to randomize the Individual in a fashion appropriate to its representation.

Another way to create an individual is to read it from a binary or text stream. `ec.Species` provides two methods for this:

#### **ec.Species Methods**

---

```
public Individual newIndividual(EvolutionState state, LineNumberReader reader) throws IOException
    Produces a new individual read from the stream.

public Individual newIndividual(EvolutionState state, DataInput input) throws IOException
    Produces a new individual read from the given DataInput.
```

---

These methods create Individuals by cloning the prototype, then calling the equivalent `readIndividual(...)` method in `ec.Individual`. See Section 3.2.4 for more information on those methods.

### 3.2.3 Reading and Writing Populations and Subpopulations

Populations and Subpopulations have certain predefined methods for reading and writing, which you should know how to use. If you subclass `Population` or `Subpopulation` (relatively rare) you may need to reimplement these methods. `Population`'s methods are:

```
public void printPopulationForHumans(EvolutionState state, int log);
public void printPopulation(EvolutionState state, int log);
public void printPopulation(EvolutionState state, PrintWriter writer);
public void readPopulation(EvolutionState state, LineNumberReader reader)
    throws IOException;
public void writePopulation(EvolutionState state, DataOutput output)
    throws IOException;
public void readPopulation(EvolutionState state, DataInput input)
    throws IOException;
```

`Subpopulation`'s methods are nearly identical:

#### In Subpopulation:

```
public void printSubopulationForHumans(EvolutionState state, int log);
public void printSubopulation(EvolutionState state, int log);
public void printSubopulation(EvolutionState state, PrintWriter writer);
public void readSubopulation(EvolutionState state, LineNumberReader reader)
    throws IOException;
public void writeSubopulation(EvolutionState state, DataOutput output)
    throws IOException;
public void readSubopulation(EvolutionState state, DataInput input)
    throws IOException;
```

These methods employ similar methods in `ec.Individual` to print out, or read, Individuals. Those methods are discussed next in Section 3.2.4.

The first `Population` method, `printPopulationForHumans(...)`, prints an entire population to a log in a form pleasing to the human eye. It begins by printing out the number of subpopulations, then prints each `Subpopulation` index and calls `printSubpopulationForHumans(...)` on each `Subpopulation` in turn. `printPopulationForHumans(...)` then prints out the number of individuals, then for each `Individual` it prints the `Individual` index, then calls `printIndividualForHumans` to print the `Individual`. Overall, it looks along these lines:

```

Number of Subpopulations: 1
Subpopulation Number: 0
Number of Individuals: 1000
Individual Number: 0
Evaluated: T
Fitness: 0.234
-4.97551104730313 -1.7220830524609632 1.7908415218297096
2.3277606156190496 3.5616099573877404 -3.8002895023118617
Individual Number: 1
Evaluated: T
Fitness: 4.91235
3.1033182498148575 -3.613847679151146 -0.562978505270439
-2.860926011046968 1.9007479097991151 -3.051348823625001
...

```

The next two Population methods, both named `printPopulation(...)`, print an entire population to a log in a form that can be (barely) read by humans but can also be read back in perfectly by ECJ, resulting in identical Populations. These operate similarly to `printPopulationForHumans(...)`, except that various data types are emitted using `ec.util.Code` (Section 2.2.2).

```

Number of Subpopulations: i1|
Subpopulation Number: i0|
Number of Individuals: i1000|
Individual Number: i0|
Evaluated: F
Fitness: f0|0.0|
i6|d4600627607395240880|0.3861348728170766|d4616510324226321041|4.284844300646584|
d4614576621171274054|3.2836854885228233|d4616394543356495435|4.182010230653371|
Individual Number: i1|
Evaluated: F
Fitness: f0|0.0|
i6|d4603775819114015296|0.6217914592919627|d4612464338011645914|2.345643329183969|
d-4606767824441912859|-4.368233761797886|d4616007477858046134|3.919113503960115|
...

```

The Population method `readPopulation(..., LineNumberReader)` can read in this mess to produce a Population. It in turn does its magic by calling the equivalent method in Subpopulation.

The last two methods, `writePopulation(...)` and `readPopulation(..., DataInput)`, read and write Populations (or Subpopulations) to binary files.

### 3.2.4 About Individuals

Individuals have four basic parts:

- The Individual's fitness.  

```
public Fitness fitness;
```
- The Individual's species.  

```
public Species species;
```



- Whether the individual has been evaluated and had its Fitness set to a legal value yet.<sup>4</sup>

```
public boolean evaluated;
```

- The representation of the Individual. This could be anything from an array to a tree structure—representations of course vary and are defined by subclasses. We’ll talk about them later.

### 3.2.4.1 Implementing an Individual

For many purposes you can just use one of the standard “off-the-rack” individuals — vector individuals, genetic programming tree individuals, ruleset individuals — but if you need to implement one yourself, here are some methods you need to be aware of. First off, Individuals are Prototypes and must override the `clone()` method to deep-clone themselves, including deep-cloning their representation and their Fitness, but not their Species (which is just pointer-copied). Individuals must also implement the `setup(...)`, and `defaultBase()` methods. Additionally, Individuals have a number of methods which either should or must be overridden. Let’s start with the “must override” ones:

```
public abstract int hashCode();
public abstract boolean equals(Object individual);
```

These two standard Java methods enable hashing by value, which allows Subpopulations to remove duplicate Individuals. `hashCode()` must return a hashcode for an individual based on value of its representation. `equals(...)` must return true if the Individual is identical to the other object (which in ECJ will always be another Individual).

The next two methods are optional and may not be appropriate depending on your representation:

```
public long size();
public double distanceTo(Individual other);
```

`size()` returns an estimate of the size of the individual. The only hard-and-fast rule is that 0 is the smallest possible size (and the default returned by the method). Size information is largely used by the `ec.parsimony` package (Section 5.2.12) to apply one of several parsimony pressure techniques.

`distanceTo(...)` returns an estimate of the distance, in some metric space, of the Individual to some other Individual of the same type. In the future this method may be used for various crowding or niching methods. At present no package uses it, though all vector individuals implement it. The default implementation returns 0 if the other Individual is identical, else `Double.POSITIVE_INFINITY`.

Last come a host of functions whose purpose is to read and write individuals. You’ve seen this pattern before in Section 3.2.3. Some of these are important to implement; others can wait if you’re in a hurry to get your custom Individual up and running.

```
public void printIndividualForHumans(EvolutionState state, int log);
public void printIndividual(EvolutionState state, int log);
public void printIndividual(EvolutionState state, PrintWriter writer);
public void readIndividual(EvolutionState state, LineNumberReader reader)
                                                                    throws IOException;
public void writeIndividual(EvolutionState state, DataOutput output)
                                                                    throws IOException;
public void readIndividual(EvolutionState state, DataInput input)
                                                                    throws IOException;
```

These six methods only need to be overridden in certain situations, and in each case there’s another method which is typically overridden instead. Here’s what they do:

---

<sup>4</sup>Why isn’t this in the Fitness object? Another excellent question.

- `printlnIndividualForHumans(...)` prints an individual, whether it's been evaluated, and its fitness, out a log in a way that's pleasing and useful for real people to read. Rather than override this method, you probably should instead override this method:

```
public String genotypeToStringForHumans();
```

... which should return the representation of the individual in a human-pleasing fashion. Or, since `genotypeToStringForHumans()` by default just calls `toString()`, you can just override:

```
public String toString();
```

Overriding one or both of these methods is pretty important: otherwise `Statistics` objects will largely be printing your individuals as gibberish. Here's a typical output of these methods:

```
Evaluated: T
Fitness: 0.234
-4.97551104730313 -1.7220830524609632 1.7908415218297096
2.3277606156190496 3.5616099573877404 -3.8002895023118617
```

- Both `printlnIndividual(...)` methods print an individual, and its fitness, out in a way that can be perfectly read back in again with `readIndividual(...)`, but which can also be parsed by humans with some effort. Rather than override this method, you probably should instead override this method:

```
public String genotypeToString();
```

This method is important to implement only if you intend to write individuals out to files in such a way that you can load them back in later. If you don't implement it, `toString()` will be used, which probably won't be as helpful. This returns a `String` which can be parsed in again in the next method. Note that you need to write an individual out so that it can *perfectly* be read back in again as an identical individual. How do you do this? ECJ's classes by default all use the aging and idiosyncratic package `ec.util.Code` package developed long ago for this purpose, but which still works well. See Section 2.2.2. Here's a typical output of these methods (note the use of `ec.util.Code`):

```
Evaluated: F
Fitness: f0|0.0|
i6|d4600627607395240880|0.3861348728170766|d4616510324226321041|4.284844300646584|
d4614576621171274054|3.2836854885228233|d4616394543356495435|4.182010230653371|
```

- `readIndividual(..., LineNumberReader)` reads an individual, and its fitness, in from a `LineNumberReader`. The stream of text being read is assumed to have been generated by `printlnIndividual(...)`. Rather than override this method, you probably should instead override this method:

```
protected void parseGenotype(EvolutionState state, LineNumberReader reader)
    throws IOException;
```

This returns a `String` which can be parsed in again in the next method. Note that you need to write an individual out so that it can *perfectly* be read back in again as an identical individual. How do you do this? ECJ's classes by default all use the aging and idiosyncratic package `ec.util.Code` package developed long ago for this purpose, but which still works well. See Section 2.2.2. You only need to override this method if you plan on reading individuals in from files (by default the method just throws an error).

- The last two methods (`writeIndividual(...)` and `readIndividual(..., DataInput)`) read and write an individual, including its representation, fitness and evaluated flag, in a purely binary fashion to a stream. Don't write the Species. It's probably best instead to override the following methods to just read and write the genotype:

```
public void writeGenotype(EvolutionState state, DataOutput output)
    throws IOException;

public void readGenotype(EvolutionState state, DataInput input)
    throws IOException;
```

These methods are probably only important if you plan on using ECJ's distributed facilities (distributed evaluator, island models). The default implementations of these methods throw exceptions.

### 3.2.5 About Fitnesses

Fitnesses are separate from Individuals, and various Fitnesses can be used depending on the demands of the evolutionary algorithm. The most common Fitness is `ec.simple.SimpleFitness`, which represents fitness as a single number from negative infinity to positive infinity, where larger values are "fitter". Certain selection methods (notably fitness proportionate selection) require that the fitness be non-negative; and ideally between 0 and 1 inclusive.

There are other Fitness objects. For example, there are various multiobjective fitnesses (see Section 7.4), in which the fitness value is not one but some  $N$  numbers, and either higher or lower may be better depending on the algorithm. Other Fitnesses, like the one used in genetic programming (Section 5.2), maintain a primary Fitness statistic and certain auxiliary ones.

You probably won't need to implement a Fitness object. But you may need to use some of the methods below. Fitnesses are Prototypes and so must implement the `clone()` (as a deep-clone), `setup(...)`, and `defaultBase()` methods. Fitness has four additional required methods:

```
public abstract float fitness();
public abstract boolean isIdealFitness();
public abstract boolean equivalentTo(Fitness other);
public abstract boolean betterThan(Fitness other);
```

The first method, `fitness()`, should return the fitness cast into a value from negative infinity to positive infinity, where higher values are better. This is used largely for fitness-proportionate and similar selection methods. If there is no appropriate mechanism for this, you'll need to fake it. For example, multiobjective fitnesses might return the maximum or sum over their various objectives.

The second method, `isIdealFitness()`, returns true if the fitness in question is the best possible. This is largely used to determine if it's okay to quit. It's fine for this method to always return false if you so desire.

The third and fourth methods compare against another fitness object, of the same type. The first returns true if the two Fitnesses are in the same equivalence class: that is, neither is fitter than the other. For simple fitnesses, this is just equality. For multiobjective fitnesses this is Pareto-nondomination of one another. The second method returns true if the Fitness is superior to the one provided in the method. For simple fitnesses, this just means fitter. For multiobjective fitnesses this implies Pareto domination.

Fitnesses also have similar printing facilities to Individuals:<sup>5</sup>

---

<sup>5</sup>Starting to get redundant? Sorry about that.

```

public void printFitnessForHumans(EvolutionState state, int log);
public void printFitness(EvolutionState state, int log);
public void printFitness(EvolutionState state, PrintWriter writer);
public void readFitness(EvolutionState state, LineNumberReader reader)
                                                                    throws IOException;

public void writeFitness(EvolutionState state, DataOutput output)
                                                                    throws IOException;

public void readFitness(EvolutionState state, DataInput input)
                                                                    throws IOException;

```

As usual: the first method, `printFitnessForHumans(...)`, prints a Fitness in a way pleasing for humans to read. It simply prints out the result of the following method (which you should override instead if you ever need to):

```
public String fitnessToStringForHumans();
```

The default implementation of `fitnessToString()` simply calls:

```
public String toString();
```

The next two methods, both named `printFitness(...)`, prints a Fitness in a way that can be (barely) read by humans, and can be read by ECJ to produce an identical Fitness to the original. These methods just print out the result of the following method (which you should override instead if you ever need to):

```
public String fitnessToString();
```

The default implementation of this method calls `toString()`, which is almost certainly wrong. But all the standard Fitness subclasses implement it appropriately using the `ec.util.Code` tools (Section 2.2.2).

The method `readFitness(..., LineNumberReader)` reads into ECJ a Fitness written by these last two printers. Finally, the last two methods, `writeFitness(...)` and `readFitness(..., DataInput)`, read and write the Fitness in a binary fashion. The default implementation of these methods throws an error, but all standard subclasses of Fitness implement them properly.

Fitnesses have two auxiliary variables:

```

public ArrayList trials = null;
public Individual[] context = null;

```

These variables are used by coevolutionary processes (see Section 7.1) to keep track of the number of trials (in the form of `java.lang.Double` used to compute the Fitness value, and to maintain the context (other collaborating Individuals) which produced the best result represented by the Fitness. Outside of coevolution they're presently unused: leave them null and they won't be printed.

### 3.3 Initializers and Finishers

The **Initializer** is called at the beginning of an evolutionary run to create the initial population. The **Finisher** is called at the end of a run to clean up. In fact, it's very rare to use any Finisher other than `ec.simple.Finisher`, which does nothing at all. So nearly always you'll have this:

```
finish = ec.simple.SimpleFinisher
```

Initializers vary largely based on representation, but not for the reason you think. Initializers generally don't need to know anything about the representation of an individual in order to construct it. Instead,

certain representations require a lot of pieces which need to be in a central repository (they're Cliques). For example, the genetic programming facility (Section 5.2) has various *types*, *function sets*, *tree constraints*, *node constraints*, etc. It's not in ECJ's style to store these things as static variables because of the difficulty it presents for serialization. Instead ECJ needed a global object to hold them, and Initializers were chosen for that task. It's probably not been the smartest of decisions: Finishers (which have historically had little purpose) could have been recruited to the job, or some generic type repository perhaps. As it stands, Initializers aren't an optimal location, but there it is.<sup>6</sup>

Unless you're doing genetic programming (`ec.gp`) or using the `ec.rule` package, you'll probably use a `ec.simple.SimpleInitializer`:

```
init = ec.simple.SimpleInitializer
```

ECJ's generational<sup>7</sup> initialization procedure goes like this:

1. The `EvolutionState` asks the `Initializer` to build a `Population` by calling:

```
population = state.initializer.initialPopulation(state, 0);
```

The 0 is thread index 0: this portion of the code is single-threaded.

2. The `Initializer` then creates and sets up a `Population` by calling the following on itself. It then tells the `Population` to populate itself with individuals:

```
Population pop = setupPopulation(state, 0);  
pop.populate(state, 0);
```

Why break this out? Because there are a few `EvolutionState` subclasses which don't want to populate the population immediately or at all — they just want to set it up. For example, steady state evolution sets up a `Population` but may only gradually fill it with initial population members. In this case, the steady state system will just call `setupPopulation(...)` directly, bypassing `initialPopulation(...)`.

3. The `Population`'s `populate(...)` method is straightforward: it calls `populate(...)` in turn on each `Subpopulation` in the `Population`'s subpopulation array.
4. A `Subpopulation`'s `populate(...)` method usually works like this. First, it determines if it should create new individuals from scratch or if it should fill its array by reading `Individuals` from a file. This is determined by (as usual) a parameter! If `Subpopulation 0` should be read in from the file `/tmp/population.in`, the parameter setting would be:

```
pop.subpop.0.file = /tmp/population.in
```

`Subpopulation` will try to read individuals from this file using their `readIndividual(..., LineNumberReader)` method. However if the number of individuals in the file doesn't match the size of the `Subpopulation`, the `Subpopulation` will be resized to match the file, deleting the original `Individuals`. `Subpopulation` will then read individuals using the `newIndividual(..., LineNumberReader)` method in `ec.Species` (slightly less efficient). (See Section 3.2.2).

If no file is provided, the `Subpopulation` assumes that it is generating individuals from scratch. This is by far the most common case. To do this, `Subpopulation` generates new individuals using the standard `newIndividual(...)` method in `ec.Species` (see Section 3.2.2). ECJ can also check to make sure that the `Subpopulation` does not produce duplicate individuals, if you set the following parameter (in this case, in `Subpopulation 0`):

---

<sup>6</sup>This makes it problematic to have both a "rule" representation and a genetic programming representation in the same run without a little hacking, since both require their own `Initializer`. Perhaps this might be remedied in the future.

<sup>7</sup>ECJ's Steady State evolution mechanism has a different initialization procedure. See Section 4.2 for more information.

```
pop.subpop.0.duplicate-retries = 100
```

The default value is no retries.

This says that if the Subpopulation creates a duplicate individual, it will try up to 100 times to replace it with a new, original individual. After that it will give up and use the duplicate individual.

### 3.4 Evaluators and Problems

ECJ evaluates (assesses the fitness of) Individuals in a Population by passing it to an `ec.Evaluator`. Various evolutionary algorithms and other stochastic search algorithms have their own special kinds of Evaluators. Evaluators perform this fitness assessment by cloning one or more **Problems**, discussed in the next Section, and asking these Problems to evaluate the individuals on their behalf. Evaluators hold the prototypical Problem here:

```
public Problem p_problem;
```

This problem is loaded from parameters. For example, to specify that we will use the Artificial Ant Problem to test our genetic programming Individuals, we'd say:

```
eval.problem = ec.app.ant.Ant
```

The basic Evaluator is `ec.simple.SimpleEvaluator`. This class evaluates a Population first by determining how many threads to use. To use four threads (for example), we say:

```
evalthreads = 4
```

The default value is a single thread.

Recall from Section 2.4 that this will require at least four random number generator seeds, for example:

```
seed.0 = 1234  
seed.1 = -503812  
seed.2 = 992341  
seed.3 = -16723
```

When evaluating a Population, `ec.simple.SimpleEvaluator` will construct  $N$  Problems cloned from the Problem prototype. Then, for each Subpopulation, the Evaluator will break the Subpopulation into  $N$  chunks (one per thread), and assign each chunk to a different Problem instance. This enables the Population to be evaluated in parallel. Most commonly, we just set  $N = 1$ .

Certain Evaluator methods are required. The primary method an Evaluator must implement is

```
public abstract void evaluatePopulation(EvolutionState state);
```

This method must take the Population (that is, `state.population`) and evaluate all the individuals in it in the fashion expected by the stochastic search algorithm being employed. Additionally, an Evaluator must implement the method

```
public abstract boolean runComplete(EvolutionState state);
```

... which returns true if the Evaluator believes the process has reached a terminating state. Typically this is done by scanning through the Population and determining if any of the Individuals have ideal fitnesses. If you don't want to be bothered, it's fine to have this method always return false.

### 3.4.1 Problems

Evaluators assess the fitness of individuals typically by creating one or more Problems and handing them chunks of Subpopulations to evaluate. There are two ways that an Evaluator can ask a Problem to perform evaluation:

- For each Individual, the Evaluator can call the Problem’s *evaluation* method. This method varies depending on the kind of Problem. Problems which adhere to `ec.simple.SimpleProblemForm` — by far the most common situation — use the following method:

```
public void evaluate(EvolutionState state, Individual ind,
                    int subpopulation, int threadnum);
```

When this approach is taken, the Problem must assign a fitness immediately during the `evaluate(...)` method. In practice, ECJ doesn’t do this all that much.

- The more common approach allows a Problem to perform fitness evaluation in bulk. In this approach, the Evaluator will first call the following method once:

```
public void prepareToEvaluate(EvolutionState state, int thread);
```

This signals to the Problem that it must prepare itself to begin evaluating a series of Individuals, and then afterwards assign fitness to all of them. Next the Evaluator calls the Problem’s *evaluation* method for each Individual, typically using the method `evaluate(...)` as before. Finally, the Evaluator calls this method:

```
public void finishEvaluating(EvolutionState state, int thread);
```

Using this approach, the Problem is permitted to delay assigning fitness to Individuals until `finishEvaluating(...)` is called.

When ECJ is preparing to exit various Statistics objects sometimes construct a Problem in order to re-evaluate the fittest Individual of the run, solely to have such evaluation print out useful information to tell the user how the Individual operates. This special version of evaluation is done with the following `ec.simple.SimpleProblemForm` method:

```
public void describe(EvolutionState state, Individual ind, int subpopulation,
                    int threadnum, int log);
```

When this method is called, the expectation is that the individual will be evaluated for the purpose of writing out interesting descriptive information to the log. For example, a fit Artificial Ant agent might show the map of the trail it produces as it wanders about eating pellets of food. If you prefer you don’t have to implement this method: and in fact many Problems don’t. The default version (in `ec.Problem`) does nothing at all.

Problem is a Prototype, and so it must implement the `clone()` (as a deep-clone), `setup(...)`, and `defaultBase()` methods: although in truth the default base is rarely used. Problem’s “default” default base is `problem`, which is very rarely used.

### 3.4.2 Implementing a Problem

Commonly the only method a Problem needs to implement is the `evaluate(...)` method. For example, let’s imagine that our Individuals are of the class `ec.vector.IntegerVectorIndividual`, discussed in Section 5.1. The genotype for `IntegerVectorIndividual` is little more than an array of integers. Let us presume that the fitness of these individuals is defined as the product of their integers.

The example below does five basic things:

1. If the individual has already been evaluated, we don't bother evaluating it again. It's possible you'd might want to evaluate it anyway (perhaps if you had a dynamically changing fitness function, for example).
2. We do a sanity check: if the individual is of the wrong type, we issue an error.
3. We compute product of the values in the genome.
4. We set the fitness to that product, and test to see if the fitness is optimal (in this case, if it's equal to `Double.POSITIVE_INFINITY`).
5. We set the individual's evaluated flag.

The implementation is pretty straightforward:

```
package ec.app.myapplication;
import ec.*;
import ec.simple.*;
import ec.vector.*;
public class MyProblem extends Problem implements SimpleProblemForm
{
    public void evaluate(EvolutionState state, Individual ind,
                        int subpopulation, int thread)
    {
        if (ind.evaluated) return;

        if (!(ind instanceof IntegerVectorIndividual))
            state.output.fatal("Whoa! It's not an IntegerVectorIndividual!!!");

        int[] genome = ((IntegerVectorIndividual)ind).genome;
        double product = 1.0;

        for(int x=0; x<genome.length; x++)
            product = product * genome[x];

        ((SimpleFitness)ind.fitness).setFitness(state, product,
            product == Double.POSITIVE_INFINITY);
        ind.evaluated = true;
    }
}
```

## 3.5 Breeders

Individuals are selected bred to create new Individuals using a subclass of `ec.Breeder`. Because this is so central to the differences among various evolutionary algorithms, many such algorithms implement their own Breeder subclasses. A Breeder consists of a single method:

```
public abstract Population breedPopulation(EvolutionState state);
```

This method is required to take the current Population, found here...

```
state.population
```

... and return a Population to be used for the next generation, consisting of individuals selected and bred from the previous Population in a manner appropriate for the algorithm being used. The Population returned



can be the original Population, or it can be an entirely new Population cloned from the original (Population is a Group, recall — see Section 3.1.5).

The most common Breeder is `ec.simple.SimpleBreeder`, which implements a basic form of generational breeding common to the Genetic Algorithm and to Genetic Programming, among others. `SimpleBreeder` has facilities for multithreaded breeding and a simple form of elitism, and works as follows:

1. For each Subpopulation in the Population,
  - (a) Determine the  $N$  fittest Individuals in the Subpopulation.
  - (b) Create a new Subpopulation.
  - (c) Load these  $N$  individuals (the “elites”) into last (highest) slots of the new Subpopulation’s individuals array.
  - (d) Break the remaining unfilled region of this individuals array into  $M$  chunks, one chunk per thread.
  - (e) For each of the  $M$  threads (in parallel),
    - i. Construct a new Breeding Pipeline.
    - ii. Use this Breeding Pipeline to populate the thread’s chunk with newly-bred Individuals.
2. Assemble all the new Subpopulations into a new Population and return it.

The number of elites ( $N$ ) in each Subpopulation is a parameter. To set 10 elites for Subpopulation 0 (for example), you’d say:

```
breed.elite.0 = 10
```

The default value is no elites.

You don’t have to specify the elitism on a per-Subpopulation basis. If you specified a default subpopulation using the parameter `pop.default-subpop` (see Section 3.2.1), `SimpleBreeder` will try this instead when it can’t find an elitism parameter, and will issue a warning letting you know it’s doing so. For example, if you declared:

```
pop.default-subpop = 0 breed.elite.0 = 10
```

... then when `SimpleBreeder` must determine elitism for (say) Subpopulation 4, and that parameter doesn’t exist, it’ll use 10 instead (the value for Subpopulation 0).

Ordinarily elites never have their fitness reevaluated. But if you have a dynamic fitness function, you may wish to reevaluate their fitness each generation to see if it’s still the same. To do this for Subpopulation 0, you say:

```
breed.reevaluate-elites.0 = true
```

The default value is false.

Again, you don’t have to specify the elitism on a per-Subpopulation basis, if you specified a default subpopulation using the parameter `pop.default-subpop`. `SimpleBreeder` will use the default value and issue a warning.

The number of threads ( $M$ ) is also a parameter. To set it to 4, you’d say:

```
breedthreads = 4
```

The default value is a single thread.

As was the case for the `evalthreads` parameter (for `Evaluator`), recall from Section 2.4 that this will require at least four random number generator seeds, one per thread. For example:



Figure 3.3 Two example Breeding Pipelines.

```

seed.0 = 1234
seed.1 = -503812
seed.2 = 992341
seed.3 = -16723

```

Last but not least, certain Breeders allow you to change the subpopulation breeding order to a *sequential* one. Let's say you have 3 subpopulations. If you turn on the following parameter:

```
breed.sequential = true
```

... then SimpleBreeder will only breed subpopulation 0 the first generation, then only subpopulation 1 the second generation, then only subpopulation 2 the second generation, then only subpopulation 0 the third generation, and so on. Obviously if you only have one subpopulation, this parameter has no effect. Why would you want this? ECJ only uses it for a common kind of coevolution, discussed later in Section 7.1.4.1. Otherwise you should keep this set to `false`. The only reason it's mentioned here is so that if you make a Breeder subclass, you can check for this parameter and signal an error if it is `true`, indicating that your Breeder subclass doesn't support it.

All that remains is the breeding procedure itself, for which SimpleBreeder (and many Breeders) constructs a **Breeding Pipeline**.

### 3.5.1 Breeding Pipelines and BreedingSources

A Breeding Pipeline is a chain of selection and breeding operators whose function is to draw from Individuals in an old Subpopulation to produce individuals in a new Subpopulation.

Breeding Pipelines consist of two kinds of objects. First there are **Selection Methods**, which select

Individuals from the old Subpopulation and return them. Then there are **Breeding Pipelines** (what would have better been called *Breeding Operators*), which take Individuals from Selection Methods or from other Breeding Pipelines, modify them in some way, and return them.

The Breeding Pipeline structure isn't actually a *pipeline*: it's really a tree (or in some situations, a directed acyclic graph). The leaf nodes in the graph tree are the Selection Methods (subclasses of `ec.SelectionMethod`), and the nonleaf nodes are the Breeding Pipeline objects (subclasses of `ec.BreedingPipeline`).

Each Breeding Pipeline object can have some  $N$  **sources** (children) from which it draws Individuals. Both `ec.SelectionMethod` and `ec.BreedingPipeline` are subclasses of the abstract superclass `ec.BreedingSource`, and so can function as sources for Breeding Pipelines. Selection Methods do not have sources: rather, they draw Individuals directly from the old Subpopulation.

Breeding Sources (and Breeding Pipeline, and Selection Methods) are Prototypes, and so must implement the `clone()`, `defaultBase()`, and `setup(...)` methods. Breeding Sources also implement three special methods which perform the actual selection and breeding, which we describe here. When a Breeder wishes to produce a series of new Individuals from an old Subpopulation, it begins by calling the method

```
public abstract void prepareToProduce(EvolutionState state, int subpopulation,
                                     int thread);
```

This instructs the BreedingSource to prepare for a number of requests for Individuals drawn from subpopulation number subpopulation. During this method the BreedingSource will, at a minimum, call `prepareToProduce(...)` on each of its sources.

Next, the Breeder calls the following zero or more times to actually produce the Individuals:

```
public abstract int produce(int min, int max, int start, int subpopulation,
                           Individual[] inds, EvolutionState state, int thread);
```

This instructs the BreedingSource to produce between min and max Individuals drawn from subpopulation number subpopulation. The Individuals are to be placed in the inds array at slots `inds[start]`, `inds[start+1]`, ... Finally the method returns the actual number of Individuals produced.

Last, the Breeder calls the following method to give the BreedingSource an opportunity to clean up. The BreedingSource in turn, at a minimum, will call the same method on each of its sources.

```
public abstract void finishProducing(EvolutionState state, int subpopulation,
                                     int thread);
```

Additionally, Breeding Sources implement three other methods. The first:

```
public abstract int typicalIndsProduced();
```

This method returns the number of individuals a BreedingSource would produce by default if not constrained by min and max. The method can return any number  $> 0$ .

The next method:

```
public abstract boolean produces(EvolutionState state, Population newpop,
                                int subpopulation, int thread);
```

... returns true if the BreedingSource believes it can validly produce Individuals of the type described by the given Species, that is, by `newpop.subpops[subpopulation].species`. This is basically a sanity check. At the minimum, the BreedingSource should call this method on each of its sources and return false if any of them return false.

Last, we have the hook...

```
public void preparePipeline(Object hook);
```

You don't have to implement this at all. ECJ does not call this method nor implement it in any of its `BreedingSources` beyond the default implementation (which in `BreedingPipeline` calls the method in turn on each of its sources). This method simply exists in the case that you need a way to communicate with all the methods of a `BreedingPipeline` at some unusual time.

### 3.5.2 SelectionMethods

Selection Methods by default implement the `typicalIndsProduced()` method to return `SelectionMethod.INDS_PRODUCED` (that is, 1).

Furthermore, the default implementation of the `produces` method,

```
public abstract boolean produces(EvolutionState state, Population newpop,
                                int subpopulation, int thread);
```

...just returns true. But you may wish to use this method to check to make sure that your `SelectionMethod` knows how to work with the kind of `Fitnesses` found in the given subpopulation, that is, `state.population[subpopulation].f.prototype`.

The default implementations of `prepareToProduce(...)` and `finishProducing(...)` do nothing at all; though some kinds of `SelectionMethods`, such as `Fitness Proportionate Selection` (`ec.select.FitProportionateSelection`), use `prepareToProduce(...)` to prepare probability distributions based on the Subpopulation in order to select properly.

`SelectionMethods` are sometimes called upon not to produce an `Individual` but to provide an *index* into a subpopulation where the `Individual` is located — perhaps to kill that `Individual` and place another `Individual` in its stead. To this end, `SelectionMethods` have an alternative form of the `produce(...)` method:

```
public abstract int produce(int subpopulation, EvolutionState state,
                             int thread);
```

This method must return the index of the selected individual in the Subpopulation given by

```
state.population.subpops[subpopulation].individuals;
```

#### 3.5.2.1 Implementing a Simple SelectionMethod

Implementing a `SelectionMethod` can be as simple as overriding the “alternative” form of the `produce(...)` method. You don't have to implement the “standard” form of `produce(...)` because its default implementation calls the alternative form and handles the rest of the work for you.

To select an individual at random from a Subpopulation, you could simply implement the “alternative” form to return a random number between 0 and the size of the subpopulation in question:

```
public int produce(int subpopulation, EvolutionState state, int thread)
{
    return state.random[thread.next(
        state.population.subpops[subpopulation].individuals.length)];
}
```

You'll want to always implement the alternative form of `produce(...)`. But in some cases you may wish to also reimplement the “standard” form of `produce(...)` for some special reason.<sup>8</sup> It's a bit more involved but not too hard. We start by determining how many individuals we'll produce, defaulting with 1:

---

<sup>8</sup>Many ECJ selection methods implement the “standard” form in addition to the “alternative” simply to be pedantic — it's very slightly faster — but that's not a good reason nowadays.

```

public int produce(int min, int max, int start, int subpopulation, Individual[] inds,
                  EvolutionState state, int thread)
{
    int n = 1;
    if (n>max) n = max;
    if (n<min) n = min;

```

Next we select  $n$  individuals from the old subpopulation and place them into slots `inds[start] ... inds[start+n-1]`. Here we're do it randomly:

```

    for(int q = 0; q < n; q++)
    {
        Individual[] oldinds = state.population.subpops[subpopulation].individuals;
        inds[start+q] = oldinds[state.random[thread].nextInt(
            state.population.subpops[subpopulation].individuals.length)];
    }
    return n;
}

```

### 3.5.2.2 Standard Classes

There are a number of standard SelectionMethods available in ECJ, all found in the `ec.select` package.

- `ec.select.FirstSelection` always returns the first individual in the Subpopulation. This is largely used for testing purposes.
- `ec.select.RandomSelection` returns an Individual chosen uniformly at random. Basically it does the same thing as the example given above.
- `ec.select.FitProportionateSelection`<sup>9</sup> uses **Fitness-Proportionate Selection**, sometimes called **Roulette Selection**, to pick individuals. Thus `ec.select.BestSelection` requires that all fitnesses be non-negative.
- `ec.select.SUSSelection` selects individuals using **Stochastic Universal Sampling**, a low-variance version of Fitness-Proportionate selection in which highly fit individuals are unlikely to *never* be chosen. Every new generation, and  $M$  selection events thereafter, it shuffles the Subpopulation, then computes the next  $M$  individuals to be selected in the future. ECJ assumes that  $M$  is the size of the Subpopulation. Fitnesses must be non-negative. You have the option of whether or not to shuffle the Subpopulation first:

```
base.shuffle = true
```

The default value is false.

- `ec.select.SigmaScalingSelection` (written by Jack Compton, a former undergraduate at GMU) is another low-variance version of Fitness-Proportionate Selection, in which modified versions of the Individuals' fitnesses are used to reduce the variance among them. This is done by first computing the mean  $\mu$  and standard deviation  $\sigma$  among the fitnesses. If  $\sigma = 0$  no change is made. Otherwise each modified fitness  $g$  is then treated as  $g \leftarrow 1 + \frac{f - \mu}{2\sigma}$ . This can result in negative modified fitnesses, so we introduce a fitness floor: modified fitnesses are bounded to be no less than the floor. Original fitnesses must be non-negative To set this floor to 0.1 (a common value), you'd say:

```
base.scaled-fitness-floor = 0.1
```

0.1 is the default value already, so this is redundant.

SigmaScalingSelection default base is `select.sigma-scaling`.

---

<sup>9</sup>It's called `FitProportionateSelection` rather than `FitnessProportionateSelection` for a historical reason: MacOS 9 didn't allow filenames longer than 32 characters, and `FitnessProportionateSelection.class` is 35 characters long.

- `ec.select.BestSelection` picks among the best or worst  $N$  individuals in the population, depending on a parameter. The selection procedure among these  $N$  is fitness-proportionate selection. Thus `ec.select.BestSelection` requires that all fitnesses be non-negative. To make `BestSelection` pick among the best 5 Individuals, you could say:

```
base.pick-worst = false
base.n = 5
```

The default value for `pick-worst` is false; and 1 for `n`.

`BestSelection`'s default base is `select.best`.

- `ec.select.BoltzmanSelection` (also written by Jack Compton) works like Fitness-Proportionate Selection, but uses modified fitness values according to a Boltzman (Simulated-Annealing-style) cooling schedule. Initially `BoltzmanSelection` has a high *temperature*  $T$ , and for each successive generation it decreases  $T$  by a *cooling rate*  $R$  as  $T \leftarrow T * R$ . Each modified fitness  $g$  is computed as  $g \leftarrow e^{f/T}$ , where  $f$  is the original fitness. Fitnesses must be non-negative. When the temperature reaches 1.0, `BoltzmanSelection` reverts to `FitnessProportionateSelection`. To set the initial temperature to 1000 and the cooling rate to 0.99, you'd say:

```
base.starting-temperature = 1000
base.cooling-rate = 0.99
```

The default temperature is 1.0; and the default cooling rate is 0.0, which causes `BoltzmanSelection` to behave exactly like `FitProportionateSelection`.

`BoltzmanSelection`'s default base is `select.boltzman`.

- `ec.select.GreedyOverselection` is a variation of Fitness-Proportionate Selection which was common in the early genetic programming community (see Section 5.2), but no longer. The Individuals are sorted and divided into the "fitter" and "less fit" groups. With a certain probability the "fitter" individuals will be selected (using Fitness-Proportionate Selection), else the "less fit" individuals will be selected (also using Fitness-Proportionate Selection). Fitnesses must be non-negative. To specify that the "fitter" group is 25% of the Subpopulation, and that individuals are chosen from it 40% of the time, you'd say:

```
base.top = 0.25
base.gets = 0.40
```

`GreedyOverselection`'s default base is `select.greedy`.

- `ec.select.TournamentSelection` first chooses  $N$  individuals entirely at random with replacement (thus the same Individual may be chosen more than once). Then from among those  $N$  it returns the fittest (or least fit, a parameter setting) Individual, breaking ties randomly.  $N$  is often an integer but in fact it doesn't have to be: it can be any real-valued number  $N > 0$ . If  $N$  isn't an integer, it's interpreted as follows: with probability  $N - \lfloor N \rfloor$  we choose  $\lceil N \rceil$  individuals at random, else we choose  $\lfloor N \rfloor$  individuals at random. Fitnesses must be non-negative. The most common setting for  $N$  is 2. To use 2, and return the fittest individual rather than the least-fit one, say:

```
base.size = 2
base.pick-worst = false
```

By default, `pick-worst` is false, so the second parameter is redundant here.

`TournamentSelection`'s default base is `select.tournament`.

- Finally, `ec.select.MultiSelection` is a special version of a `SelectionMethod` with  $N$  other `SelectionMethods` as sources. Each time it must produce an individual, it picks one of these `SelectionMethods` at random

(using certain probabilities) and has it produce the Individual instead. To set up MultiSelection with two sources, TournamentSelection (chosen 60% of the time) and FitnessProportionateSelection (chosen 40% of the time), you'd say:

```
base.num-selects = 2
base.select.0 = ec.select.TournamentSelection
base.select.0.prob = 0.60
base.select.1 = ec.select.FitnessProportionateSelection
base.select.1.prob = 0.40
```

MultiSelection's default base is `select.multiselect`.

### 3.5.3 BreedingPipelines

BreedingPipelines (`ec.BreedingPipeline`) take Individuals from sources, typically modify them in some way, and hand them off. Some BreedingPipelines are mutation or crossover operators; others are more mundane utility pipelines. BreedingPipelines specify the required number of sources they use with the following method:

```
public abstract int numSources();
```

This method must return a value  $> 0$ , or it can return the value `BreedingPipeline.DYNAMIC_SOURCES`, which indicates that the BreedingPipeline can vary its number of sources, and that the user must specify the number of sources with the parameter like this:

```
base.num-sources = 3
```

You specify each source with a parameter. For example, to stipulate sources 0, 1, and 2, you might say:

```
base.source.0 = ec.select.TournamentSelection
base.source.1 = ec.select.TournamentSelection
base.source.2 = ec.select.GreedyOverselection
```

One trick available to you is to state that a source is the same source as a previous one using a special value called `same`. For example, in the example above *two* TournamentSelection operators are created. But if you said the following instead:

```
base.source.0 = ec.select.TournamentSelection
base.source.1 = same
base.source.2 = ec.select.GreedyOverselection
```

...then sources 0 and 1 will be the exact same object. At any rate, the sources are then stored in the following instance variable:

```
public BreedingSource[] sources;
```

Unlike SelectionMethods, BreedingPipelines guarantee a **copy-forward protocol**: any Individual produced by a BreedingPipeline will be unique to that thread. The protocol is simple: if a BreedingPipeline requests an Individual from a source, and that source is a SelectionMethod, the BreedingPipeline will copy the Individual and modify and hand off the copy. But if the source is another BreedingPipeline, the BreedingPipeline will not copy the Individual but instead will just modify it directly and hand it off. What's the point of this? It enables multiple BreedingPipelines, one per thread, to be attached to an old Population and have *all* of them selecting Individuals out of that Population, modifying them, and generating new Individuals without the need for any locking.

Some BreedingPipelines, like crossover pipelines, have a very specific number of children they produce by default (the value returned by `typicalIndsProduced()`). However many others (mutation operators, etc.) simply return whatever Individuals they receive from their sources. For these, BreedingPipeline has a default implementation of `typicalIndsProduced()` which should work fine: it simply calls `typicalIndsProduced()` on all of its sources, and returns the minimum. This computation is done via a simple utility function, `minChildProduction()`, one of two such methods which might be useful to you:

```
public int minChildProduction();
public int maxChildProduction();
```

BreedingPipeline has default implementations of the `produces(...)`, `prepareToProduce(...)`, `finishProducing(...)`, and `preparePipeline(...)` methods, all of which call the same methods on the BreedingPipeline's children.

One final option common to most BreedingPipelines which make modifications (mutation, crossover): you can specify the probability that the pipeline will operate at all, or if Individuals will simply be passed through. For example, let's say you're using a crossover pipeline of some sort, which creates two children from its sources, then crosses them over and returns them. If you state:

```
base.likelihood = 0.8
```

...then with an 0.8 probability crossover will occur as normal. But with a 0.2 probability two Individuals from the sources will be simply copied and returned, with no crossover occurring.

### 3.5.3.1 Implementing a Simple BreedingPipeline

To implement a BreedingPipeline, at a minimum, you'll need to override two methods: `numSources()` and `produce(...)`. `numSources()` is easy. Just return the number of sources your BreedingPipeline requires, or `BreedingPipeline.DYNAMIC.SOURCES` if the number can be any size specified by the user. For example, to make a mutation pipeline, we probably want a single source, which we'll extract Individuals from and mutate:

```
public int numSources() { return 1;}
```

Now we need to implement the `produce(...)` method. Most mutation procedures ask their sources to produce some number of Individuals for them (however many the source prefers), and then mutate those Individuals and return them. We can ask our source to produce sources like this:

```
public int produce(int min, int max, int start, int subpopulation, Individual[] inds,
                  EvolutionState state, int thread)
{
    int n = sources[0].produce(min,max,start,subpopulation,inds,state,thread);
```

The source has taken the liberty of filling slots `inds[start] ... inds[start+n-1]` with  $\min \leq n \leq \max$  Individuals. Next we need to decide whether to bother mutating at all, based on the `likelihood` parameter. We'll use the following utility method to help us:

---

#### ec.BreedingPipeline Methods

```
public int reproduce(int n, int start, int subpopulation, Individual[] inds, EvolutionState state, int thread,
                    boolean produceChildrenFromSource)
    If produceChildrenFromSource is true, extracts n Individuals from Source 0 of the BreedingPipeline, and places them
    in locations inds[start] ... inds[start+n-1], else works with the existing Individuals in those slots. If Source 0 is a
    SelectionMethod, the individuals in inds[start] ... inds[start+n-1] are replaced with clones. Then n is returned.
```

---



We do a coin-flip to determine whether or not to use this method to just clone some individuals from the Source or to go ahead and mutate them. If the former, since our Source has already produced the children, we call this method, passing in *false* for the last argument:

```
// should we bother mutating at all, or just reproduce?
if (!state.random[thread].nextBoolean(likelihood))
    return reproduce(n, start, subpopulation, inds, state, thread, false);
```

At this point we're committed to mutating the  $n$  Individuals. First we need to clone them if Source 0 (our only Source) was a SelectionMethod, since it doesn't clone them for us:

```
// clone the individuals if necessary
if (!(sources[0] instanceof BreedingPipeline))
    for(int q=start;q<n+start;q++)
        inds[q] = (Individual)(inds[q].clone());
```

Now we can mutate the Individuals. How mutation occurs depends on the representation of the Individual of course, so we'll fake it with a comment:

```
for(int q=start;q<n+start;q++)
{
    // modify inds[q] somehow
}
return n;
}
```

### 3.5.3.2 Standard Classes

Most BreedingPipelines are custom for your representation: vectors and trees etc. all have their own special ways of being crossed over or mutated. However there are some utility BreedingPipelines you should be aware of, all stored in the `ec.breed` package:

- `ec.breed.ReproductionPipeline` is by the most common utility BreedingPipeline. In response to a request for  $N$  individuals, `ReproductionPipeline` requests the same number from its single source, then simply returns them (copying if necessary). `ReproductionPipeline` has one rarely-used parameter, which indicates if it *must* copy the individuals even if it's not required to maintain the copy-forward protocol:

```
base.must-clone = true
```

By default, `must-clone` is false.

- Also common is `ec.breed.MultiBreedingPipeline`, which takes some  $M$  sources—determined by the user—and when asked to produce Individuals, chooses randomly among its sources to produce the Individuals for it. It then returns those Individuals. This is a common BreedingPipeline used in genetic programming (Section 5.2). Recall that to stipulate the number of sources, you say:

```
base.num-sources = 2
```

Each source can be accompanied with a probability that this source will be chosen. For example, to state that the first Source is a `ReproductionPipeline`, chosen 10% of the time, and that the second is a `VectorCrossoverPipeline`, chosen 90% of the time, we'd say something like:

```
base.source.0 = ec.vector.breed.VectorCrossoverPipeline
base.source.0.prob = 0.90
base.source.1 = ec.breed.ReproductionPipeline
base.source.1.prob = 0.10
```

You can also state that the number of Individuals returned by any source must be exactly the same — specifically, the maximum that any one of them would return in response to a given request. For example, if you had a Crossover pipeline (which normally returns 2 Individuals) and a Reproduction pipeline (which normally returns 1 Individual), you could force both of them to return 2 Individuals if called on. This is done by saying:

```
base.generate-max = true
```

By default, `generate-max` is true, so this is redundant.

- `ec.breed.BufferedBreedingPipeline` buffers up Individual requests and then hands them out one by one. When you first call `produce()` on a `BufferedBreedingPipeline`, regardless of the number of Individuals requested, it will in turn demand some  $N$  children from its single source. It then stores them in a buffer and hands them to this and later `produce()` requests until they are depleted, at which time it requests  $N$  more, and so on. This value of  $N$  is set like this:

```
base.num-inds = 10
```

Why would you want to do this? Primarily tricks like the following. Let's say you want to create a crossover operator which produces two children, which are then fed into *another* different crossover operator and thus are crossed over *again*. Ordinarily you'd think you could do it along these lines:

```
pop.subpop.0.pipe.0 = ec.app.myCrossover
pop.subpop.0.pipe.0.source.0 = ec.app.myOtherCrossover
pop.subpop.0.pipe.0.source.1 = same
```

Looks good, right? Not so fast. The `myCrossover` class will request exactly one individual from each of its sources. First it'll request from source 0, which will cross over two children, return one, *and throw away the other*. Then it'll request from source 1, which will do exactly the same thing again even though it's the same object. As a result you're not crossing over two individuals twice. You're crossing over individuals which are the result of separate earlier crossovers. But if you did it instead like this:

```
pop.subpop.0.pipe.0 = ec.app.myCrossover
pop.subpop.0.pipe.0.source.0 = ec.select.BufferedBreedingPipeline
pop.subpop.0.pipe.0.source.1 = same
pop.subpop.0.pipe.0.source.0.num-inds = 2
pop.subpop.0.pipe.0.source.0.source.0 = ec.app.myOtherCrossover
pop.subpop.0.pipe.0.source.0.source.1 = same
```

Now `myCrossover` requests one child from `BufferedBreedingPipeline`, which in turn demands *two* children from `myOtherCrossover`, which crosses over two Individuals and returns them. `BufferedBreedingPipeline` returns one of the Individuals. Then `myOtherCrossover` requests the second child, and `BufferedBreedingPipeline` returns the other Individual out of its buffer. Problem solved.

- `ec.breed.ForceBreedingPipeline` takes a single source. In response for a request for  $N$  individuals, `ForceBreedingPipeline` in turn requests up to some  $M$  Individuals from its source. If  $M > N$ , then `ForceBreedingPipeline` requests exactly  $N$  individuals. If  $M < N$ , then `ForceBreedingPipeline` repeatedly requests  $M$  individuals to fill the  $N$ , until possibly the last request, where it requests the remainder. For example, if  $N = 15$  and  $M = 4$ , then `ForceBreedingPipeline` will request 4, then 4, then 4, then 3 Individuals and return them all.

This trick is sort of the counterpart to `BufferedBreedingPipeline`: it gives you a way of demanding a certain number of individuals from a pipeline (like a `CrossoverPipeline`) which doesn't normally return that number. It's only occasionally useful in practice though.

- Last, `ec.breed.GenerationSwitchPipeline` takes two sources. In response to a request for individuals, for generations 1 through  $N - 1$  `GenerationSwitchPipeline` will request Individuals from source 0. For



Figure 3.4 Two example Breeding Pipelines (Repeat of Figure 3.3).

generations  $N$  and on, `GenerationSwitchPipeline` will request `Individuals` from source 1. You specify the switch-generation  $N$  as:

```
base.switch-at = 15
```

Like `ReproductionPipeline`, `GenerationSwitchPipeline` can guarantee that both of its sources always return the same number of individuals (the maximum of the two) with:

```
base.generate-max = true
```

By default, `generate-max` is true, so this example is redundant.

### 3.5.4 Setting up a Pipeline

Setting up a pipeline using parameters sometimes isn't entirely obvious. Let's do the two examples shown in Figure 3.4, in both cases setting up a pipeline for Subpopulation 0.

#### 3.5.4.1 A Genetic Algorithm Pipeline

The left figure is a Breeding Pipeline common to the Genetic Algorithm: two `Individuals` are selected from the old Subpopulation, then copied and crossed over, and the two children are then mutated and added to the new Subpopulation. To build the pipeline, we work our way backwards: first defining the mutator as the top element, then the crossover as its source, then the selector as the crossover's two sources.

First the mutator. We'll use `ec.vector.VectorMutationPipeline`, a common mutator for vector individuals. Subpopulation 0's species holds the prototypical pipeline:

```
pop.subpop.0.species.pipe = ec.vector.VectorMutationPipeline
```

Next we define its sole source: the crossover operator (`ec.vector.VectorCrossoverPipeline`)

```
pop.subpop.0.species.pipe.source.0 = ec.vector.VectorCrossoverPipeline
```

Back to building the Pipeline. Crossover has two sources. We'd like them to both be the Tournament Selector (they could be different Tournament Selectors; it doesn't really matter):

```
pop.subpop.0.species.pipe.source.0.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.0.source.1 = same
```

Tournament Selection has a tournament size operator. Since we're using the same selector for both sources, we only need to set it once:

```
pop.subpop.0.species.pipe.source.0.source.0.size = 2
```

We could also just set the default parameter for all tournament selectors:

```
select.tournament.size = 2
```

Perhaps we'd like the second source to use a tournament size of 4. To do this we'd need to use a separate selector, so we could do this:

```
pop.subpop.0.species.pipe.source.0.source.1 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.0.source.1.size = 4
```

This would also override the default we just set, so it'd work whether or not we used the default-setting approach for source 0.

`VectorCrossoverPipeline` and `VectorMutationPipeline`, discussed later, have various parameters; to simplify the pipeline-building procedure, the `ec.vector` package (Section 5.1) puts these parameters in the `Species`, not the pipeline. For completeness sake, let's include some of them here:

```
pop.subpop.0.species = ec.vector.FloatVectorSpecies
pop.subpop.0.species.crossover-type = one
pop.subpop.0.species.mutation-prob = 1.0
pop.subpop.0.species.mutation-type = gauss
pop.subpop.0.species.mutation-stdev = 0.01
```

### 3.5.4.2 A Genetic Programming Pipeline

The right figure is a typical Genetic Programming pipeline (see Section 5.2). We begin with the root:

```
pop.subpop.0.species.pipe = ec.breed.MultiBreedingPipeline
```

As discussed earlier, `MultiBreedingPipeline` can take any number of sources, so we have to specify it (to 2 here). We also need to state the sources and the probabilities for each source. We'll do 10% Reproduction for the first source and 90% Genetic Programming Crossover for the second. We won't require the two sources to produce the same number of individuals:

```

pop.subpop.0.species.pipe.num-sources = 2
pop.subpop.0.species.pipe.generate-max = false
pop.subpop.0.species.pipe.source.0 = ec.breed.ReproductionPipeline
pop.subpop.0.species.pipe.source.0.prob = 0.10
pop.subpop.0.species.pipe.source.1 = ec.gp.koza.CrossoverPipeline
pop.subpop.0.species.pipe.source.1.prob = 0.90

```

In Genetic Programming normally we'd use Tournament Selection for all selectors. But we'll do various selectors as shown in the Figure. First the Fitness Proportionate Selection source for the ReproductionPipeline:

```

pop.subpop.0.species.pipe.source.0.source.0 = ec.select.FitProportionateSelection

```

Next TournamentSelection (tournament size 7) as Crossover's first source:

```

pop.subpop.0.species.pipe.source.1.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.1.source.0.size = 7

```

Last, Sigma Scaling Selection as Crossover's second source:

```

pop.subpop.0.species.pipe.source.1.source.1 = ec.select.SigmaScalingSelection
pop.subpop.0.species.pipe.source.1.source.1.scaled-fitness-floor = 0.1

```

(The default setting for `pop.subpop.0.species.pipe.source.1.source.1.scaled-fitness-floor` is already 0.1, so it doesn't really need to be set.)

## 3.6 Exchangers

An Exchanger is a subclass of `ec.Exchanger`, and is called both before and after breeding. Exchangers form the basis of **Island Models** in ECJ and will be discussed in-depth in Section 6.2.

Besides `setup(...)`, Exchangers have three basic functions, called at different times in the evolutionary cycle:

```

public abstract Population preBreedingExchangePopulation(EvolutionState state);
public abstract Population postBreedingExchangePopulation(
                                                    EvolutionState state);

public abstract String runComplete(EvolutionState state);

```

The first method is called prior to breeding a population. It's largely available for Island Models to ship off members of the Population to remote ECJ processes. The second method is called immediately *after* breeding a Population, and enables Island Models to import members from remote ECJ processes, possibly displacing newly-bred individuals. The third method is called after `preBreedingExchangePopulation(...)` to determine whether or not the Exchanger thinks ECJ should shut down its process because some *other* process has found the optimal individual. To cause ECJ to shutdown, return a String with a shutdown message of some sort (which will get printed out); otherwise return null.

Unless you're doing Island Models, almost certainly you'll use a default Exchanger called `ec.simple.SimpleExchanger` which does nothing to the Population at all. To wit:

```

exch = ec.simple.SimpleExchanger

```

## 3.7 Statistics

ECJ provides a large number of **statistics hooks**, places where ECJ will call arbitrary methods on a Statistics object throughout the process. Statistics objects are subclasses of `ec.Statistics` and often follow one or more **statistics forms** which provide alternate Statistics hooks. For example, `ec.steadystate.SteadyStateStatisticsForm`, discussed later in Section 4.2, stipulates hooks for Statistics in steady-state evolution. The Statistics class `ec.simple.SimpleStatistics` and has hooks for both regular generational evolution (defined in `ec.Statistics`) and `SteadyStateStatisticsForm`. Another basic Statistics class, `ec.simple.SimpleShortStatistics`, has hooks only for generational evolution. Other Statistics classes exist, such as those found in Genetic Programming (Section 5.2).

You can have as many Statistics objects as you want, but one Statistics object (usually arbitrarily chosen) must be the **statistics root**. To define an `ec.simple.SimpleStatistics` as the root, you say:

```
stat = ec.simple.SimpleStatistics
```

Statistics objects usually, but not always, have a **file** which they log their statistics results out to. It's common to stipulate that file as:

```
stat.file = $out.stat
```

This tells `SimpleStatistics` to write to a file called `out.stat`, located right where the user launched ECJ (for a reminder on the meaning of the "\$", see Section 2.1.2). If you are running with multiple jobs (Section 2.5), ECJ will automatically append the prefix "`jobs.n.`" to this filename, where *n* is the job number. Thus the statistics file for job number 5 will be "`jobs.5.out.stat`". If no file is provided, `SimpleStatistics` will simply print out to the screen.

If ECJ is restarted from a checkpoint, `SimpleStatistics` will append to existing files rather than overwriting them. `SimpleStatistics` also has an option to compress the file using GZIP (and thus add a ".gz" suffix at the very end, as in "`jobs.5.out.stat.gz`". Note that if this option is used, `SimpleStatistics` will simply overwrite the file if restarted from a checkpoint. The parameter is:

```
stat.zip = true
```

For each generation, the `SimpleStatistics` object prints out the best Individual of the generation using the `printIndividualForHumans(...)` method. For example, generation 0 might have:

```
Generation: 0
Best Individual:
Subpopulation 0:
Evaluated: T
Fitness: -1503.8322
2.4502202187677815 0.9879236448665667 0.7631586426217085 0.6854305126240172
```

At the end of the run, the `SimpleStatistics` object prints out the best Individual of the run:

```
Best Individual of Run:
Subpopulation 0:
Evaluated: T
Fitness: -185.78166
-1.0393115193403102 -2.006026366200021 -0.03642166362331428 -1.1196984643947918
```

If your Problem implements the `describe(...)` method, `SimpleStatistics` will also ask the best-of-run Individual to describe itself at the very end of the file.

### 3.7.1 Creating a Statistics Chain

So how do you add additional Statistics objects? As children of the root, or of one another. Any Statistics object can have some  $n$  children. Children are called the same hooks as their parents are. To add another Statistics object (say, `ec.simple.SimpleShortStatistics`), we might add a child to the root:

```
stat.num-children = 1
stat.child.0 = ec.simple.SimpleShortStatistics
stat.child.0.file = $out2.stat
```

Notice that the file has changed: you don't want both Statistics objects writing to the same file! If we wanted to add a third Statistics object (say, another `ec.simple.SimpleShortStatistics`), we could do it this way:

```
stat.num-children = 2
stat.child.0 = ec.simple.SimpleShortStatistics
stat.child.0.file = $out2.stat
stat.child.1 = ec.simple.SimpleShortStatistics
stat.child.1.file = $out3.stat
```

...or we could do it this way:

```
stat.num-children = 1
stat.child.0 = ec.simple.SimpleShortStatistics
stat.child.0.file = $out2.stat
stat.child.0.num-children = 1
stat.child.0.child.0 = ec.simple.SimpleShortStatistics
stat.child.0.child.0.file = $out3.stat
```

The point is, you can hang a Statistics object as a child of any other Statistics object. Pick your poison.

### 3.7.2 Tabular Statistics

`SimpleShortStatistics` writes out a different kind of statistics from `SimpleStatistics`. In its basic form, for each generation it writes out a line of the following values, each separated by a space.

1. The generation number
2. The mean fitness of the entire population for this generation
3. The best fitness of the entire population for this generation
4. The best fitness of the entire population so far in the run

For example, we might have values like this...

```
0 -1851.9916400146485 -1559.68 -1559.68
1 -1801.2400487060547 -1557.7627 -1557.7627
2 -1758.2322434082032 -1513.4955 -1513.4955
3 -1715.5276463623047 -1420.0074 -1420.0074
4 -1675.379030883789 -1459.842 -1420.0074
5 -1637.332774291992 -1426.798 -1420.0074
...
```

You can add sizing information as well. If you set:

```
stat.child.0.do-size = true
```

... you'll get

1. The generation number
2. (If `do-size` is true) The average size of an individual this generation
3. (If `do-size` is true) The average size of an individual so far in the run
4. (If `do-size` is true) The size of the best individual this generation
5. (If `do-size` is true) The size of the best individual so far in the run
6. The mean fitness of the entire population for this generation
7. The best fitness of the entire population for this generation
8. The best fitness of the entire population so far in the run

You can also turn on timing information. If you set:

```
stat.child.0.do-size = true stat.child.0.do-time = true
```

You'll additionally get:

1. The generation number
2. (If `do-time` is true) How long initialization took in milliseconds, or how long the previous generation took to breed to form this generation
3. (If `do-time` is true) How long evaluation took in milliseconds this generation
4. (If `do-size` is true) The average size of an individual this generation
5. (If `do-size` is true) The average size of an individual so far in the run
6. (If `do-size` is true) The size of the best individual this generation
7. (If `do-size` is true) The size of the best individual so far in the run
8. The mean fitness of the entire population for this generation
9. The best fitness of the entire population for this generation
10. The best fitness of the entire population so far in the run

Finally, you can add in per-subpopulation information. If you set:

```
stat.child.0.do-size = true stat.child.0.do-time = true stat.child.0.do-subpops = true
```

you get the whole shebang:

1. The generation number
2. (If `do-time` is true) How long initialization took in milliseconds, or how long the previous generation took to breed to form this generation
3. (If `do-time` is true) How long evaluation took in milliseconds this generation
4. *Once for each subpopulation...*
  - (a) (If `do-size` is true) The average size of an individual this generation for this subpopulation
  - (b) (If `do-size` is true) The average size of an individual so far in the run for this subpopulation



- (c) (If do-size is true) The size of the best individual this generation for this subpopulation
  - (d) (If do-size is true) The size of the best individual so far in the run for this subpopulation
  - (e) The mean fitness of the subpopulation for this generation
  - (f) The best fitness of the subpopulation for this generation
  - (g) The best fitness of the subpopulation so far in the run
5. (If do-size is true) The average size of an individual this generation
  6. (If do-size is true) The average size of an individual so far in the run
  7. (If do-size is true) The size of the best individual this generation
  8. (If do-size is true) The size of the best individual so far in the run
  9. The mean fitness of the entire population for this generation
  10. The best fitness of the entire population for this generation
  11. The best fitness of the entire population so far in the run

**Restricting Rows with a Modulus** If you’ve got a lot of subpopulations, or very long runs, there’s another way you can reduce the size of your file: only output results ever  $n$  generations. For example, if we had set

```
stat.child.0.modulus = 2
```

.. then the aforementioned statistics would look like this:

```
0 -1851.9916400146485 -1559.68 -1559.68
2 -1758.2322434082032 -1513.4955 -1513.4955
4 -1675.379030883789 -1459.842 -1420.0074
...
```

An important thing to keep in mind. Let’s say you want to run for 1000 generations and gather the last. Thus your generations are 0 through 999. If you did a modulus of 100 (say), your reported results would be for generations 0, 100, 200, 300, 400, 500, 600, 700, 800, and 900. That is, you’d have lost the last 99 generations, including the final generation. How might you fix this? The easiest way is to simply run for 1001 generations. Then your last reported generation will be 1000.

### 3.7.3 Implementing a Statistics Object

A basic Statistics object implements one or more of the following hooks:

```
public void preInitializationStatistics(EvolutionState state);
public void postInitializationStatistics(EvolutionState state); // Generational
public void preCheckpointStatistics(EvolutionState state);
public void postCheckpointStatistics(EvolutionState state);
public void preEvaluationStatistics(EvolutionState state); // Generational
public void postEvaluationStatistics(EvolutionState state); // Generational
public void prePreBreedingExchangeStatistics(EvolutionState state);
public void postPreBreedingExchangeStatistics(EvolutionState state);
public void preBreedingStatistics(EvolutionState state); // Generational
public void postBreedingStatistics(EvolutionState state); // Generational
public void prePostBreedingExchangeStatistics(EvolutionState state);
public void postPostBreedingExchangeStatistics(EvolutionState state);
public void finalStatistics(EvolutionState state, int result);
```

When these statistics hooks are called should be self-explanatory from the method name. Note that the methods marked `Generational` are only called by generational `EvolutionState` objects—notably the `ec.simple.SimpleEvolutionState` object. There are also some additional hooks called by the `ec.steadystate` package for steady-state evolution (see Section 4.2).

The `finalStatistics(...)` method, called at the end of an evolutionary run, contains one additional argument, `result`. This argument will be either `ec.EvolutionState.R_SUCCESS` or `ec.EvolutionState.R_FAILURE`. Success simply means that the optimal individual was discovered, and nothing more.

Whenever you override one of these methods, make certain to call `super(...)` first. Let's say that we'd like to know what the size is of the very first individual created after initialization. We might create a `Statistics` subclass which overrides this to print this size out to the screen:

```
public void postInitializationStatistics(EvolutionState state)
{
    super.postInitializationStatistics(state); // always call this
    state.output.println(state.population.subpops[0].individuals[0].size(), 0); // stdout
}
```

We could also write to a file, but to do so we'd need to determine the file name. We could do it in a manner similar to `SimpleStatistics` (ignoring the compression):

```
public static final String P_STATISTICS_FILE = "file";
public int log = 0; // 0 by default means stdout

public void setup(final EvolutionState state, final Parameter base)
{
    super.setup(state, base);
    File statisticsFile = state.parameters.getFile(base.push(P_STATISTICS_FILE), null);
    if (statisticsFile != null) try
    {
        log = state.output.addLog(statisticsFile, true, false, null, false);
    }
    catch (IOException i)
    {
        state.output.fatal("An IOException occurred trying to create the log "
            + statisticsFile + ":\n" + i);
    }
    // else we'll just keep the log at 0, which is stdout
}
```

Now we can write out to the log:

```
public void postInitializationStatistics(EvolutionState state)
{
    super.postInitializationStatistics(state); // always call this
    state.output.println(state.population.subpops[0].individuals[0].size(), log);
}
```

## 3.8 Debugging an Evolutionary Process

A hint. One helpful way to debug ECJ is via BeanShell (<http://www.beanshell.org>). This tool is essentially a command-line for Java. Using it, we can create an ECJ process and get it going quite easily. Let's say that the parameter file is called "foo.params". We can type the following into BeanShell:

```

show();
import ec.*;
args = new String[] { "-file", "foo.params" };
database = Evolve.loadParameterDatabase(args);
state = Evolve.initialize(database, 0);
state.run(EvolutionState.C_STARTED_FRESH);
Evolve.cleanup(state);

```

This is basically a variation of the main code found in Section 2.5. When `state.run(...)` is executed, ECJ will go through the entire evolutionary run. It turns out that `state.run(...)` is really just a cover for three methods (at least, if we're starting from scratch rather than from a checkpoint). They are: first calling `state.startFresh()`, then repeatedly calling `state.evolve()`, and when it returns something other than `state.R_NOTDONE`, finally calling `state.finish(...)`, passing in what was returned. In other words:

```

show();
import ec.*;
args = new String[] { "-file", "foo.params" };
database = Evolve.loadParameterDatabase(args);
state = Evolve.initialize(database, 0);
state.startFresh();
result = EvolutionState.R_NOTDONE;
while( result == EvolutionState.R_NOTDONE )
{
    result = state.evolve();
}
state.finish(result);
Evolve.cleanup(state);

```

So how can we use this for debugging? Well, BeanShell gives us full access to Java. Let's try it on the `ecsuite.params` file. We begin with the preliminaries:

```

show();
args = new String[] { "-file", "ecsuite.params" };

```

BeanShell responds with:

```
<[Ljava.lang.String;@39617189>
```

This is the String array. We continue with:

```
database = Evolve.loadParameterDatabase(args);
```

BeanShell responds with the ParameterDatabase:

```
<{} : ({} : ({eval=ec.simple.SimpleEvaluator, pop.subpop.0=ec.Subpopulation,
quit-on-run-complete=true, generations=1000,
pop.subpop.0.species.pipe.source.0=ec.vector.breed.VectorCrossoverPipeline,
pop.subpop.0.species.min-gene=-5.12, eval.problem=ec.app.ecsuite.ECSuite,
state=ec.simple.SimpleEvolutionState, pop.subpop.0.species.mutation-type=gauss,
pop=ec.Population, pop.subpop.0.duplicate-retries=2, select.tournament.size=2,
pop.subpops=1, pop.subpop.0.species.mutation-stdev=0.01,
pop.subpop.0.species.pipe=ec.vector.breed.VectorMutationPipeline,
pop.subpop.0.species.max-gene=5.12, pop.subpop.0.species.pipe.source.0.source.1=same,
pop.subpop.0.species.pipe.source.0.source.0=ec.select.TournamentSelection,
pop.subpop.0.species=ec.vector.FloatVectorSpecies, breed=ec.simple.SimpleBreeder,
pop.subpop.0.species.mutation-prob=1.0, pop.subpop.0.species.genome-size=100,
pop.subpop.0.species.crossover-type=one, finish=ec.simple.SimpleFinisher,
parent.0=././ec.params, init=ec.simple.SimpleInitializer,
pop.subpop.0.species.ind=ec.vector.DoubleVectorIndividual,
pop.subpop.0.species.fitness=ec.simple.SimpleFitness, pop.subpop.0.size=1000,
eval.problem.type=rastrigin, stat=ec.simple.SimpleStatistics,
exch=ec.simple.SimpleExchanger, stat.file=$out.stat} : ({checkpoint-modulo=1,
evalthreads=1, checkpoint=false, breedthreads=1, prefix=ec, seed.0=time})))>
```

Now we initialize the EvolutionState from the database:

```
state = Evolve.initialize(database, 0);
```

This causes ECJ to start printing to the screen something along these lines:

```
| ECJ
| An evolutionary computation system (version 19)
| By Sean Luke
| Contributors: L. Panait, G. Balan, S. Paus, Z. Skolicki, R. Kicinger, E. Popovici,
|               K. Sullivan, J. Harrison, J. Bassett, R. Hubley, A. Desai, A. Chircop,
|               J. Compton, W. Haddon, S. Donnelly, B. Jamil, and J. O'Beirne
| URL: http://cs.gmu.edu/~eclab/projects/ecj/
| Mail: ecj-help@cs.gmu.edu
|       (better: join ECJ-INTEREST at URL above)
| Date: July 10, 2009
| Current Java: 1.6.0_20 / Java HotSpot(TM) 64-Bit Server VM-16.3-b01-279
| Required Minimum Java: 1.4
```

```
Threads:  breed/1 eval/1
Seed: 1853290822
<ec.simple.SimpleEvolutionState@1f5b0afd>
```

Notice the last line — BeanShell is returning the EvolutionState. Now we fire up the EvolutionState to initialize the first population:

```
state.startFresh();
```

... and get back:

```
Setting up
Initializing Generation 0
```

The first population has been created. Let's look at the first Individual:

```
state.population.subpops[0].individuals[0]
```

This produces:

```
<ec.vector.DoubleVectorIndividual@87740549{699618236}>
```

Not very helpful. Instead, let's have it printed:

```
state.population.subpops[0].individuals[0].printIndividualForHumans(state, 0);
```

This produces something more useful:

```
Evaluated: F
```

```
Fitness: 0.0
```

```
-4.3846934361930945 4.051323475292111 2.750742781209575 -2.1599970035296088
3.5139838195638236 -4.326431483145531 -1.5799722524229094 -4.64489169381555
-4.809825694271426 -0.6969239813124668 -4.322411553562226 4.8723307904232565
2.8978088843319947 -4.311437772193992 -1.556903048013028 2.876699531303326
-1.5461627480422133 -3.406470106152458 0.3510231690045371 -1.26870148662141
-2.9943682283832675 -1.1321325429409796 -4.780798908878881 -2.789054768098288
2.7957975471728034 -2.4529277934521363 0.06864524959557006 -2.807030901927618
-3.817734647565329 3.0018199187738803 3.893346256074625 -4.1700250768556355
-3.3035366716916714 -1.5300889532287534 -1.2924365390313826 2.6878356877535623
4.344108056131552 1.0732802812225044 1.804809997034555 0.6627493849916508
1.6556742582736854 -3.8324177646471913 0.2901815515514814 -0.5045301890375606
-2.755111883054377 -4.057309896490254 -2.097059222061862 -2.062611078568839
3.676980437590175 3.4010063830636517 5.001876654997903 2.3637174851440808
-2.3242430228722846 -0.2027490501614988 4.948796285958214 3.645393286308912
-0.9981883696957627 -2.4911201811073296 2.281601570422807 -3.0028177298996583
-0.6949487749058276 2.4115725052273005 2.2705630820859133 3.8198793397976756
3.927188087275849 3.5439728479577974 4.195897069928313 4.064291914283307
-1.6071055662352376 -0.45138576561254506 0.5382601925283925 2.2824947546503687
-0.0837300863613164 -2.4997930740673895 0.06696037058102089 1.782243737261787
-3.390249634178219 4.669336185081783 2.371290190775591 1.8743739255868377
0.13349732700681827 2.808175830805574 -4.2297879656940705 -0.5781599273148448
3.4174595199606577 2.5509508748123793 0.9574470878471297 1.181916131827328
3.3128918249657184 3.5085201808925843 -0.8921840350705308 -4.016933626993176
2.5591127486976983 1.580181276449899 -0.6102226049991097 1.0644092417475743
0.5897983455130262 2.5504671849586904 -2.230897886457403 1.8133759722806326
```

Now we're getting somewhere. You could print out all the individuals like this:

```
for(int i = 0; i < state.population.subpops[0].individuals.length; i++ )
    state.population.subpops[0].individuals[i].printIndividualForHumans(state, 0);
```

... but I won't torture you with what comes out. Instead, notice that the Individual has not been evaluated yet—just initialized with the Population. We go through one evolutionary loop—evaluating the individuals and breeding a new Population—by calling

```
state.evolve();
```

This produces:

```
Subpop 0 best fitness of generation: Fitness: -1493.534
<2>
```

The first line was printed to the screen. The second line indicates what `state.evolve()` returned. 2 is the value of `EvolutionState.R_NOTDONE`, indicating that the process needs to be evolved more.

Let's look at that first individual again:

```
state.population.subpops[0].individuals[0].printIndividualForHumans(state, 0);
```

...yielding...

Evaluated: F

Fitness: -1806.1432

```
-1.5595820609909528 -2.9941135630034292 -3.188550961391961 0.8673223056511647
2.4308132097811472 -3.6298006589453533 -4.62193495641744 1.7381186900517611
-3.2707539202577953 -2.8517369832386144 -4.701099579700639 1.1683479248633841
0.10118833856168477 2.7982137159130787 -1.3673458253800685 -4.548719487000453
-1.7852252742508177 -1.662999422245311 -4.891889992368657 2.0689413066938824
4.64815452362056 4.03620579726471 -2.6065781548997413 2.8384398494616585
-1.6231723965539844 -0.19641152832494305 -0.8025430015631594 -4.337733534634894
-4.259188069209607 0.0974585410674078 4.878006291864429 4.187577755641656
3.9507153065207605 -3.3456633008586922 3.7163666200189596 -0.7581028665673978
4.28299933455259 1.8522464455693997 4.4324032846812935 -1.3209545115697914
4.239911043319335 -2.7741200087506352 -3.181419981396656 -0.4574562816089688
3.9209870275697982 0.31049605413333237 -0.46868091240064014 -4.570530964131764
-0.9126484738704782 3.6348709305820153 -1.800821491837854 -0.8548399118205554
-1.6874962921883667 2.628667604603462 0.060377157894385663 3.194354857448187
1.2106237734207714 -3.477534436566739 1.919326547065771 -3.74880517912247
4.076653684533312 -2.9153006121227034 -2.4460232838375973 0.6128610868842217
0.7785108819209824 -1.213371979065718 3.2441049504290587 -1.352037820951835
1.151316091162472 0.3915293759690397 -0.15229424767569708 1.8192706794904545
-3.057866603248519 -3.2217378304635926 3.7963181147558447 1.9609441782591566
2.1365399986514815 -0.7608502832241196 -1.2202190662246202 -3.2592371482282956
-2.612971504172355 3.1496849987738167 -5.083084415090031 -4.243405086300351
3.8516939433487387 -4.87008846122508 1.012854792831603 3.77728764346906
2.843506550933032 4.705462097924235 1.4291349248648448 3.8398215224809875
1.1776568359195472 -4.784524531392207 2.765230136436807 -2.6521295800350555
-2.271480494878218 -2.018481022639772 -2.2536397207045686 -1.5048357519436404
```

It's a different Individual: the next generation one to be exact. Notice that although it has a "fitness", in fact it's not been evaluated: the "fitness" is just nonsense cloned from a previous Individual.

So how do you see a population with fitness-evaluated Individuals? Generational ECJ `EvolutionState` processes assess the fitness of a Population, then create a new Population and throw away the old one. You can hold onto the old Population pretty easily. Just do this:

```
p = state.population;
state.evolve();
```

Now `p` holds the old Population, filled with now-evaluated Individuals, and `state.population` holds the next-generation Population, which hasn't been evaluated yet. For example, if you say:

```
p.subpops[0].individuals[0].printIndividualForHumans(state, 0);
```

... you will get back something like this:

Evaluated: T

Fitness: -1779.4391

-1.5595820609909528 -2.9941135630034292 -3.188550961391961 0.8673223056511647  
2.4308132097811472 -3.6298006589453533 -4.62193495641744 1.7381186900517611  
-3.2707539202577953 -2.8517369832386144 -4.701099579700639 1.1683479248633841  
0.10118833856168477 2.7982137159130787 -1.3673458253800685 -4.548719487000453  
-1.7852252742508177 -1.662999422245311 -4.891889992368657 2.0689413066938824  
4.64815452362056 4.03620579726471 -2.6065781548997413 2.8384398494616585  
-1.6231723965539844 -0.19641152832494305 -0.8025430015631594 -4.337733534634894  
-4.259188069209607 0.0974585410674078 4.878006291864429 4.187577755641656  
3.9507153065207605 -3.3456633008586922 3.7163666200189596 -0.7581028665673978  
4.28299933455259 1.8522464455693997 4.4324032846812935 -1.3209545115697914  
4.239911043319335 -2.7741200087506352 -3.181419981396656 -0.4574562816089688  
3.9209870275697982 0.31049605413333237 -0.46868091240064014 -4.570530964131764  
-0.9126484738704782 3.6348709305820153 -1.800821491837854 -0.8548399118205554  
-1.6874962921883667 2.628667604603462 0.060377157894385663 3.194354857448187  
1.2106237734207714 -3.477534436566739 1.919326547065771 -3.74880517912247  
4.076653684533312 -2.9153006121227034 -2.4460232838375973 0.6128610868842217  
0.7785108819209824 -1.213371979065718 3.2441049504290587 -1.352037820951835  
1.151316091162472 0.3915293759690397 -0.15229424767569708 1.8192706794904545  
-3.057866603248519 -3.2217378304635926 3.7963181147558447 1.9609441782591566  
2.1365399986514815 -0.7608502832241196 -1.2202190662246202 -3.2592371482282956  
-2.612971504172355 3.1496849987738167 -5.083084415090031 -4.243405086300351  
3.8516939433487387 -4.87008846122508 1.012854792831603 3.77728764346906  
2.843506550933032 4.705462097924235 1.4291349248648448 3.8398215224809875  
1.1776568359195472 -4.784524531392207 2.765230136436807 -2.6521295800350555  
-2.271480494878218 -2.018481022639772 -2.2536397207045686 -1.5048357519436404

Notice that this Individual has been evaluated. Its Fitness is valid.





## Chapter 4

# Basic Evolutionary Processes

### 4.1 Generational Evolution

ECJ is most commonly used for generational evolution: where a whole Population is evaluated, then updated, at a time. There are a number of packages which use generational evolution, but the two most common are the `ec.simple` package, which does Genetic Algorithm style generational evolution, and the `ec.es` package which does Evolution Strategies.

#### 4.1.1 The Genetic Algorithm (The `ec.simple` Package)

We've pretty much covered everything in the `ec.simple` package throughout Section 3. But just a quick reminder:

- `ec.simple.SimpleEvolutionState` subclasses `ec.EvolutionState` to provide the generational top-level loop shown in Figure 4.1. Each generation the entire Population is handed to the Evaluator, then the Breeder. The class adds no new parameters beyond those defined in `EvolutionState`.
- `ec.simple.SimpleBreeder` subclasses `ec.Breeder` to provide multithreaded breeding and elitism. `SimpleBreeder` was discussed at length in Section 3.5.
- `ec.simple.SimpleEvaluator` subclasses `ec.Evaluator` to provide multithreaded evaluation. `SimpleEvaluator` adds no new parameters beyond those defined in `Evaluator`, and was discussed at length in Section 3.4.
- `ec.simple.SimpleFitness` subclasses `ec.Fitness` to provide a simple fitness consisting of a single floating-point number, where higher fitness values are preferred. `SimpleFitness` also holds a boolean flag indicating whether the fitness assigned is the optimal fitness. `SimpleFitness` adds no new parameters beyond those defined in `Fitness`, and was discussed at length in Section 3.2.5.
- `ec.simple.SimpleProblemForm` defines the kind of methods which must be implemented by Problems used by a `SimpleEvaluator`, and was discussed at length in Section 3.4.1. As a reminder, the two methods defined by `SimpleProblemForm` are `evaluate(...)`, which evaluates an individual and sets its fitness; and `describe(...)`, which evaluates an individual solely for the purpose of writing a detailed description about the individual's performance out to a stream. They look like this:

```
public void evaluate(EvolutionState state, Individual ind,
                    int subpopulation, int threadnum);
public void describe(EvolutionState state, Individual ind,
                    int subpopulation, int threadnum, int log);
```



Figure 4.1 Top-Level Loop of ECJ's SimpleEvolutionState class, used for basic generational EC algorithms. Various sub-operations are shown occurring before or after the primary operations. The full population is revised each iteration. A repeat of Figure 1.1.

- `ec.simple.SimpleInitializer` subclasses `ec.Initializer` to provide multithreaded breeding and elitism. `SimpleInitializer` adds no new parameters beyond those defined in `Initializer`, and was discussed at length in Section 3.3.
- `ec.simple.SimpleFinisher` subclasses `ec.Finisher` and does nothing at all. `SimpleFinisher` was discussed at length (so to speak) in Section 3.3.
- `ec.simple.SimpleExchanger` subclasses `ec.Exchanger` and does nothing at all. `SimpleExchanger` was mentioned in Section 3.6.
- `ec.simple.SimpleStatistics` subclasses `ec.Statistics` and outputs the best-of-generation individual each generation, plus the best-of-run individual at the end. `SimpleStatistics` was discussed at length in Section 3.7.
- `ec.simple.SimpleShortStatistics` subclasses `ec.Statistics` and gives numerical statistics about the progress of the generation. `SimpleStatistics` was also discussed at length in Section 3.7.
- `ec.simple.SimpleDefaults` implements `ec.DefaultsForm` and provides the package default parameter base.

**Example** Let's put these together to do a simple genetic algorithm. We start with the basic parameters:

```
# Threads and Seeds
evalthreads = 1
breedthreads = 1
seed.0 = time

# Checkpointing
checkpoint = false
checkpoint-modulo = 1
prefix = ec
```

Next a basic generational setup:

```
# The basic setup
state = ec.simple.SimpleEvolutionState
init = ec.simple.SimpleInitializer
finish = ec.simple.SimpleFinisher
exch = ec.simple.SimpleExchanger
breed = ec.simple.SimpleBreeder
eval = ec.simple.SimpleEvaluator
stat = ec.simple.SimpleStatistics
pop = ec.Population

# Basic parameters
generations = 200
quit-on-run-complete = true
pop.subpops = 1
pop.subpops.0 = ec.Subpopulation
pop.subpop.0.size = 1000
pop.subpop.0.duplicate-retries = 0
breed.elite.0 = 0
stat.file = $out.stat
```

We'll use Individuals of the form `ec.vector.IntegerVectorIndividual`, discussed later in Section 5.1. This is not much more than a cover for a one-dimensional array of integers:

```
# Representation
pop.subpops.0.species = ec.vector.IntegerVectorSpecies
pop.subpop.0.species.ind = ec.vector.IntegerVectorIndividual pop.subpop.0.species.genome-size
= 100
```

For fitness, we'll use SimpleFitness:

```
# Fitness
pop.subpop.0.species.fitness = ec.simple.SimpleFitness
```

In Section 3.5.4 we laid out a simple Genetic Algorithm Pipeline:

```
# Pipeline
pop.subpop.0.species.pipe = ec.vector.VectorMutationPipeline
pop.subpop.0.species.pipe.source.0 = ec.vector.VectorCrossoverPipeline
pop.subpop.0.species.pipe.source.0.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.0.source.1 = same
select.tournament.size = 2
```

Because they are so common, Vector pipelines are unusual in that they define certain probabilities in Species rather than in the Pipeline, mostly for simplicity. We haven't discussed these yet (we'll get to them in Section 5.1), but here's one possibility:

```
pop.subpop.0.species.crossover-type = one
pop.subpop.0.species.mutation-prob = 0.01
```

In Section 3.4.2 we defined a simple Problem in which the fitness of an IntegerVectorIndividual was the product of the integers in its genome. Let's use it here.

```
package ec.app.myapp;
import ec.*;
import ec.simple.*;
import ec.vector.*;
public class MyProblem extends Problem implements SimpleProblemForm
{
    public void evaluate(EvolutionState state, Individual ind,
                        int subpopulation, int thread)
    {
        if (ind.evaluated) return;

        if (!(ind instanceof IntegerVectorIndividual))
            state.output.fatal("Whoa! It's not an IntegerVectorIndividual!!!");

        int[] genome = ((IntegerVectorIndividual)ind).genome;
        double product = 1.0;

        for(int x=0; x<genome.length; x++)
            product = product * genome[x];

        ((SimpleFitness)ind.fitness).setFitness(state, product,
            product == Double.POSITIVE_INFINITY);
        ind.evaluated = true;
    }
}
```

Now we state that we're using this class:

```
eval.problem = myapp.MyProblem
```

... and we're ready to go! Save out as the file `ga.params`, we then run ECJ as:

```
java ec.Evolve -file ga.params
```

### 4.1.2 Evolution Strategies (The `ec.es` Package)

The `ec.es` package implements basic versions of the  $(\mu, \lambda)$  and  $(\mu + \lambda)$  algorithms. Since it's also generational, it largely depends on and extends the `ec.simple` package. The way these algorithms are done is via special Breeders: the aptly named `ec.es.MuCommaLambdaBreeder` and `ec.es.MuPlusLambdaBreeder`. These work in conjunction with a specially formulated Selection operator called `ec.es.ESSelection`.

Evolution Strategies differs from the Genetic Algorithm mostly in that it does **truncation selection**: an entire segment of the Population (all but the best  $\mu$ ) is simply lopped off, and breeding occurs among the remainder. ECJ does this by having `ESSelection` only select among the best  $\mu$  Individuals in the Population. In truncation selection, each remaining Individual gets the same number of opportunities to breed. `ESSelection` works with the two Breeders to guarantee this: each time the `BreedingPipeline` is pulsed for an individual, the parent has been pre-determined by the Breeder, and `ESSelection` simply returns that parent.

`ESSelection` doesn't have to be the only selection operator in your `BreedingPipeline`: you can have various other ones. But you should have at *least* one `ESSelection` operator in your pipeline, else it just isn't Evolution Strategies any more. Also note that `ESSelection` always returns the Individual pre-determined by the Breeder: thus if you have multiple `ESSelection` operators in your pipeline, they will *all* return that same Individual.

The  $\mu$  variable stipulates the number of parents left after truncation selection has lopped off the population. The  $\lambda$  variable stipulates how many children are generated by the  $\mu$  parents in total. Each Individual in the  $\mu$  gets to produce exactly  $\frac{\lambda}{\mu}$  children. Obviously  $\mu$  must divide evenly into  $\lambda$ . Often  $\lambda$  is also the initial population size: but this doesn't have to be the case. If not, it will be the population size of the second and later generations.

The `MuCommaLambdaBreeder` has two parameters, which, not surprisingly, specify the values of  $\mu$  and  $\lambda$  in a given Subpopulation. For Subpopulation 0, this will provide  $\mu = 1$  and  $\lambda = 10$ , to make a (10, 100) Evolution Strategy.

```
breed = ec.es.MuCommaLambdaBreeder
es.mu.0 = 10
es.lambda.0 = 100
```

This algorithm differs from  $(\mu, \lambda)$  in that, after creating the children, the  $\mu$  parents join the  $\lambda$  children to form the next generation Population. Thus the next generation is  $\mu + \lambda$  in size. Again, the initial population can be any size (traditionally I think it's  $\mu + \lambda$ ). The `MuPlusLambdaBreeder` subclasses `MuCommaLambdaBreeder` and adds no new parameters, though you'd change the Breeder of course:

```
breed = ec.es.MuPlusLambdaBreeder
```

It's common in Evolution Strategies to use a mutation-only pipeline. Here's one:

```
pop.subpop.0.pipe = ec.vector.VectorMutation
pop.subpop.0.pipe.source.0 = ec.es.ESSelection
```

Last but not least, `ec.es.ESDefaults` provides the package default parameter base.

**Example** We build off of the example shown in Section 4.1.1, so let's use that file:

```
parent.0 = ga.params
```

Next, let's override some parameters to use Evolution Strategies:

```
breed = ec.es.MuCommaLambdaBreeder
es.mu.0 = 10
es.lambda.0 = 100
pop.subpop.0.pipe = ec.vector.VectorMutation
pop.subpop.0.pipe.source.0 = ec.es.ESSelection
```

Evolution Strategies also often uses a floating-point array representation. The Genetic Algorithm example in Section 4.1.1 used an integer array representation. We could change it to an array of Doubles like this:

```
pop.subpop.0.species = ec.vector.FloatVectorSpecies
pop.subpop.0.species.ind = ec.vector.DoubleVectorIndividual
```

IntegerVectorIndividual has a simple default mutator: randomizing the integers. This is why the mutation-prob is set low (to 0.01). Since we're using floating-point values, let's change the mutation type to gaussian mutation with a standard deviation of 0.01, happening 100% of the time:

```
pop.subpop.0.species.mutation-type = gauss
pop.subpop.0.species.mutation-stdev = 0.01
pop.subpop.0.species.mutation-prob = 1.0
```

Since we're using a different representation, we need to change our Problem a bit:

```
package ec.app.myapp;
import ec.*;
import ec.simple.*;
import ec.vector.*;
public class MySecondProblem extends Problem implements SimpleProblemForm
{
    public void evaluate(EvolutionState state, Individual ind,
                        int subpopulation, int thread)
    {
        if (ind.evaluated) return;

        if (!(ind instanceof DoubleVectorIndividual))
            state.output.fatal("Whoa! It's not an DoubleVectorIndividual!!!");

        double[] genome = ((DoubleVectorIndividual)ind).genome;
        double product = 1.0;

        for(int x=0; x<genome.length; x++)
            product = product * genome[x];

        ((SimpleFitness)ind.fitness).setFitness(state, product,
            product == Double.POSITIVE_INFINITY);
        ind.evaluated = true;
    }
}
```

Finally, we set the problem:

```
eval.problem = ec.app.myapp.MySecondProblem
```

... and we're ready to go! Save out the file as `es.params`, and then run ECJ as:

```
java ec.Evolve -file es.params
```

## 4.2 Steady-State Evolution (The `ec.steadystate` Package)

In Steady-state Evolution, each iteration some  $n$  individuals are bred from the existing Population, have their fitnesses assessed, and then are stuck back in the Population, displacing  $n$  others. ECJ's implementation of Steady-State Evolution sets  $n = 1$ , which is the most common case.

To implement Steady-State Evolution ECJ requires special versions of: an `EvolutionState`, a `Breeder`, and an `Evaluator`:

```
state = ec.steadystate.SteadyStateEvolutionState
breed = ec.steadystate.SteadyStateBreeder
eval = ec.steadystate.SteadyStateEvaluator
```

Additionally Breeding Sources must adhere to certain rules, as must Exchangers and Statistics.

ECJ's top-level Steady-State Evolution loop, implemented in the class `ec.steadystate.SteadyStateEvolutionState`, is shown in Figure 4.2. Compare to Figure 4.1, which shows a generational loop. At this abstract level, the two are very similar: the primary differences lie in the Statistics methods called. But inside the `Breeder` and `Evaluator` there are some significant differences. One important detail: Steady-State Evolution *must* at present be single-threaded:

```
evalthreads = 1
breedthreads = 1
```

`SteadyStateEvolutionState` can either run until some  $n$  evaluations have been processed, or until some  $m$  "generations" have been processed. Normally `SteadyStateEvolutionState` uses generations. If you'd prefer to work in evaluations, you can say something like this:

```
evaluations = 10000
```

What does a "generation" mean in this context? ECJ sums up the sizes of all of the subpopulations to form the *total population size*  $p$ . Whenever  $p$  individuals have been evaluated, that's a new generation.

Steady-State evolution in ECJ works in two stages:

- **Stage 1: Initialization.** The Subpopulation begins completely empty. One-by-one, individuals are created at random and then sent off to be evaluated. In Steady-State evolution, this evaluation happens immediately and each individual is returned immediately. However in a distributed version called Asynchronous Evolution (discussed in Section 6.1), individuals are shipped off to remote sites and may not come back for a while. In this case evolution does not wait for them, but continues generating new individuals.

When an individual returns with its fitness assessed, it then is added to the Subpopulation until the Subpopulation is full, at which time the system advances to...

- **Stage 2: The Steady State.** One-by-one, individuals are *bred* and then sent off to be evaluated. Again, in Steady-State evolution, this evaluation happens immediately and each individual is returned immediately. However in Asynchronous Evolution individuals are shipped off to remote sites and may not come back for a while. In this case evolution does not wait for them, but continues breeding new individuals.

When an individual returns with its fitness assessed, an existing Subpopulation member is marked for death, and is then replaced with the new individual. Note that if using Asynchronous Evolution, some of these new individuals may be stragglers created during Initialization.

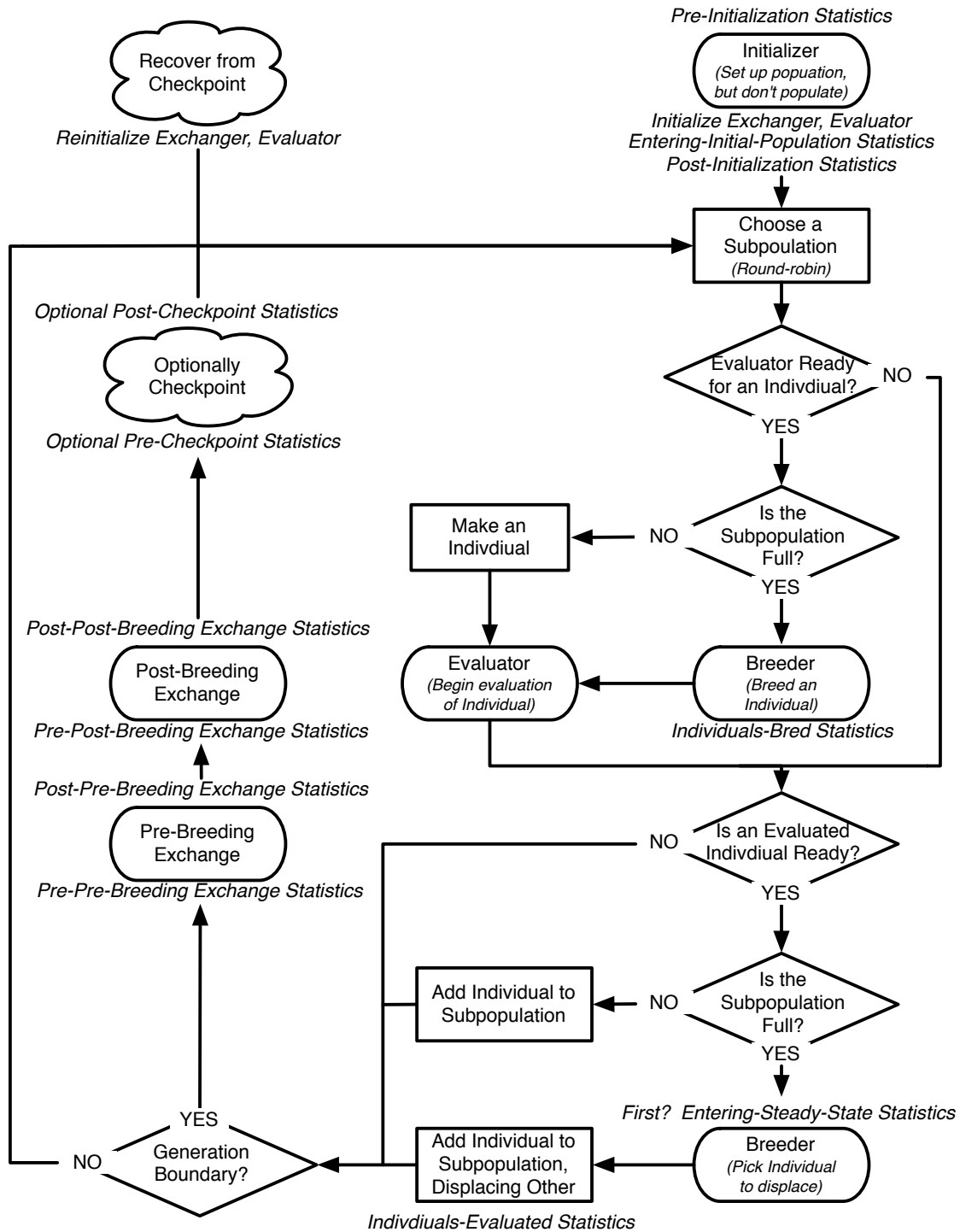


Figure 4.2 Top-Level Loop of ECJ's SteadyStateEvolutionState class, used for simple steady-state EC and Asynchronous Evolution algorithms. "First?" means to perform the Statistics whenever the Subpopulation in question is picking an Individual to displace for the very first time. (Each Subpopulation will do it once, but possibly at different times).



These stages aren't immediately obvious from Figure 4.2: they're distinguished by both "Is the Subpopulation Full?" branches.

The Evaluator in question is `ec.steadystate.SteadyStateEvaluator`, which has no special parameters.

The Breeder in question is `ec.steadystate.SteadyStateBreeder`. It contains the `SelectionMethod` which is used to select individuals for death. Typically this `SelectionMethod` is set up to select *unfit* or *random* individuals rather than *fit* ones (which wouldn't make much sense). To pick an unfit individual using Tournament Selection for Subpopulation 0, for example, we might say:

```
steady.deselector.0 = ec.select.TournamentSelection
steady.deselector.0.size = 2
steady.deselector.0.pick-worst = true
```

To pick a random individual, you could change this to

```
steady.deselector.0.size = 1
```

...or you could just use...

```
steady.deselector.0 = ec.select.RandomSelection
```

All `BreedingSources` in your breeding pipeline, and in particular `SelectionMethods`, must implement `ec.steadystate.SteadyStateBSourceForm`. This interface contains a method called to update the `BreedingSource` that a new individual has entered into the population, displacing an old one. This is important because some `SelectionMethods`, such as `FitProportionateSelection`, rely on precomputed statistics which will be no longer correct. This method is:

```
public void individualReplaced(SteadyStateEvolutionState state,
                              int subpopulation, int thread, int individual);
```

At the very least, a `BreedingSource` must implement this method to call the same method on *its* sources. This means that they must be of `SteadyStateBSourceForm` as well. To guarantee this, the following check method must be implemented:

```
public void sourcesAreProperForm(SteadyStateEvolutionState state);
```

This method should likewise call its namesake on all sources. If the sources are not of `SteadyStateBSourceForm`, it should then issue an ordinary Output "error" (not a fatal error) indicating this problem.

`SteadyStateBSourceForm` is implemented by `TournamentSelection` and its variations; and also by `FirstSelection` (of course), and by `RandomSelection`. You can implement `SteadyStateBSourceForm` for `FitProportionateSelection` and its ilk but it's not implemented by default because they'd be so inefficient. I strongly suggest sticking with `TournamentSelection`.

Exchangers are called every *generation* worth of evaluations. Since Exchangers also modify individuals in the population, they also need to update the `BreedingSources` of this fact. This is done by having them call the following method in the Breeder, which lets all the relevant `BreedingSources` know:

```
public void individualReplaced(SteadyStateEvolutionState state,
                              int subpopulation, int thread, int individual);
```

(Yes, it's the same name). If an Exchanger makes this promise, it gets to implement the `ec.steadystate.SteadyStateExchangerForm` interface as a badge of honor.

The steady state nature of evolution requires a different set of Statistics hooks. These hooks are defined in the interface `ec.steadystate.SteadyStateStatisticsForm`. Note that no ECJ Statistics class presently implements these hooks; but `SteadyStateEvolutionState` can handle non-`SteadyStateStatisticsForm` Statistics objects gracefully.

Several hooks you’ve already seen in Section 3.7:

```
public void preCheckpointStatistics(EvolutionState state);
public void postCheckpointStatistics(EvolutionState state);
public void prePreBreedingExchangeStatistics(EvolutionState state);
public void postPreBreedingExchangeStatistics(EvolutionState state);
public void prePostBreedingExchangeStatistics(EvolutionState state);
public void postPostBreedingExchangeStatistics(EvolutionState state);
public void finalStatistics(EvolutionState state, int result);
```

Other hooks are special to steady-state evolution. First, we have:

```
public void enteringInitialPopulationStatistics(
    SteadyStateEvolutionState state);
public void enteringSteadyStateStatistics(int subpop,
    SteadyStateEvolutionState state);
public void generationBoundaryStatistics(EvolutionState state);
```

The first method is called immediately after the Population has been constructed but before any Individuals have been created. The second method is called when a Subpopulation enters its “steady state”, that is, it has been filled with Individuals. The third method is called whenever a generation boundary is reached, that is, when a Population’s worth of individuals has been evaluated.

Last, we have two hooks which are called whenever an individual is bred or evaluated:

```
public void individualsBredStatistics(SteadyStateEvolutionState state,
    Individual[] individuals);
public void individualsEvaluatedStatistics(SteadyStateEvolutionState state,
    Individual[] newIndividuals, Individual[] oldIndividuals,
    int[] subpopulations, int[] indices);
```

The first method is called when one or more Individuals has just been bred (at present only one individual will be bred at a time). The second method is called when one or more Individuals have just been evaluated and inserted into the Population, dispersing other Individuals (at present, it’ll only be one individual evaluated). The Subpopulations and Individual indices in the Subpopulations where the event occurred are provided as well.

Last but not least, `ec.steady.SteadyStateDefaults` provides the package default parameter base.

**Example** We again build off of the example shown in Section 4.1.1, and continue to use generation boundaries rather than evaluations to keep things simple. So let’s use that file:

```
parent.0 = ga.params
```

Steady-State evolution doesn’t allow multiple threads:

```
evalthreads = 1
breedthreads = 1
```

Next, let’s override some parameters to use Steady-State Evolution:

```
state = ec.steadystate.SteadyStateEvolutionState
breed = ec.steadystate.SteadyStateBreeder
eval = ec.steadystate.SteadyStateEvaluator
```

Traditionally, steady-state evolutionary algorithms try hard not to produce duplicates during initialization:

```
steady.duplicate-retries = 100
```

We'll need to specify a deselector for each supopulation too. Let's do random selection:

```
steady.deselector.0 = ec.select.ec.select.RandomSelection
```

...and we're done! We don't bother changing the Statistics object. SimpleStatistics will do in a pinch when we're observing generation boundaries.



## Chapter 5

# Representations

### 5.1 Vector and List Representations (The `ec.vector` Package)

The `ec.vector` package defines one of the most common representations used in evolutionary algorithms: one-dimensional arrays of numbers or objects. In Genetic Algorithms such things have often been referred to as **chromosomes**. ECJ supports Individuals consisting of a single vector of the following types:

Array Type	Individual	Species
boolean	<code>ec.vector.BitVectorIndividual</code>	<code>ec.vector.VectorSpecies</code>
byte	<code>ec.vector.ByteVectorIndividual</code>	<code>ec.vector.IntegerVectorSpecies</code>
short	<code>ec.vector.ShortVectorIndividual</code>	<code>ec.vector.IntegerVectorSpecies</code>
int	<code>ec.vector.IntegerVectorIndividual</code>	<code>ec.vector.IntegerVectorSpecies</code>
long	<code>ec.vector.LongVectorIndividual</code>	<code>ec.vector.IntegerVectorSpecies</code>
float	<code>ec.vector.DoubleVectorIndividual</code>	<code>ec.vector.FloatVectorSpecies</code>
double	<code>ec.vector.DoubleVectorIndividual</code>	<code>ec.vector.FloatVectorSpecies</code>
Subclasses of <code>ec.vector.VectorGene</code>	<code>ec.vector.GeneVectorIndividual</code>	<code>ec.vector.GeneVectorSpecies</code>

Note that all the integer-type Individuals (byte, short, int, long) share the same species: `ec.vector.IntegerVectorSpecies`. Likewise, all the float-type Individuals (float, double) share `ec.vector.FloatVectorSpecies`. Last, `GeneVectorIndividual` holds an array of subclasses of the abstract class `ec.vector.VectorGene`, which you can subclass to do or hold any data type. All these Individuals are subclasses of the abstract class `ec.vector.VectorIndividual`; and all Species are subclasses of `ec.vector.VectorSpecies`.

Each of these Individuals has the same crossover operators and support for vector manipulation. They differ in the default initialization and mutation mechanisms available to them: obviously Gaussian mutation makes little sense on integers, for example.

To make matters simple, ECJ has a single set of BreedingPipelines for all of these data types. To do this, the BreedingPipelines call crossover or mutation functions in the Individuals themselves (or more properly, in their Species). This is different from other representations and results in some breeding parameters (crossover type, mutation probability) being found in the Species and not in the BreedingPipeline in question. But as it makes the package much smaller and simpler, so be it.

ECJ does not use generics in this package: each of the data types above has an Individual subclass all its own. The reason for this is straightforward: generics would be *exceptionally* slow. These are arrays of basic data types, and Java's generics would box and unbox them. So no generics it is.

Every vector or list has a single array which holds the genes proper. This array is:

```
public type[] genome;
```

... where *type* is the array type of the Vector (review the table above). You are welcome to read this array as you like, and modify it if you need to (don't replace it or change its length unless you're doing a List, and even then you should use the cover functions discussed in Section 5.1.2).

Many Vector genomes have minimum and maximum acceptable gene values. These values may be defined for the whole genome, for certain sections of it, or on a per-gene basis. For this reason, the easiest way to get this information is with the `minGene()` and `maxGene()` methods defined in the Vector's Species:

---

#### **ec.vector.IntegerVectorSpecies Methods**

```
public long minGene(int gene)
```

Returns the minimum valid gene value for this gene index.

```
public long maxGene(int gene)
```

Returns the maximum valid gene value for this gene index.

```
public boolean inNumericalTypeRange(long geneVal)
```

Returns true if the given gene value is within the numerical type range of the prototypical Individual for this Species. Note that just because the gene value lies within the type range does *not* mean that it's a valid value — you should use `minGene(...)` and `maxGene(...)` to determine this instead.

---

---

#### **ec.vector.FloatVectorSpecies Methods**

```
public double minGene(int gene)
```

Returns the minimum valid gene value for this gene index.

```
public double maxGene(int gene)
```

Returns the maximum valid gene value for this gene index.

```
public boolean inNumericalTypeRange(double geneVal)
```

Returns true if the given gene value is within the numerical type range of the prototypical Individual for this Species. Note that just because the gene value lies within the type range does *not* mean that it's a valid value — you should use `minGene(...)` and `maxGene(...)` to determine this instead.

---

### **5.1.1 Vectors**

The default use of this representation is as *fixed-length* vectors which can be crossed over and mutated. The size of the vectors is specified in the Species. For example, to define Individuals of the form of 100 bytes in Subpopulation 0, we might do this:

```
pop.subpop.0.species = ec.vector.IntegerVectorSpecies
pop.subpop.0.species.ind = ec.vector.ByteVectorIndividual
pop.subpop.0.species.genome-size = 100
```

The package `ec.vector.breed` contains Breeding Pipelines which know how to manipulate all such vectors. Specifically:

- `ec.vector.breed.VectorMutationPipeline` takes a single source and mutates Individuals from that source by calling the `mutate(...)` method on them (discussed later).
- `ec.vector.breed.VectorCrossoverPipeline` takes a two sources and draws one Individuals at a time each from these sources, then crosses them over by calling the `crossover(...)` method on one of them (discussed later).

- `ec.vector.breed.MultipleVectorCrossoverPipeline`, by Beenish Jamil, a former undergraduate at GMU, takes Individuals from a variable number of sources, then performs uniform crossover between all of them. For example, imagine if there were three Individuals,  $A$ ,  $B$ , and  $C$ . For each index  $i$ , `MultipleVectorCrossoverPipeline` randomly shuffles the values among  $A_i$ ,  $B_i$ , and  $C_i$  with a certain `crossoverProbability`, described next.

#### 5.1.1.1 Initialization

`VectorIndividuals` are initialized by being cloned from the prototypical `Individual`, and then having the following method called on them:

```
public void reset(EvolutionState state, int thread);
```

How `VectorIndividuals` implement this method depends on their type:

- **Floating Point** (float, double): each gene is set to a random value between its legal minimum and maximum values, inclusive.
- **Integer** (byte, short, int, long): each gene is set to a random value between its legal minimum and maximum values, inclusive.
- **Boolean**: each gene is set to a random value.
- **Gene**: each gene is randomized by calling the following method on its `ec.vector.VectorGene`:

```
public void reset(EvolutionState state, int thread);
```

Floating Point and Integer individuals need minimum and maximum values for their genes. These values can be set in any combination of three ways:

- Global settings for the entire `Individual`. For example:

```
pop.subpop.0.species.min-gene = -14
pop.subpop.0.species.max-gene = 92
```

- Settings for *segments* along the `Individual`. The number of segments is:

```
pop.subpop.0.species.num-segments = 4
```

Segments may be either specified by stating the *start indices* of each segment:

```
pop.subpop.0.species.segment-type = start

pop.subpop.0.species.segment.0.start = 0
pop.subpop.0.species.segment.1.start = 15
pop.subpop.0.species.segment.2.start = 50
pop.subpop.0.species.segment.3.start = 80
```

... or they may be specified by stating the *end indices* of each segment.<sup>1</sup>

```
pop.subpop.0.species.segment-type = end

pop.subpop.0.species.segment.0.end = 14
pop.subpop.0.species.segment.1.end = 49
pop.subpop.0.species.segment.2.end = 79
pop.subpop.0.species.segment.3.end = 100
```

---

<sup>1</sup>Why do this? It's a long story, with no good excuses.

Then we specify the min and max gene values for all genes in each segment:

```
pop.subpop.0.species.segment.0.min-gene = -14
pop.subpop.0.species.segment.0.max-gene = 92
pop.subpop.0.species.segment.1.min-gene = 0
pop.subpop.0.species.segment.1.max-gene = 100
pop.subpop.0.species.segment.2.min-gene = 0
pop.subpop.0.species.segment.2.max-gene = 50
pop.subpop.0.species.segment.3.min-gene = -0
pop.subpop.0.species.segment.3.max-gene = 1
```

The code for segments was provided by Rafal Kicingier, then a PhD student at GMU.

- Last, settings for *individual genes* may be stated. Not all genes have to be stated this way, just the ones you want. For example:

```
pop.subpop.0.species.min-gene.59 = 2
pop.subpop.0.species.max-gene.59 = 26
pop.subpop.0.species.min-gene.72 = 3
pop.subpop.0.species.max-gene.59 = 19
```

The rule is: individual gene settings override segment settings, which in turn override global settings.

#### 5.1.1.2 Crossover

The class `ec.vector.breed.VectorCrossoverPipeline` works by calling the following method on one of the `Individual`s, passing in the other:

```
public void defaultCrossover(EvolutionState state, int thread, VectorIndividual ind);
```

The two `Individual`s must be of the same type. You could override this method in a custom `VectorIndividual` of some sort to do your own crossover type.

All `VectorIndividual`s have default implementations of this method, which follow parameters specified in their `Species`. The first thing to know about crossover in vectors is that most forms (one-point, two-point, uniform) only occur along **chunk boundaries**. By default chunks are the size of a single gene; but you can make them larger. For example, you could specify that crossover can only occur every seven bytes in our `ByteVectorIndividual`:

```
pop.subpop.0.species.chunk-size = 7
```

Why would you do this? Mostly because you have encoded genes such that groups of seven genes together constitute a unit of some sort which shouldn't be broken up arbitrarily via crossover.

ECJ's `ec.vector.breed.VectorCrossoverPipeline` supports five types of crossover. The crossover type is specified like this:

```
pop.subpop.0.species.crossover-type = one-point
```

The types are:

- `one-point` performs standard One-Point crossover.
- `two-point` performs standard Two-Point crossover.
- `uniform` performs parameterized Uniform crossover, where every gene is crossed over independently with a certain probability. `ec.vector.breed.MultipleVectorCrossoverPipeline` also uses this probability. The probability is stated like this:



```
pop.subpop.0.species.crossover-prob = 0.25
```

(It doesn't make sense to use a probability over 0.5). It's important to note that the `crossover-probability` parameter is the *per-gene* crossover probability, not the probability that the *entire individual* will be crossed over. For that, see the `likelihood` parameter in Section 3.5.

- `line` performs Line Recombination. The two individuals are treated as points in space. A straight line is drawn through both points, and two children are created along this line. If the individuals are  $\vec{x}$  and  $\vec{y}$ , we draw two random values  $\alpha$  and  $\beta$ , each between  $-p$  and  $1 + p$  inclusive. Then the two children are defined as  $\alpha\vec{x} + (1 - \alpha)\vec{y}$  and  $\beta\vec{y} + (1 - \beta)\vec{x}$  respectively. For Integer vector individuals, the values of each index are then rounded to the nearest integer. The most common setting of  $p$  is 0.25, which allows children to be somewhat outside the hypercube formed with the parents at corners. To set the value of  $p$ , we state:

```
pop.subpop.0.species.line-extension = 0.25
```

Line Recombination cannot be performed on `BitVectorIndividuals` or `GeneVectorIndividuals`. ECJ's implementation of Line Recombination is due to William Haddon, a former student at GMU.

- `intermediate` performs Intermediate Recombination. This is very similar to Line Recombination. The two individuals are again treated as points in space. Children are created somewhere in the vicinity of the hypercube formed with the two individuals as corners. If the individuals are  $\vec{x}$  and  $\vec{y}$ , for *every index*  $i$  we draw two independent random values  $\alpha_i$  and  $\beta_i$ , each between  $-p$  and  $1 + p$  inclusive. Then genes at index value  $i$  of the two children are defined as  $\alpha_i x_i + (1 - \alpha_i) y_i$  and  $\beta_i y_i + (1 - \beta_i) x_i$  respectively. For Integer vector individuals, the values of each index are then rounded to the nearest integer. Again, the most common setting of  $p$  is 0.25, which allows children to be somewhat outside the hypercube formed with the parents at corners. Again, to set the value of  $p$ , we state:

```
pop.subpop.0.species.line-extension = 0.25
```

Intermediate Recombination cannot be performed on `BitVectorIndividuals` or `GeneVectorIndividuals`. ECJ's implementation of Intermediate Recombination is also due to William Haddon.

- `sbx` performs Simulated Binary Crossover. This crossover, by Kalyanmoy Deb, attempts to simulate the effect of binary crossover on doubles and floats as if they were encoded as bits. It relies on the Polynomial Distribution and as such has a parameter called the *crossover distribution index* which specifies the kind of distribution. The standard accepted value for this index is 20 (but it's not the ECJ default: you must set it):

```
pop.subpop.0.species.crossover-distribution-index = 20
```

You can use any integer  $\geq 0$ . There are probably several different versions of SBX out there. The implementation in ECJ is a faithful reproduction of the one which appears in the NSGA-II C code found here: <http://www.iitk.ac.in/kangal/codes/nsga/nsgaorig.tar>

`VectorCrossoverPipeline` can be set to return both crossed-over children or just one of them. This parameter isn't in the `Species` but is actually in the `Pipeline` proper. Let's say that `VectorCrossoverPipeline` was the root `Pipeline` for Subpopulation 0. To set it to return just one child, you'd say:

```
pop.subpop.0.species.pipe = ec.vector.breed.VectorCrossoverPipeline
pop.subpop.0.species.pipe.toss = true
```

Alternatively you could use the default base:

```
vector.xover.toss = true
```

The default value is false.

### 5.1.1.3 Multi-Vector Crossover

The class `ec.vector.breed.MultipleVectorCrossoverPipeline` performs uniform crossover among the Individuals in question as described earlier. It ignores the `crossover-type` parameter, but it does respect, and indeed require, the parameter:

```
pop.subpop.0.species.crossover-prob = 0.25
```

This is the probability that a given gene index will be shuffled among the various Individuals. As before, it doesn't make much sense for this to be set to any value over 0.5.

### 5.1.1.4 Mutation

Similarly, the BreedingPipeline `ec.vector.breed.VectorMutationPipeline` does its work by calling the following method on the Individual in question:

```
public void defaultMutate(EvolutionState state, int thread);
```

You could override this method in a custom `VectorIndividual` of some sort to do your own custom kind of mutation.

Every `VectorIndividual` has a default implementation of this method. Mutation parameters are also specified in the Species of each Individual. At present the default mutation procedures do not respect chunk boundaries, unlike Crossover. However, each gene is only mutated with a certain probability. This probability is defined as:

```
pop.subpop.0.species.mutation-probability = 0.1
```

Different representations have different kinds of mutations. Specifically:

- **Boolean** (boolean): the default mutation operator simply flips the boolean value with the given mutation probability.
- **Integer** (byte, short, int, long): the default mutation operator simply randomizes the integer value between the legal minimum and maximum values (with the mutation probability of course). There are almost certainly smarter approaches than this — you'll probably want to implement something in a subclass.
- **Gene**: the default mutation operator for `ec.vector.GeneVectorIndividual`, with the given mutation probability, calls the following `VectorGene` method:

```
public void mutate(EvolutionState state, int thread);
```

You are responsible for implementing this method. The default version of the method calls:

```
public void reset(EvolutionState state, int thread);
```

...on the `VectorGene`.

- **Floating Point** (float, double): allows uniform gene randomization (`reset`), Gaussian mutation (`gauss`), or Polynomial mutation (`polynomial`).

**Uniform Mutation** Uniform gene randomization simply sets the gene to a random value between its minimum and maximum legal values. To use it, you say:

```
pop.subpop.0.species.mutation-type = reset
```

**Gaussian Mutation** Gaussian mutation adds Gaussian noise to the current value. If the result is outside the bounds of minimum and maximum legal values, another Gaussian noise is tried instead, and so on, until a legal value is found. You will need to specify a standard deviation for the Gaussian random noise and also the number of times ECJ should attempt to apply Gaussian noise within the minimum and maximum bounds before giving up and simply keeping the original value. For example:

```
pop.subpop.0.species.mutation-type = gauss
pop.subpop.0.species.mutation-stdev = 0.1
pop.subpop.0.species.out-of-bounds-retries = 20
```

Note that if `out-of-bounds-retries` is 0, then ECJ behaves specially: it never stops retrying. The default value is 100.

Alternatively you can add Gaussian noise to the current values and simply allow the results to freely exceed the min/max gene bounds. No retries are attempted (and `out-of-bounds-retries` is ignored): noise is simply added and that's it. If you do this, be certain that the particular problem you're testing works properly with values outside the stated initial bounds. To lift the bounds in this way, you say:

```
pop.subpop.0.species.mutation-bounded = false
```

This does not affect the use of gene bounds in determining initialization values. Even if you declare the mutation to be bounded, you'll still need to provide `out-of-bounds-retries` (though it won't be used).

**Polynomial Mutation** Finally, Polynomial mutation is by Kalyanmoy Deb, and uses a distribution chosen from the Polynomial distribution series rather than a Gaussian distribution. All Polynomial distributions have a min and max value ranging in  $[-1, 1]$ , and Polynomial mutation adjusts this to reflect the min and max gene ranges. The Polynomial distribution used is defined by the *distribution index* and by far the most common setting (but not the ECJ default) is 20. There are at least four different versions of Polynomial mutation in the literature or various common code. We have implemented two versions: the **standard** version, which can produce values outside the min/max range, and an **alternative** version found in the NSGA-II C code among certain other places (<http://www.iitk.ac.in/kangal/codes/nsga/nsgaorig.tar>), which modifies the distribution to always mutate values within the min/max range. Note that we've found the alternative version superior for multiobjective optimization problems. In ECJ, **the alternative version is the default**.

To use the *alternative* version, you'll need to state something like this.

```
pop.subpop.0.species.mutation-type = polynomial
pop.subpop.0.species.mutation-distribution-index = 20
pop.subpop.0.species.alternative-polynomial-version = true
pop.subpop.0.species.out-of-bounds-retries = 20
```

You can use any integer  $\geq 0$  for the distribution index. Note that even though the alternative version doesn't actually need `out-of-bounds-retries`, ECJ requires it anyway, so just put something arbitrary there.

To use the *standard* version, you'll need to state something along these lines.

```

pop.subpop.0.species.mutation-type = polynomial
pop.subpop.0.species.mutation-distribution-index = 20
pop.subpop.0.species.alternative-polynomial-version = false
pop.subpop.0.species.out-of-bounds-retries = 20

```

Here you'll need to provide a rational value for out-of-bounds-retries.

Just as was the case for Gaussian noise, you have the option of lifting the bounds entirely and letting Polynomial Mutation freely exceed the min/max gene bounds. In this case, no retries are attempted (and out-of-bounds-retries is ignored): noise is simply added and that's it. Again, if you do this, be certain that the particular problem you're testing works properly with values outside the stated initial bounds. To lift the bounds in this way, you say:

```

pop.subpop.0.species.mutation-bounded = false

```

Once again, this does not affect the use of gene bounds in determining initialization values. And even if you declare the mutation to be bounded, you'll still need to provide out-of-bounds-retries (though it won't be used).

Numeric representations (integer, floating-point) provide two utility methods which may be useful to you in creating or debugging a custom mutator:

```

public boolean isInRange();
public void clamp();

```

The first method returns true if all the genes in the representation are within their minimum and maximum per-gene bounds. The second method insures that this is the case: if a gene is larger than its maximum, it is set to the maximum; and if it is smaller than the minimum, it is set to the minimum.

**Example** In the Genetic Algorithms section (Section 4.1.1), we provided an example for a pipeline for crossing over and mutating VectorIndividuals:

```

pop.subpop.0.species.pipe = ec.vector.breed.VectorMutationPipeline
pop.subpop.0.species.pipe.source.0 = ec.vector.breed.VectorCrossoverPipeline
pop.subpop.0.species.pipe.source.0.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.0.source.1 = same
select.tournament.size = 2

```

The example used IntegerVectorIndividuals. Let's change it to DoubleVectorIndividuals, and also specify some constraints.

```

pop.subpops.0.species = ec.vector.FloatVectorSpecies
pop.subpop.0.species.ind = ec.vector.DoubleVectorIndividual
pop.subpop.0.species.genome-size = 100

```

Let's also add some minimum and maximum gene information, both globally and for the first gene:

```

pop.subpops.0.species.min-gene = 0.0
pop.subpops.0.species.max-gene = 1.0
pop.subpops.0.species.min-gene.0 = 0.0
pop.subpops.0.species.max-gene.0 = 2.0

```

We'll do gaussian mutation with 100% likelihood, and some uniform crossover:

```

pop.subpop.0.species.crossover-type = uniform
pop.subpop.0.species.crossover-prob = 0.25
pop.subpop.0.species.mutation-prob = 1.0
pop.subpop.0.species.mutation-type = gauss
pop.subpop.0.species.mutation-stdev = 0.1

```

**Another Example** Bit vectors used to be very common. Let's change our Individual to one. Note that it uses `ec.vector.VectorSpecies` as its Species:

```

pop.subpops.0.species = ec.vector.VectorSpecies
pop.subpop.0.species.ind = ec.vector.BooleanVectorIndividual
pop.subpop.0.species.genome-size = 128

```

Perhaps our encoding uses each 32 bits to represent some important concept, so we'd like to prevent crossover from occurring anywhere except on 32-bit boundaries:

```

pop.subpops.0.species.chunk-size = 32

```

Booleans only have one mutation parameter: probability of bit-flip. Let's also include two-point crossover on the chunk boundaries:

```

pop.subpop.0.species.crossover-type = one
pop.subpop.0.species.mutation-prob = 0.01

```

## 5.1.2 Lists

While vectors are arrays of fixed length, Lists are arrays of arbitrary and possibly changing length. List representations are much less common in evolutionary algorithms.

With a few changes, the classes in ECJ's `ec.vector` package may be used for lists as well as for vectors. ECJ supports list representations in four ways:

- Useful utility methods for manipulating lists.
- List-oriented initialization
- List-oriented crossover procedures.
- Default vector mutation also works fine for lists.

Let's cover each in turn. But first read the previous section on Vectors (Section 5.1.1). And note that lists completely ignore the chunk facility: it's too complex.

### 5.1.2.1 Utility Methods

`VectorIndividual` provides methods for manipulating the genome without knowing the type of the object in question. First, the genome (the array) may be retrieved and set:

```

public Object getGenome();
public void setGenome(Object genome);

```

Second, the genome length can be changed. If the length is shortened, the array will be truncated. If the length is lengthened, the extra area will be set as follows: boolean array slots will be set to false. Numeric

array slots will be set to 0. VectorGene array slots will be set to a blank VectorGene cloned from the prototype. The methods for checking or changing the length are:

```
public int size();
public int genomeLength();
public void setGenomeLength(int length);
public void reset(EvolutionState state, int thread, int newSize);
```

The first two methods are synonymous. The last method resizes the individual, then resets it entirely (randomizes its values).

Third, VectorIndividual supplies methods for cutting a splicing genome arrays, which is quite useful for manipulating variable-length lists. For example, you can **split** a genome into some  $n$  pieces. To do this you provide  $n - 1$  split points. The first piece will start at index 0 and run up to, but not including, the first split point; the second piece will start with the split point and run up to, but not including, the second split point; and so on. The final piece will run to the end of the genome. The  $n$  pieces are placed into the provided Object array.

```
public void split(int[] points, Object[] pieces);
```

VectorIndividual can also **join** certain pieces — concatenate them — to form the genome, replacing the original genome.

```
public void join(Object[] pieces);
```

Last, VectorIndividual can clone genes in the arrays split off of a genome before joining them again. This is useful for concatenating genes to an Individual for example. Obviously, genomes consisting of arrays of numbers or booleans don't need to be "cloned" — but the VectorGenes found in GeneVectorIndividual need to be copied or else the concatenated gene slots will share pointers with the original genes, which you perhaps didn't want. Note that the pieces array is modified in place.

```
public void cloneGenes(Object[] pieces);
```

### 5.1.2.2 Initialization

The initial size of the genome can be hard-coded in parameters (just like it was for vectors), or you can specify one of two algorithms for picking a size. To hard-code the size, just do it like a vector:

```
pop.subpop.0.species.genome-size = 100
```

Alternatively you can ask ECJ to pick the genome size uniformly from between a minimum and maximum size inclusive:

```
pop.subpop.0.species.genome-size = uniform
pop.subpop.0.species.min-initial-size = 10
pop.subpop.0.species.max-initial-size = 80
```

Last, you can ask ECJ to pick the genome size from the geometric distribution. Here ECJ will start at the minimum size, then flip a coin of a given probability. If the coin comes up heads (true), ECJ will increase the size by one, then flip the coin again, and so on, until the coin comes up tails (false). That'll be the resulting size. For this you need to provide the probability and the minimum size:

```
pop.subpop.0.species.genome-size = geometric
pop.subpop.0.species.min-initial-size = 10
pop.subpop.0.species.geometric-prob = 0.95
```

### 5.1.2.3 Crossover

In the `ec.vector.breed` package is one special pipeline made for crossing over lists, called, not surprisingly, `ec.vector.breed.ListCrossoverPipeline`. This class was developed by Stephen Donnelly, then an undergraduate at GMU.

`ListCrossoverPipeline` performs either one-point or two-point list crossover. In one-point list crossover, we pick random indexes  $i$  and  $j$  in each of two individuals  $A$  and  $B$  (which may have different lengths). We then swap the string of genes  $A_i, \dots, A_{end}$  and  $B_j, \dots, B_{end}$ . It's possible for these strings to be all of the Individual, or entirely empty.

In two-point list crossover, for each individual we pick *two* random indexes:  $i \leq j$  for Individual  $A$  and  $k \leq l$  for Individual  $B$ . We then swap the string of genes  $A_i, \dots, A_j$  with  $B_k, \dots, B_l$ . Again, it's possible for these strings to be all of the Individual, or entirely empty.

You specify the form of crossover like you do with Vectors: in the Species.

```
pop.subpop.0.species.crossover-type = one-point
```

(or two-point, if you prefer).

`ListCrossoverPipeline`'s crossover can be constrained in several ways. First, you can stipulate that the children are no less than a certain size. Let's say that `ListCrossoverPipeline` was the root Pipeline for Subpopulation 0. To set it to never return children less than 5 in length, you might say:

```
pop.subpop.0.species.pipe = ec.vector.breed.ListCrossoverPipeline
pop.subpop.0.species.pipe.min-child-size = 5
```

The default is 0.

To set it to never remove more than a certain percentage, and no less than another percentage, of material from a parent, you could say:

```
pop.subpop.0.species.pipe.min-crossover-percent = 0.25
pop.subpop.0.species.pipe.max-crossover-percent = 0.50
```

The defaults are 0.0 and 1.0 respectively.

`ListCrossoverPipeline` will try repeatedly to find crossover points which met these constraints. If it fails, it will simply return the two parents. You set the number of times it will try like this:

```
pop.subpop.0.species.pipe.tries = 100
```

The default is 1.

Like `VectorCrossoverPipeline`, `ListCrossoverPipeline` can be set to return both crossed-over children or just one of them. To set it to return just one child, you'd say:

```
pop.subpop.0.species.pipe.toss = true
```

The default value is false.

Alternatively you could use the default base for any or all of these:

```

vector.list-xover.min-child-size = 5
vector.list-xover.min-crossover-percent = 0.25
vector.list-xover.max-crossover-percent = 0.50
vector.list-xover.tries = 100
vector.list-xover.toss = true

```

#### 5.1.2.4 Mutation

The standard vector mutation operator will work fine in list mode: it doesn't care what the length of the array is.

ECJ at present has a single mutation operators special to lists: `ec.vector.breed.GeneDuplicationPipeline`. This pipeline replicates a particular mutation operation common in Grammatical Evolution (Section 5.3). Specifically: we select two indices in the list at random such that they are not equal to one another. Then we copy the region between the two indices, and concatenate it to the end of the list. And that's it!

**Example** Let's modify the Vector example given previously. First, we'll change the crossover procedure:

```

pop.subpop.0.species.pipe = ec.vector.VectorMutationPipeline
pop.subpop.0.species.pipe.source.0 = ec.vector.ListCrossoverPipeline
pop.subpop.0.species.pipe.source.0.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.0.source.1 = same
select.tournament.size = 2

```

Let's stipulate that crossed-over individuals must be at least two genes in length if possible:

```

pop.subpop.0.species.pipe.source.0.min-child-size = 2

```

Finally, let's change the procedure for determining the initial size of an Individual:

```

pop.subpop.0.species.genome-size = geometric
pop.subpop.0.species.min-initial-size = 10
pop.subpop.0.species.geometric-prob = 0.95

```

### 5.1.3 Arbitrary Genes: `ec.vector.VectorGene`

Last, some discussion should be reserved regarding vectors of "arbitrary genes". The idea behind this is to provide you with maximum flexibility as to what you can create vectors out of. The classes involved here are:

- `ec.vector.VectorGene`, the abstract superclass of objects which can fill gene positions.
- `ec.vector.GeneVectorIndividual`, the `VectorIndividual` which contains nothing but `VectorGenes`.
- `ec.vector.GeneVectorSpecies`, the `Species` for `GeneVectorIndividuals`.

To use these classes, you'll not only need to specify the use of `GeneVectorIndividual` and `GeneVectorSpecies`, but you'll also need to state which subclass of `VectorGene` will be used to fill the `GeneVectorIndividual`. These three things are done as follows:

```

pop.subpop.0.species = ec.vector.GeneVectorSpecies
pop.subpop.0.species.ind = ec.vector.GeneVectorIndividual
pop.subpop.0.species.gene = ec.app.MySubclassOfVectorGene

```



As mentioned earlier, `GeneVectorIndividual`'s `reset(...)` method, used to initialize the `Individual`, in turn calls this method on your `VectorGene` subclass:

```
public abstract void reset(EvolutionState state, int thread);
```

Furthermore, mutation on `GeneVectorIndividuals` call this method on each `VectorGene` (with the mutation probability):

```
public abstract void mutate(EvolutionState state, int thread);
```

The default form of this method just calls `reset(...)`.

`VectorGene` is a `Prototype`, and so requires `setup(...)`, `clone(...)`, and `defaultBase()`. Default implementations are already provided for you. Except for `clone(...)` you're unlikely to need to override them.

You do need to provide two additional `VectorGene` methods:

```
public abstract int hashCode();
public abstract boolean equals(Object other);
```

The first method should provide an intelligent hash code based on the value of the `VectorGene` contents. The second method should test this `VectorGene` against another for equality and return true if they contain the same thing.

Technically that's all you need to provide. But in reality `VectorGene` is not going to be very useful unless you at least provide a way to describe the gene when it's printed to a log. The easiest way to do this is to override this method:

```
public String printGeneToStringForHumans();
```

The default version of this method simply calls `toString()` (which you could override too if you wanted). This method is in turn called by

```
public void printGeneForHumans(EvolutionState state, int verbosity, int log);
```

If you're writing `Individuals` with the intent that they be read back in again later, you'll probably want to override this method:

```
public String printGeneToString();
```

This method is called in turn by the following two methods:

```
public void printGene(EvolutionState state, int verbosity, int log);
public void printGene(EvolutionState state, PrintWriter writer);
```

The default form simply calls `toString()`, which is definitely wrong. You'll also want to override the method which reads in the gene again:

```
public void readGeneFromString(String string, EvolutionState state);
```

To write out the gene in a computer-readable fashion, I suggest using ECJ's Code package (Section 2.2.2), but it's up to you. This method is called by

```
public void readGene(EvolutionState state, LineNumberReader reader)
                                                    throws IOException;
```

Finally if you're intending to send your `Individual` over the network, either for distributed evaluation or island models, you'll need to implement these methods:

```

public void writeGene(EvolutionState state, DataOutput output)
                                                    throws IOException;
public void readGene(EvolutionState state, DataInput input) throws IOException;

```

**Example** The following VectorGene contains two doubles and mutates them by performing a random affine rotation on them (why not?). We'll implement it fully here, though keep in mind that this isn't entirely necessary. For fun, we'll use an elaborate hash code which XORs all four 32-bit segments of the two doubles together (I hope I wrote that right!).

First the parameters:

```

pop.subpop.0.species = ec.vector.GeneVectorSpecies
pop.subpop.0.species.ind = ec.vector.GeneVectorIndividual
pop.subpop.0.species.gene = ec.app.trig.TrigGene

```

Now the implementation:

```

package ec.app.trig;
import ec.util.*;
public class TrigGene extends VectorGene {
    double x;
    double y;
    public void reset(EvolutionState state, int thread) {
        double alpha = state.random[thread].nextDouble() * Math.PI * 2;
        x = Math.cos(alpha);    y = Math.sin(alpha);
    }

    public void mutate(EvolutionState state, int thread) {
        double alpha = Math.atan2(y,x);
        double dalpha = (state.random[thread].nextDouble() - 0.5) * Math.PI * 2 / 100.0;
        x = Math.cos(alpha + dalpha);    y = Math.sin(alpha + dalpha);
    }

    public int hashCode() {
        long a = Double.doubleToRawLongBits(x); long b = Double.doubleToRawLongBits(y);
        return (int) ((a & (int)-1) ^ (a >> 32) (b & (int)-1) ^ (b >> 32));
    }

    public boolean equals(Object other) {
        return (other != null && other instanceof TrigGene &&
            ((TrigGene)other).x == x && ((TrigGene)other).y == y);
    }

    public String printGeneToStringForHumans() { return ">" + x + " " + y ; }

    public String printGeneToString() {
        return ">" + Code.Encode(x) + " " + Code.Encode(y);
    }

    public void readGeneFromString(String string, EvolutionState state) {
        string = string.trim().substring(0); // get rid of the ">"
        DecodeReturn dr = new DecodeReturn(string);
        Code.decode(dr); x = dr.d;           // no error checking
        Code.decode(dr); y = dr.d;
    }

    public void writeGene(EvolutionState state, DataOutput out) throws IOException {
        out.writeDouble(x); out.writeDouble(y);
    }

    public void readGene(EvolutionState state, DataOutput in) throws IOException {
        x = in.readDouble(); y = in.readDouble();
    }
}

```

## 5.2 Genetic Programming (The ec.gp Package)

The ec.gp package is far and away the most developed and tested package in ECJ. ECJ was largely developed in order to support his package, and much of our existing literature is based on it.

ECJ's genetic programming package uses "Koza-style" tree structures [3, 4] which represent the parse trees of Lisp s-expressions. For an introduction to genetic programming, see [14]. Much of ECJ's approach

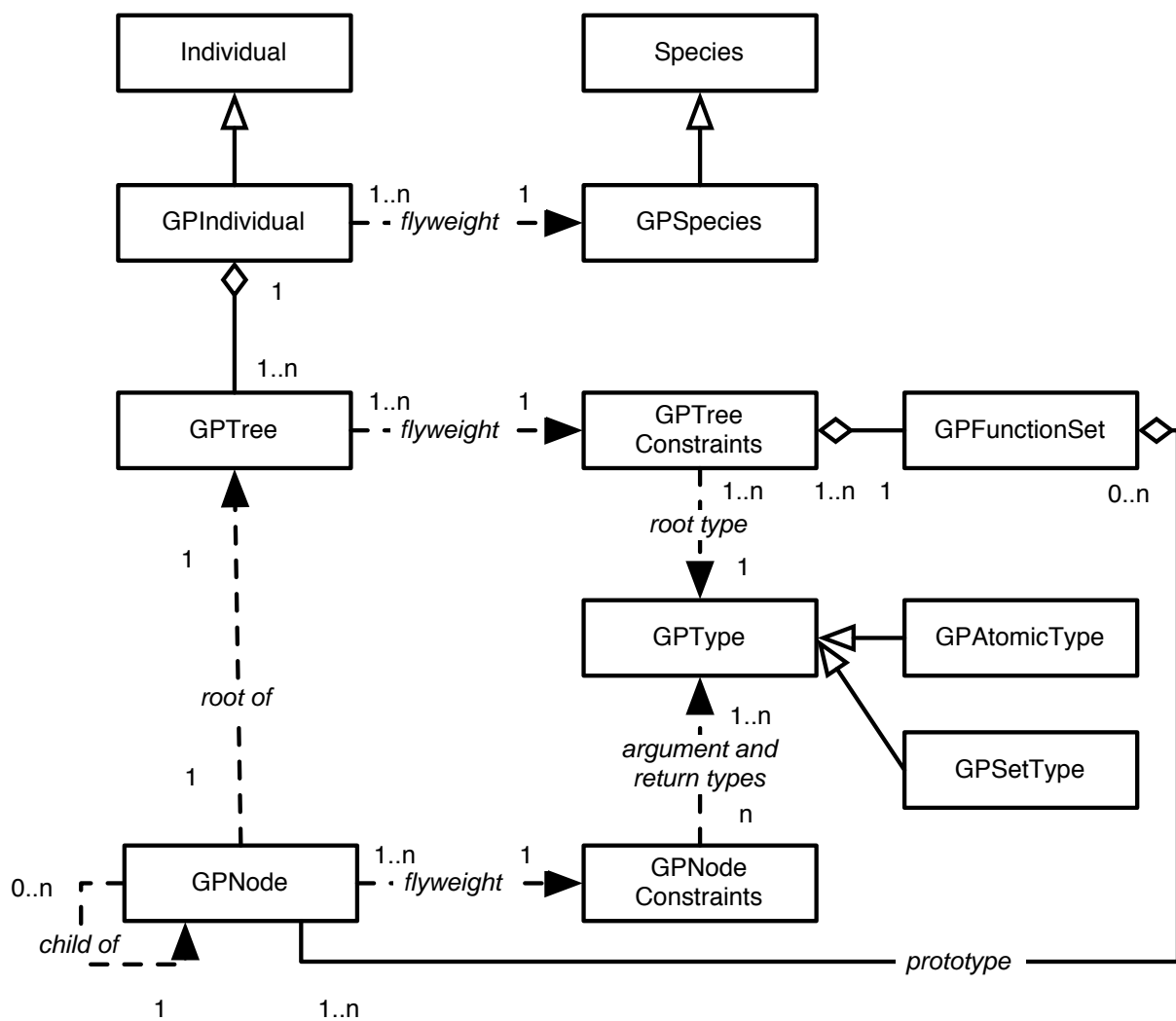


Figure 5.1 Data objects common to tree-based “Koza-style” genetic programming Individuals.

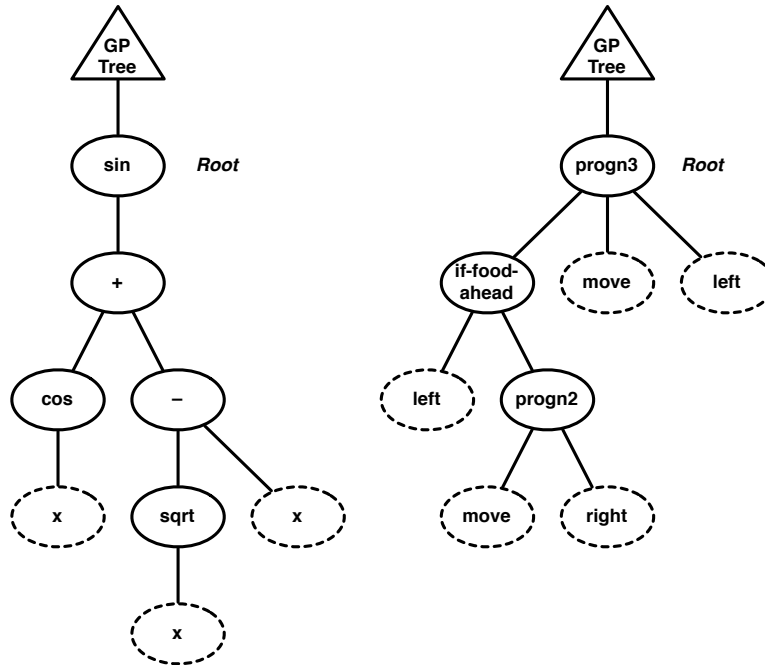


Figure 5.2 Two example genetic programming parse trees. At top is a single `ec.gp.GPTree` instance, which holds onto a single `ec.gp.GPNode` designated the *root* of the tree. `GPNodes` form the tree itself, and so have a *parent* and zero or more *children*. The parent of the root is the `GPTree` object itself. Leaf nodes, denoted with dotted ovals, are traditionally called *terminals*, and non-leaf nodes, including the root, are traditionally called *nonterminals*. Normally `GPNodes` have fixed arity. That is, all `if-food-ahead` `GPNodes` will always have two children, and all `cos` nodes will always have one child, etc.

to GP is inspired by *lil-gp* [16], an earlier C-based GP system. However *lil-gp* and many other GP systems pack the parse trees into arrays to save memory. ECJ does not: the parse trees are stored as tree structures in memory. This is much more wasteful of memory but it is faster to evaluate and far easier to manipulate.

GP's top-level class is an `Individual` called `ec.gp.GPIndividual`. `GPIndividual` holds an array of GP trees, held by `ec.gp.GPTree` objects. Each GP tree is a tree of `ec.gp.GPNode` objects. One `GPNode`, the root of the tree, is held by the `GPTree`.

`GPIndividual`, `GPTree`, and `GPNode` are all `Prototypes`, and furthermore they all adhere to the *flyweight* pattern (Section 3.1.4). `GPIndividual`'s flyweight relationship is with a `Species` (of course), called `ec.gp.GPSpecies`. `GPTrees` have a flyweight relationship with subclasses of `ec.gp.GPTreeConstraints`. `GPNodes` have a flyweight relationship with subclasses of `ec.gp.GPNodeConstraints`.

GP's tree nodes are **typed**, meaning that they can have certain constraints which specify which nodes may serve as children of other nodes. These types are defined by an abstract class called `ec.gp.GPType`, of which there are two concrete subclasses, `ec.gp.GPAtomicType` and `ec.gp.GPSetType`.

The primary function of `GPSpecies` is to build new `GPIndividuals` properly. The primary function of `GPTreeConstraints` is to hold onto the **function set** (`ec.gp.GPFunctionSet`) for a given tree. This is a set of prototypical `GPNodes`, copies of which are used to construct the tree in question. `GPTreeConstraints` also contains typing information for the tree root. The primary purpose of `GPNodeConstraints` is to provide typing and arity information for various `GPNode`.

### 5.2.1 GPNodes, GPTrees, and GPIndividuals

Figure 5.2 shows two example trees of `GPNodes` (shown as ovals). The top of each tree is a `GPTree`, and directly under it is the *root* `GPNode`. As can be seen from the figure, each `GPNode` has both a parent and

zero or more children; and each GPTree has exactly one child. Both GPNodes and GPTrees implement `ec.gp.GPNodeParent`, and can serve as parents of other GPNodes (the root has the GPTree as its parent).

### 5.2.1.1 GPNodes

A basic GPNode consists of four items:

```
public GPNodeParent parent;
public GPNode children[];
public byte argposition;
public byte constraints;
```

The parent should be self-explanatory. The `children[]` is an array holding the children to the GPNode. Leaf nodes in a GP tree (traditionally called **terminals** are permitted to either have a zero-length array or a null value for `children[]`.

The `argposition` is the position of the node in its *parent's* `children[]` array. The root's `argposition` is 0. Last, the `constraints` is a tag which refers to the GPNode's `GPNodeConstraints` object. It's a byte rather than a full pointer to save a bit of space: GPNodes make up by far the bulk of memory in a genetic programming experiment. You can get the `GPNodeConstraints` by calling the following GPNode method:

```
public final GPNodeConstraints constraints(GPInitializer initializer);
```

Why the `GPInitializer`? Because `GPNodeConstraints`, `GPTreeConstraints`, `GPTypes`, and `GPFunctionSets` are all accessible via `Initializer`, which must be a `GPInitializer`.<sup>2</sup> More on that later. Assuming you have access to the `EvolutionState` (probably called `state`) can call this function like this:

```
GPNodeConstraints constraints = myGPNode.constraints((GPInitializer)(state.initializer));
```

You will make various subclasses of GPNode to define the kinds of functions which may appear in your genetic programming tree.

### 5.2.1.2 GPTrees

Unlike GPNode, which is liberally subclassed, you'll rarely subclass GPTree. The `ec.gp.GPTree` class holds onto the root GPNode here:

```
public GPNode child;
```

Each GPTree also has a backpointer to the `GPIndividual` which holds it:

```
public GPIndividual owner;
```

GPTree also has a pointer to its `GPTreeConstraints` object. Like GPNode, GPTree uses a byte rather than a full pointer.<sup>3</sup>

```
public byte constraints;
```

Just like GPNode, you can access GPTree's constraints using this function:

```
public final GPTreeConstraints constraints(GPInitializer initializer);
```

...which is typically called like this:

---

<sup>2</sup>It wasn't a good decision to use the `Initializer` in this fashion, and one day we may change it to something else.

<sup>3</sup>This is mostly historic: GPTree doesn't fill nearly as much memory as GPNode and so doesn't really need this tight reference approach.

```
GPTreeConstraints constraints = myGPTree.constraints((GPInitializer)(state.initializer));
```

### 5.2.1.3 GPIndividual

The GPIndividual contains an array of GPTrees. In most cases, this array has a single GPTree in it:

```
public GPTree[] trees;
```

### 5.2.1.4 GPNodeConstraints

The GPNodeConstraints contains several data elements shared by various GPNodes:

```
public byte constraintNumber;
public GPType returnType;
public GPType[] childtypes;
public String name;
public float probabilityOfSelection;
public GPNode zeroChildren[] = new GPNode[0];
```

The first element is obvious: it's number of the constraints which the GPNode objects point to. The next two items hold the return type and children types of the node: more on that later. Specifically, the return type of a child in slot 0 must be compatible with the child type declared for slot 0. For now what matters is that you can determine the expected number of children to a GPNode by the length of the childtypes array.

The name variable holds the name of the GPNodeConstraints (*not* the GPNodes which refer to them): we'll define some in the next section. The probabilityOfSelection variable holds an auxiliary variable used by certain tree-building operators. Last, zeroChildren[] holds a blank, zero-length GPNode which terminals are free to use in lieu of null for their children.

```
public GPTree[] trees;
```

### 5.2.1.5 GPTreeConstraints

The GPTreeConstraints contains data elements shared by GPTrees:

```
public byte constraintNumber;
public GPType treetype;
public String name;
public GPNodeBuilder init;
public GPFunctionSet functionset;
```

The first element is again obvious. The treetype variable declares the GPType for the tree as a whole: the return type of the root must be compatible with this type. The name works similarly to the one in GPNodeConstraints.

The last two variables are critical. The init variable holds the algorithm used to generate trees or subtrees for this GPTree. We will discuss tree builders later. Last, the functionset variable holds the function set for this tree: all GPNodes appearing in this GPTree must be cloned from this function set.

### 5.2.1.6 GPFunctionSet

The GPFunctionSet contains a name (like GPNodeConstraints and GPTreeConstraints) and a set of GPNodes, clones of which may appear in the GPTree. This set is stored in various hash tables and arrays to make

lookup easy for different common queries (such as “give me all terminals”) or (“give me all nodes whose return type is *foo*”). Usually you don’t need to access this class directly: instead, we’ll set up the function set using parameters.

## 5.2.2 Basic Setup

Now let’s work towards setting a GP problem. We begin by defining the `GPIndividual` and `GPSpecies`. Usually, we’ll just use those classes directly:

```
pop.subpop.0.species = ec.gp.GPSpecies
pop.subpop.0.species.ind = ec.gp.GPIndividual
```

Let’s presume for now that we just want a single tree per `GPIndividual`. This is the usual case. Typically this class is defined by `GPTree` unless we’re doing something odd. We say:

```
pop.subpop.0.species.ind.numtrees = 1
pop.subpop.0.species.ind.tree.0 = ec.gp.GPTree
```

Different trees can have different `GPTreeConstraints` objects, or share them. This is done by defining a set of `GPTreeConstraints` (which is a `Clique`, Section 3.1.2) and giving each member of the set a unique identifier. Then `GPTrees` identify with a given `GPTreeConstraints` by using that identifier.

Since we have only one tree, we really only need to create one `GPTreeConstraints`. We’ll call it “`tc0`”.

```
gp.tc.size = 1
gp.tc.0 = ec.gp.GPTreeConstraints
gp.tc.0.name = tc0
```

Note that as a `Clique`, `GPTreeConstraints` objects all have a global parameter base of `gp.tc`. Now we assign it to the tree:

```
pop.subpop.0.species.ind.tree.0.tc = tc0
```

A `GPTreeConstraints` object in turn holds onto the `GPFunctionSet` used to construct trees which identify with it. `GPFunctionSet` is also a `clique`. We’ll call the function set “`f0`”:

```
gp.fs.size = 1
gp.fs.0 = ec.gp.GPFunctionSet
gp.fs.0.name = f0
```

Note that as a `Clique`, `GPFunctionSet` objects all have a global parameter base of `gp.fs`. We now assign this function set to our tree constraints:

```
gp.tc.0.fset = f0
```

As to types: we’ll discuss typed GP later on. For now we’ll assume that there is a single atomic type which is used universally by everyone — that is, everything can connect with everything (it’s “typeless”). This is the classic GP scenario. `GPTypes` are also a `clique`: and they have a global parameter base of `gp.type`. We define zero `GPSetTypes` and one `GPAtomicType` (which we will name, for lack of a better word, “`nil`”) like this:

```
gp.type.a.size = 1
gp.type.a.0.name = nil
gp.type.s.size = 0
```

Our `GPTreeConstraints` object needs to define the `GPType` of the tree as a whole (the “root type”). To set it to our `nil` type, we’d say:



```
gp.tc.0.returns = nil
```

Last, we need to define some `GPNodeConstraints`. A `GPNodeConstraint` object describes three things about the `GPNodes` related to them via the Flyweight pattern:

- The number of children of the `GPNode`.
- The `GPTypes` that the children of the `GPNode` must be consistent with.
- The `GPType` of that the parent of the `GPNode` must be consistent with.

More on types later. But for now we'll define a few `GPNodeConstraints` for nodes with zero, one, and two children. Since we only have one type, the types of all the children and the parent are all going to be `nil`. We'll call these `GPNodeConstraints` `nc0`, `nc1`, and `nc2`.

```
gp.nc.size = 3

gp.nc.0 = ec.gp.GPNodeConstraints
gp.nc.0.name = nc0
gp.nc.0.returns = nil
gp.nc.0.size = 0

gp.nc.1 = ec.gp.GPNodeConstraints
gp.nc.1.name = nc1
gp.nc.1.returns = nil
gp.nc.1.size = 1
gp.nc.1.child.0 = nil

gp.nc.2 = ec.gp.GPNodeConstraints
gp.nc.2.name = nc2
gp.nc.2.returns = nil
gp.nc.2.size = 2
gp.nc.2.child.0 = nil
gp.nc.2.child.1 = nil
```

### 5.2.2.1 Defining `GPNodes`

Let's imagine that we're trying to create trees that consist of the following tree nodes (which we'll create later): `ec.app.myapp.X`, `ec.app.myapp.Y`, `ec.app.myapp.Mul`, `ec.app.myapp.Sub`, `ec.app.myapp.Sin`. These nodes take 0, 0, 2, 2, and 1 children respectively and have no special types (we'll use `nil`). We could add them to the function set like this:

```
gp.fs.0 = ec.gp.GPFunctionSet
gp.fs.0.size = 5
gp.fs.0.func.0 = ec.app.myapp.X
gp.fs.0.func.0.nc = nc0
gp.fs.0.func.1 = ec.app.myapp.Y
gp.fs.0.func.1.nc = nc0
gp.fs.0.func.2 = ec.app.myapp.Mul
gp.fs.0.func.2.nc = nc2
gp.fs.0.func.3 = ec.app.myapp.Sub
gp.fs.0.func.3.nc = nc2
gp.fs.0.func.4 = ec.app.myapp.Sin
gp.fs.0.func.4.nc = nc1
```

Notice that we don't state the number of children or the types explicitly: instead we state them implicitly by assigning the appropriate `GPNodeConstraints` object.

### 5.2.3 Defining the Representation, Problem, and Statistics

GP is more complex than most other optimization procedures because of its representation. When you create a GP problem, you have two primary tasks:

- Create the `GPNodes` with which a `GPIndividual` may be constructed
- Create a `Problem` which tests the `GPIndividual`

Let's start with the first one. As an example, we'll build a simple Symbolic Regression example on two variables,  $X$  and  $Y$ . The GP tree can have `GPNodes` which subtract, multiply, and perform sine, just as was done earlier.<sup>4</sup> This means we'll need two terminals ( $X$  and  $Y$ ), and three nonterminals (subtract and multiply, each of arity 2—two children each—and cosine, with arity 1).

Recall from Section 5.2.2 that our function set would look like this:

```
gp.fs.0 = ec.gp.GPFunctionSet
gp.fs.0.size = 5
gp.fs.0.func.0 = ec.app.myapp.X
gp.fs.0.func.0.nc = nc0
gp.fs.0.func.1 = ec.app.myapp.Y
gp.fs.0.func.1.nc = nc0
gp.fs.0.func.2 = ec.app.myapp.Mul
gp.fs.0.func.2.nc = nc2
gp.fs.0.func.3 = ec.app.myapp.Sub
gp.fs.0.func.3.nc = nc2
gp.fs.0.func.4 = ec.app.myapp.Sin
gp.fs.0.func.4.nc = nc1
```

We need to make each of these classes. Each of these are `GPNodes` with a single crucial method overridden:

```
public abstract void eval(EvolutionState state, int thread, GPData input,
                        ADFStack stack, GPIndividual individual, Problem problem);
```

This method is called when the `GPNode` is being executed in the course of executing the tree. Execution proceeds depth-first like the evaluation of a standard parse tree. For example, in order to compute the expression  $\sin(x - y)$  (shown in GP form in Figure 5.3) we will call `eval(...)` on the `Sin` object, which will in turn call `eval(...)` on the `Sub` object. This will then call `eval(...)` on the  $X$  object, then on the  $Y$  object.  $X$  and  $Y$  will return their values. The `Sub` object will then subtract them and return the result, and finally `Sin` will return the sine of that.

Execution doesn't have to be just in terms of calling `eval(...)` on children, processing the results, and returning a final value. In fact, for some problems the return value may not matter at all, but simply which nodes are executed. In this case, likely the nodes themselves are doing things via side effects: moving a robot around, for example. Execution could also be tentative: an "if" node might evaluate one or another of its children and leave it at that. Execution could also be repetitive: you might make a "while" node which repeatedly evaluates a child until some test is true. Basically you can execute these nodes any way that might appear in a regular programming language.

---

<sup>4</sup>Ridiculously limited, but what did you expect? This is a demonstration example!

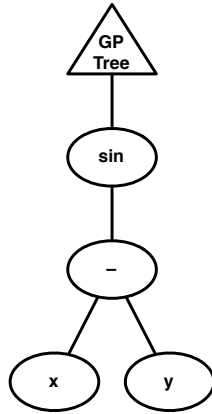


Figure 5.3 A simple GP tree representing the mathematical expression  $\sin(x - y)$ .

### 5.2.3.1 GPData

The `eval(...)` method has several arguments, only two of which should be nonobvious: `ec.gp.ADFStack` and `ec.gp.GPData`. We will discuss `ADFSStack` later in Section 5.2.10. The `GPData` object is a simple data object passed around amongst your `GPNodes` when they execute one another. It's your opportunity to pass data from node to node. In the example above, it's how the values are passed from the children to their parents: for example, it's how the `Sub` node returns its value to the `Sin` node. It's also possible that the parent needs to pass data to its child: and the `GPData` object can be used like that as well.

Typically a single `GPData` object is created and handed to the `GPNodes`, and then they hand it to one another during execution, reusing it. This avoids making lots of clones of a `GPData` object during execution.

Our `GPData` object needs to hold the return value from a child. If you are holding a simple piece of data (like a double or an int) also just need to implement a single method, `copyTo(...)`, which copies the data from your `GPData` object into another, then returns it:

```
public abstract GPData copyTo(GPData other);
```

In this case, it's simple:

```
package ec.app.myapp;
import ec.gp.*;
public class MyData extends GPData
{
    public double val;
    public void copyTo(GPData other)
    { ((MyData)other).val = val; return other; }
}
```

Now it might be the case that you need to hold more complex data. For example, what if you had an array of doubles? In this case you'd need to either clone or copy the data during the `copyTo(...)` operation. Additionally, `GPData` is a Prototype and so it needs to implement the `clone()` method as a deep clone. The default implementation just does a light clone. But with your array of doubles, you'd need to clone that. Altogether you might have something like this:

```

package ec.app.myapp;
import ec.gp.*;
public class MyData extends GPData
{
    public double[15] val = new double[15];
    public void copyTo(GPData other)
    {
        System.arraycopy(val, 0, ((MyData)other).val, 0, val.length);
        return other;
    }

    public Object clone()
    {
        MyData other = (MyData)(super.clone());
        other.val = (double[])(val.clone());
        return other;
    }
}

```

The important thing to note is that when you perform `copyTo(...)` or `clone()`, the resulting other object should not be sharing any data in common with you except constant (immutable, read-only) data. Why the two methods? Certainly most things `copyTo(...)` performs could be done with `clone()`. The reason for `copyTo(...)` is entirely for efficiency when using Automatically Defined Functions (Section 5.2.10). Perhaps in the future we might obviate the need for its use.

### 5.2.3.2 KozaFitness

You can use any fitness you like for GP. But it's common to use a particular fitness function popularized by John Koza [3]. This fitness object contains a *standardized fitness* in which 0 is the ideal result and Infinity is worse than the worst possible result. Note that this isn't yet the correct fitness according to the `ec.Fitness`. Instead, when asked for fitness, the function converts this to an *adjusted fitness* in which 1 is the ideal result and 0 is worse than the worst possible result, using the function  $\frac{1}{1 + \text{standardized}}$ . The adjusted fitness makes this a valid `Fitness` subclass. The GP fitness also has an auxiliary variable, `hits`, which originally was meant to indicate how many optimal subsolutions were discovered: it's printed out in the statistics and used for nothing else; use it as you like. This fitness is set as:

```
pop.subpop.0.species.fitness = ec.gp.koza.KozaFitness
```

The standard fitness-setting function for this class is:

```
public final void setStandardizedFitness(EvolutionState state, float value);
```

You can get (or set) the hits as:

```
int hits = myKozaFitness.hits;
```

### 5.2.3.3 GPPProblem

With all that out of our hair, let's construct the Problem. Let's attempt to create a GP tree which closely matches a set of data we've created: we'll generate the data from the function  $z = \sin(x \times y) - \sin(x) - x \times y$ , in the range  $[0, 1)$  for both  $x$  and  $y$ . What we'll do is define  $n$   $\langle x, y \rangle$  data points up front, then evaluate our individuals. For each data point, we'll set some global variables accessible by the X and Y GPNodes. The

individual will return a value, which we'll compare against the expected  $z$  result. The fitness will be the sum of squared differences.

```
package ec.app.myapp;
import ec.gp.*;
import ec.simple.*;
import ec.*;
import ec.gp.koza.*;
public class MyProblem extends GPPProblem implements SimpleProblemForm {
    final static int N = 20;
    int current;
    double[] Xs = new double[N]; // will be pointer-copied in clone(), which is okay
    double[] Ys = new double[N]; // likewise
    double[] Zs = new double[N]; // likewise
    MyData data;

    public void setup(EvolutionState state, Parameter base) {
        super.setup(state, base);

        // generate N random <x, y, z = f(x,y)> tuples
        for(int i = 0; i < N; i++) {
            double x, y;
            Xs[i] = x = state.random[0].nextDouble();
            Ys[i] = y = state.random[0].nextDouble();
            Zs[i] = Math.sin(x * y) - Math.sin(x) - x * y;
        }

        // create a GPData object
        data = (MyData) state.parameters.getInstanceForParameterEq(base.push(P_DATA),
            null, MyData.class);
        data.setup(state, base.push(P_DATA));
    }

    public Object clone() // we'll clone the GPData so each thread has its own copy
    {
        MyProblem obj = (MyProblem)(super.clone());
        obj.data = (MyData)(data.clone());
        return obj;
    }

    public void evaluate(final EvolutionState state, Individual ind,
        int subpopulation, int threadnum) {
        if (!ind.evaluated) { // don't bother reevaluating
            double sum = 0.0;

            // for each tuple, evaluate the individual. For good measure reset
            // the GPData first, though in this example it's not necessary
            data.val = 0;
            for(current = 0; current < N; current++) // note: an instance variable
                ((GPIndividual)ind).trees[0].child.eval(
                    state, threadnum, data, stack, ((GPIndividual)ind),this);
            sum += (data.val * data.val);

            // set the fitness and the evaluated flag
            KozaFitness f = (KozaFitness)(ind.fitness);
            f.setStandardizedFitness(state, sum);
            f.hits = 0; // don't bother using this
        }
    }
}
```

```

        ind.evaluated = true;
    }
}

```

#### 5.2.3.4 GPNode Subclasses

Now let's implement our five GPNode subclasses. Each will implement `toString()` to print out the node name, and also `eval(...)` discussed earlier. It's also common to implement the method `checkConstraints()` to do a final sanity-check on the node (whether it has the right number of children, etc.) but it's not necessary, and we'll omit it here.

First the terminals:

```

package ec.app.myapp;
import ec.*;
import ec.gp.*;
public class X extends GPNode {
    public String toString() { return "x" };
    public void eval(EvolutionState state, int thread, GPData input,
        ADFStack stack, GPIndividual individual, GPPProblem problem) {
        MyData data = (MyData) input;
        MyProblem prob = (MyProblem) problem;

        data.val = problem.Xs[problem.current]; // return current X value to parent
    }
}

```

... and...

```

package ec.app.myapp;
import ec.*;
import ec.gp.*;
public class Y extends GPNode {
    public String toString() { return "y" };
    public void eval(EvolutionState state, int thread, GPData input,
        ADFStack stack, GPIndividual individual, GPPProblem problem) {
        MyData data = (MyData) input;
        MyProblem prob = (MyProblem) problem;

        data.val = problem.Ys[problem.current]; // return current Y value to parent
    }
}

```

Next the Sine:

```

package ec.app.myapp;
import ec.*;
import ec.gp.*;
public class Sin extends GPNode {
    public String toString() { return "sin" };
    public void eval(EvolutionState state, int thread, GPData input,
        ADFStack stack, GPIndividual individual, GPPProblem problem) {
        MyData data = (MyData) input;
        MyProblem prob = (MyProblem) problem;

        children[0].eval(state, thread, data, stack, individual, prob);
        data.val = Math.sin(data.val);
    }
}

```

Next the Multiply and Subtract:

```

package ec.app.myapp;
import ec.*;
import ec.gp.*;
public class Mul extends GPNode {
    public String toString() { return "*" };
    public void eval(EvolutionState state, int thread, GPData input,
        ADFStack stack, GPIndividual individual, GPPProblem problem) {
        MyData data = (MyData) input;
        MyProblem prob = (MyProblem) problem;

        children[0].eval(state, thread, data, stack, individual, prob);
        double val1 = data.val;

        children[1].eval(state, thread, data, stack, individual, prob);
        data.val = val1 * data.val;
    }
}

```

```

package ec.app.myapp;
import ec.*;
import ec.gp.*;
public class Sub extends GPNode {
    public String toString() { return "-" };
    public void eval(EvolutionState state, int thread, GPData input,
        ADFStack stack, GPIndividual individual, GPPProblem problem) {
        MyData data = (MyData) input;
        MyProblem prob = (MyProblem) problem;

        children[0].eval(state, thread, data, stack, individual, prob);
        double val1 = data.val;

        children[1].eval(state, thread, data, stack, individual, prob);
        data.val = val1 - data.val;
    }
}

```

### 5.2.3.5 Statistics

The default statistics class for GP is SimpleStatistics. However, the GP package has a Statistics subclass designed for doing the same basic stuff as SimpleShortStatistics (Section 3.7.2), but with some extra GP-specific tree statistics. You can turn it on like this:

```
stat = ec.gp.koza.KozaShortStatistics
```

This statistics object has all the basic features of SimpleShortStatistics, including the do-time, do-size, do-subpops, and modulus parameters. Additionally it adds a new parameter which you can turn on like this:

```
stat.child.0.do-depth = true
```

This parameter enables tree depth information.

Beyond depth options, the full form of output differs from SimpleShortStatistics in that it reports per-tree information as well, like this:

1. The generation number
2. (If do-time is true) How long initialization took in milliseconds, or how long the previous generation took to breed to form this generation
3. (If do-time is true) How long evaluation took in milliseconds this generation
4. *Once for each subpopulation...*
  - (a) (If do-depth is true) Output of the form [a b c ...], representing the average depth of each tree *a*, *b*, *c*, etc. of an individual this generation for this subpopulation
  - (b) (If do-size is true) Output of the form [a b c ...], representing the average size of each tree *a*, *b*, *c*, etc. of an individual this generation for this subpopulation
  - (c) (If do-size is true) The average size of an individual this generation for this subpopulation
  - (d) (If do-size is true) The average size of an individual so far in the run for this subpopulation
  - (e) (If do-size is true) The size of the best individual this generation for this subpopulation
  - (f) (If do-size is true) The size of the best individual so far in the run for this subpopulation
  - (g) The mean fitness of the subpopulation for this generation
  - (h) The best fitness of the subpopulation for this generation
  - (i) The best fitness of the subpopulation so far in the run
5. (If do-depth is true) Output of the form [a b c ...], representing the average depth of each tree *a*, *b*, *c*, etc. of an individual this generation
6. (If do-size is true) Output of the form [a b c ...], representing the average size of each tree *a*, *b*, *c*, etc. of an individual this generation
7. (If do-size is true) The average size of an individual this generation
8. (If do-size is true) The average size of an individual so far in the run
9. (If do-size is true) The size of the best individual this generation
10. (If do-size is true) The size of the best individual so far in the run
11. The mean fitness of the entire population for this generation
12. The best fitness of the entire population for this generation
13. The best fitness of the entire population so far in the run



## 5.2.4 Initialization

To use GP we'll need to define the initializer as a subclass of `ec.gp.GPInitializer`:

```
init = ec.gp.GPInitializer
```

ECJ has traditionally followed the *lil-gp* default for disallowing duplicates in the initial Population: if a duplicate is created, ECJ will try 100 times to create another non-duplicate Individual in its stead. If this fails, the last duplicate created will be allowed. We say this in the standard way:

```
pop.subpop.0.duplicate-retries = 100
```

To create trees, ECJ relies on a tree-creation algorithm in the form of an `ec.gp.GPNodeBuilder`, part of the `GPTreeConstraints` object. The `GPNodeBuilder` for `GPTreeConstraints 0` is specified like this:

```
gp.tc.0.init = ec.gp.koza.HalfBuilder
```

ECJ provides quite a number of node builders in the `ec.gp.koza` and `ec.gp.build` packages. You request a tree with the following function:

```
public abstract GPNode newRootedTree(EvolutionState state, GPType type,  
                                     int thread, GPNodeParent parent, GPFunctionSet set,  
                                     int argposition, int requestedSize);
```

This method builds a tree of `GPNodes` whose root return type is compatible with `type`, attached to the given `GPNodeParent`, at position `argposition`, and built from clones of `GPNodes` in the function set `set`. The root node is returned. Several `GPNodeBuilders` also produce the tree of the requested size: others ignore this function. You can also ask the `GPNodeBuilder` to pick its own tree size from a distribution specified by the user in parameters, by passing `ec.gp.GPNodeBuilder.NOSIZEGIVEN` for the size (this is the usual thing done by most initialization procedures).

If you are using a `GPNodeBuilder` which generates trees of a certain size, and `ec.gp.GPNodeBuilder.NOSIZEGIVEN` is used (as usual), then you can specify a distribution of sizes in two ways. First, you can have the `GPNodeBuilder` pick a size uniformly from among a minimum and maximum size, for example:

```
gp.tc.0.init.min-size = 10  
gp.tc.0.init.max-size = 20
```

Alternatively you can specify the distribution of sizes manually. To stipulate probabilities sizes for 1, 2, 3, 4, and 5, you'd say:

```
gp.tc.0.init.num-sizes = 5  
gp.tc.0.init.size.0 = 0.2  
gp.tc.0.init.size.1 = 0.1  
gp.tc.0.init.size.2 = 0.2  
gp.tc.0.init.size.3 = 0.25  
gp.tc.0.init.size.4 = 0.25
```

ECJ has a whole bunch of `GPNodeBuilder` algorithms available to you. I wrote a shoot-out paper describing and comparing nearly all of these algorithms [8]. Here is the run-down:

- `ec.gp.koza.FullBuilder` generates full trees using Koza's FULL algorithm. You cannot request a size. It requires a minimum and maximum depth, for example:

```
gp.tc.0.init = ec.gp.koza.FullBuilder
gp.tc.0.init.min-depth = 2
gp.tc.0.init.max-depth = 6
```

Alternatively:

```
gp.koza.full.min-depth = 2
gp.koza.full.max-depth = 6
```

- `ec.gp.koza.GrowBuilder` generates arbitrary trees depth-first using Koza's GROW algorithm. You cannot request a size. It requires a minimum and maximum depth, for example:

```
gp.tc.0.init = ec.gp.koza.GrowBuilder
gp.tc.0.init.min-depth = 2
gp.tc.0.init.max-depth = 6
```

Alternatively:

```
gp.koza.grow.min-depth = 2
gp.koza.grow.max-depth = 6
```

- `ec.gp.koza.HalfBuilder` generates arbitrary trees depth-first using Koza's RAMPED HALF-AND-HALF algorithm. You cannot request a size. This is nothing more than flipping a coin to decide whether to use GROW or FULL. `HalfBuilder` is the default builder for creating GP trees in ECJ, but it's not particularly good. It requires a minimum and maximum depth, and the probability of doing GROW, for example:

```
gp.tc.0.init = ec.gp.koza.HalfBuilder
gp.tc.0.init.min-depth = 2
gp.tc.0.init.max-depth = 6
gp.tc.0.init.growp = 0.5
```

Alternatively:

```
gp.koza.half.min-depth = 2
gp.koza.half.max-depth = 6
gp.koza.half.growp = 0.5
```

- `ec.gp.build.PTC1` is a modification of GROW which guarantees that trees will be generated with a given mean. You cannot request a size. Additionally, each terminal and nonterminal can specify its probability of being chosen from the function set as `PTC1` constructs the tree. `PTC1` requires an expected size and a maximum depth:

```
gp.tc.0.init = ec.gp.build.PTC1
gp.tc.0.init.expected-size = 10
gp.tc.0.init.max-depth = 6
```

Alternatively:

```
gp.build.ptc1.expected-size = 10
gp.build.ptc1.max-depth = 6
```

`PTC1` requires that its function sets adhere to the interface `ec.gp.build.PTCFunctionSetForm`. This interface contains three tables of probabilities for your `GPNodes` to be selected:

```

public float[] terminalProbabilities(int type);
public float[] nonterminalProbabilities(int type);
public float[] nonterminalSelectionProbabilities(int expectedTreeSize);

```

The first function returns, for a given GPType number, a distribution of desired selection probabilities for terminals of that type. The order of the terminals is the same as the following array in GPFunctionSet:

```

public GPNode[type][] terminals;

```

The second function returns, for a given GPType number, a distribution of desired selection probabilities for nonterminals of that type. The order of the nonterminals is the same as the following array in GPFunctionSet:

```

public GPNode[type][] nonterminals;

```

The final function returns, for a given desired tree size, the probability that a nonterminal (of a given GPType return type) should be selected over a terminal of the same GPType. This is only used by PTC1, not PTC2 below.

You don't need to implement this interface: the `ec.gp.build.PTCFunctionSet` class does it for you:

```

gp.fs.size = 1
gp.fs.0 = ec.gp.build.PTCFunctionSet
gp.fs.0.name = f0

```

This function set computes all the above probabilities from user-specified probabilities as parameters. The probabilities are specified by each GPNodeConstraints object. Following the example we started in Section 5.2.2, we might state that the terminals X and Y (node constraints 0) should be picked with 0.5 probability each, and the nonterminals Mul, Sub (node constraints 2) and Cos (node constraints 1) should be picked with 0.3, 0.3, and 0.4 probability:

```

gp.nc.0.prob = 0.5
gp.nc.1.prob = 0.3
gp.nc.2.prob = 0.4

```

What if you wanted Mul and Sub to have different probabilities? You'd need to create different GPNodeConstraints. For example, we could create a new, separate GPNodeConstraints for Sub:

```

gp.nc.size = 4
gp.nc.3 = ec.gp.GPNodeConstraints
gp.nc.3.name = nc3
gp.nc.3.returns = nil
gp.nc.3.size = 2
gp.nc.3.child.0 = nil
gp.nc.3.child.1 = nil

```

Now we assign it to Sub:

```

gp.fs.0.func.3 = ec.app.myapp.Sub
gp.fs.0.func.3.nc = nc3

```

...and last change the probabilities of Sub and Mul to be different:

```
gp.nc.0.prob = 0.5
gp.nc.1.prob = 0.3
gp.nc.2.prob = 0.25
gp.nc.3.prob = 0.35
```

- `ec.gp.build.PTC2` generates trees near to a desired size (which you request) by picking randomly from the current outer edge in the tree and adding a node. When the tree is large enough, all the remaining edge slots are filled with terminals. Additionally, each terminal and nonterminal can specify its probability of being chosen from the function set as PTC2 constructs the tree. PTC2 requires a desired size and a maximum depth:

```
gp.tc.0.init = ec.gp.build.PTC2
gp.tc.0.init.expected-size = 10
gp.tc.0.init.max-depth = 6
```

Alternatively:

```
gp.build.ptc2.expected-size = 10
gp.build.ptc2.max-depth = 6
```

Like PTC1, PTC2 requires that function sets adhere to the `PTCFunctionSetForm` interface. Just use `PTCFunctionSet`.

- `ec.gp.build.RandomBranch` generates trees near to a desired size (which you request) using the `RANDOMBRANCH` algorithm. Beyond the size distributions, this algorithm has no additional parameters.
- `ec.gp.build.Uniform` generates trees near to a desired size (which you request) using the `UNIFORM` algorithm, which selects trees of any tree size. You can select sizes either using the user distribution, or according to the natural distribution of tree sizes. To do the second, you'd say:

```
gp.tc.0.init = ec.gp.build.Uniform
gp.tc.0.init.true-dist = true
```

Alternatively:

```
gp.breed.uniform.true-dist = true
```

**WARNING:** This algorithm is complex and I fear it may be suffering from bit-rot. I have been told it's not working properly any more but have not debugged it yet.

- `ec.gp.build.RandTree` (by Alexander Chircop) generates trees near to a desired size (which you request) using the `RAND.TREE` algorithm, which selects trees distributed uniformly using Dyck words. No extra parameters are needed beyond the tree size selection. **WARNING:** I suspect this algorithm may have some bugs.

## 5.2.5 Breeding

ECJ has a large number of breeding pipeline operators for GP trees. This includes the most common operators used in GP (`ec.gp.koza.Crossover`, `ec.gp.koza.Mutation`), and several more found in the `ec.gp.breed` package.

Pipelines generally pick a single `GPTree` in a given `GPindividual` in which to do mutation or crossover. In most cases you can lock down the specific `GPTree`, or let the pipeline choose it at random.

Once they've picked a `GPTree`, GP breeding operators often need to choose `GPNodes` in the tree in which to perform crossover, mutation, etc. To do this, they make use of a `ec.gp.GPNodeSelector`. A `GPNodeSelector` is a simple interface for picking nodes, consisting of the following methods:

```

public abstract void reset();
public abstract GPNode pickNode(EvolutionState s, int subpopulation,
                                int thread, GPIndividual ind, GPTree tree);

```

When a breeding pipeline needs to pick a node in a particular GPTree of a particular GPIndividual, it first will call the `reset()` to get the GPNodeSelector to ready itself, then it will call `pickNode(...)` to select a node. If the breeding pipeline needs another node in the same tree, it can call `pickNode(...)` again as many times as necessary.

The standard GPNodeSelector is `ec.gp.koza.KozaNodeSelector`, which picks certain kinds of nodes with different probabilities. The kinds of nodes you can state probabilities for are: the root, nonterminals, terminals, and all nodes. The most common settings are (here as default parameters):

```

gp.koza.ns.terminals = 0.1
gp.koza.ns.nonterminals = 0.9
gp.koza.ns.root = 0.0

```

This says to pick terminals 10% of the time, nonterminals 90% of the time, the root (specifically) 0% of the time and any arbitrary node 0% of the time. The arbitrary-node percentage is whatever is left over from the other three percentages. (The root could still be picked, since it's a nonterminal or a terminal—but it won't be *special*ly picked).

Why might the breeding pipeline need to call `pickNode(...)` repeatedly? Most likely because the chosen GPNode has type constraint problems. For example, in order to do crossover between the subtrees rooted by two GPNodes, the nodes need to be type-compatible with one another's parent nodes: otherwise the tree locations wouldn't be valid. Pipelines with these issues will try some  $n$  times to pick compatible nodes; if they fail all  $n$  times, the parents are returned rather than the generated children.

Here are the breeding pipelines that come with ECJ. In each case, let's presume that we're placing the pipeline as the root pipeline of Subpopulation 0, parameter-wise:

- `ec.gp.koza.CrossoverPipeline` performs standard subtree crossover: it requests a GPIndividual from each of its two sources; then a tree is selected from each GPIndividual, then a node is selected in each tree, and finally the two subtrees rooted by those nodes are swapped. CrossoverPipeline has several parameters. The first three:

```

pop.subpop.0.species.pipe.tries = 1
pop.subpop.0.species.pipe.maxdepth = 17
pop.subpop.0.species.pipe.toss = false

```

This tells CrossoverPipeline that children may not have a depth which exceeds 17 (the common value). The pipeline will try just one times to find type-valid and depth-legal crossover points before giving up and just returning the parents instead. This is the most common setting in Genetic Programming. If `toss=true` then only one child is returned—the other is thrown away. The default value for `toss` is `false`.

```

pop.subpop.0.species.pipe.tree.0 = 0
pop.subpop.0.species.pipe.tree.1 = 0

```

This tells CrossoverPipeline that it should pick GPNodes in GPTree 0 of each individual. If either of these parameters is missing entirely, then CrossoverPipeline will pick that tree at random. At any rate, the GPTrees chosen must have the same GPTreeConstraints. Finally we have:

```

pop.subpop.0.species.pipe.ns.0 = ec.gp.koza.KozaNodeSelector
pop.subpop.0.species.pipe.ns.1 = same

```

This states that the GPNodeSelector for both GPIndividuals should be a `KozaNodeSelector`. You can state them independently or node selector 1 can be `same`.

The default parameter base versions for all of these would be:

```
gp.koza.xover.tries = 1
gp.koza.xover.maxdepth = 17
gp.koza.xover.toss = false
gp.koza.xover.tree.0
gp.koza.xover.tree.1
gp.koza.xover.ns = ec.gp.koza.KozaNodeSelector
```

**Important note:** the default version of the parameter for node selectors is just `ns`. There's no `ns.0` or `ns.1`.

- `ec.gp.koza.MutationPipeline` performs standard subtree mutation: it requests a `GPIndividual` from a single source; then a tree is selected; then a node is selected in that tree; and finally the subtree rooted by that node is replaced in its entirety by a randomly-generated tree.

`MutationPipeline` has many parameters similar to `CrossoverPipeline`:

```
pop.subpop.0.species.pipe.tries = 1
pop.subpop.0.species.pipe.maxdepth = 17
pop.subpop.0.species.pipe.tree.0 = 0
pop.subpop.0.species.pipe.ns = ec.gp.koza.KozaNodeSelector
```

Note that the node selector is just `ns`, not `ns.0`.

The replacing subtree is generated using a `GPNodeBuilder`. The standard `GPNodeBuilder` is a `GrowBuilder` with the following default values:

```
gp.koza.grow.min-depth = 5
gp.koza.grow.max-depth = 5
```

These are strange default values, but that's common GP original settings. You stipulate the `GPNodeBuilder` as:

```
pop.subpop.0.species.pipe.build.0 = ec.gp.koza.GrowBuilder
```

Though `GrowBuilder` ignores size demands, if you replaced with another builder such as `PTC2`, you can also optionally stipulate that the replacing subtree must be about the same size as the original subtree. Here's the parameter:

```
pop.subpop.0.species.pipe.equal = true
```

The default setting is `false`.

The default parameter base versions for all of these would be:

```
gp.koza.mutate.tries = 1
gp.koza.mutate.maxdepth = 17
gp.koza.mutate.tree.0 = 0
gp.koza.mutate.pipe.ns = ec.gp.koza.KozaNodeSelector
gp.koza.mutate.build.0 = ec.gp.koza.GrowBuilder
gp.koza.mutate.equal = true
```

- `ec.gp.breed.InternalCrossoverPipeline` selects two `GPNodes` in the same `GPIndividual`, such that neither `GPNode` is in the subtree rooted by the other. The `GPNodes` may be in different `GPTrees`, or they may be in the same `GPTree`. It then swaps the two subtrees.

`InternalCrossoverPipeline`'s parameters are essentially identical to those in `CrossoverPipeline`. For example:

```

pop.subpop.0.species.pipe.tries = 1
pop.subpop.0.species.pipe.maxdepth = 17
pop.subpop.0.species.pipe.tree.0 = 0
pop.subpop.0.species.pipe.tree.1 = 0
pop.subpop.0.species.pipe.ns.0 = ec.gp.koza.KozaNodeSelector
pop.subpop.0.species.pipe.ns.1 = same

```

The default parameter base versions for all of these would be:

```

gp.breed.internal-xover.tries = 1
gp.breed.internal-xover.maxdepth = 17
gp.breed.internal-xover.toss = false
gp.breed.internal-xover.tree.0
gp.breed.internal-xover.tree.1
gp.breed.internal-xover.ns = ec.gp.koza.KozaNodeSelector

```

**Important note:** just as is the case for CrossoverPipeline, the default version of the parameter for node selectors is just `ns`. There's no `ns.0` or `ns.1`.

- `ec.gp.breed.MutatePromotePipeline` selects a `GPNode`, other than the root, and replaces its parent (and its parent's subtree) with the `GPNode` and its subtree. This was called the *PromoteNode* algorithm in [1] and is similar to the *Deletion* algorithm in [11].

`MutatePromotePipeline`'s parameters are pretty simple. Because its constraints are tighter, it doesn't use a `GPNodeSelector`: instead it searches among all nodes in the tree to find one which is type-compatible with its parent. thus its parameters are simply the number of times it tries before giving up, and returning the original tree. Like previous methods, if the tree parameter doesn't exist, a tree is picked at random (which is usually what'd you'd want anyway).

```

pop.subpop.0.species.pipe.tries = 1
pop.subpop.0.species.pipe.tree.0 = 0

```

The default parameter base versions:

```

gp.breed.mutate-promote.tries = 1
gp.breed.mutate-promote.tree.0

```

- `ec.gp.breed.MutateDemotePipeline` selects a `GPNode`, then replaces the node with a new nonterminal. The old Node becomes a child of the new node at a random argument location, and the remaining child slots are filled with terminals. This was called the *DemoteNode* algorithm in [1] and is similar to the *Insertion* algorithm in [11].

`MutatePromotePipeline` is similar to `MutatePromotePipeline`: it doesn't use a `GPNodeSelector`, and tries a certain number of times to find valid node points before giving up and returning the tree. parameters are pretty simple. Because its constraints are tighter, it doesn't use a `NodeSelector`: instead it searches among all nodes in the tree to find one which is type-compatible with its parent, and which wouldn't create a tree deeper than a maximum legal value. Like previous methods, if the tree parameter doesn't exist, a tree is picked at random (which is usually what'd you'd want anyway).

```

pop.subpop.0.species.pipe.tries = 1
pop.subpop.0.species.pipe.maxdepth = 17
pop.subpop.0.species.pipe.tree.0 = 0

```

The default parameter base versions:

```
gp.breed.mutate-demote.tries = 1
gp.breed.mutate-demote.maxdepth = 17
gp.breed.mutate-demote.tree.0 = 0
```

- `ec.gp.breed.MutateSwapPipeline` selects a `GPNode` with at least two children, then selects two children of that node such that each is type-compatible with the other. Then it swaps the two subtrees rooted by those children.

`MutateSwap`'s parameters are simple because it doesn't use a `GPNodeSelector` (the constraints are too complex). You simply specify the tree (or have one picked at random if none is specified) and the number of tries:

```
pop.subpop.0.species.pipe.tries = 1
pop.subpop.0.species.pipe.tree.0 = 0
```

The default parameter base versions for all of these would be:

```
gp.breed.mutate-swap.tries = 1
gp.breed.mutate-swap.tree.0
```

- `ec.gp.breed.MutateOneNodePipeline` selects a `GPNode`, then replaces that node with a different node of the same arity and type constraints. This was called the *OneNode* algorithm in [1].

`MutateOneNodePipeline` uses a `GPNodeSelector` to pick the node. You also specify the tree number: or if you don't specify anything, one will be picked at random (which is usually what'd you'd want).

```
pop.subpop.0.species.pipe.ns.0 = ec.gp.koza.KozaNodeSelector
pop.subpop.0.species.pipe.tree.0 = 0
```

The default parameter base versions:

```
gp.breed.mutate-one-node.ns.0 = ec.gp.koza.KozaNodeSelector
gp.breed.mutate-one-node.tree.0 = 0
```

- `ec.gp.breed.MutateAllNodesPipeline` selects a `GPNode`, then for *every* node in the subtree rooted by the `GPNode`, it replaces each node with a different node of the same arity and type constraints. This highly destructive operator was called the *AllNodes* algorithm in [1].

`MutateAllNodesPipeline` uses a `GPNodeSelector` to pick the `GPNode`. You also specify the tree number: or if you don't specify anything, one will be picked at random (which is usually what'd you'd want).

```
pop.subpop.0.species.pipe.ns.0 = ec.gp.koza.KozaNodeSelector
pop.subpop.0.species.pipe.tree.0 = 0
```

The default parameter base versions:

```
gp.breed.mutate-one-node.ns.0 = ec.gp.koza.KozaNodeSelector
gp.breed.mutate-one-node.tree.0 = 0
```

- `ec.gp.breed.RehangPipeline` is an oddball mutator of my own design meant to be highly destructive. It selects a nonterminal other than the root, and designates it the "new root". It then picks a child subtree of this new root, which is disconnected from its parent. The new root becomes the root of the tree. The original parent of the new root becomes the new root's child, filling the spot vacated by the disconnected subtree. The grandparent then fills the spot vacated by the parent, and so on, clear up to the root. Then finally the disconnected subtree fills remaining spot. Figure 5.4 shows this procedure. There are two parameters, as usual:



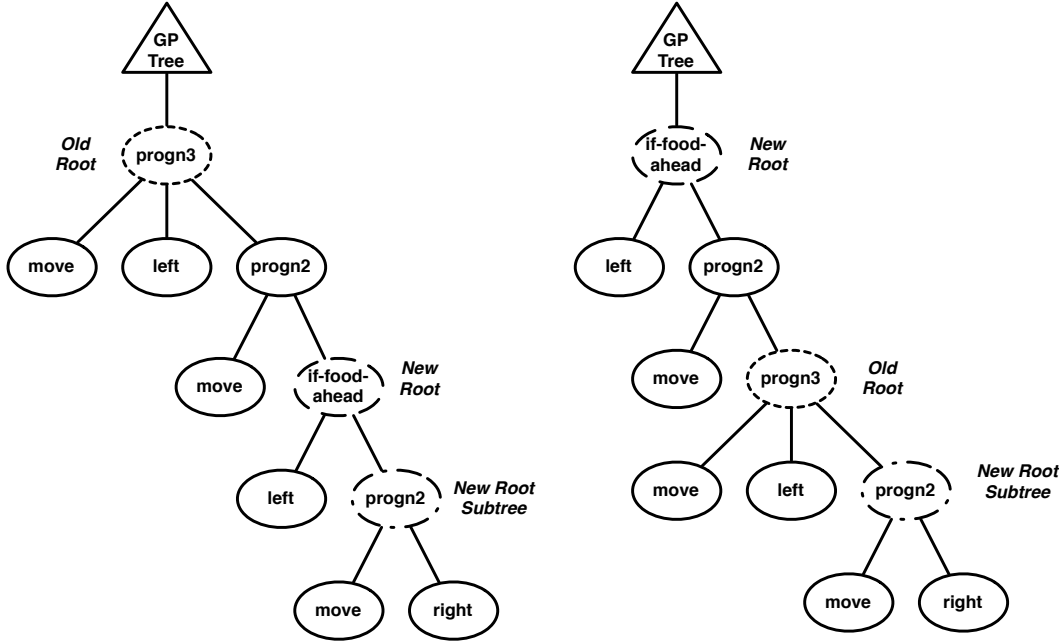


Figure 5.4 Rehangging a tree. A new root is chosen at random from among the nonterminals except for the original root. Then a subtree of that new root is chosen at random and disconnected. The tree is then rehung as shown: the parent of the new root becomes its child; the grandparent becomes the parent's child, and so on up to the root. The disconnected subtree then fills the remaining spot.

```
pop.subpop.0.species.pipe.tries = 1
pop.subpop.0.species.pipe.tree.0 = 0
```

The default parameter base versions for all of these would be:

```
gp.breed.rehang.tries = 1
gp.breed.rehang.tree.0
```

**Warning:** Because of the complexity of its rehangging process, RehangPipeline ignores all typing information.

- `ec.gp.breed.MutateERCPipeline` works similarly to the *Gaussian* algorithm in [1]. The algorithm picks a random node in a random tree in the `GPIndividual`, then for every Ephemeral Random Constant (ERC) in the subtree rooted by that node, it calls `mutateERC()` on that ERC. ERCs are discussed later in Section 5.2.9. As usual, the two common parameters:

```
pop.subpop.0.species.pipe.tries = 1
pop.subpop.0.species.pipe.tree.0 = 0
```

The default parameter base versions for all of these would be:

```
gp.breed.mutate-erc.tries = 1
gp.breed.mutate-erc.tree.0
```

If you wished to mutate the ERCs in the *entire tree*, you could set the node selector parameters like this:

```

gp.breed.mutate-erc.ns.0.terminals = 0.0
gp.breed.mutate-erc.ns.0.nonterminals = 0.0
gp.breed.mutate-erc.ns.0.root = 1.0

```

## 5.2.6 A Complete Example

Much of these initial parameters could have been entered simply by including the parameter file `ec/gp/koza/koza.params`. But we'll go through it in detail. First some basic generational parameters:

```

# Threads and Seeds
evalthreads = 1
breedthreads = 1
seed.0 = time

# Checkpointing
checkpoint = false
checkpoint-modulo = 1
prefix = ec

# The basic setup
state = ec.simple.SimpleEvolutionState
finish = ec.simple.SimpleFinisher
exch = ec.simple.SimpleExchanger
breed = ec.simple.SimpleBreeder
eval = ec.simple.SimpleEvaluator
stat = ec.simple.SimpleStatistics
pop = ec.Population
pop.subpops = 1
pop.subpops.0 = ec.Subpopulation
pop.subpop.0.duplicate-retries = 0
pop.subpop.0.size = 1024
breed.elite.0 = 0
stat.file = $out.stat
quit-on-run-complete = true

```

Genetic programming typically doesn't run very long, and (for the time being) require their own Initializer. We'll also use `KozaFitness`. Following the *lil-gp* example, we'll set the duplicate retries to 100:

```

init = ec.gp.GPInitializer
generations = 51
pop.subpop.0.species.fitness = ec.gp.koza.KozaFitness
pop.subpop.0.duplicate-retries = 100

```

For good measure, let's attach `KozaShortStatistics` to the statistics chain. This isn't standard in the `koza.params` file, but what the heck.

```

stat.num-children = 1
stat.child.0 = ec.gp.koza.KozaShortStatistics
stat.child.0.gather-full = true
stat.child.0.file = $out2.stat

```

Our initializer will work by using `HalfBuilder` to build trees. We define its parameters here:

```

# HalfBuilder
gp.koza.half.min-depth = 2
gp.koza.half.max-depth = 6
gp.koza.half.growp = 0.5

```

We begin by defining the tree constraints, node constraints, types, and function sets for the problem:

```

# Types
gp.type.a.size = 1
gp.type.a.0.name = nil
gp.type.s.size = 0

# Basic Function Set Parameters (more later)
gp.fs.size = 1
gp.fs.0 = ec.gp.GPFunctionSet
gp.fs.0.name = f0

# Tree Constraints
gp.tc.size = 1
gp.tc.0 = ec.gp.GPTreeConstraints
gp.tc.0.name = tc0
gp.tc.0.fset = f0
gp.tc.0.returns = nil
gp.tc.0.init = ec.gp.koza.HalfBuilder

# Node Constraints
gp.nc.size = 3

gp.nc.0 = ec.gp.GPNodeConstraints
gp.nc.0.name = nc0
gp.nc.0.returns = nil
gp.nc.0.size = 0

gp.nc.1 = ec.gp.GPNodeConstraints
gp.nc.1.name = nc1
gp.nc.1.returns = nil
gp.nc.1.size = 1
gp.nc.1.child.0 = nil

gp.nc.2 = ec.gp.GPNodeConstraints
gp.nc.2.name = nc2
gp.nc.2.returns = nil
gp.nc.2.size = 2
gp.nc.2.child.0 = nil
gp.nc.2.child.1 = nil

```

Now we define the GP elements of the Species and the Individual:

```

# Representation
pop.subpop.0.species = ec.gp.GPSpecies
pop.subpop.0.species.ind = ec.gp.GPIndividual
pop.subpop.0.species.ind.numtrees = 1
pop.subpop.0.species.ind.tree.0 = ec.gp.GPTree
pop.subpop.0.species.ind.tree.0.tc = tc0

```

Here's a basic GP breeding pipeline:

```

# Pipeline
pop.subpop.0.species.pipe = ec.breed.MultiBreedingPipeline
pop.subpop.0.species.pipe.generate-max = false
pop.subpop.0.species.pipe.num-sources = 2
pop.subpop.0.species.pipe.source.0 = ec.gp.koza.CrossoverPipeline
pop.subpop.0.species.pipe.source.0.prob = 0.9
pop.subpop.0.species.pipe.source.1 = ec.breed.ReproductionPipeline
pop.subpop.0.species.pipe.source.1.prob = 0.1

```

For no good reason, we'll define the selection methods, and various other parameters using the default parameter bases for CrossoverPipeline and ReproductionPipeline:

```

# Reproduction
breed.reproduce.source.0 = ec.select.TournamentSelection

# Crossover
gp.koza.xover.source.0 = ec.select.TournamentSelection
gp.koza.xover.source.1 = same
gp.koza.xover.ns.0 = ec.gp.koza.KozaNodeSelector
gp.koza.xover.ns.1 = same
gp.koza.xover.maxdepth = 17
gp.koza.xover.tries = 1

# Selection
select.tournament.size = 7

```

Since Crossover is using a node selector, let's define some parameters for that:

```

# Node Selectors
gp.koza.ns.terminals = 0.1
gp.koza.ns.nonterminals = 0.9
gp.koza.ns.root = 0.0

```

Let's presume that we have created the X, Y, Sin, Mul, and Sub methods described in Section 5.2.3. We'll now hook them up.

```

# Our Function Set
gp.fs.0 = ec.gp.GPFunctionSet
gp.fs.0.size = 5
gp.fs.0.func.0 = ec.app.myapp.X
gp.fs.0.func.0.nc = nc0
gp.fs.0.func.1 = ec.app.myapp.Y
gp.fs.0.func.1.nc = nc0
gp.fs.0.func.2 = ec.app.myapp.Mul
gp.fs.0.func.2.nc = nc2
gp.fs.0.func.3 = ec.app.myapp.Sub
gp.fs.0.func.3.nc = nc2
gp.fs.0.func.4 = ec.app.myapp.Sin
gp.fs.0.func.4.nc = nc1

```

Also we're using the MyProblem and MyData classes defined in Section 5.2.3 as our Problem. Even though we're not using ADFs (see Section 5.2.10), we need to define a few items here as well.

```
# Our Problem eval.problem = ec.app.myapp.MyProblem
eval.problem.data = ec.app.myapp.MyData
gp.problem.stack = ec.gp.ADFStack
gp.adf-stack.context = ec.gp.ADFContext
```

Phew! That was a lot of parameters. Thankfully nearly all of them are already defined for you in `ec/gp/koza/koza.params`.

## 5.2.7 GPNodes in Depth

GPNode has a gazillion utility methods to assist various crossover, mutation, statistics, and tree-building operators in their tasks of making, breaking, printing, reading, writing and examining trees of GPNodes. Let's look at some of them here, and divvy up the rest in later sections where they're more appropriate.

First, the two abstract methods which you *must* override:

```
public String toString();
public abstract void eval(EvolutionState state, int thread, GPData input,
                        ADFStack stack, GPIndividual individual, Problem problem);
```

The first method prints out the node in a human-readable fashion, with no whitespace. Except for rare cases such as Ephemeral Random Constants (Section 5.2.9), this should be a single simple symbol like "cos" or "if-food-ahead". The second method we introduced in Section 5.2.3, of course.

One method which is commonly overridden by example applications in ECJ, but it's not necessary for you to do so:

```
public void checkConstraints(EvolutionState state, int tree, GPIndividual
                           typicalIndividual, Parameter individualBase)
```

This method is called after the prototypical GPNode is loaded into a function set to give it an opportunity to double-check that everything is okay: that the GPNode has the right number of children, that it *has* children, that it has a parent, etc., and issue an error if not. The primary function of this method is to allow Automatically Defined Functions (Section 5.2.10) a chance to make sure that everything is working. You don't really need to override this method (if you do, be sure to call `super(...)`) but you'll see it often implemented in the examples.

Next come some methods which are usually overridden by Ephemeral Random Constants but rarely by any other kind of GPNode (the default implementation suffices):

### ec.gp.GPNode Methods

---

```
public String name()
```

Returns the name of the GPNode. By default, this calls `toString()`.

```
public int nodeHashCode()
```

Returns a hash code appropriate to your GPNode (hash by value, not by address).

```
public boolean nodeEquals(GPNode node)
```

Returns true if this method is identical to the given node.

```
public boolean nodeEquivalentTo(GPNode node)
```

Returns true if the two nodes are the same "kind" of node—usually meaning they could have been cloned from the same prototype node. The default form of this function returns true if the two nodes are the same class, have the same length child array, and have the same constraints. Often `nodeEquals(...)` and `nodeEquivalentTo(...)` may return the same thing, but in Ephemeral Random Constants, they often return different values. For example, two ERCs that are the same class, and have the same constraints, may hold different values (2.34 vs. 3.14 say). These ERCs would be equivalent to one another but not equal to one another. You'd rarely need to override this method.

```
public String toStringForHumans()
    Writes the node to a String in a fashion readable by humans. The default version simply calls toString.

public void resetNode(EvolutionState state, int thread)
    Randomizes the node. Ephemeral Random Constants randomize their internal values; other GPNodes typically
    do nothing.
```

---

Next, some test functions:

### ec.gp.GPNode Methods

---

```
public int atDepth()
    Returns the depth of the node (the root is at depth 0).

public int depth()
    Returns the depth of the subtree rooted by the node (terminals have a subtree depth of 1).

public GPNodeParent rootParent()
    Returns the parent of the root of the tree in which the GPNode resides. Though the method returns a GPNodePar-
    ent, this returned object should always be some kind of GPTree.

public boolean contains(GPNode subnode)
    Returns true subnode exists somewhere within the subtree rooted by the GPNode.

public pathLength(int nodesearch)
    Returns the sum of all paths from all nodes in the GPNode's subtree to the GPNode itself. The nodesearch parameter
    allows us to restrict which nodes have paths included in the sum: only leaf nodes (terminals), non-leaf nodes
    (nonterminal), or all nodes:

        public static final int GPNode.NODESEARCH_ALL;
        public static final int GPNode.NODESEARCH_TERMINALS;
        public static final int GPNode.NODESEARCH_NONTERMINALS;

public int numNodes(int nodesearch)
    Returns the number of nodes in the subtree rooted by GPNode. The nodesearch parameter allows us to restrict
    which nodes are included in the total, using the constants above.

public int meanDepth(int nodesearch)
    Returns the path length divided by the number of nodes. The nodesearch parameter allows us to restrict which
    nodes are included in the total, using the constants above.
```

---

Two methods permit even more flexibility in filtering exactly which GPNodes you want to consider, by using a `ec.gp.GPNodeGatherer` object. This object contains a single method which you may override:

```
public boolean test(GPNode thisNode);
```

In a subclass, override this method to return true if the node provided is one which you'd like to include. The default form simply returns true for all nodes. Armed with a `GPNodeGatherer` you've constructed, you can then call one of the two `GPNode` methods:

### ec.gp.GPNode Methods

---

```
public int numNodes(GPNodeGatherer gatherer)
    Returns the number of nodes (filtered by the GPNodeGatherer) in the subtree rooted by the GPNode.
```

`public int nodeInPosition(int p, GPNodeGatherer gatherer, int nodesearch)`  
returns the *p*th node in the subtree, filtered either by one of the following node search constants (note there are four here, whereas there were three in the previous methods):

```
public static final int GPNode.NODESEARCH_ALL;  
public static final int GPNode.NODESEARCH_TERMINALS;  
public static final int GPNode.NODESEARCH_NONTERMINALS;  
public static final int GPNode.NODESEARCH_CUSTOM;
```

In the case of the first three constants, the GPNodeGatherer should be null. The fourth constant indicates that we should use the GPNodeGatherer instead (and it must be provided).

---

And now we come to rigamarole which should look familiar to you if you've trudged through the Vector chapter.

### ec.gp.GPNode Methods

---

`public int printNodeForHumans(EvolutionState state, int log)`  
Prints a node to a log in a fashion readable by humans. You don't want to override this method: it calls `toStringForHumans()` by default — override that instead.

`public int printNode(EvolutionState state, int log)`  
Prints a node to a log in a fashion readable by humans and also parsable by `readNode(...)`. You don't want to override this method: it calls `toString()` by default — override that instead.

`public int printNode(EvolutionState state, PrintWriter writer)`  
Prints a node to a writer in a fashion readable by humans and also parsable by `readNode(...)`. You don't want to override this method: it calls `toString()` by default — override that instead.

`public String toStringForError()`  
Writes a node to a string in a fashion useful for error messages. The default writes out the name and the tree the node is in, which works fine.

`public GPNode readNode(DecodeReturn dret)`  
Generates a GPNode from the DecodeReturn via a light clone: children and parents are not produced. The default version clones the the node, then reads a string from the DecodeReturn. This string should match `toString()` exactly. If not, returns null to indicate an error. Otherwise returns the GPNode. This default implementation should be fine in most cases: though Ephemeral Random Constants (Section 5.2.9) require a different procedure.

`public void writeNode(EvolutionState state, DataOutput output) throws IOException`  
Writes the node, but not any of its children or parents, out to *output*.

`public void readNode(EvolutionState state, DataInput input) throws IOException`  
Reads a node from *input*. Children and parents are not produced.

---

Last are a host of different ways of cloning a GPNode or tree. In most cases the default implementations work just fine:

### ec.gp.GPNode Methods

---

`public GPNode lightClone()`  
Light-clones a GPNode, including its children array, but not any children or parents.

`public Object clone()`  
Deep-clones a GPNode, except for its parent. All children are cloned as well.

`public final GPNode cloneReplacing(GPNode newSubtree, GPNode oldSubtree)`  
Deep-clones a GPNode, except that, if found within its cloned subtree, *oldSubtree* is replaced with a deep-cloned version of *newSubtree*.

```

public final GPNode cloneReplacing(GPNode[] newSubtrees, GPNode[] oldSubtrees)
    Deep-clones a GPNode, except that, if found within its cloned subtree, each of the oldSubtrees is replaced with a
    clone of the corresponding member of newSubtrees.

public final GPNode cloneReplacingNoSubclone(GPNode newSubtree, GPNode oldSubtree)
    Deep-clones a GPNode, except that, if found within its cloned subtree, oldSubtree is replaced with newSubtree (not
    a clone of newSubtree).

public final GPNode cloneReplacingAtomic(GPNode newNode, GPNode oldNode)
    Deep-clones a GPNode, except that, if found within its cloned subtree, oldNode is replaced with newNode (not a
    clone of newNode).

public final GPNode cloneReplacingAtomic(GPNode[] newNodes, GPNode[] oldNodes)
    Deep-clones a GPNode, except that, if found within its cloned subtree, each of the oldNodes is replaced with the
    corresponding member of newNodes (not a clone).

public final void replaceWith(GPNode newNode)
    Replaces the GPNode with newNode right where it lives in its GPTree.

```

---

These are the primary public methods. There are plenty of other public methods, but they're largely used internally and you'll rarely need them.

## 5.2.8 GPTrees and GPIndividuals in Depth

Unlike GPNode, there's nothing in a GPTree that you have to override or modify. It's pretty rare to subclass GPTree, though it's perfectly reasonable to do so. But there are a number of methods you should be aware of, many of which are probably very familiar by now. First, let's cover the three non-familiar ones:

### ec.gp.GPTree Methods

---

```

public int treeNumber()
    Returns the position of the GPTree in its GPIndividual's trees[] array. This is an  $O(n)$  operation—it works by
    scanning through the array until it finds the GPTree. If the tree is not found (which would indicate an error), then
    GPTree.NO_TREENUM is returned.

public final void verify(EvolutionState state)
    An auxillary debugging method which verifies many features of the structure of the GPTree and all of its GPNodes.
    This method isn't called by ECJ but has proven useful in determining errors in GPTree construction by various
    tree building or breeding algorithms.

public void buildTree(EvolutionState state, int thread)
    Builds a tree and attaches it to the GPTree, displacing the original, using the tree-generation algorithm defined for
    its GPTreeConstraints. No specific tree size is requested.

```

---

Next come cloning and tests for removing duplicates:

### ec.gp.GPTree Methods

---

```

public boolean treeEquals(GPTree tree)
    Returns true if the GPNodes which make up the GPTree are structured the same and equal in value to one another.
    Override this to provide more equality if necessary.

public int treeHashCode()
    Returns a hash code generated for the structure and makeup of the GPNodes in the GPTree. Override this to add
    additional hash information.

```



```
public GPTree lightClone()
    Performs a light clone on the GPTree: the GPNodes are not cloned but are instead the pointer to the root is simply copied.

public Object clone()
    Performs a deep clone on the GPTree, including all of its GPNodes but not its parent GPIndividual.
```

---

Last, the standard methods for printing and reading:

### ec.gp.GPTree Methods

---

```
public void printTreeForHumans(EvolutionState state, int log)
    Prints the tree in a human-readable fashion to log.

public void printTree(EvolutionState state, int log)
    Prints to log the tree in a fashion readable both by humans and also by readTree(..., LineNumberReader). By default this uses the Code package (Section 2.2.2).

public void printTree(EvolutionState state, PrintWriter writer)
    Prints to writer the tree in a fashion readable both by humans and also by readTree(..., LineNumberReader). By default this uses the Code package.

public void readTree(EvolutionState state, LineNumberReader reader) throws IOException
    Reads a tree produced by printTree(...). By default, this uses the Code package (Section 2.2.2).

public void writeTree(EvolutionState state, DataOutput output) throws IOException
    Writes a tree to output in binary fashion such that it can be read by readTree(..., DataInput).

public void readTree(EvolutionState state, DataInput input) throws IOException
    Reads a tree from input in binary fashion that had been written by writeTree(...).
```

---

#### 5.2.8.1 Pretty-Printing Trees

GPTrees have a particular gizmo that's not well known but is quite nice: you can print out GPTrees in one of (at present) four styles:

- Lisp (the default):

```
(* (+ x (- (% x x) (cos x))) (exp x))
```

- A style easily converted to C, C++, Java, or C#:

```
(x + ((x % x) - cos(x))) * exp(x)
```

To print out trees this way you'd use a parameter along these lines (notice the lower-case "c"):

```
pop.subpop.0.species.ind.tree.0.print-style = c
```

...or using the default parameter base:

```
gp.tree.print-style = c
```

Printing in C-style has two options. First, by default ECJ prints out two-child GPNodes as if they were operators "b a c" rather than as "a(b, c)". This is what's being done above. But if you're not using mathematical operators and would prefer to see 2-child GPNodes as functions, you can do it like this:

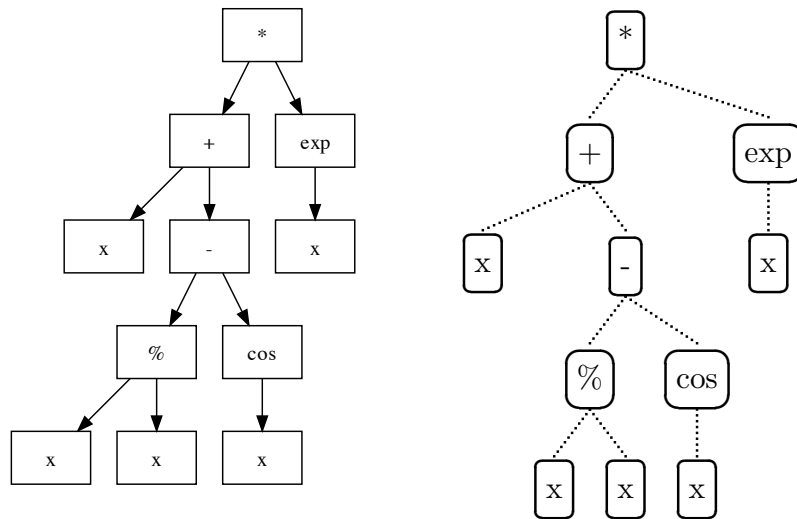


Figure 5.5 Auto-generated trees: in ".dot" format (left) and in LaTeX format (right)

```
pop.subpop.0.species.ind.tree.0.c-operators = false
```

...or using the default parameter base:

```
gp.tree.c-operators = false
```

This results in the following:

```
*(+(x, -(%(x, x), cos(x))), exp(x))
```

This doesn't seem useful for the example here (Symbolic Regression) but for other problems it's probably the right thing to do, particularly if all the GPNodes aren't operators. Additionally, by default ECJ prints out zero-child GPNodes as constants, as in "a", rather than as zero-argument functions, as in "a()". If you'd prefer zero-argument functions, you might say:

```
pop.subpop.0.species.ind.tree.0.c-variables = false
```

...or using the default parameter base:

```
gp.tree.c-variables = false
```

This results in the following:

```
x() + ((x() % x()) - cos(x())) * exp(x())
```

Again, whether this will be useful to you is based on exactly what kind of problem you're emitting. ECJ does not at present have support for converting if-statements (such as the "if-food-ahead" node in Artificial Ant) into a brace format appropriate to C. But hopefully these options will help you get most of the ugly parsing work out the way.

- .dot format: used by GraphViz to produce high-quality trees and graphs. The code below produces the tree shown at left in Figure 5.5:

```

digraph g {
node [shape=rectangle];
n[label = "*"];
n0[label = "+"];
n00[label = "x"];
n0 -> n00;
n01[label = "-"];
n010[label = "%"];
n0100[label = "x"];
n010 -> n0100;
n0101[label = "x"];
n010 -> n0101;
n01 -> n010;
n011[label = "cos"];
n0110[label = "x"];
n011 -> n0110;
n01 -> n011;
n0 -> n01;
n -> n0;
n1[label = "exp"];
n10[label = "x"];
n1 -> n10;
n -> n1;
}

```

To generate trees in .dot format, you'd say:

```
pop.subpop.0.species.ind.tree.0.print-style = dot
```

...or using the default parameter base:

```
gp.tree.print-style = dot
```

- $\text{\LaTeX}$  format, which emits the following code:

```

\begin{bundle}{\gpbox{*}}\chunk{\begin{bundle}{\gpbox{+}}\chunk{\gpbox{x}}
\chunk{\begin{bundle}{\gpbox{-}}\chunk{\begin{bundle}{\gpbox{%}}
\chunk{\gpbox{x}}\chunk{\gpbox{x}}\end{bundle}}\chunk{\begin{bundle}
{\gpbox{cos}}\chunk{\gpbox{x}}\end{bundle}}\end{bundle}}\end{bundle}}
\chunk{\begin{bundle}{\gpbox{exp}}\chunk{\gpbox{x}}\end{bundle}}\end{bundle}

```

To generate trees in  $\text{\LaTeX}$  format, you'd say:

```
pop.subpop.0.species.ind.tree.0.print-style = latex
```

...or using the default parameter base:

```
gp.tree.print-style = latex
```

This code works with the  $\text{\LaTeX}$  ecltree and fancybox packages to produce a tree. Note that you'll have to replace the “%” with “\%” to make it legal  $\text{\LaTeX}$ . The code works with the following boilerplate to produce the tree shown at right in Figure 5.5.

```

\documentclass[]{article}
\usepackage{epic}      % required by ecltree and fancybox packages
\usepackage{ecltree}   % to draw the GP trees
\usepackage{fancybox}  % required by \Ovalbox

\begin{document}

% minimum distance between nodes on the same line
\setlength{\GapWidth}{1em}

% draw with a thick dashed line, very nice looking
\thicklines \drawwith{\dottedline{2}}

% draw an oval and center it with the rule. You may want to fool with the
% rule values, though these seem to work quite well for me. If you make the
% rule smaller than the text height, then the GP nodes may not line up with
% each other horizontally quite right, so watch out.
\newcommand{\gpbox}[1]{\Ovalbox{#1}\rule[-.7ex]{0ex}{2.7ex}}

% And now the code
\begin{bundle}{\gpbox{*}}\chunk{\begin{bundle}{\gpbox{+}}\chunk{\gpbox{x}}
\chunk{\begin{bundle}{\gpbox{-}}\chunk{\begin{bundle}{\gpbox{\%}}
\chunk{\gpbox{x}}\chunk{\gpbox{x}}\end{bundle}}\chunk{\begin{bundle}
{\gpbox{cos}}\chunk{\gpbox{x}}\end{bundle}}\end{bundle}}\end{bundle}}
\chunk{\begin{bundle}{\gpbox{exp}}\chunk{\gpbox{x}}\end{bundle}}\end{bundle}

% Finally end the document
\end{document}

```

### 5.2.8.2 GPIndividuals

Okay, there's not much in-depth here. GPIndividual implements all the standard Individual methods discussed in Section 3.2. Note that the distanceTo(...) method is *not* implemented. Two methods you might want to be aware of:

#### ec.gp.GPIndividual Methods

---

```
public final void verify(EvolutionState state)
```

An auxillary debugging method which verifies many features of the structure of the GPIndividual and ll of its GPTrees (and their GPNodes). This method isn't called by ECJ but has proven useful in determining errors in GPtree construction by various tree building or breeding algorithms.

```
public long size()
```

By default, returns the number of nodes in all the trees held by the GPIndividual.

---

### 5.2.9 Ephemeral Random Constants

An **Ephemeral Random Constant** or ERC [3] is a special GPNode, usually a terminal, which represents a constant such as 3.14159 or true or 724 or complex:  $3.24 + 5i$  or "aegu". Usually ERCs are used to add constants to programs which rely on math (such as Symbolic Regression).

An ERC needs to be able to do three things:

- Set itself to a random value when first created.
- Mutate to a new value when asked to do so.

- Stay fixed at that value all other times (as a constant).

In ECJ, ERCs are usually subclasses of `ec.gp.ERC`, an abstract superclass which provides basic functionality. For example, if we were building a subclass of `ERC` which represents a floating point numerical constant, we might add a single instance variable to hold its value:

```
package ec.app.myapp;
import ec.gp.*;
public class MyERC extends ERC {
    public double value;
    // other methods go here...
}
```

We'll also probably need to implement some or most of the following methods to modify, read, write, or compare the `ERC`.

```
public String name();
public boolean nodeEquals(GPNode node);
public int nodeHashCode();
public void resetNode(EvolutionState state, int thread);
public void mutateERC(EvolutionState state, int thread);
public String toStringForHumans();
public String encode();
public boolean decode(DecodeReturn ret);
public void readNode(EvolutionState state, DataInput input) throws IOException
public void writeNode(EvolutionState state, DataOutput output)
                                                                    throws IOException
public abstract void eval(EvolutionState state, int thread, GPData input,
                        ADFStack stack, GPIndividual individual, Problem problem);
```

Let's go through these in turn:

```
public String name();
```

This is not the default `GPNode` implementation. When an `ERC` prints itself out in various ways, it writes its name, then its value. By default the name is simply "ERC", and if you have only one `ERC` type, you don't have to override it. For example, an `ERC` printed out might be `ERC[3.14159]`. However if you have more than one class of `ERCs`, holding different kinds of values, you'll want to distinguish them both for humans and for ECJ to read back in again. To do this, override `name()` to return a unique symbol for each kind of `ERC`. For example, you might just return "ERC1" versus "ERC2", resulting in `ERC1[3.14159]` versus `ERC2[921]`.

```
public boolean nodeEquals(GPNode node);
public int nodeHashCode();
```

Override the first method to test for equality with the second node: it's the same kind of `ERC`, the same class, has the same values, etc. Override the second method to provide a hash code for the `ERC` based on its type and the values it contains. By default you can avoid implementing this second method, and just implement the `encode()` method, discussed later. The default version of `nodeHashCode()` calls `encode()` and then hashes the `String`.

```
public void resetNode(EvolutionState state, int thread);
public void mutateERC(EvolutionState state, int thread);
```

Override the first method to entirely randomize the value of the `ERC`. Override the second method to mutate the value of the `ERC` when called to do so by the `MutateERCPipeline`. By default, `mutateERC(...)` just

calls `resetNode(...)`, which is probably not what you want. Instead the mutation will likely need to be a small deviation from the current value.

```
public String toStringForHumans();
public String encode();
public boolean decode(DecodeReturn ret);
public void readNode(EvolutionState state, DataInput input) throws IOException
public void writeNode(EvolutionState state, DataOutput output) throws IOException
```

As usual, override `toStringForHumans()` to provide a pretty version of the ERC for human consumption, used by `GPNode`'s printer functions. The default version just calls `toString()`. You probably want to write something prettier. The `encode()` and `decode(...)` methods are supposed to use the `Code` package (Section 2.2.2) to encode and decode the ERC in a reasonable fashion.

Finally, the `readNode(...)` and `writeNode(...)` methods, as usual, read and write the ERC in binary fashion. You only need to implement these methods if you're planning on writing over the network (such as using distributed evaluation or island models). But they're easy so why not? And of course, there's the method to actually execute the ERC as code. This typically returns the ERC's internal value (it's a constant after all):

```
public abstract void eval(EvolutionState state, int thread, GPData input,
                        ADFStack stack, GPIndividual individual, Problem problem);
```

**Example** Let's create an ERC and add it to our existing example from Section 5.2.6. We'll make an ERC which represents constants between 0.0 and 1.0, not including 1.0. Our mutator will add a little Gaussian noise to the node. Here's a full class:

```
package ec.app.myapplication;
import ec.gp.*;
public class MyERC extends ERC {
    public double value;
    public String toStringForHumans() { return "" + value; }
    public String encode() { return Code.encode(value); }

    public boolean decode(DecodeReturn ret) {
        int pos = dret.pos;
        String data = dret.data;
        Code.decode(dret);
        if (dret.type != DecodeReturn.T_DOUBLE) // uh oh! Restore and signal error.
            { dret.data = data; dret.pos = pos; return false; }
        value = dret.d
        return true;
    }

    public boolean nodeEquals(GPNode node)
        { return (node.getClass() == this.getClass() && ((MyERC)node).value == value); }

    public void readNode(EvolutionState state, DataInput input) throws IOException
        { value = dataInput.readDouble(); }

    public void writeNode(EvolutionState state, DataOutput output) throws IOException
        { dataOutput.writeDouble(value); }

    public void resetNode(EvolutionState state, int thread)
        { val = state.random[thread].nextDouble(); }

    public void mutateNode(EvolutionState state, int thread) {
```

```

double v;
do v = value + state.random[thread].nextGaussian() * 0.01;
while( v < 0.0 || v >= 1.0 );
value = v;
}

public void eval(EvolutionState state, int thread, GPData input, ADFStack stack,
                GPIndividual individual, Problem Problem)
{ ((MyData)data).val = value; }
}

```

Now let's set up the parameters to use it. We'll change the function set. Our ERC is a terminal so it takes no arguments: we'll use nc0 as its constraints.

```

# Our Function Set
gp.fs.0 = ec.gp.GPFunctionSet
gp.fs.0.size = 6
gp.fs.0.func.0 = ec.app.myapp.X
gp.fs.0.func.0.nc = nc0
gp.fs.0.func.1 = ec.app.myapp.Y
gp.fs.0.func.1.nc = nc0
gp.fs.0.func.2 = ec.app.myapp.Mul
gp.fs.0.func.2.nc = nc2
gp.fs.0.func.3 = ec.app.myapp.Sub
gp.fs.0.func.3.nc = nc2
gp.fs.0.func.4 = ec.app.myapp.Sin
gp.fs.0.func.4.nc = nc1
gp.fs.0.func.5 = ec.app.myapp.MyERC
gp.fs.0.func.5.nc = nc0

```

... and we're done.

## 5.2.10 Automatically Defined Functions and Macros

**Automatically Defined Functions** (ADFs) [4] are a standard way of creating some modularism in Genetic Programming. They define multiple trees in the GPIndividual and essentially define a function calling structure where certain trees can call other trees as subfunctions. ADFs are the primary reason why GPIndividual has multiple GPTrees.

The simplest kind of ADF is found in Figure 5.6. Here each ADF function is a terminal, and when it is evaluated, it simply evaluates the corresponding ADF tree, then returns the tree's return value. The ADF function is a GPNode, an instance of ec.gp.ADF. It's very rare to further specialize this class. Notice that calling is nested — an ADF can call another ADF and so on. However it's not very common to have *recursive* calls because you'll need to construct some kind of stopping base-case criterion to avoid infinite recursive loops.

ADFs add a two more parameters to the standard GPNode suite. Let's say we're adding a zero-argument ADF. Beyond the node constraints, we also need to specify which tree will be called when the ADF is executed; and also a simple name for the ADF to distinguish it from other ADFs and GPNodes. Ideally this name should only have lowercase letters, numbers, and hyphens (that is, "Lisp-style"):

```

gp.fs.0.func.5 = ec.gp.ADF
gp.fs.0.func.5.nc = nc0
gp.fs.0.func.5.tree = 1
gp.fs.0.func.5.name = ADF0

```



Figure 5.6 A GP Individual with two no-argument (terminal) ADFs. The primary GP Tree has functions in its function set that can call the other two trees; in turn the ADF 1 tree has a function in its function set that can call the ADF 2 tree. The return values of the various ADF trees become the return values of their respective calling functions.

**It's traditional that the first tree be called ADF0, the second ADF1, and so on.** Since typically the first GP Tree in the array is the "main" tree, and the second tree is ADF0, this means that ADF0's associated tree number is usually 1, and ADF1's associated tree number is usually 2. If you don't specify a name, ECJ will maintain this tradition by setting the name to "ADF" + (tree number - 1), which is usually right anyway. It'll also issue a warning.

The name of the ADF (in this case "ADF1") and its associated GP tree (in this case, tree 1) are stored in the `ec.gp.ADF` class like this:

```
public int associatedTree;
public String name;
```

The `name()` method returns the value of the name variable.

We'll show how to set up the ADF tree itself in the Example below.

ADFs can also have arguments to the functions. In Figure 5.7, we have an ADF with two arguments. The way this works is as follows: when an ADF function is called, we first evaluate its children, then hold their return values in storage. We then call the corresponding ADF tree. In that tree there may be one or more special terminal GPNodes, two different kinds of instances of `ec.gp.ADFArgument`. One group of instances, when evaluated, will return the value of the first child. The second group return the value of the second child. This enables the ADF tree to use arguments in its "function call" so to speak.

ADFAArguments add one additional parameter: the child number associated with the argument. For example:





Figure 5.7 A GPIndividual with one 2-argument ADF. The primary GP Tree has a function in its function set that can call the ADF tree. This function first evaluates its children, then executes the ADF tree. In the ADF tree there are two terminal functions (ARG 1 and ARG 2) which, when evaluated, return the values of the two children respectively. The return value of the ADF tree becomes the return value of the ADF function.

```
gp.fs.1.func.6 = ec.gp.ADFArgument
gp.fs.1.func.6.nc = nc0
gp.fs.1.func.5.arg = 0
gp.fs.1.func.5.name = ARG0
```

If you don't specify a name, ECJ will set the name to "ARG" + arg number, which is usually right anyway. It'll also issue a warning.

The ADFArgument's name and argument number are stored in the `ec.gp.ADFArgument` class as:

```
public int argument; public String name;
```

Again, the `name()` method returns the value of the name variable.

ECJ also supports **Automatically-Defined Macros** (or ADMs), described in [18]. These differ from ADFs only in when the children are evaluated. When an ADM node is evaluated, its children are not evaluated first; rather the ADM immediately calls its associated tree. When an argument node in that tree (again, a terminal) is evaluated, we teleport back to the associated child and evaluate it right then and there, then return its value. Note that this means that children may never be evaluated; or can be evaluated multiple times, as shown in Figure 5.8.

ADMs are just like ADFs in their parameters:

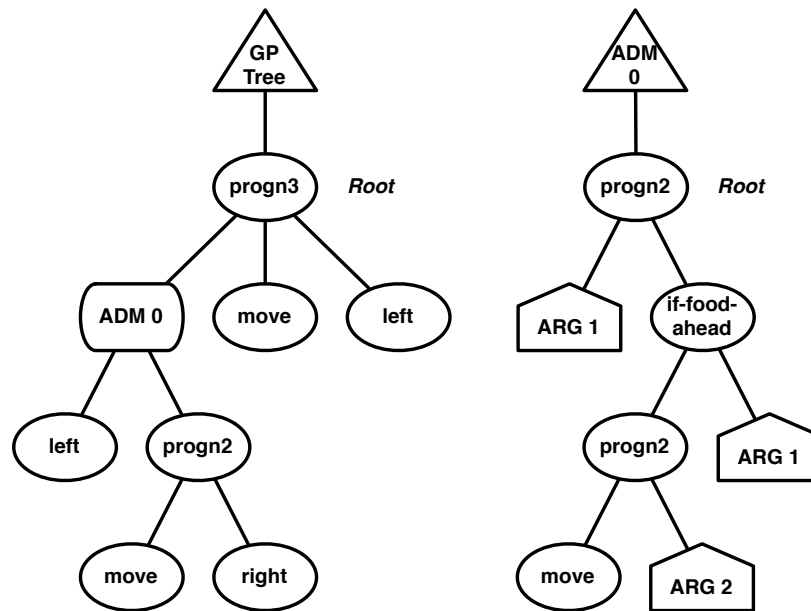


Figure 5.8 A GPIndividual with one 2-argument ADM. The primary GP Tree has a function in its function set that can call the ADM tree. This function delays the evaluation of its children, and immediately executes the ADM tree. In the ADM tree there are two terminal functions (ARG 1 and ARG 2) which, when evaluated, evaluate (or, as necessary, re-evaluate) the original children to the ADM function and then return their values. Notice that in this example child # 1 may be evaluated *twice*, and child #2, depending on whether there's food ahead, may be *never evaluated*.

```
gp.fs.0.func.6 = ec.gp.ADM
gp.fs.0.func.6.nc = nc2
gp.fs.0.func.6.tree = 1
# This will be called "ADM1"
gp.fs.0.func.6.name = ADM1
```

If an ADF or ADM tree has arguments, it probably will require its own separate GPTreeConstraints, because it needs to have its own GPFunctionSet with those arguments defined. See the Example below.

### 5.2.10.1 About ADF Stacks

ADFs and ADMs are a bit complex. To do their magic, they need a special object called an `ec.gp.ADFStack`. This is actually two stacks of `ec.gp.ADFContext` objects which store the location and current return values of various children. These classes are almost never overridden: here's the standard (default) parameters for them:

```
gp.problem.stack = ec.gp.ADFStack
gp.adf-stack.context = ec.gp.ADFContext
```

You *could* have defined it this way...

```
eval.problem.stack = ec.gp.ADFStack
eval.problem.stack.context = ec.gp.ADFContext
```

... but there are advantages in using the default parameters, particularly when getting to Grammatical Evolution (Section 5.3).

**Example** ADFs and ADMs are fairly straightforward to implement but they can require a fair number of parameters. Continuing with the example in started in Section 5.2.6 and extended in Section 5.2.9, let's add a 2-argument ADF to the individual. This will require adding a second GPTree and its own GPTreeConstraints.

Let's begin by modifying the GPFunctionSet of the original tree to include this ADF:

```
# Our Function Set
gp.fs.0 = ec.gp.GPFunctionSet
gp.fs.0.size = 7
gp.fs.0.func.0 = ec.app.myapp.X
gp.fs.0.func.0.nc = nc0
gp.fs.0.func.1 = ec.app.myapp.Y
gp.fs.0.func.1.nc = nc0
gp.fs.0.func.2 = ec.app.myapp.Mul
gp.fs.0.func.2.nc = nc2
gp.fs.0.func.3 = ec.app.myapp.Sub
gp.fs.0.func.3.nc = nc2
gp.fs.0.func.4 = ec.app.myapp.Sin
gp.fs.0.func.4.nc = nc1
gp.fs.0.func.5 = ec.app.myapp.MyERC
gp.fs.0.func.5.nc = nc0
gp.fs.0.func.6 = ec.gp.ADF
gp.fs.0.func.6.nc = nc2
gp.fs.0.func.6.tree = 1
gp.fs.0.func.6.name = ADF1
```

Let's create a second function set for our second (ADF) tree. This set will have all the same functions as the main tree, except for the ADF function (we don't want to call ourselves recursively!) Instead we'll add two ADFArgument nodes to represent the two children.

```
gp.fs.size = 2

# Our Second Function Set
gp.fs.1 = ec.gp.GPFunctionSet
gp.fs.1.name = f1
gp.fs.1.size = 8
gp.fs.1.func.0 = ec.app.myapp.X
gp.fs.1.func.0.nc = nc0
gp.fs.1.func.1 = ec.app.myapp.Y
gp.fs.1.func.1.nc = nc0
gp.fs.1.func.2 = ec.app.myapp.Mul
gp.fs.1.func.2.nc = nc2
gp.fs.1.func.3 = ec.app.myapp.Sub
gp.fs.1.func.3.nc = nc2
gp.fs.1.func.4 = ec.app.myapp.Sin
gp.fs.1.func.4.nc = nc1
gp.fs.1.func.5 = ec.app.myapp.MyERC
gp.fs.1.func.5.nc = nc0
gp.fs.1.func.6 = ec.gp.ADFArgument
gp.fs.1.func.6.nc = nc0
gp.fs.1.func.6.arg = 0
gp.fs.1.func.6.name = ARG0
gp.fs.1.func.7 = ec.gp.ADFArgument
gp.fs.1.func.7.nc = nc0
gp.fs.1.func.7.arg = 1 gp.fs.1.func.7.name = ARG1
```

Now we create a new GPTreeConstraints which uses this function set:

```
gp.tc.size = 2

# Our Second Tree Constraints
gp.tc.1 = ec.gp.GPTreeConstraints
gp.tc.1.name = tc1
gp.tc.1.fset = f1
gp.tc.1.returns = nil
gp.tc.1.init = ec.gp.koza.HalfBuilder
```

Next we add the second tree to the GPIndividual:

```
pop.subpop.0.species.ind.numtrees = 2
pop.subpop.0.species.ind.tree.1 = ec.gp.GPTree
pop.subpop.0.species.ind.tree.1.tc = tc1
```

The ADF stack and ADF context were already defined in the previous examples, but we'll do it again here for clarity:

```
gp.problem.stack = ec.gp.ADFStack
gp.adf-stack.context = ec.gp.ADFContext
```

...and we're done!

An important note: because the main GP Tree and the ADF have different function sets and thus different GPTreeConstraints, standard GP Crossover (see Section 5.2.5) won't cross over the ADF tree with a main tree of some other individual or vice versa. But if, for example, your GPIndividual had two ADF trees that had the *same* GPTreeConstraints, they could get crossed over with arbitrary other ADF trees in another GPIndividuals.

## 5.2.11 Strongly Typed Genetic Programming

Sometimes it's useful to constrain which GPNodes may serve as children of other GPNodes, or the root of the GPTree. For example, consider a GPNode called (if *test then else*). This node evaluates *test*, and based on its result (true or false) it either evaluates and returns *then* or *else*. Let's presume that *then* and *else* (and *if*) return doubles. On the other hand, *test* is intended to return a boolean. So you'll need to have some GPNodes in your function set which return doubles and others which return booleans; the (if ...) node itself returns a double.

The problem is not that you have nodes which return different values — this is easily handled by hacking your GPData object. The problem is that you now have constraints on valid tree structures: you can't plug a node which returns a double (say, (sin ...)) into the *test* slot of your (if ...) node, which is expecting a boolean.<sup>5</sup> This is where strong typing comes in.

ECJ's typing system is simple but sufficient for many common uses. It's not as sophisticated as a full polymorphic typing system but it also doesn't have the hair-raising complexity that such a system requires. ECJ is complex enough as it is thank you very much!

ECJ's system is based on **type objects**, subclasses of the abstract superclass `ec.gp.GPType`, and nodes are allowed to connect as parent and child if their corresponding type objects are **type compatible**. ECJ's type objects of two kinds: atomic types and set types. An atomic type is just a single object (in fact, it's theoretically just a symbol, or in some sense, an integer). A set type is a set of atomic types. Type compatibility is as follows:

- Two atomic types are compatible if they are the same.

---

<sup>5</sup>Well, you *could* if you assumed that 0.0 was false and anything else was true. But this is a hack. The right way to do it is to constrain things properly

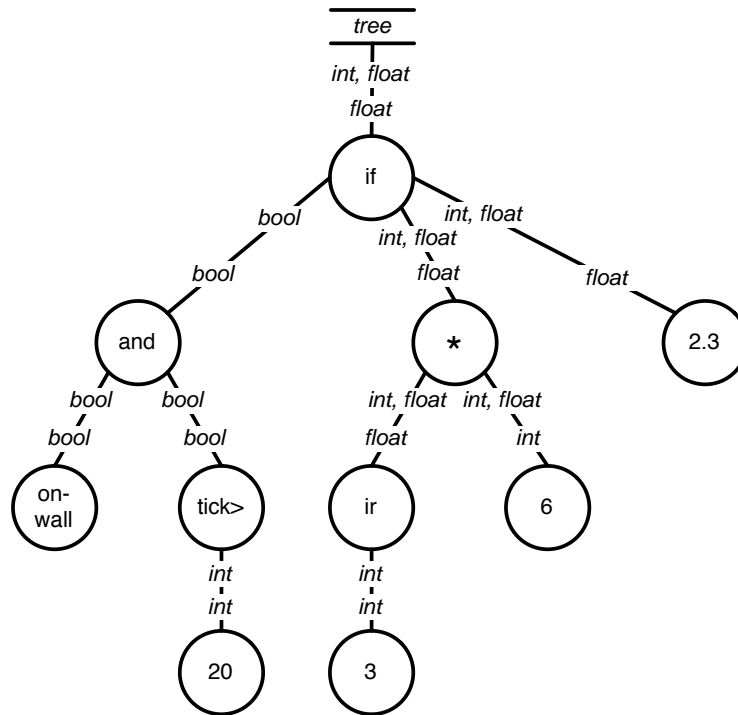


Figure 5.9 A typed genetic programming parse tree. Types of the form “int, float” are set types. All others are atomic types. A repeat of Figure 1.4.

- A set type is compatible with an atomic type if it contains the atomic type in its set.
- Two set types are compatible if their intersection is nonempty.

Every GPNode is assigned a type object to represent the “return type” of the GPNode. Furthermore every nonterminal GPNode is assigned a type object for each of its children: this is called the “child type” or “argument type” or that particular child slot. Last, the GPTree itself is assigned “root type”: a type for the root of the tree. Each GPTree in a GPIndividual can have a different root type. Here’s what must be true about any given GPTree.

- For any parent and child in a tree, with the child in slot *C*, the return type of the child must be compatible with the child type of the parent for slot *C*.
- The return type of the root GPNode must be compatible with the GPTree’s root type. This ensures that the tree is returning

You can see an example of a GPTree with type constraints listed in Figure 5.9.

Every GPNodeBuilder and GP Breeding Pipeline must maintain the constraints guaranteed by typing. The issue is guaranteeing that if you replace one GPNode with another, that the second GPNode will be legal. This is done primarily with the following two utility functions:

#### ec.gp.GPNode Methods

public GPType parentType(GPInitializer initializer)

If the GPNode’s parent is another GPNode, this returns the type of the parent’s child slot presently filled by the GPNode. If the GPNode’s parent is a GPTree, this returns the type of the tree’s root.

```
public final boolean swapCompatibleWith(GPInitializer initializer, GPNode node)
    Returns true swapping the GPNode into the slot presently occupied by node is acceptable, type-wise.
```

---

Before you can assign types, you'll need to define them. Each type is given a unique symbol. As an example, let's begin by creating two atomic types, called (uninterestingly) "boolean" and "nil" (we say "nil" instead of "double" so we don't have to redefine the GPTreeConstraints and GPNodeConstraints we defined earlier, which all use "nil" as their types). We'd say:

```
gp.type.a.size = 2
gp.type.a.0.name = boolean
gp.type.a.1.name = nil
```

We might also define a set type or two. For fun, let's create a set type which contains both booleans and doubles. We'll also have to stipulate the atomic types encompassed by the set type:

```
gp.type.s.size = 1
gp.type.s.0.name = boolean-or-nil
gp.type.s.0.size = 2
gp.type.s.0.member.0 = boolean
gp.type.s.0.member.1 = nil
```

What's the point of set types, you might ask. Why not just atomic types? Set types are particularly useful for describing situations where a given GPNode can adapt to several different kinds of children in a given slot. This is particularly useful for simulating notions of subtyping or subclassing. For example, a GPNode like (sin ...) might declare that its child can either be an "integer" or a "double", by having its child type defined as a set type of "number" which encapsulates both integers and doubles.

What this typing facility *cannot* do is dynamically change types as necessary. For example, you cannot say that a GPNode like (+ ... ...) returns a double if either of its children is of type double, but if both of its children are of type integer, then it returns an integer.

You are also restricted to a finite number of types. For example, consider a GPNode called (matrix-multiply ... ...) which takes two children which return matrices. You cannot say that if the left child is an  $M \times N$  matrix, and your right child is an  $N \times P$  matrix, that the return type will be an  $M \times P$  matrix. This is partly because you can't define dynamic typing, but it's also because  $M$ ,  $N$ , and  $P$  can be any of an infinite number of numbers, resulting in an infinite number of types: that would be hard on your fingers. There exist polymorphic typing systems for genetic programming but they're fairly bleeding-edge. If you need things like this, I suggest instead to look at Grammatical Evolution (Section 5.3).

**Example** Let's add some typing to the example we started in Section 5.2.6 and continued in Sections 5.2.9 and 5.2.10. We'll add a boolean type as before, and a few functions which rely on it.

First the boolean type:

```
gp.type.a.size = 2
gp.type.a.0.name = boolean
gp.type.a.1.name = nil gp.type.s.size = 1
gp.type.s.0.name = boolean-or-nil
gp.type.s.0.size = 2
gp.type.s.0.member.0 = boolean
gp.type.s.0.member.1 = nil
```

Next we'll need to modify the node constraints. We'll not add a new tree constraints because we don't want any trees which return booleans.

For the node constraints, let's add three new GPNodeConstraints:

- A function which returns a double (nil), has three children, and the first child needs to be a boolean. The other two children are nil. This would be for things like the (if ... ..) node.
- A function which takes two booleans and returns a boolean. This would be for functions like (nand ... ..).
- A function which takes two doubles and returns a boolean. This would be for functions like (> ... ..).

These three new GPNodeConstraints would be:

```
gp.nc.size = 6

# ... first come the original node constraints, then these:

# Example: (if ... ..)
gp.nc.3 = ec.gp.GPNodeConstraints
gp.nc.3.name = nc3
gp.nc.3.returns = nil
gp.nc.3.size = 3
gp.nc.3.child.0 = boolean
gp.nc.3.child.1 = nil
gp.nc.3.child.2 = nil

# Example: (nand ... ..)
gp.nc.4 = ec.gp.GPNodeConstraints
gp.nc.4.name = nc2
gp.nc.4.returns = boolean
gp.nc.4.size = 2
gp.nc.4.child.0 = boolean
gp.nc.4.child.1 = boolean

# Example: (> ... ..)
gp.nc.5 = ec.gp.GPNodeConstraints
gp.nc.5.name = nc2
gp.nc.5.returns = boolean
gp.nc.5.size = 2
gp.nc.5.child.0 = nil
gp.nc.5.child.1 = nil
```

Recall that our GPData looks like this:

```
package ec.app.myapp;
import ec.gp.*;
public class MyData extends GPData
{
    public double val;
    public GPData copyTo(GPData other)
    { ((MyData)other).val = val; return other; }
}
```

We could modify this to include a boolean data type as well, but we'll just use the val variable to store both boolean and real-valued data. Before you go about that, re-read Section 5.2.3.1 (on GPData) and its discussion about clone() and copyTo(...).

Now let's define our three new GPNodes. First our If-statement:

```

package ec.app.myapp;
import ec.*;
import ec.gp.*;
public class If extends GPNode {
    public String toString() { return "if" };
    public void eval(EvolutionState state, int thread, GPData input,
        ADFStack stack, GPIndividual individual, GPPProblem problem) {
        MyData data = (MyData) input;
        children[0].eval(state, thread, data, stack, individual, prob);
        if (data.val != 0.0) // true
            children[1].eval(state, thread, data, stack, individual, prob);
        else
            children[2].eval(state, thread, data, stack, individual, prob);
        // the result be stored in data
    }
}

```

Next, the nand node:

```

package ec.app.myapp;
import ec.*;
import ec.gp.*;
public class Nand extends GPNode {
    public String toString() { return "nand" };
    public void eval(EvolutionState state, int thread, GPData input,
        ADFStack stack, GPIndividual individual, GPPProblem problem) {
        MyData data = (MyData) input;
        children[0].eval(state, thread, data, stack, individual, prob);
        boolean left = (data.val != 0.0);
        children[1].eval(state, thread, data, stack, individual, prob);
        boolean right = (data.val != 0.0);
        data.val = !(left && right) ? 1.0 : 0.0;
    }
}

```

Next, the > node:

```

package ec.app.myapp;
import ec.*;
import ec.gp.*;
public class GreaterThan extends GPNode {
    public String toString() { return ">" };
    public void eval(EvolutionState state, int thread, GPData input,
        ADFStack stack, GPIndividual individual, GPPProblem problem) {
        MyData data = (MyData) input;
        children[0].eval(state, thread, data, stack, individual, prob);
        double left = data.val;
        children[1].eval(state, thread, data, stack, individual, prob);
        double right = data.val;
        data.val = (left > right) ? 1.0 : 0.0;
    }
}

```

Notice that these functions hijack the double value to store boolean information. This is okay because we *know* that the recipient of this information will understand it. How do we know? Because the typing constraints have made it impossible to be otherwise.



So let's add these functions to the function set of our main GP Tree:

```
# Our Main Tree Function Set
gp.fs.0 = ec.gp.GPFunctionSet
gp.fs.0.size = 10
gp.fs.0.func.0 = ec.app.myapp.X
gp.fs.0.func.0.nc = nc0
gp.fs.0.func.1 = ec.app.myapp.Y
gp.fs.0.func.1.nc = nc0
gp.fs.0.func.2 = ec.app.myapp.Mul
gp.fs.0.func.2.nc = nc2
gp.fs.0.func.3 = ec.app.myapp.Sub
gp.fs.0.func.3.nc = nc2
gp.fs.0.func.4 = ec.app.myapp.Sin
gp.fs.0.func.4.nc = nc1
gp.fs.0.func.5 = ec.app.myapp.MyERC
gp.fs.0.func.5.nc = nc0
gp.fs.0.func.6 = ec.gp.ADF
gp.fs.0.func.6.nc = nc2
gp.fs.0.func.6.tree = 1
gp.fs.0.func.6.name = ADF1
gp.fs.0.func.7 = ec.app.myapp.If
gp.fs.0.func.7.nc = nc3
gp.fs.0.func.8 = ec.app.myapp.Nand
gp.fs.0.func.8.nc = nc4
gp.fs.0.func.9 = ec.app.myapp.GreaterThan
gp.fs.0.func.9.nc = nc5
```

...and we're done!

**Mixing ADF and ADMs and Typed GP** A quick note. The return type of an ADF node must match the root type of its corresponding ADF tree. Additionally, the child type of a certain slot in an ADF node must match the return type of the corresponding ADFArgument.

### 5.2.11.1 Inside GPTypes

If you want to create a GPNodeBuilder or a GP Breeding Pipeline, you ought to go in more detail about GPTypes.

ec.gp.GPType is an abstract superclass of ec.gp.GPAtomicType and ec.gp.GPSetType, which define the atomic and set types respectively. The two basic data elements in a GPType are:

```
public String name;
public int type;
```

The first variable, like GPFunctionSet, GPTreeConstraints, and GPNodeConstraints, holds the name of the type (as defined in the parameters). The second variable holds a uniquely-assigned integer for this type. The important feature for types is to determine whether they are **type-compatible** with one another. The compatibility function is this:

#### ec.gp.GPType Methods

---

```
public boolean compatibleWith(GPInitializer initializer, GPType type)
```

Returns true if the return type of this GPNode is type-compatible with the given type.

---

GPAAtomicTypes are simple: they are compatible with one another if their type integer is the same. A GPSetType instead is a set of GPAAtomicTypes, stored in different ways for query convenience:

```
public Hashtable types_h;  
public int[] types_packed;  
public boolean[] types_sparse;
```

The first is the GPAAtomicTypes in the set stored in a Hashtable. The second is an array of the GPAAtomicTypes. And the third is an array of booleans, one for each GPAAtomicType number, which is true if that GPAAtomicType is a member of the set.

A GPSetType is compatible with a GPAAtomicType if the GPSetType contains the GPAAtomicType as an element. Two GPSetTypes are compatible with one another if their intersection is nonempty.

### 5.2.12 Parsimony Pressure (The ec.parsimony Package)

Genetic programming has a serious bloat problem: as evolution progresses, the size of the trees inside the population tend to grow without bound. This is a problem that exists for various arbitrary-length representations (lists, graphs, rule sets, etc.) but genetic programming has studied it the most.

The most common simple way of keeping trees down is to make it illegal to produce a tree larger than a certain depth. For example, Koza's standard rules, adhered to by the basic parameters in ECJ, stipulate that crossover and mutation operators may not produce a child which is deeper than 17 nodes [3], for example:

```
gp.koza.xover.tries = 1  
gp.koza.xover.maxdepth = 17
```

Here if the crossover operation produces a child greater than 17, it is not forwarded on; but rather its (presumably smaller) parent is forwarded on in its stead. This is a fairly crude approach to parsimony pressure, but it's fairly effective. Another approach — which can be done at the same time — is to modify the selection operator to favor smaller individuals. This notion is called **parsimony pressure**.

In the ec.parsimony package, ECJ has several SelectionMethods which select both based on fitness and on size (smaller size being preferred). These methods compute size based on GPNode's size() function. Many of these selection methods were compared and discussed at length in [9]. Here they are:

- ec.parsimony.LexicographicTournamentSelection is a straightforward TournamentSelection operator, except that the fitter Individual is preferred (as usual), but when both Individuals are the same fitness, the smaller Individual is preferred. Parameters for this operator are basically the same as for TournamentSelection.

Let us presume that the SelectionMethod is the first source of the pipeline of Subpopulation 0. Then the basic parameters are the same as in TournamentSelection:

```
pop.subpop.0.species.pipe.source.0 = ec.parsimony.LexicographicTournamentSelection  
pop.subpop.0.species.pipe.source.0.size = 7  
pop.subpop.0.species.pipe.source.0.pick-worst = false
```

Or using the default parameters:

```
pop.subpop.0.species.pipe.source.0 = ec.parsimony.LexicographicTournamentSelection  
select.lexicographic-tournament.size = 7  
select.lexicographic-tournament.pick-worst = false
```

The problem with this method is that bloat control only comes into effect for problems where lots of fitness ties occur. This problem lead us to two modifications of the basic idea:

- ec.parsimony.BucketTournamentSelection is like LexicographicTournamentSelection, except that first individuals are placed into  $N$  classes ("buckets") based on fitness. The subpopulation is first sorted

by fitness. Then the bottom  $\frac{\text{subpopulation size}}{N}$  individuals are placed in the worst bucket, plus any individuals remaining in the subpopulation with the same fitness as the best individual in that bucket. Next the bottom remaining  $\frac{\text{subpopulation size}}{N}$  are placed in the second worst bucket, plus any individuals remaining in the population with the same fitness as the best individual in that bucket. This continues until all the individuals are exhausted. BucketTournamentSelection then works like LexicographicTournamentSelection except that instead of comparing based on fitness, it compares based on the bucket the individual is in. The number of buckets is defined by num-buckets:

```
pop.subpop.0.species.pipe.source.0 = ec.parsimony.BucketTournamentSelection
pop.subpop.0.species.pipe.source.0.size = 7
pop.subpop.0.species.pipe.source.0.pick-worst = false
pop.subpop.0.species.pipe.source.0.num-buckets = 10
```

Or using the default parameters:

```
pop.subpop.0.species.pipe.source.0 = ec.parsimony.BucketTournamentSelection
select.bucket-tournament.size = 7
select.bucket-tournament.pick-worst = false
select.bucket-tournament.num-buckets = 10
```

- ec.parsimony.ProportionalTournamentSelection is like TournamentSelection, except that it *either* selects based on fitness *or* selects based on size. It determines which one to do by flipping a coin of a certain probability that *fitness* will be used. The parameters are:

```
pop.subpop.0.species.pipe.source.0 = ec.parsimony.ProportionalTournamentSelection
pop.subpop.0.species.pipe.source.0.size = 7
pop.subpop.0.species.pipe.source.0.pick-worst = false
pop.subpop.0.species.pipe.source.0.fitness-prob = 0.9
```

Or using the default parameters:

```
pop.subpop.0.species.pipe.source.0 = ec.parsimony.ProportionalTournamentSelection
select.bucket-tournament.size = 7
select.bucket-tournament.pick-worst = false
select.bucket-tournament.fitness-prob = 0.9
```

- ec.parsimony.DoubleTournamentSelection is actually two TournamentSelections in a row. In short, we do a TournamentSelection of tournament size  $N$  based on fitness: but the entrants to that tournament are not chosen uniformly at random from the subpopulation, but rather are the winners of  $N$  *other* tournament selections, each performed based on size. Alternatively, we can first do tournament selections on fitness, then have a final tournament on size.

Thus there are roughly twice as many parameters: ones describing the final tournament, and ones describing the initial (“qualifying”) tournaments

```

pop.subpop.0.species.pipe.source.0 = ec.parsimony.DoubleTournamentSelection

# Final tournament
pop.subpop.0.species.pipe.source.0.size = 2
pop.subpop.0.species.pipe.source.0.pick-worst = false

# Qualifying tournaments
pop.subpop.0.species.pipe.source.0.size2 = 2
pop.subpop.0.species.pipe.source.0.pick-worst2 = false

# Make the qualifying tournament based on size
pop.subpop.0.species.pipe.source.0.do-length-first = true

```

Or using the default parameters:

```

pop.subpop.0.species.pipe.source.0 = ec.parsimony.DoubleTournamentSelection

# Final tournament
select.double-tournament.size = 7
select.double-tournament.pick-worst = false

# Qualifying tournaments
select.double-tournament.size2 = 7
select.double-tournament.pick-worst2 = false

# Make the qualifying tournament based on size
select.double-tournament.do-length-first = true

```

- `ec.parsimony.TarpeianStatistics` implements the “Tarpeian” parsimony pressure method [13]. This method identifies the individuals in the subpopulation with above-average size. Notice that this may not be half the subpopulation: it could be a very small number if they are very large and the others are very small. Then a certain proportion of these individuals, picked randomly, are assigned a very bad fitness, and their evaluated flags are set. This happens *before* evaluation, so the evaluation procedure doesn’t bother to evaluate those individuals further.

The Tarpeian method isn’t a selection procedure: it’s a fitness assignment procedure. As such it’s not implemented as a `SelectionMethod` but rather as a `Statistics` subclass which hooks into the evolutionary loop prior to evaluation.

Let’s say that `TarpeianStatistics` is the only child of our primary `Statistics` object. The parameters would look like this:

```

stat.num-children = 1
stat.child.0 = ec.parsimony.TarpeianStatistics
stat.child.0.kill-proportion = 0.2

```

## 5.3 Grammatical Evolution (The `ec.gp.ge` Package)

Grammatical Evolution (GE) [17] is an approach to building genetic programming trees using a grammar interpreted by a list representation. The trees are then evaluated and tested. In ECJ the procedure for evaluating Individuals in GE works roughly like this.

- The representation is an arbitrarily long list of bytes (See Section 5.1.2), implemented as a special subclass of `ByteVectorIndividual` called `ec.gp.ge.GEIndividual`.

- The Fitness is typically a KozaFitness.
- The GEIndividual's species is a `ec.gp.ge.GESpecies`.
- The GESpecies holds a **grammar**, loaded from a file and produced via an `ec.gp.ge.GrammarParser`, which we will interpret according to the bytes in the Individual.
- To assess the fitness of a GEIndividual, we hand it to the GESpecies which interprets it according to the grammar, producing a GPIIndividual.
- The GPIIndividual is then evaluated in the normal ECJ fashion and its Fitness is set.
- We then transfer the Fitness to the GEIndividual.

We'd like to use plain-old GP test problems to make this as easy as possible. To pull this off we have to do a few minor hacks. First, we must insert the GE conversion process in-between the Evaluator and the GPPProblem that the user wrote. To do this we have a special Problem class called `ec.gp.ge.GEProblem`, which is assigned as the Evaluator's problem. The GPPProblem is then set up as a subsidiary of the GEProblem. When the Evaluator wishes to evaluate a `ByteVectorIndividual`, it calls the GEProblem, which converts it to a `GPIIndividual` and then hands the `GPIIndividual` to the GPPProblem to evaluate. This is done as follows:

```
eval.problem = ec.gp.ge.GEProblem
eval.problem.problem = ec.app.myproblem.MyGPPProblem
```

Note that GPPProblems usually require auxiliary parameters (not the least of which is their `GPData`), so we'll need to also say things like...

```
eval.problem.problem.data = ec.app.myproblem.MyGPData
```

... etc. You don't need to define the `ADFStack` or `ADFContext` specially because the default parameters for them suffice:

```
gp.problem.stack = ec.gp.ADFStack
gp.adf-stack.context = ec.gp.ADFContext
```

### 5.3.1 GEIndividuals, GESpecies, and Grammars

The reason we use `GEIndividual` instead of just `ByteVectorIndividual` is simple: when we print out the `GEIndividual` we wish to print out the equivalent `GPIIndividual` as well. That's the *only* thing `GEIndividual` does beyond being just a `ByteVectorIndividual`.

The `GEProblem` doesn't actually do the translation from list to tree. Instead, it calls on the `GESpecies` to do this dirty work. First, let's set the `GESpecies`, `GEIndividual`, `GrammarParser`, and `Fitness` of Species 0:

```
pop.subpop.0.species = ec.gp.ge.GESpecies
pop.subpop.0.species.parser = ec.gp.ge.GrammarParser
pop.subpop.0.species.ind = ec.gp.ge.GEIndividual
pop.subpop.0.species.fitness = ec.gp.koza.KozaFitness
```

`GESpecies` requires a **grammar file** for each tree in the `GPIIndividual`. If your `GPIIndividual` has two trees (say), you'll need two grammar files, which are specified like this:

```
pop.subpop.0.species.file.0 = foo.grammar
pop.subpop.0.species.file.1 = bar.grammar
```

Alternatively you can use the default parameter base:

```
ge.species.file.0 = foo.grammar
ge.species.file.1 = bar.grammar
```

Grammar files are text files consisting of lines. Each line can be blank whitespace (which is ignored), a comment (which starts with a #, just like parameter files), or a **grammar rule**. A grammar rule has a **head**, followed by whitespace, followed by the string " ::= ", then more whitespace, and finally a **body**. Here is an example rule:

```
<foo> ::= (hello <bar>) | (yo) | (whoa <bar> <baz>)
```

In this example, <foo> is the head, and (hello <bar>) | (yo) | (whoa <bar> <baz>) is the body. The head is always a simple, usually lower-case, symbol surrounded in angle brackets. The body is a collection of **clauses** separated by the pipe symbol ( | ). Each clause is a single GP node in Lisp s-expression form. For example, (hello <bar>) defines a GP node named hello, which has a single child defined by the symbol <bar>.

Terminal (leaf-node) GPNodes are defined as S-expressions with no children. For example, (yo) is a terminal GPNode whose name is yo. And a GPNode can have more than one child, such as (whoa <bar> <baz>).

For each symbol that appears in the body of a rule in angle brackets, there must be at least one such symbol in a head of a rule. However you can have multiple such heads:

```
<foo> ::= (hello <bar>)
<foo> ::= (yo) | (whoa <bar> <baz>)
```

This says the exact same thing as the previous rule example. You are free to use pipe symbols or create separate rules with the same head, or mix and match the two. You can also have plain angle-bracket symbols in the body like this:

```
<foo> ::= (blah <bar>) | <okay>
```

This says that <foo> can expand to either the GPNode blah or it can expand to whatever <okay> expands to. You'd probably want to watch out for recursive definitions. For example, these wouldn't be great rules:

```
<foo> ::= (blah <bar>) | <okay>
<okay> ::= <foo> | (yuck)
```

The **head of the first rule in the file** is the entry point of the grammar and represents the root node of the GPTree. In all of our examples we have called it <start>, but it doesn't have to be that. For example, here is a complete grammar for a very simple tree which produces the same kinds of GPIndividuals as described in the example in Section 5.2.3:

```
<start> ::= <op>
<op> ::= (x) | (y)
<op> ::= (sin <op>) | (* <op> <op>) | (- <op> <op>)
```

### 5.3.1.1 Strong Typing

At this point we don't see a lot of power. But Grammatical Evolution can define all sorts of typing requirements by specifying which symbols appear where in rules. For example, we could extend this grammar to handle the Strongly-Typed Genetic Programming example shown in Section 5.2.11:

```

<start> ::= <op>
<op> ::= (x) | (y) | (sin <op>) | (* <op> <op>) | (- <op> <op>)
<op> ::= (if <bool> <op> <op>)
<bool> ::= (nand <bool> <bool>)
<bool> ::= (> <op> <op>)

```

If you have a grammar like this, you’d imagine it would need to be accompanied by GPNodes that have been appropriately typed in a strongly typed context. But it’s not true. Recall that strong typing is intended to provide constraints for building trees, mutating them, and crossing them over using traditional GP methods. But since Grammatical Evolution is doing all of this based on a list of bytes, GPTypes serve no function.

As a result, you should just have a single GPType even in a “strongly typed” example such as the one above. You can have more if you like, but it serves no purpose and may trip you up.

### 5.3.1.2 ADFs and ERCs

Each GPNode in the grammar is looked up according to its name. Thus if you have a grammar element called (*sin...*), you’ll need to have a GPNode whose name() method returns sin. This goes for ERCs and ADFs as well. Typically the name() of an ERC is simply ERC, the name() of an ADF is something like ADF1, and the name() of an ADFArgument is often something like ARG0. Thus you might have a grammar that looks like:

```

<start> ::= <op>
<op> ::= (x) | (y) | (sin <op>) | (* <op> <op>) | (- <op> <op>)
<op> ::= (ERC)
<op> ::= (ADF1 <op> <op>)

```

## 5.3.2 Translation and Evaluation

Translation is done as follows. GESpecies is handed a GEIndividual, which is little more than an arbitrary-length ByteVectorIndividual. You should set this GEIndividual to have maximal min-gene and max-gene values, so it can have any of all 256 possible settings per gene:

```

pop.subpop.0.species.min-gene = -128
pop.subpop.0.species.max-gene = 127

```

GESpecies starts working through the byte array and the grammar, depth-first. Each time GESpecies comes across a point in the grammar where multiple expansions are possible, it consults the next byte in the byte array. Let’s say there are four possible expansions of the head <foo>, and the next byte says 145. We compute  $145 \bmod 4 = 1$  and so pick expansion 1 (the second expansion — the first one is expansion 0).

This continues until one of two things happens. Either the tree is completed, or we run out of bytes. In the first case, the tree is finished and the GEProblem hands it to the GPProblem to be evaluated. In the second case, the fitness is simply set to the worst possible Koza Standardized Fitness (Float.MAX\_VALUE).

If the GEIndividual is valid and the GPIndividual is then set to the GPProblem to be evaluated, after evaluation, the GEIndividual’s fitness is set to the GPIndividual’s fitness, the GEIndividual’s evaluated flag is set to the GPIndividual’s flag, and evaluation of the GEIndividual is completed.

To do the translation, GESpecies relies on certain methods that may be useful to you:

### ec.gp.ge.GESpecies Methods

---

```

public int makeTree(EvolutionState state, GEIndividual ind, GPTree tree, int position, int treeNum,
int threadnum)

```

Builds a GPTree from a GEIndividual. The tree is stored in *tree*. The tree number (and thus the particular grammar to be used) is defined by *treeNum*. Bytes are read from the GEIndividual starting at *position*. When the tree has

been generated, the first unread byte position is returned. If the tree could not be built because there were not enough bytes, then `ec.gp.ge.GESpecies.BIG_TREE_ERROR` is returned instead.

```
public int makeTrees(EvolutionState state, GEIndividual ind, GPTree[] trees, int threadnum)
```

Builds an entire array of GPTrees, sufficient to create an entire GPIndividual, from a GEIndividual. The trees are stored in *trees*. Bytes are read from the GEIndividual starting at position 0. When the tree has been generated, the first unread byte position is returned. If the trees could not be built because there were not enough bytes, then `ec.gp.ge.GESpecies.BIG_TREE_ERROR` is returned instead.

```
public int consumed(EvolutionState state, GEIndividual ind, int threadnum)
```

Computes and returns the number of bytes that would be consumed to produce a GPIndividual from the given GEIndividual. This is done by actually building an individual, then throwing it away. If the GPIndividual could not be built because there were not enough bytes, then `ec.gp.ge.GESpecies.BIG_TREE_ERROR` is returned instead.

---

**Handling ERCs** When the GESpecies needs to produce an ERC from the grammar, it consults the next byte in the array. If the byte is 192, it looks up 192 in a special hash table stored in GESpecies. If it finds that 192 has hashed to an existing ERC, it clones this ERC and uses the clone. Else, it creates a new ERC, calls `reset()` on it, stores it in the hash table with 192, clones it, and uses the clone. This way all ERCs whose byte is 192 are the same value.

**Handling ADFs and Multiple Trees** Recall that the GESpecies maintains a separate grammar for each tree in the GPIndividual (and thus for each separate ADF). However there's only a single byte array in the GEIndividual. This is handled straightforwardly the array is used to build the first tree; then the unused remainder of the array is used to build the second tree; and so on. If not all trees were able to be built, evaluation of the GPIndividual is bypassed and the fitness is set to the `(Float.MAX_VALUE)`.

**Grammatical Evolution Does Not Support GroupedProblemForm** GroupedProblemForm (Section 7.1.2) evaluates several Individuals together. What if at least one of them is a GEIndividual, and the GEIndividual cannot generate a valid GPIndividual? This creates a number of hassles. For example, if we're doing cooperative coevolution, and one Individual can't be generated, what's the fitness of the group? Is it fair to penalize the group thusly? Likewise if we're doing competitive coevolution, and one of the competitors can't be generated, what's the resulting fitness of the other? For this reason, at present Grammatical Evolution does not support GroupedProblemForm.

### 5.3.3 Initialization and Breeding

Since we're doing lists, we'll need to define the parameters for creating new lists in the first place. The default parameters in `ge.params` do geometric size distribution as follows:

```
pop.subpop.0.species.genome-size = geometric
pop.subpop.0.species.geometric-prob = 0.85
pop.subpop.0.species.min-initial-size = 5
```

Change this to your heart's content.

While we could use things like `ListCrossoverPipeline`, GE has two idiosyncratic list breeding operators:

- `ec.vector.breed.GeneDuplicationPipeline` picks two random indices in the list, copies the byte sequence between them, then tacks the copy to the end of the list. Because this doesn't rely on anything special to GE, it's located in `ec.breed.vector`.
- `ec.gp.ge.breed.GETruncationPipeline` determines how many bytes were consumed in the production of the GPTree. It then truncates the list to remove the unused bytes.



The default pipeline in `ge.params` is a `MultiBreedingPipeline` which with 0.9 probability performs `GETruncation`, followed by crossover; and with 0.05 probability performs `GETruncation` followed by `GeneDuplication`; and 0.05 probability simply does plain `VectorMutation`. In all cases we use `Tournament Selection` with size of 7.

```
pop.subpop.0.species.pipe = ec.breed.MultiBreedingPipeline
pop.subpop.0.species.pipe.num-sources = 3

pop.subpop.0.species.pipe.source.0 = ec.vector.breed.ListCrossoverPipeline
pop.subpop.0.species.pipe.source.0.source.0 = ec.gp.ge.breed.GETruncationPipeline
pop.subpop.0.species.pipe.source.0.source.0.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.0.source.1 = same
pop.subpop.0.species.pipe.source.0.prob = 0.9

pop.subpop.0.species.pipe.source.1 = ec.vector.breed.GeneDuplicationPipeline
pop.subpop.0.species.pipe.source.1.source.0 = ec.gp.ge.breed.GETruncationPipeline
pop.subpop.0.species.pipe.source.1.source.0.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.1.prob = 0.05

pop.subpop.0.species.pipe.source.2 = ec.vector.breed.VectorMutationPipeline
pop.subpop.0.species.pipe.source.2.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.2.prob = 0.05

select.tournament.size = 7
```

You are of course welcome to change this any way you like.

The default pipeline also defines the mutation probability for `VectorMutationPipeline`, and includes a crossover type even though it's unused (to quiet complaints from ECJ):

```
pop.subpop.0.species.mutation-prob = 0.01
# This isn't used at all but we include it here to quiet a warning from ECJ
pop.subpop.0.species.crossover-type = one
```

### 5.3.4 Dealing with GP

Once you've created your grammar and set up your `GEIndividual` and `GESpecies`, you'll still need to define all the same `GPNodes`, `GPFunctionSets`, `GPNodeConstraints`, `GPTreeConstraints`, etc. as usual. All GE is doing is giving you a new way to breed and create trees: but the tree information must still remain intact. However, certain items must be redefined because the `GPSpecies` is no longer in the normal parameter base (`pop.subpop.0.species`) but rather is the subsidiary to `GESpecies`, and thus at the parameter base (`pop.subpop.0.species.gp-species`). The simplest way to do this is to include `ec/gp/koza/koza.params` to define all the basic GP stuff, then create all of your `GPNodes` etc., and then override certain GP parameters, namely:

```

# We define a dummy KozaFitness here, and set the number of trees to 1.
# If you're doing ADFs, you'll need to add some more trees.
pop.subpop.0.species.gp-species = ec.gp.GPSpecies
pop.subpop.0.species.gp-species.fitness = ec.gp.koza.KozaFitness
pop.subpop.0.species.gp-species.ind = ec.gp.GPIndividual
pop.subpop.0.species.gp-species.ind.numtrees = 1
pop.subpop.0.species.gp-species.ind.tree.0 = ec.gp.GPTree
pop.subpop.0.species.gp-species.ind.tree.0.tc = tc0

# We also need a simple dummy breeding pipeline for GP, which will never
# be used, but if it's not here GP will complain. We'll just use Reproduction.
pop.subpop.0.species.gp-species.pipe = ec.breed.ReproductionPipeline
pop.subpop.0.species.gp-species.pipe.num-sources = 1
pop.subpop.0.species.gp-species.pipe.source.0 = ec.select.TournamentSelection

```

This is enough to convince the GP system to go along with our bizarre plans.

### 5.3.5 A Complete Example

We will continue the example given in Section 5.2.11. We don't include the parameters and Java files specified so far in that example, but you'll of course need them.

That example showed how to do a strongly-typed Individual with various functions, plus an ADF and an ERC. The ADF points to a tree that contains the same basic functions, plus another ERC and two ADFArguments. Keep in mind that strong typing just gets in our way, but in this example, it's fairly harmless. We use all the existing parameters, but change a few. First, let's do the most of the ones we've discussed so far:

```

# Basic parameters that we redefine
eval.problem = ec.gp.ge.GEProblem
pop.subpop.0.species = ec.gp.ge.GESpecies
pop.subpop.0.species.parser = ec.gp.ge.GrammarParser
pop.subpop.0.species.gp-species = ec.gp.GPSpecies
pop.subpop.0.species.fitness = ec.gp.koza.KozaFitness
pop.subpop.0.species.ind = ec.gp.ge.GEIndividual
pop.subpop.0.species.min-gene = -128
pop.subpop.0.species.max-gene = 127
pop.subpop.0.species.mutation-prob = 0.01
pop.subpop.0.species.crossover-type = one
pop.subpop.0.species.genome-size = geometric
pop.subpop.0.species.geometric-prob = 0.85
pop.subpop.0.species.min-initial-size = 5

# The pipeline
pop.subpop.0.species.pipe = ec.breed.MultiBreedingPipeline
pop.subpop.0.species.pipe.num-sources = 3
pop.subpop.0.species.pipe.source.0 = ec.vector.breed.ListCrossoverPipeline
pop.subpop.0.species.pipe.source.0.source.0 = ec.gp.ge.breed.GETruncationPipeline
pop.subpop.0.species.pipe.source.0.source.0.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.0.source.1 = same
pop.subpop.0.species.pipe.source.0.prob = 0.9
pop.subpop.0.species.pipe.source.1 = ec.vector.breed.GeneDuplicationPipeline
pop.subpop.0.species.pipe.source.1.source.0 = ec.gp.ge.breed.GETruncationPipeline
pop.subpop.0.species.pipe.source.1.source.0.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.1.prob = 0.05
pop.subpop.0.species.pipe.source.2 = ec.vector.breed.VectorMutationPipeline
pop.subpop.0.species.pipe.source.2.source.0 = ec.select.TournamentSelection
pop.subpop.0.species.pipe.source.2.prob = 0.05
select.tournament.size = 7

# GP hacks
pop.subpop.0.species.gp-species.fitness = ec.gp.koza.KozaFitness
pop.subpop.0.species.gp-species.ind = ec.gp.GPIndividual
pop.subpop.0.species.gp-species.pipe = ec.breed.ReproductionPipeline
pop.subpop.0.species.gp-species.pipe.num-sources = 1
pop.subpop.0.species.gp-species.pipe.source.0 = ec.select.TournamentSelection

```

You don't have to specify all this: just include `ge.params` as a parent of your parameter file. Since we're doing ADFs, we'll need **two trees** rather than just one. `ge.params` by default defines a single tree. We do two here:

```

# More GP hacks to handle an ADF
pop.subpop.0.species.gp-species.ind.numtrees = 2
pop.subpop.0.species.gp-species.ind.tree.0 = ec.gp.GPTree
pop.subpop.0.species.gp-species.ind.tree.0.tc = tc0
pop.subpop.0.species.gp-species.ind.tree.1 = ec.gp.GPTree
pop.subpop.0.species.gp-species.ind.tree.1.tc = tc1

```

Next let's hook up the Problem, which was originally called `ec.app.myapp.MyProblem`:

```
# The problem
eval.problem.problem = ec.app.myapp.MyProblem
eval.problem.problem.data = ec.app.myapp.MyData
```

Now for each tree (the main tree and the ADF) we need to define the grammar files:

```
# The grammars
ge.species.file.0 = myproblem.grammar
ge.species.file.1 = adf0.grammar
```

### 5.3.5.1 Grammar Files

The myproblem.grammar file defines our basic functions (both boolean and real-valued) and also our ADF and ERC.

```
# This is the grammar file for the main tree
<start> ::= <op>
<op> ::= (x) | (y) | (sin <op>) | (* <op> <op>) | (- <op> <op>)
<op> ::= (ERC)
<op> ::= (ADF <op> <op>)
<op> ::= (if <bool> <op> <op>)
<bool> ::= (nand <bool> <bool>)
<bool> ::= (> <op> <op>)
```

The adf0.grammar file uses just the real-valued functions, plus an ERC and two ADFArguments:

```
# This is the grammar file for ADF0, which requires two arguments
<start> ::= <op>
<op> ::= (x) | (y) | (sin <op>) | (* <op> <op>) | (- <op> <op>)
<op> ::= (ERC)
<op> ::= (ARG0) | (ARG1)
```

## 5.3.6 How Parsing is Done

GESpecies parses the grammar files by running them through an ec.gp.ge.GrammarParser, a special class which converts the files into parse graphs. The parse graphs consist of two kinds of nodes: a ec.gp.ge.GrammarFunctionNode and a ec.gp.ge.GrammarRuleNode, both subclasses of the abstract class ec.gp.ge.GrammarNode. You can replace ec.gp.ge.GrammarParser with your own subclass if you like, via:

```
pop.subpop.0.species.parser = ec.app.MySpecialGrammarParser
```

The parse graph produced by a GrammarParser is rooted with a GrammarRuleNode which indicates the initial rule. For example, reconsider the grammar below.

```
<start> ::= <op>
<op> ::= (x) | (y) | (sin <op>) | (* <op> <op>) | (- <op> <op>)
<op> ::= (if <bool> <op> <op>)
<bool> ::= (nand <bool> <bool>)
<bool> ::= (> <op> <op>)
```

The parse graph for this grammar is shown in Figure 5.10. Once the parse graph has been generated, the GESpecies wanders through this graph as follows:



Figure 5.10 Parse graph (orderings among children not shown). The entry point is `<start>`. GrammarRuleNodes are shown with angle brackets. GrammarFunctionNodes are shown without brackets.

- If on a GrammarRuleNode, the GESpecies selects exactly one child to traverse, using the next number in the GEIndividual array.
- If on a GrammarFunctionNode, the GESpecies traverses each and every child of the node in order. These become children to the GrammarFunctionNode's GPNode.

GrammarFunctionNodes hold GPNode prototypes and represent them. Their children are arguments to the node in question. GrammarRuleNodes hold choice points in the rule, and their children are the choices (only one choice will be made).

## 5.4 Rulesets and Collections (The `ec.rule` Package)

Let's get one thing out of the way right now. Though we had rulesets in mind when we developed it, the `ec.rule` package isn't *really* for rulesets. Not only can the package be used for things other than rules, but it's not even sets: it's *collections* (or "bags" or "multisets") of arbitrary objects.

The representation defined by this package is fairly straightforward: an `ec.rule.RuleIndividual` contains one or more `ec.rule.RuleSets`, each of which contain zero or more `ec.rule.Rules`. A Rule is an abstract superclass which can contain anything you want. And that's about it.

Problem domains for which the `ec.rule` package is appropriate are often also good candidates for the variable-length lists found in the `ec.vector` package. You'll need to think about which is a better choice for you. Also beware that of the various representation packages in ECJ, `ec.rule` is definitely the least used and least tested. So its facilities are somewhat cruder than the others and it's possible you may see bugs.

Each level of the `ec.rule` package (individual, ruleset, rule) is a Prototype and has a Flyweight relationship with a central object special to that level (for a reminder on Flyweights, see Section 3.1.4). Specifically:

Object	In Flyweight Relationship With
<code>ec.rule.RuleIndividual</code>	<code>ec.rule.RuleSpecies</code>
<code>ec.rule.RuleSet</code>	<code>ec.rule.RuleSetConstraints</code>
<code>ec.rule.Rule</code>	<code>ec.rule.RuleConstraints</code>

The `ec.rule` package follows the same approach as the `ec.vector` package does when it comes to breeding: two basic breeding operators are provided (`ec.rule.breed.RuleCrossoverPipeline` and `ec.rule.breed.RuleMutationPipeline`) which simply call default mutation and crossover functions in the `Individuals` themselves. Thus to do more sophisticated breeding you have the choice of either overriding these functions or creating new breeding pipelines which perform more detailed operations on their own.

### 5.4.1 RuleIndividuals and RuleSpecies

`RuleIndividuals` and `RuleSpecies` are specified in parameters in the standard way:

```
pop.subpop.0.species = ec.rule.RuleSpecies
pop.subpop.0.species.ind = ec.rule.RuleIndividual
```

A `RuleIndividual` is a subclass of `Individual` which simply consists of an array of `RuleSets`:

```
public RuleSet[] rulesets;
```

Each `RuleSet` can be a different class. You'd think that the number and class of `RuleSets` was specified in the `RuleSpecies` (like in `ec.vector`). But for no good reason it's not the case: you specify them in the parameters for the prototypical individual, along these lines:

```
pop.subpop.0.species.ind.num-rulesets = 2
pop.subpop.0.species.ind.ruleset.0 = ec.rule.RuleSet
pop.subpop.0.species.ind.ruleset.1 = ec.app.myapp.MyRuleSet
```

Alternatively, you can use the `RuleIndividual`'s default parameter base:

```
rule.individual.num-rulesets = 2
rule.individual.ruleset.0 = ec.rule.RuleSet
rule.individual.ruleset.1 = ec.app.myapp.MyRuleSet
```

Though for many applications you will probably just have a single `RuleSet`, and it'll probably just be an `ec.rule.RuleSet`:

```
pop.subpop.0.species.ind.num-rulesets = 1
pop.subpop.0.species.ind.ruleset.0 = ec.rule.RuleSet
```

### 5.4.2 RuleSets and RuleSetConstraints

A `RuleSet` contains an arbitrary number of `Rules` (`ec.rule.Rule`), anywhere between zero and up. It's largely your job to customize the breeding and initialization procedures appropriate to your problem to constraint the number and type of rules. They're defined here:

```
public Rule[] rules;
public int numRules;
```

Notice that the number of rules in the array may be *less* than the array size, that is, `numRules`  $\leq$  `rules.length`. The rules themselves run from `rules[0]` ... `rules[numRules - 1]`. This is done because, like `ArrayList` etc., `rules` is variable in size and can grow and shrink. `RuleSet` contains a number of utility methods for manipulating the order and number of these rules:

#### **ec.rule.RuleSet Methods**

---

```
public int numRules()
    Returns the number of rules in the RuleSet.
```

```

public void randomizeRulesOrder(EvolutionState state, int thread)
    An auxillary debugging method which verifies many features of the structure of the GPTree and all of its GPNodes.
    This method isn't called by ECJ but has proven useful in determining errors in GPTree construction by various
    tree building or breeding algorithms.

public void addRule(Rule rule)
    Adds the rule to the end of the ruleset, increasing the length of the RuleSet array as necessary.

public void addRandomRule(EvolutionState state, int thread)
    Produces a new randomly-generated rule and adds it to the RuleSet. The Rule is created by cloning the prototypical
    Rule from the RuleSet's RuleSetConstraints, then calling reset(...) on it.

public Rule removeRule(int index)
    Removes a rule located at the given index from the RuleSet and returns it. All rules are shifted down to fill the
    void.

public Rule removeRandomRule(EvolutionState state, int thread)
    Removes a random rule returns it. All rules are shifted down to fill the void.

public RuleSet[] split(int[] points, RuleSet[] sets)
    Breaks the RuleSet into n disjoint groups, then clones the rules in those groups and adds them to the respective
    sets, which must be provided in the given array. The first group of rules starts at 0 and ends below points[0]: this
    goes into sets[0]. Intermediate groups, which go into sets[i], start at points[i] and end below points[i+1]. The final
    group, which goes into sets[points.length], starts at points[points.length - 1] and continues to the end of the rule
    array. If points.length = 0, then all rules simply get put into sets[0]. Note that the size of sets must be one more
    than the size of points. The sets are returned.

public RuleSet[] split(EvolutionState state, int thread, RuleSet[] sets)
    For each rule in the RuleSet, clones the rule and adds the clone to a randomly chosen RuleSet from sets. Returns
    sets.

public RuleSet[] splitIntoTwo(EvolutionState state, int thread, RuleSet[] sets, float probability)
    For each rule in the RuleSet, clones the rule and, with the given probability, adds the clone to sets[0], else sets[1].
    Note that sets must be two in length.

public void join(RuleSet other)
    Copies the rules in the other RuleSet, then adds them to the end of this RuleSet.

public void copyNoClone(RuleSet other)
    Deletes all the rules in the RuleSet. Then places all the rules from the other RuleSet into this RuleSet. No cloning is
    done: both RuleSets now have pointers to the same rules.

```

---

Groups of RuleSets have a flyweight relationship with a RuleSetConstraints object. RuleSetConstraints is a Clique. You specify the number and class of RuleSetConstraints and assign each a unique name. Let's say you need two RuleSetConstraints objects, both instances of RuleSetConstraints itself. You'd write this:

```

rule.rsc.size = 2
rule.rsc.0 = ec.rule.RuleSetConstraints
rule.rsc.0.name = rsc1
rule.rsc.1 = ec.rule.RuleSetConstraints
rule.rsc.0.name = rsc2

```

RuleSetConstraints specify a number of constraints which guide the initialization and mutation of RuleSets, specifically:

- A distribution for choosing the number of Rules an initial RuleSet will have. This guides how the RuleSet's reset(...) method operates. The distribution can either be uniform with a minimum and

maximum, or you can specify a histogram of possible size probabilities. We'll do the first case for Ruleset 0 and the second case for Ruleset 1 below:

```
# RuleSetConstraints 0 will have between 5 and 10 rules inclusive
rule.rsc.0.reset-min-size = 5
rule.rsc.0.reset-max-size = 10

# RuleSetConstraints 1 will have 0 to 4 rules with these probabilities...
rule.rsc.1.reset-num-sizes = 5
rule.rsc.1.size.0 = 0.1
rule.rsc.1.size.1 = 0.2
rule.rsc.1.size.2 = 0.2
rule.rsc.1.size.3 = 0.3
rule.rsc.1.size.4 = 0.4
```

- The probability of adding, deleting, and rearranging rules, when the RuleSet's `mutate(...)` method is called, typically by the BreedingPipeline `ec.rule.RuleMutationPipeline`. When this method is called, the Ruleset first mutates all of its rules by calling `mutate(...)` on them. Then it repeatedly flips a coin of a given probability: each time the coin comes up true, or until the number of rules is equal to the minimum number of initial rules as specified above, one rule is deleted. Afterwards it repeatedly flips a coin of another probability: each time the coin comes up true, or until the number of rules is equal to the maximum number of initial rules, one new rule is added at random. Finally, with a certain probability the rule ordering is shuffled. Here's some examples of specifying these probabilities:

```
rule.rsc.0.p-add = 0.1
rule.rsc.0.p-del = 0.1
rule.rsc.0.rand-order = 0.25

rule.rsc.1.p-add = 0.5
rule.rsc.1.p-del = 0.6
rule.rsc.1.rand-order = 0.0
```

Once you've specified a RuleSetConstraints, you then attach one to each RuleSet. For example:

```
pop.subpop.0.species.ind.ruleset.0.constraints = rsc2
pop.subpop.0.species.ind.ruleset.1.constraints = rsc1
```

... or alternatively use the default parameter base...

```
rule.individual.constraints = rsc2
```

Once set, you can access the constraints with the following method:

#### **ec.rule.RuleSet Methods**

---

```
public final RuleSetConstraints ruleSetConstraints(RuleInitializer initializer)
    Returns the RuleSet's RuleSetConstraints
```

---

RuleSetConstraints one method for choosing random initial values under the constraints above:

#### **ec.rule.RuleSetConstraints Methods**

---

```
public int numRulesForReset(RuleSet ruleset, EvolutionState state, int thread)
    Returns a random value from the initial (reset(...)) distribution, to use as the number of rules to initialize or reset the RuleSet.
```



---

Additionally, the various addition, deletion, and randomization probabilities can be accessed like this:

```
RuleSetConstraints rsc = myRuleSet.ruleSetConstraints((RuleInitializer)(state.init));
float addition = rsc.p_add;
float deletion = rsc.p_del;
float shuffling = rsc.p_randorder;
```

RuleSets also contain all the standard reading and writing methods, none of which you'll need to override unless you're making a custom RuleSet.

---

### ec.rule.RuleSet Methods

```
public void printRuleSetForHumans(EvolutionState state, int log)
    Writes a RuleSet to a log in a fashion easy for humans to read.

public void printRuleSet(EvolutionState state, int log)
    Writes a RuleSet to a log in a fashion that can be read back in via readRule(...), typically by using the Code package.

public void printRuleSet(EvolutionState state, PrintWriter writer)
    Writes a RuleSet to a Writer in a fashion that can be read back in via readRule(...), typically by using the code package.

public void readRuleSet(EvolutionState state, LineNumberReader reader) throws IOException
    Reads a RuleSet written by printRuleSet(...) or printRuleSetToString(...), typically using the Code package.

public void writeRuleSet(EvolutionState state, DataOutput output) throws IOException
    Writes a RuleSet in binary fashion to the given output.

public void readRuleSet(EvolutionState state, DataInput input) throws IOException
    Reads a RuleSet in binary fashion from the given input.
```

---

### 5.4.3 Rules and RuleConstraints

RuleSetConstraints also contain the **prototypical Rule** for RuleSets adhering to a given constraints. RuleSets will clone this Rule to create Rules to fill themselves with. `ec.rule.Rule` is an abstract superclass which doesn't do anything by itself: you're required to subclass it to make the Rule into the kind of thing you want to create a collection of in your Ruleset.

The prototypical Rule is specified like this:

```
pop.subpop.0.species.ind.ruleset.0.rule = ec.app.MyRule
pop.subpop.0.species.ind.ruleset.1.rule = ec.app.MyOtherRule
```

You can get the prototypical rule like this:

```
RuleSetConstraints rsc = myRuleSet.ruleSetConstraints((RuleInitializer)(state.init));
Rule prototype = rsc.rulePrototype;
```

Each Rule has a flyweight-related RuleConstraints object, which is defined similarly to RuleSetConstraints (it's also a Clique). For example, to create a single RuleConstraints in the clique, you might say:

```
rule.rc.size = 1
rule.rc.0 = ec.rule.RuleConstraints
rule.rc.0.name = rc1
```

RuleConstraints are essentially blank: they define no special parameters or variables. You can use them however you see fit. If you don't really care, you can just make a single RuleConstraints object as above and assign it to your prototypical rules, such as:

```
pop.subpop.0.species.ind.ruleset.0.rule.constraints = rc1
pop.subpop.0.species.ind.ruleset.1.rule.constraints = rc1
```

...or use the default parameter base:

```
rule.rule.constraints = rc1
```

A Rule is abstract, and so has certain abstract methods which must be overridden, as well as others which *ought* to be overridden. First the required ones:

---

#### ec.rule.Rule Methods

```
public abstract int hashCode()
```

Returns a hash code for the rule, based on value, suitable for weeding out duplicates.

```
public abstract int compareTo(Object other)
```

Returns 0 if this Rule is identical in value to *other* (which will also be a Rule), -1 if this Rule is "less" than the other rule in sorting order, and 1 if the Rule is "greater" than the other rule in sorting order.

```
public abstract void reset(EvolutionState state, int thread)
```

Randomizes the value of the rule.

---

Rules are Prototypes and so implement the clone(), setup(...), and defaultBase() methods. You'll most likely need to override the clone() and setup(...) methods as usual. Additionally, you may want to override:

---

#### ec.rule.Rule Methods

```
public void mutate(EvolutionState state, int thread)
```

Mutates the Rule in some way and with some probability. The default implementation simply calls reset(...), which is probably much too harsh.

```
public abstract int compareTo(Object other)
```

Returns 0 if this Rule is identical in value to *other* (which will also be a Rule), -1 if this Rule is "less" than the other rule in sorting order, and 1 if the Rule is "greater" than the other rule in sorting order.

```
public abstract void reset(EvolutionState state, int thread)
```

Randomizes the value of the rule.

---

Then there are the standard printing and reading methods. You'll need to override at least printRuleToStringForHumans(), and probably will want to override toString(). The others you can optionally override depending on the kind of experiments you're doing.

---

#### ec.rule.Rule Methods

```
public String toString()
```

Writes the Rule to a String, typically in a fashion that can be read back in via readRule(...). You'll want to override this method or printRuleToString(). You probably want to use the Code package to write the rule out. You only really need to implement this method if you expect to write Individuals to files that will be read back in later.

```
public String printRuleToStringForHumans()
```

Writes the Rule to a String in a fashion easy for humans to read. The default implementation of this method simply calls toString(). You'll probably want to override this method.

```
public void printRuleForHumans(EvolutionState state, int log)
    Writes a Rule to a log in a fashion easy for humans to read. The default implementation of this method calls
    printRuleToStringForHumans(), which you should probably override instead.
```

```
public String printRuleToString()
    Writes the Rule to a String, typically in a fashion that can be read back in via readRule(...). The default implemen-
    tation of this method simply calls toString(). You'll want to override this method or toString(). You probably want
    to use the Code package to write the rule out. You only need to implement this method if you expect to write
    Individuals to files that will be read back in later.
```

```
public void printRule(EvolutionState state, int log)
    Writes a Rule to a log in a fashion that can be read back in via readRule(...). The default implementation of this
    method calls printRuleToString(), which you should probably override instead.
```

```
public void printRule(EvolutionState state, PrintWriter writer)
    Writes a Rule to a Writer in a fashion that can be read back in via readRule(...). The default implementation of this
    method calls printRuleToString(), which you should probably override instead.
```

```
public void readRule(EvolutionState state, LineNumberReader reader) throws IOException
    Reads a Rule written by printRule(...) or printRuleToString(...), typically using the Code package. The default does
    nothing. You only need to implement this method if you expect to read Individuals from files.
```

```
public void writeRule(EvolutionState state, DataOutput output) throws IOException
    Writes a Rule in binary fashion to the given output. The default does nothing. You only need to implement this
    method if you expect to read and write Rules over a network (such as the distributed evaluation or island models).
```

```
public void readRule(EvolutionState state, DataInput input) throws IOException
    Reads a Rule in binary fashion from the given input. The default signals an error. You only need to implement this
    method if you expect to read and write Rules over a network (such as the distributed evaluation or island models).
```

---

## 5.4.4 Initialization

Basic Initialization works as follows:

1. The RuleSpecies method `newIndividual(EvolutionState, int)` produces a RuleIndividual by calling `super.newIndividual(...)`—cloning a RuleIndividual prototype—and then calling `reset(...)` on the resultant RuleIndividual.
2. The RuleIndividual's `reset(...)` by default just calls `reset(...)` on each of the RuleSets.
3. A RuleSet's `reset(...)` method calls `numRulesForReset(...)` on the RuleSetConstraints to pick a random number of rules to generate (see Section 5.4.2). It then produces an array of that size and fills it with rules cloned from the RuleSetConstraint's prototypical Rule. Then it calls `reset(...)` on each of the Rules.
4. You are responsible for implementing a Rule's `reset(...)` method.

You can of course intervene and modify any of these methods as you see fit.

## 5.4.5 Mutation

As in the case in the `ec.vector` package, the `ec.rule.breed.RuleMutationPipeline` class doesn't mutate rules directly, but rather calls a method on them to ask them to mutate themselves. The procedure is as follows:

1. The RuleMutationPipeline calls `preprocessIndividual(...)` on the RuleIndividual.
2. The RuleIndividual's `preprocessIndividual(...)` method calls `preprocessRules(...)` on each of the RuleSets.

3. The RuleSet's preprocessRules(...) method by default does nothing: override it as you like.
4. The RuleMutationPipeline then calls mutate(...) on the RuleIndividual.
5. The RuleIndividual's mutate(...) method by default just calls mutate(...) on each of its RuleSets.
6. The RuleSet's mutate(...) method does several modifications to the rules in the RuleSet, in this order:
  - (a) All the Rules in the RuleSet have mutate(...) called on them.
  - (b) A coin is repeatedly flipped of a certain probability (p\_del), and each time it comes up true, a rule is deleted at random using removeRandomRule(...). The individual will not shrink smaller than its specified minimum size.
  - (c) A coin is repeatedly flipped of a certain probability (p\_add), and each time it comes up true, a rule is added at random using addRandomRule(...). That method clones a new Rule from the prototypical Rule, then calls reset(...) on it. The individual will not grow larger than its specified maximum size.
  - (d) With a certain probability (p\_randorder), the order of the rules is shuffled using randomizeRulesOrder(...).

The three probabilities (p\_del, p\_add, and p\_randorder), and the minimum and maximum rule sizes, are discussed in Section 5.4.2, and are determined by RuleSetConstraints parameters, also discussed in that Section.

7. A Rule's mutate(...) method by default simply calls reset(...), which is probably not what you want. You'll probably want a much more subtle mutation if any, and so will need to override this method.
8. You are responsible for implementing a Rule's reset(...) method.
9. Finally, the RuleMutationPipeline calls postprocessIndividual(...) on the RuleIndividual.
10. The RuleIndividual's postprocessIndividual(...) method calls postprocessRules(...) on each of the RuleSets.
11. The RuleSet's postprocessRules(...) method by default does nothing: override it as you like.

Often rules need to be in a carefully-constructed dance of constraints to be valid in an Individual. The intent of the preprocessIndividual(...), postprocessIndividual(...), preprocessRules(...), and postProcessRules(...) methods is to give your RuleIndividual a chance to fix RuleSets that have been broken by crossover or mutation. The default implementation of these methods doesn't do much:

---

#### **ec.rule.RuleSet Methods**

public void preprocessRules(EvolutionState state, int thread)

A hook called prior to mutation or crossover to prepare for possible breakage of Rules due to the mutation or crossover event. The default implementation does nothing.

public void postprocessRules(EvolutionState state, int thread)

A hook called after to mutation or crossover to fix possible breakage of Rules due to the mutation or crossover event. The default implementation does nothing.

---



---

#### **ec.rule.RuleIndividual Methods**

public void preprocessIndividual(EvolutionState state, int thread)

Calls preprocessRules(...) on each RuleSet in the Individual.

public void postprocessIndividual(EvolutionState state, int thread)

Calls postprocessRules(...) on each RuleSet in the Individual.

---

### 5.4.6 Crossover

Unlike RuleMutationPipeline, the `ec.rule.breed.RuleCrossoverPipeline` performs direct crossover on two RuleIndividuals. Here is the procedure:

1. The RuleCrossoverPipeline calls `preprocessIndividual(...)` on each RuleIndividual.
2. The RuleIndividual's `preprocessIndividual(...)` method calls `preprocessRules(...)` on each of the RuleSets.
3. The RuleSet's `preprocessRules(...)` method by default does nothing: override it as you like.
4. For each pair of RuleSets, one per RuleIndividual...
  - (a) Each RuleSet  $A$  and  $B$  is split into two pieces,  $A_1$  and  $A_2$  (and  $B_1$  and  $B_2$ ) by calling `splitIntoTwo(...)`
  - (b) A new RuleSet  $A'$  is formed from the union of  $A_1$  and  $B_1$ , and likewise, a new RuleSet  $B'$  is formed from the union of  $A_2$  and  $B_2$ .
  - (c) If  $A'$  and  $B'$  do not minimum and maximum size constraints (see Section 5.4.2), go to (a) and try again.
  - (d) Else  $A'$  and  $B'$  replace  $A$  and  $B$  respectively in each RuleIndividual.
5. Finally, the RuleCrossoverPipeline calls `postprocessIndividual(...)` on each RuleIndividual.
6. The RuleIndividual's `postprocessIndividual(...)` method calls `postprocessRules(...)` on each of the RuleSets.
7. The RuleSet's `postprocessRules(...)` method by default does nothing: override it as you like.

RuleCrossoverPipeline has a few parameters which guide its operation. First, any given Rule will migrate from one Individual's RuleSet to the other only with a certain probability. Second, the CrossoverPipeline can be set up to return only one child (tossing the second) rather than returning two. By default it returns both children. To set both of these parameters, let's say that the RuleCrossoverPipeline is the root pipeline for the species. We'd say:

```
pop.subpop.0.species.pipe = ec.rule.breed.RuleCrossoverPipeline
pop.subpop.0.species.pipe.crossover-prob = 0.1
pop.subpop.0.species.pipe.toss = true
```

It doesn't make any sense to have a rule crossover probability higher than 0.5. As usual, you could use the default parameter base as well:

```
rule.xover.crossover-prob = 0.1
rule.xover.toss = true
```



## Chapter 6

# Parallel Processes

ECJ has various built-in methods for parallelism, and they can be used in combination with one another:

- Multiple breeding and evaluation **threads**, already discussed in Section 2.4.
- **Distributed evaluation**: sending chunks of Individuals to remote computers to be evaluated. This is typically done in a generational fashion, but a variation of this is **asynchronous evolution**, in which Individuals are sent to multiple remote computers in a steady-state fashion. Additionally, remote computers can (given time) engage in a little evolutionary optimization of their own on the chunks they've received before sending them back. This is known as **opportunistic evolution**.<sup>1</sup>
- **Island models**: multiple parallel evolutionary processes occasionally send fit individuals to one another.<sup>2</sup>

### 6.1 Distributed Evaluation (The `ec.eval` Package)

Distributed Evaluation connects one **master** ECJ process with some  $N$  **slave** ECJ processes. The master handles the evolutionary loop, but when Individuals are evaluated, they are shipped off to the remote slaves to do this task. This way evaluation can be parallelized.

Distributed Evaluation is only useful if the amount of time you save by parallelizing evaluation exceeds the amount of time lost by shipping Individuals over the network (and sending at least Fitnesses back). Generally this means that evaluation needs to take a fair bit of time per Individual: perhaps several seconds.

There are two kinds of `EvolutionState` objects which can work with Distributed Evaluation:

- `SimpleEvolutionState`, which sends entire Populations off to be evaluated in parallel.
- `SteadyStateEvolutionState`, which sends individuals off to be evaluated one at a time, in a fashion called **asynchronous evolution**.

#### 6.1.1 The Master

To set up Distributed Evaluation, you first need to set up the Master. This is done just like a regular evolutionary computation process: but there are some additional parameters which must be defined. First,

---

<sup>1</sup>ECJ's built-in distributed evaluation is meant for clusters. However, Parabon Inc. has developed a grid-computing version, called Origin, which runs on hundreds of thousands or even millions of machines. See the ECJ main website for more information.

<sup>2</sup>ECJ's built-in island models are meant for clusters. However, a version of ECJ was ported to run on top of the DR-EA-M system, a peer-to-peer evolutionary computation network facility developed from a grant in Europe. See the ECJ main website for more information.



Figure 6.1 Layout of the Distributed Evaluation package.

we must define the **master problem**, nearly always as `ec.eval.MasterProblem`. The presence of the master problem turns on distributed evaluation:

```
eval.masterproblem = ec.eval.MasterProblem
```

The `MasterProblem` is the interface that connects the distributed evaluation system to a regular evolutionary computation loop. When it is defined, ECJ replaces the `Problem` prototype with a `MasterProblem` prototype. The `Problem` doesn't go away—it's rehung as a variable in the `MasterProblem`. Specifically, you can get it like this:

```
Problem originalProblem = ((MasterProblem)(state.evaluator.p_problem)).problem;
```

When your evolutionary computation process wishes to evaluate one or more individuals, it hands them to the `MasterProblem`, which it thinks is the `Problem` for the application. But the `MasterProblem` doesn't send them to a clone of your underlying `Problem` but rather routes them to the `Slave`.

Slaves register themselves over the network with an object in the Master process called a `ec.eval.SlaveMonitor`, which maintains one `ec.eval.SlaveConnection` per `Slave` to communicate with the remote `Slave`. The `SlaveMonitor` listens in on a specific socket port for incoming `Slaves` to register themselves. You'll need to define this port (to something over 2000). Here's what's standard:

```
eval.master.port = 15000
```

Your `MasterProblem` will submit `Individuals` to the `SlaveMonitor`, which will in turn direct them to one of the `SlaveConnections`. `SlaveConnections` don't just ship of single `Individuals` to `Slaves`—that would be far too inefficient use of network bandwidth. Instead, they often batch them up into **jobs** for the `Slave` to perform. Here's how you define the size of a job:

```
eval.masterproblem.job-size = 100
```

The idea is to make the job large enough to pack `Individuals` into a network packet without wasting space. If your `Individuals` are large, then a large job size won't have any efficiency benefit. If they're very



small, the job size will have a huge benefit.

You also need to keep the Slaves humming along, ideally by keeping the TCP/IP streams filled with waiting jobs queued up. Increasing this number can have a significant effect on performance. Here's a reasonable minimum:

```
eval.masterproblem.max-jobs-per-slave = 3
```

ECJ would prefer to compress the network streams to make them more efficient. But it can't do it without the `jzlib`/`ZLIB` library which you must install separately<sup>3</sup> (see the ECJ main webpage or <http://www.jcraft.com/jzlib>). Once it's installed, you can turn it on like this:

```
eval.compression = true
```

## 6.1.2 Slaves

Slaves are started up on separate CPUs, often in different machines from the Master. You can have as many Slaves as you like: the more the merrier. The `Slave` class replaces `ec.Evolve` to handle its own startup, so you don't start up a Slave using the standard `ec.Evolve` procedure. Instead you'd type:

```
java ec.eval.Slave -file slave.params
```

The slave parameters must include all the evolutionary parameters and also all the master parameters (in fact, you might as well say something like...)

```
parent.0 = master.params
```

Slaves set themselves up with their own nearly complete `EvolutionState` and `Evaluator` objects — enough to evaluate `Individuals` and also perform evolution if necessary. A slave distinguishes itself by setting a special internal parameter: `eval.i-am-slave = true`. You don't need to set this parameter — it's set programmatically by `ec.eval.Slave` when it's fired up. But you should be aware of it: it's used by `ec.Evaluator` to determine whether to replace the `Problem` with the `MasterProblem` (it needs to know if the process is a Master or a Slave, and since your Slave probably included the Master parameters — including the `eval.masterproblem` parameter — it looks confusingly like a Master).

The first thing a Slave needs to know is where the Master is so it can set itself up. This is done with something like this:

```
eval.master.host = 129.8.2.4
```

Remember that you'll also need the port (among other Master parameters!)

Next the Slave needs to know whether it should return entire `Individuals` or just the `Fitnesses` of those `Individuals`. `Individuals` are generally much bigger than `Fitnesses`, and if you only return `Fitnesses` you can cut your network traffic in half. The problem is that in some custom experiments your fitness evaluation procedure might modify the `Individual` (it depends on the nature of your experiment). You'll need to state whether to return entire `Individuals` or not:

```
eval.return-inds = false
```

Slaves can come and go at any time dynamically. If new slaves show up, the Master will immediately start taking advantage of them. If a Slave disappears, the `Individuals` it was responsible for will be reassigned to another Slave.

Slaves can be multithreaded: they can process multiple `Individuals` in parallel in the background as they come in. This is done with the standard parameter:

---

<sup>3</sup>Sure, Java has compression routines. Unfortunately they're entirely broken for network streams: they don't support "partial flush", a critical item for sending stuff across networks. They're only really useful for compressing files on disks.

```
evalthreads = 4
```

This will create four threads on the Slave process to evaluate Individuals. Note that it makes no sense for this parameter to exceed Master's `eval.masterproblem.job-size` parameter. Also, if you exceed the actual number of CPUs or cores allocated to your Slave process, it doesn't make much sense either. Last, Slaves are *not* multithreaded when evaluating Individuals with `GroupedProblemForm` (such as coevolved evaluation, see Section 7.1.2).

### 6.1.3 Opportunistic Evolution

Slaves have the option of doing some evolutionary computation of their own, a procedure known as **opportunistic evolution** [20]. The procedure works like this:

1. The Master sends the Slave a large Job.
2. The Slave evaluates the Individuals in the Job.
3. The Slave has a maximum allotted time to evaluate the Individuals. If it has not yet exceeded this time, it treats the Job as a Population and does some evolution on the Slaves.
4. When the time is up, the Slave returns the most recent Individuals in the Population in lieu of the original Individuals. The new Individuals replace the old ones in the Master's evolutionary process. This means that the Slave cannot just return Fitnesses, but must return whole Individuals.

This procedure is turned on with:

```
eval.run-evolve = true
```

You'll also need to specify the amount of time (in milliseconds) allotted to the Slave. Here's how you'd set it to six seconds:

```
eval.runtime = 6000
```

Last, if you're doing opportunistic evolution, you **must** return whole Individuals, not just Fitnesses. After all, you could have entirely different individuals after running on the Slave. Thus you'll need to set:

```
eval.return-inds = true
```

The procedure for evolution is entirely specified by the Slave's parameters just as if it were specified in a regular ECJ process, including breeding and evaluation threads, etc. There's absolutely no reason the Slave can't have its own evolutionary algorithm that's *different* from the Master's evolutionary algorithm — just specify it differently in the Slave's parameters. The only thing that'd be required is that the Slave and the Master have exactly the same kinds of Individuals, Fitnesses, and Species in their Subpopulations.

Note that Opportunistic Evolution won't work with coevolution or other procedures which require `GroupedProblemForm` (Section 7.1.2). Additionally, although Steady-State Evolution (via Asynchronous Evolution, see Section 6.1.4) can work with Opportunistic Evolution in *theory*, it'd be quite odd to do so.

### 6.1.4 Asynchronous Evolution

ECJ's distributed evaluation procedure works intuitively with generational methods (such as `ec.simple.SimpleEvolutionState` but it also works nicely with Steady-State evolution (`ec.simple.SteadyStateEvolutionState`). This procedure is called **asynchronous evolution**. See [20] for more information.

The procedure is similar to Steady State Evolution, as discussed in Section 4.2. The Population starts initially empty. Then the algorithm starts creating randomly-generated Individuals and shipping them off to

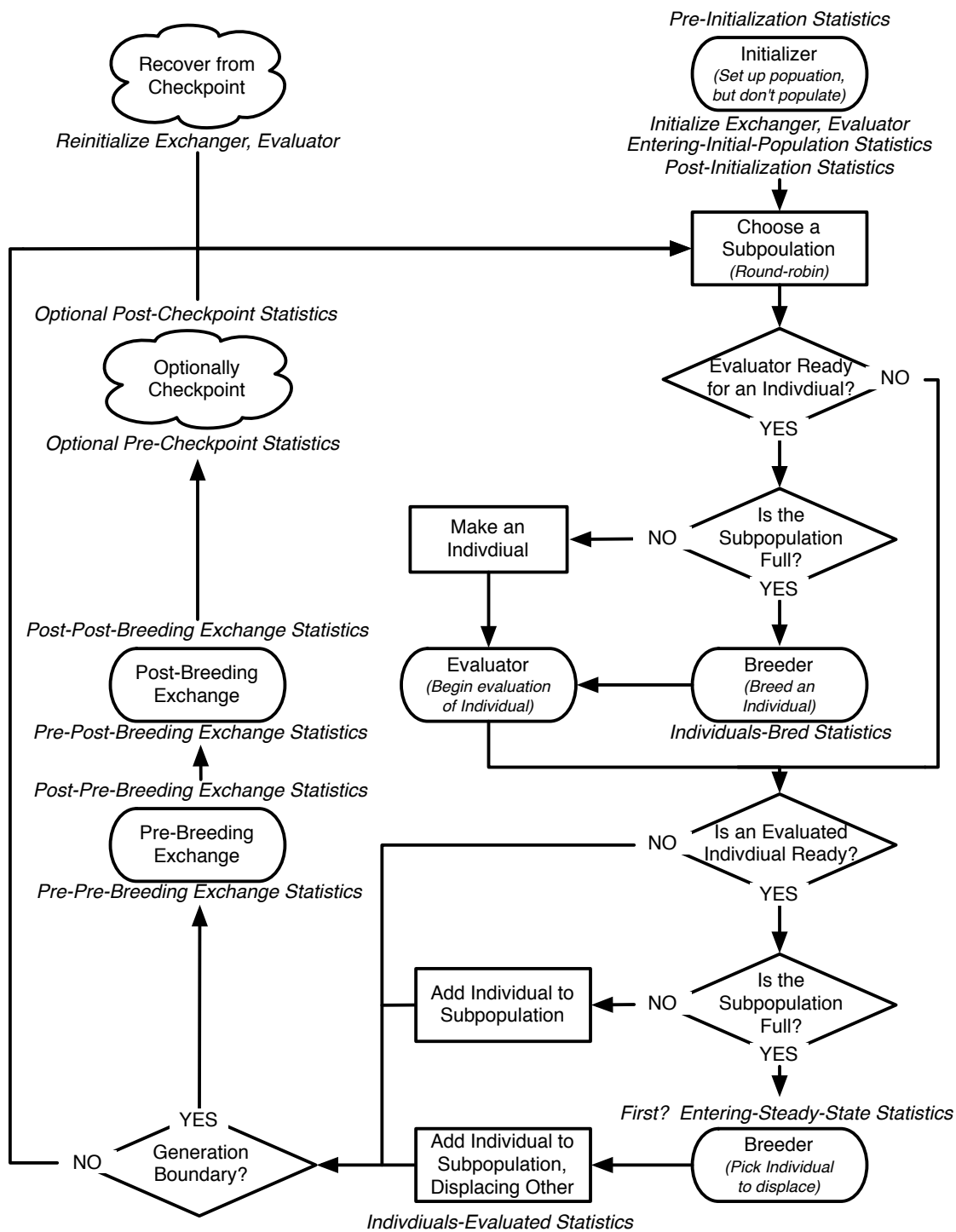


Figure 6.2 Top-Level Loop of ECJ's SteadyStateEvolutionState class, used for simple steady-state EC and Asynchronous Evolution algorithms. "First?" means to perform the Statistics whenever the Subpopulation in question is picking an Individual to displace for the very first time. (Each Subpopulation will do it once, but possibly at different times). A repeat of Figure 4.2.

remote Slaves to be evaluated. If a Slave is available, the algorithm will generate an Individual for it to work on. When a Slave has finished evaluating an Individual and has returned it (or its Fitness), the Individual is then placed into the Population.

At some point the Population will fill up. At this point the algorithm shifts to “steady state” mode. When a Slave returns an Individual, and there’s no space in the Population, the algorithm makes room by marking an existing Individual for death and replacing it with the newcomer, just like it’s done in Steady State Evolution. And when a Slave becomes available, an Individual will no longer be created at random to give to it: rather, the Individual will be bred from the existing Population.

This procedure requires some careful consideration. First, note that at the point that the algorithm shifts to “steady state” mode, there are probably a large number of Individuals being evaluated on Slaves which were not bred from the Population but were created at random. Until those Individuals have made their way into the Population, we won’t be in a true “steady state”.

Second, Steady-State Evolution assumes the production of one Individual at a time: but distributed evaluation allows more than one individual per Job. This is reconciled as follows. when Steady-State Evolution starts up, it calls `prepareToEvaluate(...)` on the Problem (the `MasterProblem`) once. Thereafter whenever an individual is sent to the Problem to be evaluated, it calls `evaluate(...)`. Recall from Section 3.4.1 that this process does not require the Problem to immediately assign Fitness—it can bulk up Individuals for evaluation and is only required to provide Fitness on or prior to a call to `finishEvaluating(...)`. However, Steady-State Evolution *never calls finishEvaluating(...)*. As a result, the distributed evaluator is free to assess Individuals in any order and any way it likes, and take as long as it wishes to assign them a Fitness. The distributed evaluator will wait for up to `job-size` worth of calls to `evaluate(...)`, then pack those Individuals together in one Job and ship them out to a remote Slave for evaluation. In “steady-state” mode, when the Individuals come back, they are placed in the Population, killing and replacing other individuals already there. Depending on the selection process for marking Individuals for death, it’s entirely possible that an Individual newly placed into the Population may be immediately marked for death and replaced with another Individual from the same Job! You can get around this by setting the `job-size` and `max-jobs-per-slave` to as low as 1:

```
eval.masterproblem.job-size = 1
eval.masterproblem.max-jobs-per-slave = 1
```

...but of course this will make the network utilization poor.

## 6.1.5 The MasterProblem

The `MasterProblem` is where much of the magic lies in the interface between ECJ and the distributed evaluator, so it’s worth mentioning some of its details.

To start, let’s discuss how `MasterProblem` handles checkpointing. Evaluators also three methods, which we didn’t discuss in Section 3.4:

```
public void initializeContacts(EvolutionState state);
public void reinitializeContacts(EvolutionState state);
public void closeContacts(EvolutionState state, int result);
```

These methods in turn call similar methods in the prototypical `Problem`:

```
public void initializeContacts(EvolutionState state);
public void reinitializeContacts(EvolutionState state);
public void closeContacts(EvolutionState state, int result);
```

result will be one of `EvolutionState.R.FAILURE` (the most common case) or `EvolutionState.R.SUCCESS` (which only happens if the Evaluator in this process or some external process found the ideal individual).

The default implementation of these `Problem` methods does nothing at all. But in `MasterProblem` these methods are used to handle reconnection of Slaves after a checkpoint recovery. The first two methods create

both a new `SlaveMonitor`. The final method shuts down the monitor cleanly.

`MasterProblem` also has special methods used only by Steady-State Evolution (and thus Asynchronous Evolution):

---

#### **ec.eval.MasterProblem Methods**

```
public boolean canEvaluate()
```

Returns true if a Slave is available to take an Individual.

```
public boolean evaluatedIndividualAvailable()
```

Returns true if a Slave has a completed Individual waiting to be introduced to the Population.

```
public QueueIndividual getNextEvaluatedIndividual()
```

Blocks until an Individual is available from a Slave. Then returns it as an `ec.steadystate.QueueIndividual`. A `QueueIndividual` is a very simple class which just contains the Individual and the Subpopulation that the Individual should be introduced into. It has the following instance variables:

```
    public Individual ind;  
    public int subpop;
```

---

`MasterProblem` also implements all the methods defined in `SimpleProblemForm` and `GroupedProblemForm` (Section 7.1.2), in addition to the following two methods normally implemented by `Evaluator` (and which are called on the `MasterProblem` from the `Evaluator`'s methods of the same names):

---

#### **ec.eval.MasterProblem Methods**

```
public void prepareToEvaluate(EvolutionState state, int threadnum)
```

Creates a new queue of Individuals who are out waiting to be sent to be processed.

```
public void finishEvaluating(EvolutionState state, int threadnum)
```

Sends all Individuals presently in the queue out in a Job. Then waits for all slaves to complete evaluation of all Individuals.

---

**Sending Initialization Data from the Master to the Slaves** `MasterProblem` has one additional feature: three methods you can override such that, whenever a Slave is started up, you are given the chance to send one-time data from the Master to the Slave. Note that this facility is untested and may be somewhat fragile: we suggest that before you use this facility you consider instead placing the data as a file that each Slave can read.

Here's how it works. You begin by declaring a special subclass of `MasterProblem`—let's call it `classec.app.myapp.MyMasterProblem`. Then you must tell ECJ that you're using it (on both sides, master and slave) with something like this:

```
eval.masterproblem = ec.app.myapp.MyMasterProblem
```

An instance of this `MasterProblem` subclass is used by the Master. When a Slave shows up, the Master will call the method `sendAdditionalData(...)` on that instance to give you the chance to send data down the stream to the Slave before any evaluation occurs. The Slave receives this data by calling the method `receiveAdditionalData(...)` on a dummy instance of this same subclass. This dummy instance will not be used for any purpose except to handle the data reception. That method must store away the transferred data, typically as instance variable in your subclass which you have created for this purpose. Finally, each time the Slave creates a new `EvolutionState` (and it may do so many times), it calls `transferAdditionalData(...)` on that dummy instance to give it a chance to modify and set up `EvolutionState` appropriately to reflect the data it had received from the Master.

`sendAdditionalData(...)` will be called once on the Master for each Slave that shows up. `receiveAdditionalData(...)` will only be called once on the Slave side. `transferAdditionalData(...)` may be called multiple times on the Slave side, every time a new `EvolutionState` is loaded and set up. It is entirely up to you to handle the protocol for sending and receiving this data: if you mess up (write more data than you read on the other side, say), it's undefined what'll happen. Probably something bad.

---

### ec.eval.MasterProblem Methods

---

`public void sendAdditionalData(EvolutionState state, DataOutputStream dataOut)`

Called from the `SlaveMonitor`'s `accept()` method to optionally send additional data to the Slave via the `dataOut` stream. By default it does nothing.

`public void receiveAdditionalData(EvolutionState state, DataInputStream dataIn)`

Called on a dummy `MasterProblem` by the Slave. You should use this method to store away received data via the `dataIn` stream for later transferring to the current `EvolutionState` via the `transferAdditionalData(...)` method. You should NOT expect this `MasterProblem` to be used for by the Slave for evolution (though it might). By default this method does nothing, which is the usual situation. The `EvolutionState` is provided solely for you to be able to output warnings and errors: do not rely on it for any other purpose (including access of the random number generator or storing any data).

`public void transferAdditionalData(EvolutionState state`

Called on a dummy `MasterProblem` by the Slave to transfer data previously loaded via `receiveAdditionalData(...)` to a running `EvolutionState` at the beginning of evolution. This method may be called multiple times if multiple `EvolutionStates` are created. By default this method does nothing, which is the usual situation. Unlike in `receiveAdditionalData(...)`, the provided `EvolutionState` is "live" and you can set it up however you like.

---

## 6.2 Island Models (The ec.exchange Package)

In addition to Distributed Evaluation, ECJ supports Island Models: separate ECJ processes ("islands") which connect over the network and occasionally hand highly-fit Individuals to one another. This facility is handled by an Exchanger called `ec.exchange.IslandExchange` which ships individuals off to other islands immediately before breeding, and immediately after breeding brings in new individuals provided it by other islands. You'll run your separate islands as ordinary processes.

Most of the issues in Island models surround the particular topology being chosen. Which islands will send Individuals to which other islands? How many at a time? How often? Are Individuals sent synchronously or asynchronously? Etc. ECJ manages all topology and connection parameters via a special ECJ process called the **island model server**. Each island will connect to and register itself with the server. When the islands have all connected, the server will tell them which islands need to hook up to which other islands and how. After the islands have hooked up, they're given the go-ahead to start their evolutionary processes. If the islands are acting synchronously, each generation will wait for the server to give them the go-ahead to continue to the next generation; this go-ahead only occurs after all islands have finished the previous generation. Finally, when an island discovers an optimal individual, it will signal the server to let the other islands know (so they can shut down).

As you can see, the server really does little more than tell the islands how to connect and act as a referee. Thus it's actually a very lightweight process. You can run the server either as its own process like this:

```
java ec.exchange.IslandExchange -file server.params -p param=value ... (etc.)
```

...or an ECJ island can also do double-duty, serving as the server. All islands, whether ordinary islands or double-duty island-server combos, are fired up in the same standard ECJ fashion:

```
java ec.Evolve -file island.params -p param=value -p param=value ... (etc.)
```

...or (as usual):

```
java ec.Evolve -checkpoint myCheckpointFile.gz
```

A double-duty island-server combo would differ from a plain island solely in the parameters it defines: it'd need server parameters in addition to client parameters.

**Mixing Island Models, Threading, and Distributed Evaluation** There's absolutely no reason you create an unholy union of Island Models and Distributed Evaluation. For example, it's perfectly reasonable to have an Island Model, where each island maintains its own pool of Slaves to do distributed evaluation. It'd be a lot of parameter files though! Island Models also work perfectly fine in a multithreaded environment.

### 6.2.1 Islands

You set up an ECJ process as an island by defining a special Exchange object for it:

```
exch = ec.exchange.IslandExchange
```

IslandExchange maintains the island's **mailbox**. Prior to breeding, the IslandExchange procedure will send some fit individuals off to mailboxes of remote islands. The procedure for selecting Individuals is defined along these lines:

```
exch.select = ec.select.TournamentSelection
```

After breeding, the IslandExchange will empty its own mailbox and introduce into the Population all of the Individuals contained therein. These Individuals will displace some of the recently-bred Individuals, which never get a chance to be selected.<sup>4</sup> The selection method for picking the Individuals to die and be displaced is defined as:

```
exch.select-to-die = ec.select.RandomSelection
```

If this parameter isn't defined, individuals are picked at random.

An Island needs to know where the Server is so it can register itself, and the socket port on which the server is listening, for example:

```
exch.server-addr = 128.2.30.4  
exch.server-port = 8999
```

When an Island registers itself with the Server, it'll tell it two things. First, it'll tell the Server the island's **name**, a String which uniquely identifies the island (and by which the Server looks up topology information for the island). Second, it'll tell the Server the socket port on which it'll receive incoming Individuals from other islands. Let's say that you're creating an island called StatenIsland. You might specify the following:

```
exch.id = StatenIsland  
exch.client-port = 9002
```

Note that the client socket port should be (1) higher than 2000 and (2) different from other client ports, and the server port, if they're running on the same machine or in the same process. You'll also probably want to have compressed network streams. Like was the case in Distributed Evaluation, this can't be done without the *jzlib*/*ZLIB* library, which you must install separately (see the ECJ main webpage or <http://www.jcraft.com/jzlib>). This is because Java's compression facilities are broken. Once this library installed, you can turn it on like this:

---

<sup>4</sup>Life's not fair.

```
exch.compression = true
```

Be certain to give your island a unique random number seed different from other islands! Don't set the seed to time, since it's possible that two islands will have the same seed because they were launched within the one millisecond of one another. I'd hard-set the seed on a per-island basis.

```
seed.0 = 5921623
```

If your islands share the same file system, you'll want to make sure they don't overwrite each other's statistics files etc. To do this, for example, StatenIsland might change its statistics file name to:

```
stat.file = $statenisland.stat
```

If you have multiple Statistics files you'll need to change all of them; the same goes for other files being written out. Also, if you are checkpointing, and your islands might overwrite each others' checkpoint files, you need to change the checkpoint prefix on a per-island basis. For example:

```
prefix = statenisland
```

Last, you can cut down on the verbosity of the islands by setting...

```
exch.chatty = false
```

## 6.2.2 The Server

The Server holds all the parameters for setting up the island topology. But first we must clue your ECJ process into realizing that it is a Server in the first place. This is done with:

```
exch.i-am-server = true
```

Next we need to state how many islands are in the island model graph:

```
exch.num-islands = 3
```

As discussed in Section 6.2.1, each island has a unique name (id). Here you will state which island in your graph has which id:

```
exch.island.0.id = StatenIsland
exch.island.1.id = ConeyIsland
exch.island.2.id = EllisIsland
```

Each island has some number of connections *to* other islands (the islands it'll send Individuals, or *migrants*, to). In this example, we'll say that StatenIsland sends migrants to ConeyIsland, which sends migrants to EllisIsland, which sends migrants to both StatenIsland and ConeyIsland:

```
exch.island.0.num-mig = 1
exch.island.0.mig.0 = ConeyIsland
exch.island.1.num-mig = 2
exch.island.1.mig.0 = EllisIsland
exch.island.2.num-mig = 2
exch.island.2.mig.0 = StatenIsland
exch.island.2.mig.1 = ConeyIsland
```

StatenIsland and ConeyIsland send 10 migrants to each of the islands they're connected to. But we want EllisIsland to send 50 migrants to each of the (two) islands it's connected to:



```

exch.island.0.size = 10
exch.island.1.size = 10
exch.island.2.size = 50

```

Alternatively you can use a default parameter base of sorts:

```

exch.size = 10

```

Each island has a maximum mailbox capacity: if there is no room, further immigrants will be dropped and disappear into the ether. You should make your mailbox big enough to accept immigrants at a reasonable rate, but not so large that in theory they could entirely overwhelm your population! I suggest a mailbox three or four times the size of the expected immigrants. How about 100 or 200?

```

exch.island.0.mailbox-capacity = 200
exch.island.1.mailbox-capacity = 200
exch.island.2.mailbox-capacity = 200

```

Alternatively you can use a default parameter base of sorts:

```

exch.mailbox-capacity = 10

```

Last you'll need to stipulate two additional parameters on a per-island basis: the **start-generation** (in which generation the island will start sending Individuals out) and **modulus** (how many generations the island will wait before it sends out another batch of Individuals). These are mostly set to maximize network utilization: perhaps you may wish the islands to send out individuals at different times so as not to clog your network, for example. Here we'll tell each island to send out individuals every three generations, but to start at different initial generations so to be somewhat staggered:

```

exch.island.0.mod = 3
exch.island.1.mod = 3
exch.island.2.mod = 3
exch.island.0.start = 1
exch.island.1.start = 2
exch.island.2.start = 3

```

Alternatively you can use a default parameter base of sorts:

```

exch.mod = 3
exch.start = 2

```

### 6.2.2.1 Synchronicity

Island models can be either **synchronous** or **asynchronous**. In a synchronous island model, islands wait until they all have reached the next generation before sending immigrants to one another. In the asynchronous island model, islands go at their own pace and send immigrants whenever they feel like it. This means that one evolutionary process on one computer may run much faster than another one (good, because it doesn't waste resources waiting for the other one to catch up) but it may overwhelm the other process with multiple generations of immigrants before the other process can get around to processing them (usually bad). Generally speaking asynchronicity is preferred — and is the default setting.

If for some reason you want to turn on synchronicity, you do this:

```

exch.sync = true

```

Note that the modulo and start-generation of islands results in a predictable behavior for synchronous

island models: but since asynchronous islands can go at their own pace, the modulo and start-generation happen when they happen for each island.

Note too that because asynchronous island models go at their own pace, and are subject to the whims of the speed of the operating system and the CPU time allotted to the process, there's no way to guarantee replicability.

### 6.2.3 Internal Island Models

ECJ's Internal Island Model facility simulates islands using separate Subpopulations: each Subpopulation is an island, and occasionally highly fit Individuals migrate from Subpopulation to Subpopulation. Like any other Exchanger, the Internal Island Model facility takes Individuals from other Subpopulations immediately before Breeding, stores them, and then introduces them into their destination Subpopulations immediately after Breeding.

There are four important things to note about this facility:

- Obviously each Subpopulation must have an identical Species and Individual and Fitness prototypes.
- Internal Island Models are always synchronous.
- Because they use the Subpopulation facility, Internal Island Models are incompatible with any other ECJ procedure which relies on Subpopulations: notably coevolution.
- Because they define an Exchanger, Internal Island Models are incompatible with any other ECJ procedure which uses an Exchanger: in particular, you can't mix Internal Island Models with regular Island Models!

Why would you use Internal Island Models? I think mostly for academic purposes: to study and simulate synchronous Island Models without having to rope together a bunch of machines. You could also use Internal Island Models to run  $N$  evolutionary processes in parallel — just set the number of immigrants to zero.

Internal Island Models depend solely on a specific Exchanger, `ec.exchange.InterPopulationExchange`. To build an Internal Island Model, you first define three subpopulations and their species, individuals, breeding pipelines, the whole works, using a standard generational algorithm. Then you define the exchanger:

```
exch = ec.exchange.InterPopulationExchange
```

Let's say you have four Subpopulations acting as islands. You'll first need to stipulate the Selection Method used to select individuals to migrate to other Subpopulations and the Selection Method used to kill Individuals to make way for incoming immigrants:

```
exch.subpop.0.select = ec.select.TournamentSelection
exch.subpop.1.select = ec.select.TournamentSelection
exch.subpop.2.select = ec.select.TournamentSelection
exch.subpop.3.select = ec.select.TournamentSelection
exch.subpop.0.select-to-die = ec.select.RandomSelection
exch.subpop.1.select-to-die = ec.select.RandomSelection
exch.subpop.2.select-to-die = ec.select.RandomSelection
exch.subpop.3.select-to-die = ec.select.RandomSelection
```

If you don't define a selection method for death, ECJ will assume you mean to select individuals randomly. Alternatively you can use the default parameter base:

```
exch.select = ec.select.TournamentSelection
exch.select-to-die = ec.select.RandomSelection
```

Next you need to state the number of immigrants sent at a time; the first generation in which they'll be sent; and the modulus (the interval, in terms of generations, between successive migrations). These are basically the same as the standard Island Model. The parameters might look like this:

```
exch.subpop.0.size = 5
exch.subpop.1.size = 5
exch.subpop.2.size = 5
exch.subpop.3.size = 15
exch.subpop.0.start = 1
exch.subpop.1.start = 2
exch.subpop.2.start = 3
exch.subpop.3.start = 4
exch.subpop.0.mod = 8
exch.subpop.1.mod = 8
exch.subpop.2.mod = 8
exch.subpop.3.mod = 8
```

It's here where you could convert these to separate independent evolutionary processes: just set the `size` parameter to 0 for all subpopulations. Anyway, you can also use default parameter bases for these:

```
exch.size = 5
exch.start = 1
exch.mod = 8
```

Now we need to define the topology. For each island we'll define the number of Subpopulations it sends migrants to, and then which ones. Imagine if Subpopulation 2 sent migrants to everyone else, but the other Subpopulations just sent migrants to Subpopulation 2. We would define it like this:

```
exch.subpop.0.num-dest = 1
exch.subpop.0.dest.0 = 2
exch.subpop.1.num-dest = 1
exch.subpop.1.dest.0 = 2
exch.subpop.2.num-dest = 3
exch.subpop.2.dest.0 = 0
exch.subpop.2.dest.1 = 1
exch.subpop.2.dest.2 = 3
exch.subpop.3.num-dest = 1
exch.subpop.3.dest.0 = 2
```

Last, Internal Island Models tend to be verbose. To make them less chatty, you can say:

```
exch.chatty = false;
```

**Reminder** It's laborious to write a zillion parameters if you have large numbers of subpopulations in your internal island model experiments. You can greatly simplify this process using the `pop.default-subpop` parameter (see Section 3.2.1).

## 6.2.4 The Exchanger

In Section 3.6 we talked about various basic Exchanger methods. But three were not discussed, mostly because they're only used for Island Models (not even Internal Island Models). They are:

```
public void initializeContacts(EvolutionState state);  
public void reinitializeContacts(EvolutionState state);  
public void closeContacts(EvolutionState state, int result);
```

If these look similar to the methods in Section 6.1.5, it's with good reason. Their function is to set up networking connections, re-establish networking connections after restarting from a checkpoint, and shut down networking connections in a clean way. `IslandExchange` implements them but not `InterPopulationExchange`.

## Chapter 7

# Additional Evolutionary Algorithms

### 7.1 Coevolution (The `ec.coevolve` Package)

The coevolution package is meant to provide support for three kinds of Coevolution:

- One-Population Competitive Coevolution
- Two-Population Competitive Coevolution
- *N*-Population Cooperative Coevolution

Coevolution differs from evolutionary methods largely in how evaluation is handled (and of course, by the fact that there are often multiple subpopulations). Thus the classes in this package are basically Problems and Evaluators. The first form of Coevolution is provided by the `ec.coevolve.CompetitiveEvaluator` class. The second two are made possible by the `ec.coevolve.MultiPopCoevolutionaryEvaluator` class.

#### 7.1.1 Coevolutionary Fitness

Coevolution is distinguished by its evaluation of Individuals not separately but in groups, where the Fitness of an Individual depends on its performance in the **context** of other Individuals (either competing with them or working with them towards a common goal). For this reason, coevolution typically involves evaluating an Individual multiple times, each time with a different set of other Individuals, and then computing the Fitness based on these multiple **trials**.

To assist in this procedure, the `ec.Fitness` class has two auxiliary variables:

```
javapublic ArrayList trials = null; public Individual[] context = null;
```

The first variable (`trials`) is used to maintain the results of each trial performed. It's only used during evaluation, and then set to null again. Most commonly each trial is stored as a `java.lang.Double`, with higher values being considered better. You can change this but if you do, be sure to read Section 7.1.5 first. **Even if your trials are not `java.lang.Double`, they must be immutable: their internals cannot be modified, so they can be pointer-copied and not just cloned.**

The second variable (`context`) is used to maintain the context of the best trial discovered for the Individual. It's typically only used by Cooperative Coevolution, where it's important to retain not only the performance of the Individual (his Fitness), but which collaborating Individuals made it possible for him to achieve that Fitness. If `context` is used, it's assumed that the slots of `context` hold each collaborator, except for the Individual himself, whose slot is set to null. Section 7.1.4.2 discusses the issue of context in more detail.

## 7.1.2 Grouped Problems

Since Coevolution involves evaluating multiple Individuals at once, it will require a new kind of Problem which takes multiple Individuals at a time. This Problem Form is defined by `ec.coevolve.GroupedProblemForm`. Evaluation in coevolution involves multiple trials, along these lines:

1. Performance scores of Individuals are cleared.
2. Individuals are tested against each other in various matches (or with one another in various collaborative problems). These trials cause trial performance scores of the Individuals to accumulate. If the trials are cooperative, the best trial found for a given Individual (its *context*) is maintained.
3. The final Fitnesses of the Individuals are set based on the performance scores over all the trials.

It's up to you to store the trial results and eventually form them into final Fitness values, as discussed later. It's also up to you to maintain the best context if you find this useful, as discussed later as well (in Section 7.1.4.2). `GroupedProblemForm` will help you by defining three methods to do these various portions of the evaluation process. You'll need to implement all three methods:

### `ec.coevolve.GroupedProblemForm` Methods

---

`public abstract void preprocessPopulation(EvolutionState state, int thread, boolean[] prepareForAssessment, boolean countVictoriesOnly)`

Called prior evaluation of a Population, mostly to clear the trials of Individuals. Only clear the trials for Individuals in Subpopulations for which *prepareForAssessment* is true. Note that although this method is not static, you should not assume that this method will be called on the same Problem as is used later for `evaluate(...)`. Thus don't use this method to set any instance variables in the Problem for later use. If *countVictoriesOnly* is true, the method being used is `SingleEliminationTournament`. Commonly you'll use this method to create a brand-new trials `ArrayList` for every Fitness of every Individual in every Subpopulation.

`void evaluate(EvolutionState state, Individual[] individuals, boolean[] updateFitness, boolean countVictoriesOnly, int[] subpops, int thread)`

Evaluates the *individuals* in a single trial, setting their performance scores for that trial. Each individual will be from a certain subpopulation, specified in *subpops*. In some versions of coevolution, only certain individuals are supposed to have their performance scores updated (the others are acting as foils). In this case, the relevant individuals will be indicated in the *updateFitness* array. Typically you'll update fitness by adding trial results to the trials `ArrayList`. For any individual for whom you're updating trials, also set the Fitness value to reflect that one trial: this allows `Single Elimination Tournament` to compare Individuals based on this notional Fitness value. In doing so, do *not* set the evaluated flag for Individuals.

If you're doing cooperative coevolution, in this method you'll also probably want to maintain the context of the trial (the collaborating Individuals) if it's produced the best results so far. More on that in Section 7.1.4.2.

`public abstract void postprocessPopulation(EvolutionState state, int thread, boolean[] assessFitness, boolean countVictoriesOnly)`

Called after evaluation of a Population to form final Fitness scores for the individuals based on the various performance scores they accumulated during trials; and then to set their evaluated flags to *true*. Only assess the Fitness and set the evaluated flags for Individuals in Subpopulations for which *assessFitness* is true. Note that although this method is not static, you should not assume that this method will be called on the same Problem as was used earlier for `evaluate(...)`. You'll probably want to set the trials variable to null to let it garbage collect.

---

Do not assume that Individuals will have the same number of trials: in several versions of Coevolution this will not be the case.

You're going to have to do some work to make sure that your Fitnesses are properly updated, and this depends on the kind of Coevolution you choose to do. So how do you keep track of trials? If you're just accumulating scores or wins, then you could use `SimpleFitness` and just increment it with each new win. But since different Individuals may have different numbers of trials, it's possible that you may need to keep

track of the number of trials, or keep track of each trial result separately. Probably the best approach is to use the auxiliary variable `trials` found in `ec.Fitness` to store all your trials in an `ArrayList`. Thus you typically use `GroupedProblemForm` like this:

1. In `preprocessPopulation(...)`, set the `Fitness`'s trials to a new `ArrayList` for all `Individuals`.
2. In `evaluate(...)`, add to trials the results of this trial for each `Individual` for whom `updateFitness` is set. Then set the `Fitness` to reflect just the immediate results of the trial (particularly if `countVictoriesOnly` is set). This allows Single Elimination Tournament—if you're using that—to determine who should advance to the next round. Make certain that whatever you add to trials is `java.io.Serializable`.  
If you're doing cooperative coevolution, determine if the new trial is superior to all the existing trials previously found in the trials array. If it is, set the context of this trial using `Fitness.setContext(...)`.
3. In `postprocessPopulation(...)`, set the `Fitness` to the final result. For example, you might set it to the average or the maximum of the various trials. Finally, set the `Fitness`'s trials to null to let it GC.

Notice that unlike in `SimpleProblemForm` (Section 3.4.1) there's no `describe(...)` method. This is because to describe an `Individual`, you'd need to do so in the context of other `Individuals`. So we left it out.

**Example** Assume we've replaced the `Fitness` with our `MyCoevolutionaryFitness` class. Let's create a `Problem` similar to the example in Section 4.1.1. In our `Problem` we take two `Individuals`, and the trial performance of an `Individual` is his vector values' product minus that of his opponent. Obviously this is a stupid example since `Individuals` are in a total ordering `Fitness`-wise: so it hardly illustrates the issues in Coevolution. But it'll suffice for the demonstration here. What we'll do is set an `Individual`'s fitness to his average score over the trials performed.

```
package ec.app.myapp;
import ec.*;
import ec.simple.*;
import ec.vector.*;
import ec.coevolve.*;
public class MyCoevolutionaryProblem extends Problem implements GroupedProblemForm {

    public void preprocessPopulation(EvolutionState state, Population pop,
                                    boolean[] prepareForAssessment, boolean countVictoriesOnly) {
        for(int i = 0; i < pop.subpops.length; i++)
            if (prepareForAssessment[i])
                for(int j = 0; j < pop.subpops[i].individuals.length; j++) {
                    SimpleFitness fit = (SimpleFitness)(pop.subpops[i].individuals[j].fitness);
                    fit.trials = new ArrayList();
                }
    }

    public void evaluate(EvolutionState state, Individual[] ind, boolean[] updateFitness,
                        boolean countVictoriesOnly, int[] subpops, int threadnum) {
        int[] genome1 = ((IntegerVectorIndividual)ind[0]).genome;
        int[] genome2 = ((IntegerVectorIndividual)ind[1]).genome;
        double product1 = 1.0;
        double product2 = 1.0;

        for(int x=0; x<genome1.length; x++) product1 = product1 * genome1[x];
        for(int x=0; x<genome2.length; x++) product2 = product2 * genome2[x];

        MyCoevolutionaryFitness fit1 = (MyCoevolutionaryFitness)(ind[0].fitness);
        MyCoevolutionaryFitness fit2 = (MyCoevolutionaryFitness)(ind[1].fitness);
```

```

        if (updateFitness[0]) {
            fit1.trials.add(new Double(product1 - product2));
            // set the fitness in case we're using Single Elimination Tournament
            fit1.setFitness(state, fit1.fitness() + product1 - product2, false);
        }
        if (updateFitness[1]) {
            fit2.trials.add(new Double(product2 - product1));
            // set the fitness in case we're using Single Elimination Tournament
            fit2.setFitness(state, fit2.fitness() + product2 - product1, false);
        }
    }

    public void postprocessPopulation(EvolutionState state, Population pop,
                                     boolean[] assessFitness, boolean countVictoriesOnly) {
        for(int i = 0; i < pop.subpops.length; i++)
            if (assessFitness[i])
                for(int j = 0; j < pop.subpops[i].individuals.length; j++) {
                    SimpleFitness fit = (SimpleFitness)(pop.subpops[i].individuals[j].fitness);

                    // Let's set the fitness to the average of the trials
                    int len = fit.trials.size();
                    double sum = 0;
                    for(int l = 0; l < len; l++)
                        sum += ((Double)(fit.trials.get(l))).doubleValue();

                    // Alternatively if we were interested in how *many* times we won rather
                    // than by how much, we might set the fitness to, say, how often the doubleValue()
                    // was positive.

                    fit.setFitness(state, (float)(sum / len), false);
                    pop.subpops[i].individuals[j].evaluated = true;
                }
    }
}

```

Now we hook it up as so:

```
eval.problem = ec.myapp.MyCoevolutionaryProblem
```

It's important to note that the trials variable doesn't have to store an ArrayList of *numbers* — it can store any object reflecting the outcome of a trial that you'd like to invent. It's available for you to use.

Generally the trials variable should never need to be cloned — it'll be set up, used during Evaluation, and then destroyed. The default implementation of the Fitness.cloneTrials() method, called by Fitness.clone(), simply throws an exception if trials is null.

### 7.1.3 One-Population Competitive Coevolution

In One-Population Competitive Coevolution, the whole population (or in ECJ parlance, a whole *Subpopulation*) competes in various games. The outcomes of those games determines the fitness of Individuals in the Subpopulation. Thus the Fitness of an Individual depends on which other Individuals it plays against.

ECJ does One-Population Competitive Coevolution with a special Evaluator called `ec.coevolve.CompetitiveEvaluator`, meant to be used in generational evolutionary methods. Building on the genetic algorithm example in Section 4.1.1, we'll replace the Evaluator:



```
eval = ec.coevolve.CompetitiveEvaluator
```

The main issue in One-Population Competitive Coevolution is the layout of the trials (competitions) between various Individuals. CompetitiveEvaluator provides the following options:<sup>1</sup>

- Round Robin. Every Individual of the Subpopulation is paired up and evaluated with every other Individual, not including the Individual itself. The Fitness of both Individuals is updated. This is stipulated as:

```
eval.style = round-robin
```

- K-Random-Opponents (One-Way). Every Individual of the Subpopulation is paired up with  $K$  random opponents other than himself, picked at random with replacement. Only the Individual's fitness is updated, not his opponents'. This is stipulated, plus the size of  $K$ , as:

```
eval.style = rand-1-way  
eval.group-size = 8
```

Our experiments have tended to suggest that  $K = 6$  to  $8$  are good values. Remember that larger group sizes result in many more evaluations.

- K-Random-Opponents (Two-Way). Every Individual of the Subpopulation is paired up with *at least*  $K$  random opponents other than himself, picked at random with replacement. Both Individuals' fitnesses are updated; thus being selected as an "opponent" allows an Individual to need fewer "regular" tests in order to fill his own  $K$  quota. Overall this results in about half the total number of tests as One-Way.

```
eval.style = rand-2-way  
eval.group-size = 8
```

A few individuals would be expected to be tested more than  $K$  times. If your fitness procedure simply cannot permit this, you can have the algorithm turn off fitness updating for those Individuals with:

```
eval.over-eval = false
```

- Single Elimination Tournament. All Individuals in the Subpopulation are paired up and evaluated. In each pair, the "winner" is the individual which winds up with the superior performance. If neither Individual is superior, then the "winner" is chosen at random. The winners advance to the next level: they're paired up and evaluated, and so on, just like in a single-elimination tournament. **Important note:** It is important that the Subpopulation size be a power of two for a proper Single-Elimination Tournament.

You stipulate Single-Elimination Tournament with:

```
eval.style = single-elim-tournament
```

After a single trial, you need to increment the Fitness of the winning Individual: this will tell Single-Elimination Tournament who to advance to the next round.<sup>2</sup> After all the trials have completed, we will treat the Fitness of an Individual simply to be the number of trials it won, and so not modify the Fitness during postprocessPopulation. You can tell Single-Elimination Tournament is being used because countVictoriesOnly is true in the GroupedProblemForm methods.

---

<sup>1</sup>Note that we used to have a number of other options at one point, many of which are still in the CompetitiveEvaluator source code of various ECJ versions, such as double-elimination tournament, world cup style, etc. But these have fallen by the wayside.

<sup>2</sup>Keep in mind that two Individuals compete with one another only because they've advanced to the same level, and so have the same Fitness initially.

One-Population Competitive Coevolution can be prohibitively expensive. So if you can use it, we suggest you *not* use Round Robin. Here's why. Round Robin is  $O(n^2)$ . K-Random-Opponents (One-Way) is  $O(kn)$ . K-Random-Opponents (Two-Way) is roughly  $O(\frac{k}{2}n)$ . But Single-Elimination Tournament is  $O(n)$ . When cast in terms of performance by number of evaluations, sometimes Single-Elimination Tournament performs best, though often K-Random-Opponents performs best with  $K$  set to somewhere between 6 and 8. We did a comparison paper on the two [12].

`ec.coevolve.CompetitiveEvaluator` is multithreaded and responds to the `evalthreads` parameter.

### 7.1.4 Multi-Population Coevolution

There are two common kinds of multi-population coevolution: 2-Population Competitive Coevolution and  $N$ -Population Cooperative Coevolution. Both are handled by the `ec.coevolve.MultiPopCoevolutionaryEvaluator` class:

```
eval = ec.coevolve.MultiPopCoevolutionaryEvaluator
```

In 2-Population Competitive Coevolution Individuals in each of two Subpopulations are tested by pitting them against Individuals in the opposing Subpopulation. The Fitness of an Individual is based on several such tests. Typically one Subpopulation is of interest; the other largely acts as a foil to challenge it.

In  $N$ -Population Cooperative Coevolution, a problem is broken into  $N$  subparts, and solutions for each subpart are optimized in separate Subpopulations. An Individual in a Subpopulation is tested by combining it with one member from each of the other Subpopulations to form a joint solution, which is then evaluated. Again, the Fitness of an Individual is based on several such tests.

When `ec.coevolve.MultiPopCoevolutionaryEvaluator` tests Individuals together, it typically selects one Individual from each Subpopulation and passes them to the `evaluate(...)` method: the order of the array of Individuals passed in should correspond to the order of the Subpopulations. Pay attention to the passed-in `updateFitness` array: only these individuals should be assessed based on a given test.

`ec.coevolve.MultiPopCoevolutionaryEvaluator` tests a given Individual in a Subpopulation in combination with Individuals selected from the other Subpopulations in various ways. The number of trials of each type that an Individual receives is determined by several associated parameters.

- Individuals selected at random, with replacement, from the current generation. The number of trials of this type is:

```
eval.num-current = 4
```

Since Individuals do not yet have their Fitnesses set, the selection procedure should be one which does not consider Fitness.<sup>3</sup> You need to specify the selection procedure for each subpopulation. For example, if we had two Subpopulations, we could say:

```
eval.subpop.0.select-current = ec.select.RandomSelection
eval.subpop.1.select-current = ec.select.RandomSelection
```

`RandomSelection` is strongly recommended. Alternatively you could use the default parameter base:

```
eval.select-current = ec.select.RandomSelection
```

- Individuals selected at random, with replacement, from the previous generation. You provide the selection method for them; unlike before, this one can be based on Fitness if you like. For generation 0, Individuals are picked from the current generation and at random. For example, to specify the number of individuals of this type, and to use `TournamentSelection`, we might do:

<sup>3</sup>What's the point then? Why bother with a `SelectionMethod` at all? Why not just pick Individuals at Random? The answer is simple: the `ec.spatial` package (Section 7.2) has a variation of Multi-Population Coevolution which uses its special version of `TournamentSelection`.

```
eval.num-prev = 6
eval.subpop.0.select = ec.select.TournamentSelection
eval.subpop.0.select.size = 2
```

Again, you could use the default parameter base, though be warned that (in this example) it'll share parameters with all TournamentSelection uses.

```
eval.num-prev = 6
eval.select-prev = ec.select.TournamentSelection select.tournament.size = 2
```

- The very fittest Individuals from the previous generation (or randomly chosen if we're at the generation 0). The number of fittest Individuals stored is the same as the number of trials:

```
eval.subpop.0.num-elites = 5
```

- Individuals in the current Population are paired up using random shuffling. For each trial, all the Subpopulations are randomly shuffled<sup>4</sup>. Then each Individual #0 of each Subpopulation is paired together for a trial; likewise each Individual #1 of each Subpopulation, each Individual #2, and so on. This produces one trial for every Individual. For the next trial, all Subpopulations are randomly shuffled again, and so on.

Note that this procedure makes it possible to guarantee  $N$  trials per Individual with many fewer total trials: you can then use your saved Evaluations towards other things, such as more trials, larger Subpopulation sizes, or longer generations. To do 4 shuffled trials per Individual, you'd say:

```
eval.num-shuffled = 4
```

For competitive coevolution, you can use the same example as in Section 7.1.2 (GroupedProblemForm). For Cooperative Coevolution it's typical to base Fitness instead on the *maximum* of tests with  $N$  other collaborators. For example, let's revise the postprocessPopulation(...) method to assess trial performance as the sum of all the products of the genomes in the various Individuals:

```
public void postprocessPopulation(EvolutionState state, Population pop,
                                boolean[] assessFitness, boolean countVictoriesOnly) {
    for(int i = 0; i < pop.subpops.length; i++)
        if (assessFitness[i])
            for(int j = 0; j < pop.subpops[i].individuals.length; j++)
                if (!pop.subpops[i].individuals[j].evaluated) {
                    SimpleFitness fit = (SimpleFitness)(pop.subpops[i].individuals[j].fitness);

                    // Let's set the fitness to the maximum of the trials
                    int len = fit.trials.size();
                    double max = Double.NEGATIVE_INFINITY;
                    for(int l = 0; l < len; l++)
                        max = Math.max(max, ((Double)(fit.trials.get(l))).doubleValue());

                    fit.setFitness(state, (float)(max), false);
                    pop.subpops[i].individuals[j].evaluated = true;
                }
    }
```

ec.coevolve.MultiPopCoevolutionaryEvaluator is not at present multithreaded. It does not respond to the evalthreads parameter.

---

<sup>4</sup>Not actually shuffled — just shuffled internal index arrays to them: the effect is the same.

**Reminder** It's laborious to write a zillion parameters if you have large numbers of subpopulations in your coevolution experiments. You can greatly simplify this process using the `pop.default-subpop` parameter (see Section 3.2.1).

#### 7.1.4.1 Parallel and Sequential Coevolution

Note that the options above enable the "Parallel" and "Parallel Previous" coevolutionary techniques [6], but not the "Serial" (or "Sequential") technique whereby each Subpopulation is evaluated in turn. For Sequential coevolution, there's an additional parameter option in `ec.simple.SimpleBreeder`:

```
breed.sequential = true
```

Ordinarily `SimpleBreeder` breeds every Subpopulation every generation. But if this parameter is set, then `SimpleBreeder` will instead breed only one Subpopulation each generation. The chosen Subpopulation's index will be the generation number, modulo the number of Subpopulations.<sup>5</sup>

This parameter is very rarely used outside of sequential cooperative coevolution. `MultiPopCoevolutionaryEvaluator` responds to `SimpleBreeder`'s parameter as well to match the sequential breeding with sequential evaluation. That is, `MultiPopCoevolutionaryEvaluator` will only evaluate a single Subpopulation at a time (the same one that's later bred by `SimpleBreeder`).

Even though the evaluator and breeder are only updating one Subpopulation a generation, the Statistics classes are still doing statistics on *every* Subpopulation. However two basic Statistics subclasses (`ec.simple.SimpleStatistics` and `ec.gp.koza.KozaShortStatistics`, have some tweaks in them to print out the proper information. Specifically, when these Statistics classes print to screen the Fitnesses of the best individuals of each Subpopulation, they only do so if the individuals in question have their "evaluated" flags set. Otherwise it'll say "not evaluated". Recall that in `ec.coevolve.GroupedProblemForm` this only happens in `postProcessPopulation`, and only to individuals for whom the `assessFitness[]` slot is set. In sequential cooperative coevolution, that slot is set only for the subpopulation which is being updated that generation.

Okay fine, but what about the Statistics' logs? They'll still have nonsensical "best of generation" Individuals right? Yes, but you can tell which ones are valid simply by looking at their evaluated flags. Only the best individual of the currently-being-updated Subpopulation will have its evaluated flag set to true.

#### 7.1.4.2 Maintaining Context

When you do Cooperative Coevolution, you'll probably want to retain information not only on how the Individual performed, but which collaborating Individuals helped him do his best. It's up to you to maintain this information, but ECJ's Fitness object provides a variable which can help:

```
public Individual[] context;
```

In this variable store each collaborator, one per Subpopulation, plus one slot set to null to represent the Individual himself. The Fitness object will detect if the context exists (by default it's null) and print out the context information when `printFitnessForHumans(...)` is called. Fitness will also handle cloning this array (it calls `newFitness.setContext(oldFitness.getContext(context))`) when the Fitness is cloned, etc. Fitness has some methods you can use rather than playing with the context directly:

---

#### ec.Fitness Methods

```
public void setContext(Individual[] cont, int index)
```

Sets the context to the Individuals stored in context, under the assumption that the slot for the primary Individual is located at `cont[index]`. Clones the Individuals before adding to them to the context.

```
public void setContext(Individual[] cont)
```

Sets the context to the Individuals stored in context, under the assumption that the slot for the primary Individual can be determined because it has been set to null. Clones the Individuals before adding to them to the context.

---

<sup>5</sup>Note that though `SimpleBreeder` implements this, a number of subclasses of `SimpleBreeder` explicitly do *not*.

```
public Individual[] getContext()
    Returns the context.
```

---

You'll want to update the context during the evaluate(...) method. Here's an example Cooperative Coevolution evaluation method:

```
public void evaluate(EvolutionState state, Individual[] ind, boolean[] updateFitness,
                    boolean countVictoriesOnly, int[] subpops, int threadnum)
{
    // First compute the value of this trial.  For example, imagine if each Individual was a
    // DoubleVectorIndividual, and the fitness function was simply their sum.
    double sum = 0;
    for(int i = 0; i < ind.length; i++)
    {
        DoubleVectorIndividual dind = (DoubleVectorIndividual)(ind[i]);
        for(int j = 0; j < dind.genome.length; j++)
            sum += dind.genome[j];
    }

    // next update each Individual to reflect this trial result
    for(int i = 0; i < ind.length; i++)
    {
        if (updateFitness[i]) // need to include this trial for this Individual
        {
            // We presume that the best trial found so far has been stored at
            // position 0 of the trials array.  This makes it O(1) to determine if
            // we're the new "best trial".  Heeere we go!

            ArrayList tr = ind[i].fitness.trials;
            if (tr.size() == 0) // we're the first trial to be performed so far
            {
                tr.add(new Double(sum)); // just dump us in
                ind[i].fitness.setContext(ind, i); // set the context to us
            }
            else if (((Double)(tr.get(0))).doubleValue < sum) // we're the new best
            {
                // swap us to the 0th position
                Double t = (Double)(tr.get(0));
                tr.set(0, new Double(sum)); // I'm at 0
                tr.add(t); // move the old guy to the end of the line
                ind[i].fitness.setContext(ind, i); // set the context to us
            }

            // don't bother setting the fitness here
        }
    }
}
```

**Warning** Maintaining context in this way is expensive. If you have  $N$  individuals involved in a trial, you'll likely be updating the context on all  $N$  of them, and each context array is  $N$  size, so you're looking at an  $O(N^2)$  operation for every trial! If you're doing coevolution with 1000 subpopulations, that's a million array slots for *each* trial of *each* individual. Furthermore, if you are using `ec.simple.SimpleStatistics`, large contexts will be printed to disk.

In this situation, you probably want to simply avoid setting context: it's just too expensive.

## 7.1.5 Performing Distributed Evaluation with Coevolution

Coevolution introduces its own wrinkle to Distributed Evaluation, because it has multiple trials, and each of those trials is a separate job potentially processed on a separate machine. We need a way to merge all those trials together to compute the final fitness. Furthermore we may need a way to determine which context (see Section 7.1.4.2) among those merged trials should be retained. If your trial results are `java.lang.Double`, with higher values considered better, you probably don't need to do anything. Otherwise you'll need to make a tweak. Here's some details on how the mechanism works.

When ECJ sends off a group of Individuals to a remote site to be evaluated using a `ec.coevolve.GroupedProblemForm`, the first thing it does (in `ec.eval.Job`) is make clones of them, then strip the clones of both trials and context. When the Individuals come back, they need to have their trials and context reintegrated with the original Individuals. This is done by calling the `merge(...)` method on the Individuals.

The `Individual.merge(...)` method isn't particularly efficient but it gets the job done. First, it calls `Fitness.merge(...)` to handle merging trials and context (actual fitness value isn't bothered with — at this point it doesn't matter). However this merging isn't in the direction you think: the *old* Fitness is merged into the new Fitness. Then it copies the entire new Individual, including the merged Fitness, into the old Individual, displacing it. This way the merged Fitness retains the new fitness value, plus merged trials and context.

How are trials and context merged? Keep in mind that when the Individual was cloned and sent to the remote site to be evaluated, the clone's trials and context were stripped. When the cloned Individual returns and must be merged with the original Individual, the cloned Individual will have a new context and one or more new trials. `Fitness.merge(...)` who has the better trial: accordingly, this is the context that is retained. Then the new trials are appended to the original trials.

**The crucial item.** To determine which group of trials contains a superior trial, `Fitness.merge(...)` calls a method called `Fitness.contextIsBetterThan(Fitness other)`. This method returns true if the given Fitness's trials are superior to the "other" Fitness's trials. **This method assumes that trials are `java.lang.Double`.** If they are not, you'll need to override this method and replace it. Recall again that **even if your trials are not `java.lang.Double`, they must be immutable: their internals cannot be modified, so they can be pointer-copied and not just cloned.**

Here are the relevant methods of interest:

---

### ec.Individual Methods

```
public void merge(EvolutionState state, Individual other)
```

Merges the other Individual into the original Individual, including Fitnesses. After the merge, (1) the Individual should be a copy of other (2) the Fitness's value should be a copy of other (3) the Fitness's trials should be a merger of the original trials and those of other (4) the higher-quality context of the two Individuals should be retained. By default this method does this by calling `other.fitness.merge(fitness)` (note the backwards merge), then copying back the entirety of other into the original Individual<sup>6</sup> You'll rarely need to override this method.

---

---

### ec.Fitness Methods

```
public void merge(EvolutionState state, Fitness other)
```

Merges the other Fitness into the original Fitness. The only thing that needs to be merged is trials and context. If other has better trials than the original Individual, its context replaces the original context. Then the trials are concatenated together. Calls `contextIsBetterThan(...)` to determine which Fitness has the superior trials, and thus context. You'll not often need to override this method, but may need to override `contextIsBetterThan(...)`.

```
public boolean contextIsBetterThan(Fitness other)
```

Returns true if the my context is better than the other Fitness's context. This is computed by comparing our trials:

---

<sup>6</sup>How is this magic copying done? By use of a `ec.util.DataPipe`: we first call `other.writeIndividual(...)` into the pipe, then call `readIndividual(...)` from the pipe. It's a hack but it works, and avoids the need for a custom version of `Individual.merge` for every possible genotype.

if I have trials and the other does not, or if one of my trials is superior to all of the other's trials, then returns true. This method assumes that trials are `java.lang.Double`, and that higher values are better. If this is not the case, you'll need to override this method to customize it.

---

## 7.2 Spatially Embedded Evolutionary Algorithms (The `ec.spatial` Package)

The `ec.spatial` package helps in designing evolutionary algorithms in which the Individuals in a Subpopulation are embedded in a metric space of some kind, and their decisions about breeding and/or selection are based on their location in that space and their relationship with other Individuals. For example, an Individual chosen to breed might be restricted to only do so with Individuals near to himself.

In order for a Subpopulation to be used in a spatially-embedded way, it must implement the `ec.spatial.Space` interface. This interface enables the Breeder to set an *index* for a **target** Individual in the Space, then later request a neighbor, chosen at random, within a given *distance* in the Space of that target Individual.

### `ec.spatial.Space` Methods

---

```
public void setIndex(int thread, int index)
```

Set the index for a certain target Individual in the Space. (Each thread can have its own index).

```
public int getIndex(int thread)
```

Return the index for the current target Individual in the Space. (Each thread can have its own index).

```
public int getIndexRandomNeighbor(EvolutionState state, int thread, int distance)
```

Return the index of a neighbor in the embedded space, chosen at random, and who is no more than *distance* away from the current target individual.

---

This seems convoluted. Why store indices? Why you might ask, can't we just pass in the index into the `getIndexRandomNeighbor` method? The answer is: because at the time this method is called (usually during a `TournamentSelection` method call), the index isn't available. Another option would be to store the index not in the Space but in a Breeder,<sup>7</sup> but it turns out that Spaces can be used not only for breeding but also for Evaluation. So we're storing the indices in the only logical location left: the Space itself.

### 7.2.1 Implementing a Space

ECJ has one common space implemented for you already: a toroidal 1D ring. For fun, let's implement another space: a 2D *non*-toroidal 10x10 space.<sup>8</sup> We'll start with the easy part: storing the per-thread index array.

```
public class 2DSubpopulation extends Subpopulation implements Space {
    int[] indices;

    public void setup(EvolutionState state, Parameter base) {
        super.setup(state, base);
        indices = new int[state.breedthreads];
    }

    public Group emptyClone() {
```

---

<sup>7</sup>This is what's done in `MuCommaLambdaBreeder` (Section 4.1.2) to enable the magic of `ESSelection` always picking a certain individual.

<sup>8</sup>I didn't say it'd be useful: just an easy demonstration!

```

    2DSubpopulation other = (2DSubpopulation)(super.emptyClone());
    other.indices = new int[indices.length];
    return other;
}

```

```

public void setIndex(int thread, int index) { indices[thread] = index; }

```

```

public int getIndex(int thread) { return indices[thread]; }

```

Now we need to implement the `getIndexRandomNeighbor` method. We'll define a random neighbor as one which is within the rectangular box up to *distance* away from the target index.

```

public int getIndexRandomNeighbor(EvolutionState state, int thread, int distance) {
    if (individuals.length != 100)
        state.output.fatal("This ridiculous example requires that the Subpopulation" +
            "be exactly 100 in size.")

    int target = getIndex(thread);
    int targetX = target / 10; // integer division
    int minX = targetX - distance < 0 ? 0 : targetX - distance;
    int maxX = targetX + distance > 9 ? 9 : targetX + distance;

    int targetY = target % 10;
    int minY = targetY - distance < 0 ? 0 : targetY - distance;
    int maxY = targetY + distance > 9 ? 9 : targetY + distance;

    int x = minX + state.random(thread).nextInt(maxX - minX + 1);
    int y = minY + state.random(thread).nextInt(maxY - minY + 1);

    return x * 10 + y;
}
}

```

## 7.2.2 Spatial Breeding

To breed, you can use the Space ECJ has already done for you: a 1D Space. `ec.spatial.Spatial1DSubpopulation` can be set up to be either toroidal (the default) or non-toroidal. To use it as toroidal, you might say:

```

pop.subpop.0 = ec.spatial.Spatial1DSubpopulation
pop.subpop.0.toroidal = true

```

You'll also need a selection procedure which understands how to pick using Spaces. ECJ provides one such procedure: `ec.spatial.SpatialTournamentSelection`. This selection operator picks its tournament members randomly from the Subpopulation but constrained so that they're within *distance* of the target index. That is: it uses `getIndexRandomNeighbor(...)` to pick them.

`SpatialTournamentSelection` has three parameter options above and beyond the standard `TournamentSelection` operators. Let's say that `SpatialTournamentSelection` is our pipeline's first source:

```

pop.subpop.0.species.pipe.source.0 = ec.spatial.SpatialTournamentSelection
pop.subpop.0.species.pipe.source.0.size = 2
pop.subpop.0.species.pipe.source.0.pick-worst = false
pop.subpop.0.species.pipe.source.0.neighborhood-size = 3
pop.subpop.0.species.pipe.source.0.ind-competes = false
pop.subpop.0.species.pipe.source.0.type = uniform

```



The size parameter should be obvious: it's TournamentSelection's tournament size. Likewise the pick-worst parameter determines whether we're picking the fittest or the least fit of the tournament. The remaining three parameters are as follows. neighborhood-size defines the distance from the target index. If ind-competes is true, then at least one member of the tournament is guaranteed to be the target Individual itself. Last, if type is uniform, then an individual is picked simply using getIndexRandomNeighbor(...). However you also have the option of doing a **random walk** through the space. The walk will be neighborhood-size steps and each step will move to a neighbor immediately bordering your current position (that is, one chosen when you pass in 1 as the distance to getIndexRandomNeighbor. Random walks of this kind approximate a Gaussian distribution centered on the target Individual. You can choose to do a random walk rather than uniform selection with:

```
pop.subpop.0.species.pipe.source.0.type = random-walk
```

You can of course use the default parameters for this kind of stuff. Note that since the default parameter base is different, you need to specify all the standard Tournament Selection stuff under this default base, not under Tournament Selection's own default base (select.tournament):

```
spatial.tournament.size = 2
spatial.tournament.pick-worst = flase
spatial.tournament.neighborhood-size = 3
spatial.tournament.ind-competes = false
spatial.tournament.type = uniform
```

In order for SpatialTournamentSelection to work, you'll have to have the target Individual set. This is the job of ec.spatial.SpatialBreeder, a simple extension of SimpleBreeder. The procedure works like this. For each Individual  $i$  in the new Subpopulation to fill:

1. SpatialBreeder sets the target index to  $i$ .
2. SpatialBreeder requests one Individual from its Pipeline
3. SpatialTournamentSelection methods in that Pipeline use this index to compute distances and pick Individuals

**Warning:** Note that although SpatialBreeder is enumerating over the *new Subpopulation*, the SpatialTournamentSelection operator is using the target index to pick Individuals the *old Subpopulation*. This means that Subpopulations must not change in size from generation to generation.

SpatialBreeder is very simple: there are no parameters to set up at all.

### 7.2.3 Coevolutionary Spatial Evaluation

The spatial package also enables one of many<sup>9</sup> possible approaches to doing coevolution in a spatially embedded Subpopulation. Recall from Section 7.1.4 that the ec.coevolve.MultiPopCoevolutionEvaluator class performs multi-population coevolution: each Individual is tested  $N$  times, each time by grouping it with one Individual from each of the other Subpopulations. These other Individuals are called **collaborators**. The ec.spatial.SpatialMultiPopCoevolutionaryEvaluator extends this by allowing you to pick those collaborators in a spatial manner: their locations in their respective Subpopulations are near to the position of the target Individual in his Subpopulation:

```
eval = ec.spatial.SpatialMultiPopCoevolutionaryEvaluator
```

---

<sup>9</sup>The other obvious approach would be a variant of ec.coevolve.CompetitiveEvaluator where competitors to test Individuals in a population are selected based on nearness to the Individual. Perhaps we might put this together one day.

The way this is done is by using the `SpatialTournamentSelection` class to pick collaborators, based on the target Individual's index. Two kinds of collaborators can be picked with a Selection Method in `MultiPopCoevolutionEvaluator`: members of the current Population, and members of the previous generation's Population. You need to be careful with the first situation, since the members don't have their Fitnesses set, and so you can't use a Selection Method which is based on Fitness. In the examples in Section 7.1.4, we used `RandomSelection`. But for spatial coevolution, we can do a "random" selection which simply picks a collaborator based on the neighborhood from the target Individual. We do this by setting the size parameter in `SpatialTournamentSelection` — the tournament size — to 1:

```
eval.subpop.0.select-current = ec.spatial.SpatialTournamentSelection

# It's important that the size be 1 ---
# this causes 'random' selection, so Fitness is not considered
eval.subpop.0.select-current.size = 1
eval.subpop.0.select-current.pick-worst = false
eval.subpop.0.select-current.neighborhood-size = 3

# It's also important that this value be false ---
# otherwise the target index will always be selected
eval.subpop.0.select-current.ind-competes = false
eval.subpop.0.select-current.type = uniform

eval.subpop.1.select-current = ec.spatial.SpatialTournamentSelection
eval.subpop.1.select-current.size = 1
eval.subpop.1.select-current.pick-worst = false
eval.subpop.1.select-current.neighborhood-size = 3
eval.subpop.1.select-current.ind-competes = false
eval.subpop.1.select-current.type = random-walk
```

You can also select collaborators from the previous generation based on spatial distance. In this case, you can have the Selection Method pick based on Fitness too (though you don't have to if you don't want to). For example:

```
eval.subpop.0.select-prev = ec.spatial.SpatialTournamentSelection
eval.subpop.0.select-prev.size = 2
eval.subpop.0.select-prev.pick-worst = false
eval.subpop.0.select-prev.neighborhood-size = 3
eval.subpop.0.select-prev.ind-competes = false
eval.subpop.0.select-prev.type = uniform

eval.subpop.1.select-prev = ec.spatial.SpatialTournamentSelection
eval.subpop.1.select-prev.size = 7
eval.subpop.1.select-prev.pick-worst = false
eval.subpop.1.select-prev.neighborhood-size = 4
eval.subpop.1.select-prev.ind-competes = true
eval.subpop.1.select-prev.type = random-walk
```

The remaining parameters are basically just like those in `MultiPopCoevolutionaryEvaluator` (see Section 7.1.4), for example:

```
eval.num-current = 4
eval.num-prev = 6
eval.subpop.0.num-elites = 5
```

**Warning:** Just like was the case for Spatial Breeding, Spatial Multi-Population Coevolution requires that Subpopulations not change in size from generation to generation.

## 7.3 Differential Evolution (The ec.de Package)

The ec.de package provides a Differential Evolution framework and some basic DE breeding operators. Differential Evolution is notable for its wide range of operator options: ECJ only has a few of them defined, but others are not hard to create yourself.

Differential Evolution does not have a traditional selection mechanism: instead, children are created entirely at random but then must compete with their parents for survival. Additionally, Differential Evolution breeding operators are somewhat opaque and not particularly amenable to mixing and matching. Because of this, we have eschewed the Breeding Pipeline approach for doing breeding. Instead, Differential Evolution simply has different Breeders for each of its breeding operator approaches. Differential Evolution operators are defined by creating a subclass of ec.de.DEBreeder. We have three such operators implemented, including the default in ec.de.DEBreeder itself: more on them in a moment.

### 7.3.0.1 The Evaluator

We implement DE's unusual selection procedure through a custom SimpleEvaluator called ec.de.DEEvaluator. It turns out this code could have been done in ec.de.DEBreeder (making the package simpler) but then various Statistics objects wouldn't report the right values. So we made a simple Evaluator to do the job. DEEvaluator is very simple subclass of SimpleEvaluator and works just like it, including multithreaded evaluation. You specify it as follows:

```
eval = ec.de.DEEvaluator
```

### 7.3.1 Breeding

DEBreeder serves both as the superclass of all DE breeding operators, and also implements the basic "classic" Differential Evolution breeding operator, known as DE/rand/1/bin. We begin by talking about the top-level elements of DEBreeder, then we discuss DEBreeder's specific default operator, followed by other operators.

Most Differential Evolution breeding operators work as follows. For each Individual  $\vec{d}$  in the Population (recall that in Differential Evolution all Individuals are DoubleVectorIndividuals, so we'll treat them as vectors), we will create a single child  $\vec{c}$ . After this is done, we replace each parent with its child if the child is as good or better than the parent (this last part is done in DEEvaluator).

Children are usually, but not always created by first generating a child  $\vec{d}$  through a combination, of sorts, of Individuals than  $\vec{d}$ . We then perform a crossover between  $d$  and  $a$ , essentially replacing parts of  $d$  with  $a$ , which results in the final child  $c$ . Thus there are often two parts to creating children: first building  $d$ , then crossing over  $d$  with  $a$ .

The whole creation process is handled by a single method:

#### ec.de.DEBreeder Methods

---

```
public DoubleVectorIndividual createIndividual(EvolutionState state, int subpop, int index, int thread)
    Creates and returns a new Individual derived from the parent Individual found in
    state.population.subpops[subpop].individuals[index], using other individuals chosen from
    state.population.subpops[subpop] as necessary.
```

---

The default implementation performs the DE/rand/1/bin operation, discussed later. If you want to create a whole new breeding operator, you'll largely need to override this individual-creation method.

Because crossover is a common operation, ECJ has broken it out into a separate method. You can override just this method, if you like, to customize how crossover is done.

#### ec.de.DEBreeder Methods

---

```
public DoubleVectorIndividual crossover(EvolutionState state, DoubleVectorIndividual target,  
                                     DoubleVectorIndividual child, int thread)  
    Crosses over child (earlier referred to as  $\vec{d}$ ) with target (earlier referred to as  $\vec{a}$ ), modifying child. child is then  
    returned.
```

---

The default implementation performs uniform crossover: for each gene, with independent probability  $Cr$ , the child's gene will be replaced with the target's gene. One child gene, chosen at random, is guaranteed to *not* be replaced.

The  $Cr$  value, which must be between 0.0 and 1.0, is set by a parameter:

```
breed.cr = 0.5
```

And it's stored in the following DEBreeder instance variable:

```
public double Cr;
```

You can leave this parameter unspecified: as mentioned before, some DEBreeder subclasses do no crossover and so don't need it. But if crossover *is* performed and the parameter is unspecified (it's set to the value `ec.DEBreeder.CR_UNSPECIFIED`), it'll be assumed to be 0.5 (probably not a good choice) and a warning will be issued.

The combination of various Individuals to form  $\vec{d}$  in the first place is usually a mathematical vector operation involving certain variables. The most common variable, used by all the operators in this package, is a scaling factor  $F$ , which ranges from 0.0 to 1.0. We define  $F$  like this:

```
breed.f = 0.6
```

The  $F$  parameter is then stored in the following DEBreeder instance variable:

```
public double F;
```

Unlike  $Cr$ ,  $F$  is required as a parameter.

Some DE breeding operators need to know the fittest Individual in the Subpopulation. Thus prior to constructing individuals, DEBreeder first computes this whether it's used or not. The location of the fittest Individuals, indexed by Subpopulation, are stored in the following DEBreeder variable:

```
public int[] bestSoFarIndex;
```

For example, you can get the fittest Individual in Subpopulation 0 as:

```
DoubleVectorIndividual bestForSubpopZero = (DoubleVectorIndividual)  
    (state.population.subpops[0].individuals[bestSoFarIndex[0]]);
```

Last but not least, DEBreeder must store away the old Population before it is overwritten by the new child Individuals. This is because in DEEvaluator the original parents in the old Population get a chance to displace the children if the children are not sufficiently fit. So we need to keep the parents around until then. The parents are stored in DEBreeder as:

```
public Population previousPopulation;
```

This value is initially null, since there are no parents in the first generation. But in successive generations it's set to the parents.

A final note: though evaluation is multithreaded, breeding at present is not. The `breedthreads` parameter has no effect.

### 7.3.1.1 The DE/rand/1/bin Operator

The “classic” DE breeding operator, DE/rand/1/bin, is the default operator implemented by DEBreeder itself. The implementation follows that found on page 140 of the text *Differential Evolution* [15]. It works as follows. For each Individual  $\vec{a}$  in the Subpopulation, we select three other Individuals at random. These Individuals must be different from one another and from  $\vec{a}$ . We’ll call them  $\vec{r0}$ ,  $\vec{r1}$ , and  $\vec{r2}$ . We create a child  $\vec{d}$  whose values are defined as:

$$d_i = r0_i + F \times (r1_i - r2_i)$$

We then cross over  $d$  with  $a$ , producing a final child  $c$ , which is then placed in the new Subpopulation. Note the use of the  $F$  parameter.

To use this operator, you’ll need to specify DEBreeder as the breeder, and of course also set the  $F$  and  $Cr$  parameters as well. For example:

```
breed = ec.de.DEBreeder
breed.f = 0.6
breed.cr = 0.5
```

This operator can produce values which are outside the min/max gene range. You’ll need to specify whether or not you wish to force the operator to only produce values bounded to within that range. For example, to turn off bounding and allow values to be anything, you would say (on a per-Species basis):

```
pop.subpop.0.species.mutation-bounded = false
```

Otherwise you’d say:

```
pop.subpop.0.species.mutation-bounded = true
```

If you bound the operator, then it will try repeatedly with new values of  $\vec{r0}$ ,  $\vec{r1}$ , and  $\vec{r2}$ , until it produces a valid individual  $d_i$ .

### 7.3.1.2 The DE/best/1/bin Operator

The class `ec.de.Best1BinDEBreeder`, a subclass of `DEBreeder`, implements the DE/best/1/bin operator, plus “random jitter”, as found on page 140 of the text *Differential Evolution* [15]. It works as follows. For each Individual  $\vec{a}$  in the Subpopulation, we first identify the *fittest* Individual  $\vec{b}$  in the Subpopulation, and also select two other Individuals at random. These Individuals must be different from one another, from  $\vec{b}$ , and from  $\vec{a}$ . We’ll call them  $\vec{r1}$  and  $\vec{r2}$ . We create a child  $\vec{d}$  whose values are defined as:

$$d_i = \vec{b} + jitter() \times (r1_i - r2_i)$$

We then cross over  $d$  with  $a$ , producing a final child  $c$ , which is then placed in the new Population. `jitter()` produces different random numbers for each  $i$ . It’s defined as:

$$jitter() = F + FNOISE \times (\text{random}(0,1) - 0.5)$$

The `random(0,1)` function returns a random number between 0.0 and 1.0 inclusive. Again, note the use of the  $F$  parameter. The `FNOISE` constant is typically 0.001. It’s defined by the parameter `breed.f-noise` and is stored in the `Best1BinDEBreeder` instance variable:

```
public double F_NOISE;
```

To use this breeding operator, you’ll need to specify the breeder and the various parameters:

```

breed = ec.de.Best1BinDEBreeder
breed.f = 0.6
breed.cr = 0.5
breed.f-noise = 0.001

```

This operator can produce values which are outside the min/max gene range. You'll need to specify whether or not you wish to force the operator to only produce values bounded to within that range. For example, to turn off bounding and allow values to be anything, you would say (on a per-Species basis):

```
pop.subpop.0.species.mutation-bounded = false
```

Otherwise you'd say:

```
pop.subpop.0.species.mutation-bounded = true
```

If you bound the operator, then it will try repeatedly with new values of  $\vec{r1}$  and  $\vec{r2}$ , until it produces a valid individual  $d_i$ .

### 7.3.1.3 The DE/rand/1/either-or Operator

The class `ec.de.Rand1EitherOrDEBreeder`, again a subclass of `DEBreeder`, implements the DE/rand/1/either-or operator, as found on page 141 of the text *Differential Evolution* [15]. It works as follows. For each Individual  $\vec{a}$  in the Subpopulation, we select three other Individuals at random. These Individuals must be different from one another and from  $\vec{a}$ . We'll call them  $\vec{r0}$ ,  $\vec{r1}$ , and  $\vec{r2}$ . Then with probability  $PF$ , we create a child  $\vec{d}$  defined in the same way as the DE/rand/1/bin operator:

$$d_i = r0_i + F \times (r1_i - r2_i)$$

... else we create the child like this:

$$d_i = r0_i + 0.5 \times (F + 1) \times (r1_i + r2_i - 2 \times r0_i)$$

We do *not* cross over  $d$  with  $a$ : we simply return  $d$  as the child.

The  $PF$  probability value, which must be between 0.0 and 1.0, is defined by the parameter `breed.pf` and is stored in the `Rand1EitherOrDEBreeder` instance variable:

```
public double PF;
```

To use this breeding operator, you'll need to specify the breeder and the various parameters:

```

breed = ec.de.Rand1EitherOrDEBreeder
breed.f = 0.6
breed.pf = 0.5

```

Note that the  $Cr$  parameter is not used, since no crossover is performed.

This operator can produce values which are outside the min/max gene range. You'll need to specify whether or not you wish to force the operator to only produce values bounded to within that range. For example, to turn off bounding and allow values to be anything, you would say (on a per-Species basis):

```
pop.subpop.0.species.mutation-bounded = false
```

Otherwise you'd say:

```
pop.subpop.0.species.mutation-bounded = true
```

If you bound the operator, then it will try repeatedly with new values of  $\vec{r_0}$ ,  $\vec{r_1}$ , and  $\vec{r_2}$ , until it produces a valid individual  $d_i$ .

## 7.4 Multiobjective Optimization (The `ec.multiobjective` Package)

ECJ has three packages which handle multiobjective optimization: the `ec.multiobjective` package, and two concrete implementations (SPEA2 and NSGA-II): the `ec.multiobjective.spea2` and `ec.multiobjective.nsga2` packages respectively.

### 7.4.0.4 The `MultiObjectiveFitness` class

The `ec.multiobjective` package contains a single new kind of Fitness. Multiobjective optimization differs from other optimization algorithms in that the Fitness of an Individual is not a single value but rather consists of some  $N$  **objectives**, separate values describing the quality of the Individual on various aspects of the Problem. Thus the primary class overridden by multiobjective algorithms is a special version of Fitness called `ec.multiobjective.MultiObjectiveFitness`. Internally in this class the objective results are stored in an array of floats:

```
public float[] objectives;
```

The number of objectives is the length of this array.

You create a `MultiObjectiveFitness` and define its objectives along these lines:

```
pop.subpop.0.species.fitness = ec.multiobjective.MultiObjectiveFitness
pop.subpop.0.species.fitness.num-objectives = 3
```

Though ECJ tends to assume that higher Fitness values are better, many multiobjective algorithms assume the opposite. Thus `MultiObjectiveFitness` has the option of doing either case. ECJ assumes lower objective values are better when you state:

```
pop.subpop.0.species.fitness.maximize = false
```

`true` is the default value. This value is stored in the variable:

```
public boolean maximize;
```

You can also set the minimum and maximum objective values on a per-objective basis:

```
pop.subpop.0.species.fitness.min.0 = 0.0
pop.subpop.0.species.fitness.max.0 = 2.0
pop.subpop.0.species.fitness.min.1 = 1.0
pop.subpop.0.species.fitness.max.1 = 3.5
pop.subpop.0.species.fitness.min.2 = -10
pop.subpop.0.species.fitness.max.2 = 0
```

The default minimum is 0.0 and the default maximum is 1.0. If you like you can set global minimum and maximum values instead:

```
pop.subpop.0.species.fitness.min = 0.0
pop.subpop.0.species.fitness.max = 2.0
```

Local min/max values, if set, override the global values. The resulting minimum and maximum values are stored in the following arrays:

```
public float[] minObjectives;
public float[] maxObjectives;
```

Note that these min/max arrays are *shared*<sup>10</sup> among all Fitness objects cloned from the same Prototype. So you should treat them as read-only.

You can get and set objectives via functions, which have the added benefit of double-checking their min/max validity:

---

#### ec.multiobjective.MultiObjectiveFitness Methods

```
public int getNumObjectives()
    Returns the number of objectives.

public float[] getObjectives()
    Returns all the current objective values.

public float getObjective(int index)
    Returns a given objective value.

public void setObjectives(EvolutionState state, float[] objectives)
    Sets the objective values to the ones provided, double-checking that they are within valid minimum and maximum ranges (discussed in a moment)

public double sumSquaredObjectiveDistance(MultiObjectiveFitness other)
    Returns the sum squared distance in objective space between this Fitness and the other. That is, if for a given objective  $i$ , this fitness has value  $A_i$  and the other has value  $B_i$ , this function will return  $\sum_i (A_i - B_i)^2$ .

public double manhattanObjectiveDistance(MultiObjectiveFitness other)
    Returns the Manhattan distance in objective space between this Fitness and the other. That is, if for a given objective  $i$ , this fitness has value  $A_i$  and the other has value  $B_i$ , this function will return  $\sum_i ||(A_i - B_i)||$ .
```

---

Since MultiObjectiveFitness is a Prototype, of course you can define these parameters using a default parameter base. For example:

```
pop.subpop.0.species.fitness = ec.multiobjective.MultiObjectiveFitness
multi.fitness.num-objectives = 3

# Global objectives
multi.fitness.min = 0.0
multi.fitness.max = 2.0

# Local Overrides (heck, why not?)
multi.fitness.min.1 = 1.0
multi.fitness.max.1 = 3.5
multi.fitness.min.2 = -10
multi.fitness.max.2 = 0
```

#### 7.4.0.5 The MultiObjectiveStatistics class

To help output useful statistics about the Pareto Front, ECJ has an optional but strongly encouraged Statistics class, ec.multiobjective.MultiObjectiveStatistics. This class largely overrides the finalStatistics method to output the Pareto Front in three different ways:

---

<sup>10</sup>Why doesn't Fitness store them in the Species or something? Because Fitness, for historical and not particularly good reasons, does not have a Flyweight relationship with any object.



- A summary of the Front is printed to the screen.
- Every Individual in the Front is written to the statistics file.
- A whitespace-delimited table of the objective values of each member of the Front, one per line, is written to a special file. If the Front is 2-objective, and the file is called `front.stat`, you can easily view the Front results by firing up GNUPLOT and entering:

```
plot front.stat
```

The `MultiObjectiveStatistics` class is used, and the Front file defined, as follows:

```
stat = ec.multiobjective.MultiObjectiveStatistics
stat.front = $front.stat
```

Various multiobjective optimization algorithms subclass from `MultiObjectiveFitness` to add auxiliary fitness values. We'd like those values to be included as columns in the Front summary when it is outputted to the screen by `MultiObjectiveStatistics`. To do this, `MultiObjectiveFitness` has two special methods which are overridden by subclasses:

---

#### **ec.multiobjective.MultiObjectiveFitness Methods**

`public String[] getAuxilliaryFitnessNames()`

Returns an array of names, one per auxiliary fitness element. These will appear as headers to their respective columns.

`public double[] getAuxilliaryFitnessValues()`

Returns an array of doubles, one per auxiliary fitness element, of the current values of those elements.

---

### **7.4.1 Selecting with Multiple Objectives**

ECJ's primary multiobjective algorithms don't use `MultiObjectiveFitness` directly, but rather subclass it to add per-algorithm gizmos. But if you're not using these algorithms, you can still use `MultiObjectiveFitness` in a more "traditional" generational algorithm. This requires some thought: at the end of the day, ECJ needs to select based on this Fitness mechanism. But if there is more than one objective value, how is this done?

You could just sum the objectives to form a final "fitness" value of course. If your `SelectionMethod` uses the raw `fitness()` value (for example, `FitProportionateSelection` from Section 3.5.2) this is exactly what `MultiObjectiveFitness` returns.

But the most common approach is instead to use some form of **Pareto domination**. Pareto domination works like this. Individual *A* *Pareto dominates* Individual *B* if and only if *A* is *at least as good* as *B* in all objectives, and is *superior* to *B* in at least one objective. Notice that there are three cases where *A* might **not** Pareto-dominate *B*:

- *B* Pareto-dominates *A*
- *A* and *B* have exactly the same objective values for all objectives.
- *A* is superior to *B* in some objectives, but *B* is superior to *A* in other objectives.

`MultiObjectiveFitness` implements the `betterThan` and `equivalentTo`<sup>11</sup> methods to return pareto domination results: `betterThan` is true if the Individual Pareto-dominates the one passed in. `equivalentTo` returns true if neither Pareto-dominates the other (either of the last two cases above).

---

<sup>11</sup>Perhaps now it might make sense why the method is called `equivalentTo` and not `equalTo`.

If your `SelectionMethod` relies entirely on the `betterThan` and `equivalentTo` methods (such as `TournamentSelection`), it'll use Pareto domination to sort `Individuals`. Note however that this method has problems, not the least of which is that there are cycles in Pareto domination. That is, it's easy to construct `Individuals` *A*, *B*, and *C* such that *A* Pareto-dominates *B*, which Pareto-dominates *C*, which Pareto-dominates *A*!

Various subclasses of `MultiObjectiveFitness` for different kinds of algorithms override the `betterThan` and `equivalentTo` methods to measure fitness differently. However you can still determine Pareto Domination with the following method:

---

#### **ec.multiobjective.MultiObjectiveFitness Methods**

```
public boolean paretoDominates(MultiObjectiveFitness other)
    Returns true if this fitness Pareto-dominates other.
```

---

##### **7.4.1.1 Pareto Ranking**

An alternative method, widely used in multiobjective algorithms but not implemented in `MultiObjectiveFitness` proper, is to perform **Pareto ranking** (sometimes called *non-dominated sorting* [19], which works as follows. The **Pareto non-dominated front** (or simply **Pareto front**) is the set of `Individuals` in a Subpopulation who are dominated by *no one else*, including one another. All members of a Pareto front of a Subpopulation receive a rank of 0. We then remove those members from consideration in the Subpopulation and form the Pareto front among the remaining members. These members receive a rank of 1. We then remove *those* members from consideration as well and repeat. Ultimately every member of the Subpopulation receives a rank. These ranks become fitness values, where lower ranks are better than higher ranks. Pareto ranking plays a major part in the NSGA-II algorithm (Section 7.4.2).

Here are methods for extracting the Pareto Front and various Pareto Front Ranks:

---

#### **ec.multiobjective.MultiObjectiveFitness Methods**

```
public static ArrayList partitionIntoParetoFront(Individual[] inds, ArrayList front, ArrayList nonFront)
    Partitions inds into the front and the non-front Individuals. If front is provided, the front Individuals are placed there and returned. If front is not provided, a new ArrayList is created, the front is placed there, and it is returned. If nonFront is provided, the non-front Individuals are placed in it, else they are discarded.

public static ArrayList partitionIntoRanks(Individual[] inds)
    Partitions inds into Pareto Front ranks. Each rank is an ArrayList of Individuals. The ranks are then placed into an ArrayList in increasing order (worsening rank) starting with rank zero. This ArrayList is returned.
```

---

##### **7.4.1.2 Archives**

The most common — and often most effective — multiobjective optimization algorithms are very strongly elitist, essentially versions of the  $(\mu + \lambda)$  evolution strategy (Section 4.1.2). Specifically, they split the Population into two groups: the primary Population, and an **archive** consisting largely of the Pareto front of nondominated `Individuals` discovered so far. Any time a new `Individual` is discovered which dominates members of this front, those members are removed from the elitist archive and the `Individual` is introduced to it. ECJ has two Pareto archive multiobjective algorithms: NSGA-II and SPEA2. We discuss them next.

## **7.4.2 NSGA-II (The ec.multiobjective.nsga2 Package)**

The Non-dominated Sorting Genetic Algorithm Version II (or NSGA-II) [2] is essentially a version of the  $(\mu + \mu)$  evolution strategy using non-dominated sorting. It maintains and updates an archive (half the population) of the current best individuals (essentially the current estimate of the Pareto Front), and breeds the remaining population from the archive.

NSGA-II requires a Breeder to maintain the archive and handle the algorithm's custom breeding; and an Evaluator to compute the non-dominated sorting. These classes are defined by the `ec.multiobjective.nsga2.NSGA2Breeder` and `ec.multiobjective.nsga2.NSGA2Evaluator` classes respectively. The non-dominated sorting information is included in the fitness, and so NSGA-II uses a subclass of `MultiObjectiveFitness` called `ec.multiobjective.nsga2.NSGA2MultiObjectiveFitness`. To output the Pareto Front for the user, we also include a special Statistics subclass called `ec.multiobjective.nsga2.NSGA2Statistics`.

Simply replace `MultiObjectiveFitness` with this class:

```
pop.subpop.0.species.fitness = ec.multiobjective.nsga2.NSGA2MultiObjectiveFitness
```

You'll need to set the number of objectives, and min/max objectives, etc., as usual, something like:

```
multi.fitness.num-objectives = 3
multi.fitness.min = 0.0
multi.fitness.max = 2.0
multi.fitness.min.1 = 1.0
multi.fitness.max.1 = 3.5
multi.fitness.min.2 = -10
multi.fitness.max.2 = 0
```

This class contains two special fitness measures: the **rank** and **sparsity** of the Individual. The rank is computed as the Pareto Rank of the individual. The sparsity is a measure of distance of the individual to others on the same rank. We like sparse individuals because we don't want individuals all clustered in one area of the front. The fitness is simple: an Individual is superior to another if its rank is lower (better). If the same, an Individual is superior if its sparsity is higher. These two measures are stored in `NSGA2MultiObjectiveFitness` as the instance variables:

```
public int rank;
public double sparsity;
```

The `NSGA2Evaluator` computes and sets these values. You define it in the obvious way:

```
eval = ec.multiobjective.nsga2.NSGA2Evaluator
```

The `NSGA2Breeder` performs breeding using the archive. Again, defined in the obvious way:

```
breed = ec.multiobjective.nsga2.NSGA2Breeder
```

You're responsible for setting up breeding pipelines like you see fit.

Note that although `NSGA2Breeder` is a subclass of `SimpleBreeder`, it cannot be used with elitism and will complain if you attempt to do so.

**Where to get examples** The package `ec.app.moosuite` has NSGA-II implementations of a number of famous multiobjective test problems. Be sure to read the `moosuite.params` file, where you can specify that you want to use NSGA-II for the problems.

### 7.4.3 SPEA2 (The `ec.multiobjective.spea2` Package)

The Strength Pareto Evolutionary Algorithm 2 (SPEA2) [22] splits the Subpopulation in two two parts: the archive and the "regular" Subpopulation. Unlike the NSGA-II algorithm, this archive can vary in size. The size of the archive is a parameter in `ec.multiobjective.spea2.SPEA2Subpopulation`:

```

pop.subpop.0 = ec.multiobjective.spea2.SPEA2Subpopulation
pop.subpop.0.size = 100
pop.subpop.0.archive-size = 50

```

Each iteration SPEA2 updates the archive to include the Pareto front, plus (if there is room) additional fit individuals using a special domination-based fitness measure called *strength*. If there are too many Individuals to fit in the archive, the archive is trimmed by removing Individuals which are too close to one another. It then uses its special fitness measure to breed individuals from the Archive and place them into the “regular” Subpopulation.

To do this, SPEA2 needs to augment the MultiObjectiveFitness with a few additional values:

```

public double strength;
public double kthNNDistance;
public double fitness;

```

The first measure is the *strength* of an Individual, defined as the number of Individuals whom it dominates. The second measure (“distance”) is an inverted measure of how far the Individual is from other Individuals in the population. The final measure (the actual *fitness*) is the sum of the so-called SPEA2 “raw fitness” and the distance measure: higher fitness values are worse.

This version of MultiObjectiveFitness is called `ec.multiobjective.spea2.SPEA2MultiObjectiveFitness`, and we include it as well, for example,

```

pop.subpop.0.species.fitness = ec.multiobjective.spea2.SPEA2MultiObjectiveFitness
multi.fitness.num-objectives = 2
multi.fitness.min.0 = 1.0
multi.fitness.max.0 = 3.5
multi.fitness.min.1 = -10
multi.fitness.max.1 = 0

```

SPEA2 also requires a special Breeder and a special Evaluator:

```

eval = ec.multiobjective.spea2.SPEA2Evaluator
breed = ec.multiobjective.spea2.SPEA2Breeder

```

When breeding, SPEA2 has a special version of TournamentSelection which only selects among archive members. We’d include it in various places instead of other selection methods, for example something like:

```

pop.subpop.0.species.pipe.source.0 = ec.multiobjective.spea2.SPEA2TournamentSelection

```

**Where to get examples** The package `ec.app.moosuite` has SPEA2 implementations of a number of famous multiobjective test problems. Be sure to read the `moosuite.params` file, where you can specify that you want to use SPEA2 for the problems.

## 7.5 Other Methods

### 7.5.1 Resets (The `ec.evolve` Package)

The `ec.evolve` package presently contains a single class, `ec.evolve.RandomRestarts`. This is a subclass of `Statistics` which performs random or timed restarts. `RandomRestarts` reinitializes your entire Population either once every fixed  $N$  generations or once every  $M$  generations where  $M$  is a random integer between 1 and  $N$  inclusive. The value of  $M$  is randomized each reset.

To use `RandomRestarts`, include it as part of your `Statistics` chain. For example, if you have a single `Statistics` object at present, you could say:

```
stat.num-children = 1
stat.child.0 = ec.evolve.RandomRestarts
```

Next you need to specify  $N$ . Here we set it to 20.

```
stat.child.0.restart-upper-bound = 20
```

Last you'll need to state whether we reset exactly every  $N$  generations in some random number of generations between 1 and  $N$  inclusive (this random value changes every reset). The options are `fixed` and `random`. Here we set the reset type to `fixed`:

```
stat.child.0.restart-type = fixed
```

When it happens, population reinitialization occurs immediately after the checkpoint is incremented and checkpointing happens. In generational statistics, the hook for this is `Statistics's preEvaluationStatistics(...)` method call. In steady state statistics, the hook is `generationBoundaryStatistics(...)`.

Random Restarts is due to James O'Beirne, a student at GMU.



# Bibliography

- [1] Kumar Chellapilla. A preliminary investigation into evolving modular programs without subtree crossover. In John R. Koza, Wolfgang Banzhaf, Kumar Chellapilla, Kalyanmoy Deb, Marco Dorigo, David B. Fogel, Max H. Garzon, David E. Goldberg, Hitoshi Iba, and Rick Riolo, editors, *Genetic Programming 1998: Proceedings of the Third Annual Conference*, pages 23–31, University of Wisconsin, Madison, Wisconsin, USA, 22–25 July 1998. Morgan Kaufmann.
- [2] Kalyanmoy Deb, Amrit Pratap, Sameer Agarwal, and T. Meyarivan. A fast elitist non-dominated sorting genetic algorithm for multi-objective optimization: NSGA-II. In Marc Schoenauer, Kalyanmoy Deb, Günther Rudolph, Xin Yao, Evelyne Lutton, Juan Julian Merelo, and Hans-Paul Schwefel, editors, *Parallel Problem Solving from Nature (PPSN VI)*, pages 849–858. Springer, 2000.
- [3] John R. Koza. *Genetic Programming: On the Programming of Computers by Means of Natural Selection*. MIT Press, 1992.
- [4] John R. Koza. *Genetic Programming II: Automatic Discovery of Reusable Programs*. MIT Press, 1994.
- [5] Sean Luke. Genetic programming produced competitive soccer softbot teams for robocup97. In John R. Koza, Wolfgang Banzhaf, Kumar Chellapilla, Kalyanmoy Deb, Marco Dorigo, David B. Fogel, Max H. Garzon, David E. Goldberg, Hitoshi Iba, and Rick Riolo, editors, *Genetic Programming 1998: Proceedings of the Third Annual Conference*, pages 214–222, University of Wisconsin, Madison, Wisconsin, USA, 1998. Morgan Kaufmann.
- [6] Sean Luke. *Essentials of Metaheuristics*. 2009. Available at <http://cs.gmu.edu/~sean/book/metaheuristics/>.
- [7] Sean Luke, Cladio Cioffi-Revilla, Liviu Panait, Keith Sullivan, and Gabriel Balan. MASON: A multiagent simulation environment. *Simulation*, 81(7):517–527, July 2005.
- [8] Sean Luke and Liviu Panait. A survey and comparison of tree generation algorithms. In Lee Spector, Erik D. Goodman, Annie Wu, W. B. Langdon, Hans-Michael Voigt, Mitsuo Gen, Sandip Sen, Marco Dorigo, Shahram Pezeshk, Max H. Garzon, and Edmund Burke, editors, *Proceedings of the Genetic and Evolutionary Computation Conference (GECCO-2001)*, pages 81–88, San Francisco, California, USA, 7–11 July 2001. Morgan Kaufmann.
- [9] Sean Luke and Liviu Panait. A comparison of bloat control methods for genetic programming. *Evolutionary Computation*, 14(3):309–344, Fall 2006.
- [10] Makoto Matsumoto and Takuji Nishimura. Mersenne twister: a 623-dimensionally equidistributed uniform pseudo-random number generator. *ACM Transactions on Modeling and Computer Simulation*, 8(1):3–30, 1998.
- [11] Una-May O’Reilly. *An Analysis of Genetic Programming*. PhD thesis, Carleton University, Ottawa-Carleton Institute for Computer Science, Ottawa, Ontario, Canada, 22 September 1995.

- [12] Liviu Panait. A comparison of two competitive fitness functions. In *GECCO 2002: Proceedings of the Genetic and Evolutionary Computation Conference*, pages 503–511. Morgan Kaufmann Publishers, 2002.
- [13] Ricardo Poli. A simple but theoretically-motivated method to control bloat in genetic programming. In *Genetic Programming, Proceedings of EuroGP'2003*, pages 204–217. Springer, 14-16 April 2003.
- [14] Riccardo Poli, William B. Langdon, and Nicholas Freitag McPhee. *A Field Guide to Genetic Programming*. Available in print from lulu.com, 2008.
- [15] Kenneth Price, Rainer Storn, and Jouni Lampinen. *Differential Evolution: A Practical Approach to Global Optimization*. Springer, 2005.
- [16] Bill Punch and Douglas Zongker. lil-gp 1.1. A genetic programming system. Available at <http://garage.cse.msu.edu/software/lil-gp/>, 1998.
- [17] Conor Ryan, J. J. Collins, and Michael O'Neill. Grammatical evolution: Evolving programs for an arbitrary language. In *EuroGP 1998*, pages 83–96, 1998.
- [18] Lee Spector. Simultaneous evolution of programs and their control structures. In Peter J. Angeline and K. E. Kinnear, Jr., editors, *Advances in Genetic Programming 2*, chapter 7, pages 137–154. MIT Press, 1996.
- [19] N. Srinivas and Kalyanmoy Deb. Multiobjective optimization using nondominated sorting in genetic algorithms. *Evolutionary Computation*, 2:221–248, 1994.
- [20] Keith Sullivan, Sean Luke, Curt Larock, Sean Cier, and Steven Armentrout. Opportunistic evolution: efficient evolutionary computation on large-scale computational grids. In *GECCO '08: Proceedings of the 2008 GECCO conference companion on Genetic and evolutionary computation*, pages 2227–2232, New York, NY, USA, 2008. ACM.
- [21] Seth Tisue and Uri Wilensky. Netlogo: A simple environment for modeling complexity. In *International Conference on Complex Systems*, pages 16–21, 2004.
- [22] Eckart Zitzler, Marco Laumanns, and Lothar Thiele. SPEA2: Improving the strength pareto evolutionary algorithm for multiobjective optimization. In K. Giannakoglou, D. Tshalis, J. Periaux, K. Papailiou, and T. Fogarty, editors, *Evolutionary Methods for Design, Optimization, and Control*, pages 19–26, 2002.