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# A Species-Conserving Genetic Algorithm for Multimodal Optimization

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A Species-Conserving Genetic Algorithm for  
Multimodal Optimization

by

Michael Scott Brown

A dissertation submitted in partial fulfillment of the requirements  
for the degree of Doctor of Philosophy  
in  
Computer Science

Graduate School of Computer and Information Sciences  
Nova Southeastern University

2010

We hereby certify that this dissertation, submitted by Michael Scott Brown, conforms to acceptable standards and is fully adequate in scope and quality to fulfill the dissertation requirements for the degree of Doctor of Philosophy.

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Graduate School of Computer and Information Sciences  
Nova Southeastern University

2010

An Abstract of a Dissertation Submitted to Nova Southeastern University  
in Partial Fulfillment of the Requirements for the Degree of Doctor of Philosophy

## A Species-Conserving Genetic Algorithm for Multimodal Optimization

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The problem of multimodal functional optimization has been addressed by much research producing many different search techniques. Niche Genetic Algorithms is one area that has attempted to solve this problem. Many Niche Genetic Algorithms use some type of radius. When multiple optima occur within the radius, these algorithms have a difficult time locating them. Problems that have arbitrarily close optima create a greater problem. This paper presents a new Niche Genetic Algorithm framework called Dynamic-radius Species-conserving Genetic Algorithm. This new framework extends existing Genetic Algorithm research.

This new framework enhances an existing Niche Genetic Algorithm in two ways. As the name implies the radius of the algorithm varies during execution. A uniform radius can cause issues if it is not set correctly during initialization. A dynamic radius compensates for these issues. The framework does not attempt to locate all of the optima in a single pass. It attempts to find some optima and then uses a tabu list to exclude those areas of the domain for future iterations. To exclude these previously located optima, the framework uses a fitness sharing approach and a seed exclusion approach. This new framework addresses many areas of difficulty in current multimodal functional optimization research.

This research used the experimental research methodology. A series of classic benchmark functional optimization problems were used to compare this framework to other algorithms. These other algorithms represented classic and current Niche Genetic Algorithms.

Results from this research show that this new framework does very well in locating optima in a variety of benchmark functions. In functions that have arbitrarily close optima, the framework outperforms other algorithms. Compared to other Niche Genetic Algorithms the framework does equally well in locating optima that are not arbitrarily close. Results indicate that varying the radius during execution and the use of a tabu list assists in solving functional optimization problems for continuous functions that have arbitrarily close optima.

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This dissertation is dedicated to the three ladies in my life, Truocmai Dinh, Jessica Maikhanh Brown and Tiffany Maitam Brown, my wife and two daughters. All that I do, I do for you.

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## **Chapter 1**

### **Introduction**

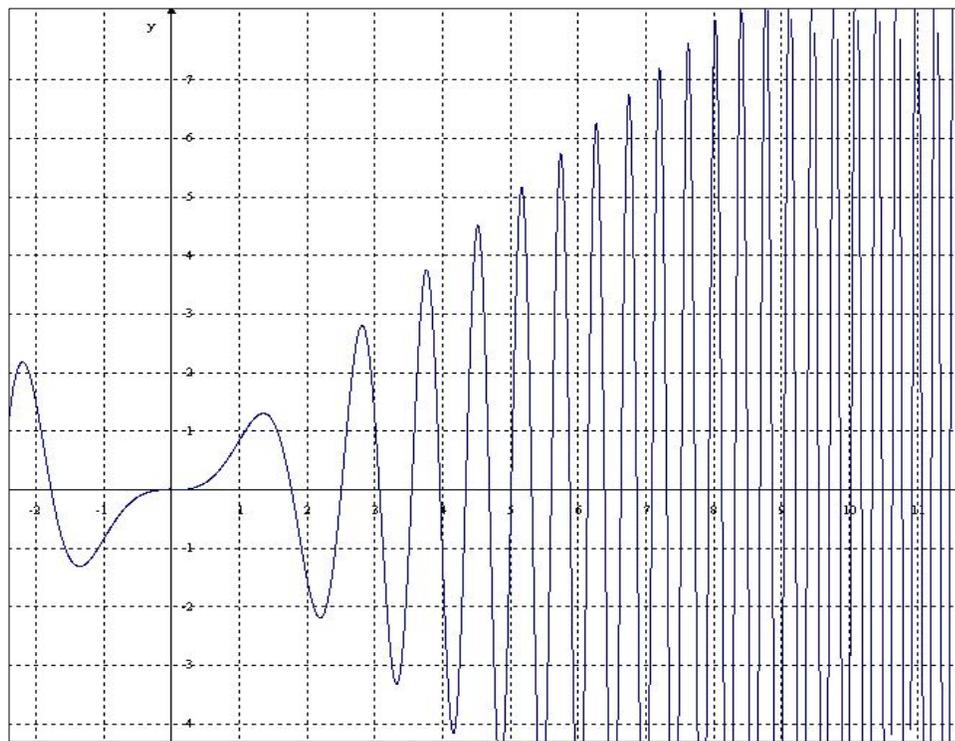
This introductory chapter is organized into seven sections. The first section defines the problem that this research addressed. The second section describes the research goal. The third section provides an overview of the approach. The fourth section explains the relevance and significance. The fifth section describes the barriers and issues that need to be overcome by this research. The sixth section provides definitions of terms used in the dissertation. The final section is a summary of this chapter.

### **Problem Statement**

Genetic Algorithms (GA) have a difficult time solving problems with multiple correct answers. When traditional GAs attempt to solve multimodal problems they often converge to only one of the possible correct or good solutions. A current area of research in GAs is called Niche Genetic Algorithms (NGA), which hopes to address this problem. NGAs can be used to solve problems that seek local optima where multiple exist.

Currently there are many NGAs. Two prominent approaches to developing NGAs are crowding and sharing (Deb & Goldberg, 1989). In crowding algorithms, members of one population coexist with members of the next population. Older individuals of the population are selected for removal based on how similar they are to newer members. A variety of NGAs use some type of crowding scheme (Cavicchio, 1970; De Jong, 1975; Jelasity & Dombi 1998; Li, Balazs, Parks & Clarkson, 2002; Ling, Wa, Yang & Wang,

2008; Raghuwanshi & Kakde, 2007). The second approach is through fitness sharing. In sharing schemes, the fitness of an individual is dependent on its distance to other individuals in the population. This increases the chance that species will form around niches by rewarding genetic isolation. There are many methods based on sharing schemes (Beasley, Bull & Martin, 1993b; Bernier, 1996; Goldberg & Richardson, 1987; Holland, 1975). While not every NGA is a crowding or sharing method, most fall into one of these two categories.



**Figure 1.** Graph of  $y = x \sin(x^2)$

Both of these approaches use some form of distance in determining what individuals perform crossover or which individuals are promoted into the next generation. There are some domains where distance is not a good indicator of niches (Ando & Kobayashi, 2005). Clearly these approaches are good if there are significant distances between niches. What is unclear is the effectiveness of these approaches when niches become

arbitrarily close. Consider functional optimization problems for continuous functions, specifically, classes of functions that have local optima that become arbitrarily close. No matter what distance a NGA expects between niches there exists an area in the domain in which niches are smaller than this distance. Figure 1 shows a graph of one such function,  $y = x \sin(x^2)$ . As  $x$  gets larger the local maximums or optima become arbitrarily close together.

### **Research Goal**

The goal of this research was to develop a new NGA that can address these types of problems for functional optimization of continuous functions. This NGA is a framework based on an existing NGA, but allows different components to be used in combination to create different algorithms. This new approach is not dependent on a static niche radius parameter that could provide poor results if selected wrong. As part of this goal the algorithm should be flexible enough to solve other types of problems that current traditional NGAs solve. This research goal will expand our current understanding of NGAs.

This dissertation addressed of few areas that make up the goal. To accomplish the goal a new NGA was developed and tested against existing NGAs. The areas that this research addressed are as follows:

- To develop a new NGA that will solve for arbitrarily close optima
- To compare this new NGA to existing NGAs to determine its effectiveness

These two research goals complement the goals defined in this chapter.

There were two hypotheses to this research. The first hypothesis was that locating optima in phases increases a NGA's ability to find optima. In the first phase the NGA finds some optima. Once some optima are located, they are used to encourage exploration in other areas of the domain. Multiple attempts to find optima are performed. Each attempt leverages already located optima, which should make it easier to locate other optima because there is less of the domain space to search. This process continues until all of the optima are located.

The second hypothesis was that many NGAs miss optima, especially arbitrarily close optima, because they use a static radius. Many current NGAs have a parameter that is used to determine if individuals are within a neighborhood. If the distance between optima is smaller than this radius, the NGA has a difficult time locating all of the optima. The value that this parameter is set to greatly affects the results of the NGA. The second hypothesis was that a dynamic radius could compensate for poor radius choices. Allowing the radius to change as the algorithm runs may allow it to adapt to conditions and find more optima.

These two hypotheses complement each other. The research hypothesis was that better results can be obtained by allowing dynamic radius and restricting areas of the domain where optima have been located. This allows investigation across the entire domain which should produce better results. Both hypotheses should increase the number of optima discovered by the NGA.

## **Approach**

The new approach for solving multimodal optimization problems used existing methods in combination to enhance a traditional GA. The new approach is a NGA framework that is based on Species Conserving Genetic Algorithm (SCGA) (Li et al., 2002). The framework is presented in a modular form and has different components that can be used in combination to create different variations of the algorithm.

SCGA enhances the traditional GA with seed selection and seed conservation steps. Seeds are identified as the fittest member within a given radius. Seeds are conserved into the next generation through a seed conservation step. Seed conservation has the seed replace the weakest individual in the new generation within the radius of the seed. This ensures that these strong individuals that are tracking different optima are preserved.

The new framework does not attempt to find all of the optima within a single pass of the algorithm. Traditional GAs perform a loop with each iteration creating a new generation of the population. Environmental pressures force the population to converge. Multimodal optimization creates a difficult problem for GAs. The new approach generates a certain number of generations in hopes of finding some optima. These optima are recorded and the environmental parameters change. These changes alter how the fitness of individuals is determined and what individuals can be seeds. This encourages future generations to avoid these optima. This allows future generations to explore other areas of the domain and locate other optima. The process of locating some optima and then changing the fitness is performed multiple times. There is an outer loop and an inner nested loop. The inner loop performs a typical SCGA algorithm. Once it

completes, seeds and optima are recorded and changes are made to the algorithm parameters. The steps consisting of the inner loop, recording and changing of the algorithm's parameters are performed in an outer loop. While traditional GAs are a single loop, this new approach uses nested loops to find some of the optima that enhances its ability to locate all of the optima.

The new framework also changes the radius used to define species. In SCGA a radius is defined and used throughout the algorithm. As other research has shown there is a limitation to algorithms of this type (Ando & Kobayashi, 2005). Poor choices for the radius produce poor results. The new approach changes the radius as the algorithm runs. The algorithm attempts to compensate for poor radius choices. After each inner loop of the algorithm completes, adjustments are made to the radius. Varying the radius mitigates the issue of incorrectly set radius.

The final difference between SCGA and this new framework is the use of a tabu list. A tabu list comes from the tabu search method and is a list that contains previously evaluated areas of the domain. The tabu list is used by the algorithm to avoid these areas in the future and concentrate on areas of the domain that have not been searched. After each completion of the inner loop, the seeds and optima from that pass are recorded on a tabu list. This list is used in future loops of the algorithm to encourage exploration by avoiding these areas.

As optima are located, the algorithm adjusts to encourage exploration into other areas of the domain. This is done in two ways. The algorithm can determine that a potential individual is too close to a member of the tabu list. This will disqualify the individual from becoming a seed. The algorithm may also adjust the fitness to individuals relative

to how close they are to individuals on the tabu list. This will decrease their chances of becoming seeds. Both of these tactics are possible within the new framework.

To validate the new approach, it was compared against other NGAs using a set of well established benchmarks. The benchmarks came from a variety of NGA research. All of the benchmarks were minimization and maximization functional optimization problems. A new benchmark was also presented. These benchmarks were used to compare the new approach to other NGAs.

The new approach was compared to other NGAs in solving the benchmarks that were defined. A variety of performance criteria were used in this comparison including proportions of peaks located and average fitness of the last 50 generations. For each benchmark the new approach was compared to multiple other NGAs. These other NGAs had been selected to cover a wide range of NGA research from modern methods to early algorithms. In many cases the new approach was compared against previously published results in other NGA research. In other cases NGAs were implemented to obtain test results. The performance criteria allowed the new approach to be evaluated against other NGA methods.

The new framework leveraged a variety of existing methods to introduce a new combination of concepts to create a NGA framework. The use of a tabu list in a NGA has been used before (McLoughlin & Cedeno, 2005; Ting & Ko, 2008; Tsai, Tseng, Chiang, & Yang, 2009). The use of a dynamic radius has also been used in other algorithms (Jelasity & Dombi, 1998). The combination of a tabu list and dynamic radius applied to the SCGA algorithm is new. This new framework was compared against other NGAs using well defined benchmarks and criteria. Comparing it against many well

established NGAs showed that this new approach can solve multimodal optimization problems.

### **Relevance and Significance**

The study of GAs is important because GAs are very useful search techniques. They have been used in almost every field of study. Much literature has been dedicated to outlining uses for GAs (Coello, 2000; Dianati, Song & Treiber, 2002; Sheikh, Raghuwanshi & Jaiswal, 2008). For example, GAs and NGAs have been used in Electrical Engineering to design electromagnetic systems (Cioffi, Formisano & Martone, 2000). In the field of Knowledge Discovery they have been used as a classifier (Pozo & Hasse, 2000). For pattern matching, NGAs can be used to match handwriting (Oliveira, Sabourin, Bortolozzi & Suen, 2002; Stefano, Cioppa & Marcelli, 1999). NGA research is useful to many fields of study.

Because GAs and NGAs are applicable to many fields of study, research in the subject has continued uninterrupted for many decades. Early researchers developed simple algorithms for multimodal optimization (Deb & Goldberg, 1989; Goldberg & Richardson, 1987; Mauldin, 1984). These algorithms solved many multimodal optimization problems. A second generation of algorithms were developed that addressed limitations of the previous algorithms (Li et al., 2002; Ling et al., 2008; Raghuwanshi & Kakde, 2007). Some research addressed the limitation that many algorithms require tuning parameters (Bernier, 1995; Fonseca & Fleming, 1993). Researchers continue to investigate NGAs.

NGA research continues to this day and is an important area of research in Artificial Intelligence. Experts in the field believe that developing new NGAs is useful and justification exists for continued investigation. These new NGAs can be used in other areas of research to solve multimodal optimization and search problems.

### **Barriers and Issues**

There were two barriers to this research. The first barrier was premature convergence which traditional GAs exhibit when attempting to locate multiple optimum. The second barrier was optimum location and preservation. Let us consider how each of these areas was addressed by this research.

#### *Premature Convergence*

One barrier to developing a NGA is to prevent global convergence. A GA naturally converges to a local optimum. This is appropriate for many types of problems, but there are problems that have multiple optima. Traditional GAs will converge to a single optimum, ignoring the other ones. The key to develop a NGA is to overcome this pressure to converge. The NGA needs to allow local convergence within niches. De Jong (1975) calls this premature convergence.

Two forces act on the generations of a traditional GA. Crossover of individuals puts pressure on the population to converge through different individuals having different probabilities of reproducing (De Jong, 1975). The algorithm exploits fit individuals in the creation of each generation. An opposite force works against this exploitation. Mutation alters individuals, which allow exploration of new areas of the domain

(Beasley, Bull & Martin, 1993a). NGAs need to balance the two forces of exploration and exploitation. Too much exploration will decrease the performance of the search, turning it into a random search (Holland, 1992). Limiting exploration too much, in favor of exploitation, leads to premature convergence (De Jong, 1975). Successful NGAs balance exploration and exploitation to locate multiple optima.

As a GA generates individuals using crossover, the amount of the domain that is being searched decreases. This is referred to as genetic drift and removes areas of the search space so greatly that even mutation cannot put them back (De Jong, 1975). This eliminates other possible solutions to the problem. In problems that have a single correct answer this convergence helps in solving the problem by eliminating areas of the domain in which the correct answer does not exist. But in multi-objective problems, it eliminates other optima.

There are two methods in GA research that could address this problem. One is to have a very large population size  $N$ . If  $N$  is very large, the GA has much more time to locate other optima before the genetic drift closes the search space. However, this leads to performance problems (De Jong, 1975). A second approach is to have a very high mutation rate. This would allow the expanding of the search space when genetic drift happens. The problem with very high mutation rates is that it prevents convergence, which is the ultimate goal of the GA and the way that the GA finds the solution (De Jong, 1975). This is why De Jong, Holland, Goldberg and other researchers believe that traditional GAs will not solve multimodal problems (De Jong, 1975; Holland, 1975; Goldberg & Richardson, 1987).

### *Optimum Preservation*

The second barrier to this research was optimum preservation. This barrier can be thought of in two parts. The first part is optimum location, identifying the areas of the domain worth preserving. Because the domain has not fully been searched when optimum location is applied, this is difficult for the NGAs. The second part is how to preserve or conserve these areas. Because of crossover and mutation, there is little guarantee that these areas of interest will be represented in the next generation. Optimum preservation is essential in NGAs.

Optimum location attempts to identify individuals within a population that can eventually lead to an optimum (De Jong, 1975). These individuals are normally individuals at, or close to, an optimum. Selection in traditional GAs focuses on the fittest members. But in multimodal optimization problems, it is possible that less fit members are also tracking a local optimum. A single generation of a population represents a very small part of the domain. Locating these individuals makes optimum location challenging.

Optimum preservation is used to ensure that the optima located are not eliminated in the population through convergence. Convergence pressures of GAs can eliminate optimum after they are discovered. The method to ensure that an optimum is preserved can be direct or subtle (Li et al., 2002). There are direct approaches like promoting an individual of interest into the next generation. More subtle approaches can be to adjust the individual's fitness to increase its chances of being selected for crossover. Regardless of the method, these located optimum need to be preserved into the next generation.

Optimum preservation is a barrier that every successful NGA needs to overcome. Some part of the algorithm needs to identify interesting individuals in the population and allow their representation in future generations (Li et al., 2002). There are a variety of methods that can be employed for optimum preservation. These methods will be described in the literature review of Chapter 2.

### **Definition of Terms**

This dissertation uses a variety of terms. Many terms are generally known in the GA field. The following list of terms should provide an overview of terms used in this dissertation.

*Baldwin Effect:* The Baldwin Effect is a biological theory that the fitness of an individual can be changed by environmental factors (Baldwin, 1896).

*Cluster:* A cluster is a set of items that share something in common. Within a cluster items should have commonality and items in different clusters should have differences (Sheikh, Raghuwanshi & Jaiswal, 2008).

*Convergence:* Convergence is a process in which new generations of a population have decreased genetic diversity. This typically occurs around an optimum.

*Crossover:* Crossover is a genetic operation that takes two individuals of a population and by interchanging genes between the two individuals creates two new individuals.

*Evolutionary Algorithms:* Evolutionary algorithms are a classification of algorithms based upon natural evolution. There are four subclasses of evolutionary algorithms:

Evolutionary Programming, Evolutionary Strategies, Genetic Algorithms and Genetic Programming (Dianati, Song & Treiber, 2002).

*Fitness:* Fitness is a measurement assigned to an individual of a population that relates to how well the individual copes with environmental pressure (De Jong, 1975).

*Genetic Algorithm:* A Genetic Algorithm is a specific type of search method that was developed by Bremermann (1958). The algorithm models the domain as a series of gene values. An initial generation of the population is created, normally randomly, of different combinations of these gene values. Genetic operations are applied to the generation to create a new generation. Over time the population converges to the optimum of the domain.

*Genetic Drift:* Genetic Drift is the change in probability or frequency that a certain gene value appears in a population (De Jong, 1975). As populations evolve certain gene values become more prevalent.

*Inversion:* Inversion is a genetic operation in which the ordering of the genes change (Holland, 1975).

*Mutation:* Mutation is a genetic operation in which a gene value is randomly changed based upon the mutation rate.

*Niche Genetic Algorithm:* A Niche Genetic Algorithm is a specific type of Genetic Algorithm that promotes genetic diversity (Mahfoud, 1995).

*Pareto Front:* The Pareto Front is the set of non-dominant optimal values for a multi-objective optimization problem (Alba, Dorronsoro, Luna, Nebro & Bouvry, 2005).

*Particle Swarm Optimization:* Particle Swarm Optimization is a specific search technique that simulates swarm intelligence.

*Premature Convergence:* Premature convergence is when a Genetic Algorithm converges at such a rate that optima are removed from the search space (De Jong, 1975).

*Seed:* A Seed is a dominant individual within a certain area of the domain in a population (Li et al., 2002).

*Selection:* Selection is a process of selecting individuals of a population to reproduce.

*Species:* The term species has different definitions in NGA research. In this research the Li et al. (2002) definition will be used. Species are individuals within a population, whose distance is less than some, pre-define parameter (Li et al., 2002).

This section contains definitions of terms used in this dissertation.

## **Summary**

This research had a few specific goals. The research produced a new NGA framework. This framework allows for the creation of multiple variations of the NGA algorithm. The problem that the new NGA addresses is functional optimization for continuous functions. Within this area, the goal is to solve for functions that have arbitrarily close optima. These types of functions are especially difficult for NGAs. A secondary goal was for the algorithm to solve other types of optimization problems equally well as other NGAs. The new NGA framework was developed to accomplish these goals.

The approach that was taken created a new NGA framework that applies existing techniques to NGA research. The NGA uses multiple passes in an effort to locate some optima and uses those optima in locating the other ones. The algorithm varies the radius

used in determining seeds. A tabu list is used to store located optima and seeds, so these areas of the domain are not revisited.

## Chapter 2

### Review of the Literature

This literature review covers the history and current state of NGA research. It is organized into five sections. The first section describes research that is relevant to this research, but that is not an NGA. NGA research is organized in the next three sections. There are sharing methods, crowding methods and a section for methods that do not fall into either of these groups. Finally, there is a summary section.

#### Relevant Research Other than NGAs

The framework presented in this paper leverages many other concepts in existing NGA research. However, some concepts of the framework come from other research areas. This section provides a literature review of other research that was influential in developing this framework.

#### *Tabu Search*

The new algorithm presented in this paper uses aspects of the tabu search. A tabu search is an optimization technique used to avoid local optima (Glover, 1989; Glover, 1990a). It has been used to solve several optimization problems (Glover, 1990b; Hansen, 1997). The tabu search has an associated memory structure that is used to store previous moves in the optimization process. This list is used to prevent the algorithm from returning to previously obtained optima.

A tabu search begins with a possible solution to the optimization problem. Each iteration of the algorithm will apply an operation that will move from one solution state to a new one. Table 1 shows the part of the tabu search algorithm that determines if a move should be made.

Table 1. Tabu Search Decision

Line Number	Pseudocode
1	Select a move
2	If the move is on the tabu list then
3	If the move satisfies the aspiration condition
4	Make the move
5	Else
6	Select another move
7	Else
8	Make the move

The operation is added to the tabu list. Future iterations of the algorithm prevent the operation from being applied, unless an aspiration condition is met. The aspiration condition determines if the move is superior to the current solution. By using the tabu list, the tabu search avoids local optima and locates the optimal solution to the problem.

The tabu search uses a short-term memory structure to track previous moves (Glover, 1990b). This memory structure is used to prevent the algorithm from revisiting previously visited states. The tabu list can be finite or infinite in length. A finite tabu list only stores a certain number of previous moves. When the list is full the oldest move will be removed when a new move is added. The algorithm prevents these moves from being made in the future. If the tabu list is finite, then the move can only be made after the previous move is purged from the tabu list. The tabu list encourages exploration.

The aspiration condition is used in a tabu search to override the tabu list (Glover, 1990b). If a move is on the tabu list, it is normally prohibited. But before the potential

move is eliminated, it is compared to an aspiration condition. If the move meets the aspiration condition, the move is performed. This allows superior moves from being eliminated because they are tabu.

The tabu search is a useful search technique. Previous research has shown that combining the tabu search with evolutionary algorithms can increase its accuracy. Tabu searches have been combined with GAs (McLoughlin & Cedeno, 2005; Ting & Ko, 2008; Tsai, et al., 2009). They have also been combined with Evolutionary Programming algorithms (Rajan & Mohan, 2004). The tabu search can provide valuable insight into solving multimodal optimization problems.

### **Fitness Sharing Methods**

A common approach to solving multimodal optimization problems is through fitness sharing. While methods for fitness sharing vary, they all alter the fitness function in some way to encourage genetic diversity. In multimodal functional optimization problems, fitness is normally directly related to the objective function. In Sharing Methods distance to other individuals is incorporated into the fitness function to encourage exploration. This prevents a single optimum from dominating the population. Some of the earliest approaches for NGA algorithms are sharing methods.

*Holland, J. H.*

Holland (1975) provides a formal framework for GA research. While it does not provide a specific NGA algorithm, it does describe some of the earlier fitness sharing concepts. Holland describes a two-armed bandit to represent the problem that can be

solved by NGAs. A two-armed bandit is a slot machine that has two handles, instead of one. Each handle has a different payout. Players may elect to pull the left or right handle. Ideally, every player would select to pull the handle with the highest payout, but there is a catch. For a given turn all of the players that select a given handle must share the payout. With this new rule the problem is not obvious which handle the players should select. Each handle is a niche and by dividing the payout, or fitness, between all individuals within a niche allows the GA to solve the problem. By defining the fitness function in such a way as to reflect other individuals in the niche, allows a traditional GA to solve for multimodal optimization problems. This is some of the earliest research in fitness sharing.

#### *Goldberg and Richardson*

Another seminal work in NGA is Goldberg and Richardson (1987). This NGA introduces a sharing function. In traditional GAs fitness functions determine the probability a member of a population will reproduce. In a multimodal problem once a traditional GA discovers a niche, it converges on it, ignoring other possible niches. A sharing function is used to reduce this convergence by using the shared fitness to determine the probability that a member will reproduce. Shared fitness penalizes individuals that are close to other individuals in the population and rewards isolated individuals. This allows the NGA to locate other niches.

In the Goldberg and Richardson (1987) method the algorithm is the same as a traditional GA, except for determining the fitness function. The algorithm uses a *shared fitness function* that accepts the distance between two members as an input parameter.

These functions must conform to three properties. The function's output must be between zero and one. When the distance is zero, the output must be one. When the distance approaches infinity, the output must be zero. When shared fitness is computed for individual  $p$ , a *niche count* is calculated by summing the sharing function of all the other members of the population with respect to  $p$ . The shared fitness is the individual's raw fitness divided by the niche count.

While the Goldberg and Richardson (1987) fitness sharing algorithm can take many forms, their research presents an example of the algorithm. The example attempts to locate the five local optima of the function  $f(x) = \sin^6(5.1 \pi x + 0.5)$ , where  $x$  is between 0 and 1. The shared fitness function selected was the power law function, which is shown below.

$$sh(d) = \begin{cases} = 1 - \left( \frac{d}{\sigma_{share}} \right)^\alpha & \text{if } d < \sigma_{share} \\ = 0 & \text{otherwise} \end{cases}$$

The parameters  $\sigma_{share}$  and  $\alpha$  are set to 0.1 and 1 respectively. The niche count,  $m_i$ , for individual  $i$  is represented by the following function.

$$m_i = \sum_{j=1}^N sh(d(x_i, x_j))$$

In the niche count  $x_i$  is individual  $i$  and  $x_j$  spans all individuals in the population  $N$ . The shared fitness of an individual is simply  $f_i' = f_i / m_i$ , where  $f_i'$  is the shared fitness,  $f_i$  is the raw fitness and  $m_i$  is the niche count. The research results showed that traditional GAs only found one optima of the function. The sharing fitness algorithm found all five

optima and had an equal number of individuals at each optimum (Goldberg & Richardson, 1987).

### *Sequential Niche Technique*

The Sequential Niche Technique (SNT) is a search technique that can be applied to a GA (Beasley, Bull & Martin, 1993b). It attempts to locate one optimum at a time. Once an optimum is located the technique adjusts the search algorithm to locate another optimum. The technique is successful because it reduces the search problem into locating a single optimum.

When applying SNT to a GA, the traditional steps of a GA are used. The fitness function, which typically is the objective function, is modified. This modified fitness function is used in the algorithm. After the algorithm runs, the best individual is recorded on a list. The modified fitness function is changed by adding a *derating function* for the fittest individual that was located. The *derating function* can take many forms, but its affect is to decrease the fitness around the located individual. This excludes this area of the domain as a place for likely optima. Future runs of the GA seek out other optima. SNT also has a *solution threshold*. If the fittest individual after each run is more fit than the *solution threshold*, it is considered an optimum. The algorithm for SNT is shown in Table 2.

SNT allows search algorithms to use previous knowledge about the problem to simplify it. This approach is attributed to other functional optimization research (Ackley, 1987). It is a useful technique that allows search algorithms to take complex problems

like multimodal functional optimization and break them into a series of much simpler problems of single functional optimization.

Table 2. Sequential Niche Technique Algorithm

Line Number	Pseudocode
1	Assign <i>modified fitness function</i> to objective function
2	While not termination condition
3	Run traditional GA using <i>modified fitness function</i>
4	After GA runs record the optimum that it finds
5	Depress optimum area in the <i>modified fitness function</i>
6	If optimum in step 4 is larger than <i>solution threshold</i> , display it as a solution
7	End loop

#### *Bernier's BDM and BPM*

The Bernier (1996) method uses a Minimum Spanning Tree (MST) for fitness sharing. It is used in each iteration of the NGA to adjust the fitness of individuals. There are two algorithms for the method: Biggest Different Method (BDM) and Biggest Proportion Method (BPM). BDM and BPM use Prim's MST algorithm although any MST algorithm might be used.

The Prim's MST algorithm is used to determine a tree,  $T$ , with minimum total weight from a graph,  $G$ . Graphs have vertices,  $V$ , and edges,  $E$ , that connect two vertices. Every edge has an associated weight,  $W$ . Because an edge connects two vertices, we can represent it as  $(u, v)$  where  $u$  and  $v$  are vertices. Prim's algorithm begins by randomly selecting a vertex for the tree  $T$ . Then it computes the weight from every vertex in  $T$  to every vertex not in  $T$  and selects the one with the minimum weight. The selected vertex and associated edge are added to  $T$ . Prim grows the minimum tree, starting with a single vertex, into a MST.

Bernier adapts MST to GAs. Individuals are represented by vertices. The weight of the edge between two individuals is defined as their Euclidean distance. Bernier's hypothesis is that by removing some number of the largest edges of the MST, what is left will be trees around each niche. Bernier offers two methods to determine what edges should be removed. The BDM looks at the longest 15% of the edges in the MST. It sorts these edges in descending order according to their weight,  $w(e_1), w(e_2) \dots w(e_n)$ . The algorithm computes the weight difference between consecutive edges, so  $\Delta_1 = w(e_1) - w(e_2)$ ,  $\Delta_2 = w(e_2) - w(e_3)$ , ...  $\Delta_{(n-1)} = w(e_{(n-1)}) - w(e_n)$ . Finally the algorithm locates the largest  $\Delta$ ,  $\Delta_x$ . All of the edges,  $w(e_1)$  through  $w(e_x)$  are removed leaving  $x + 1$  trees. Each tree corresponds to a niche. The BPM is very similar to the BDM. Instead of comparing differences between edges, it compares proportions. The top 15% of edges are sorted in descending order. Proportions are computed by dividing consecutive edges,  $p_1 = w(e_1) / w(e_2)$ ,  $p_2 = w(e_2) / w(e_3)$ , ...  $p_{(n-1)} = w(e_{(n-1)}) / w(e_n)$ . The edges with the largest proportion are removed. BDM and BPM adjust the fitness of the individuals around niches using standard fitness sharing techniques.

Results from Bernier's algorithm are very impressive. In six benchmark functions BDM and BPM located nearly 90% of the optima. A goal of this research was to develop a NGA that does not need parameters. What is unclear is how the MST parameter of 15% affects the final results. It would seem that if there were more optima than 15% of the number of individuals in the population, this algorithm would have difficulties. If there were more optima than 15% of the population, then some optima would not have their fitness adjusted through fitness sharing. Another case could be a situation where there were relatively few optima but a very large population size. Considering so many

edges could lead the algorithm into forming too many niches. Perhaps this percentage should be a parameter.

### *Fitness Sharing Summary*

This section describes many fitness sharing NGAs. Some very important concepts come out of fitness sharing methods. All of the algorithms use the idea of altering the fitness function to guide the direction that the next generation will take. SNT introduces a concept of locating a single optimum and then using the fitness function to exclude it in future generations. Fitness sharing is a useful technique to encourage exploration across the domain space.

### **Crowding Methods**

Crowding methods are another common approach to developing NGAs. Crowding methods replace members of one generation with members of a previous generation based upon their similarity. They promote genetically diverse individuals and encourage exploration across the domain space. A variety of crowding methods have been successful with multimodal functional optimization problems.

### *Cavicchio*

Cavicchio's (1970) research looks at selection schemes to solve multimodal optimization problems. This research is some of the earliest work in the NGA area. Cavicchio introduces a series of selection schemes. In Cavicchio's NGA a certain number of the fittest individuals are carried over into the next generation. The number of

individuals carried over into the next generation is a parameter of the algorithm.

Offspring also have to compete to be placed into the next generation.

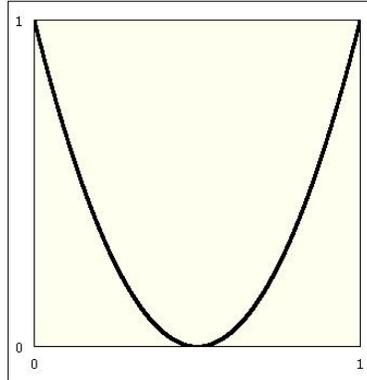
Cavicchio (1970) introduces three *Preselection Schemes*. The first scheme is based on an observation that many offspring are fit enough to be placed into the next generation, but not more fit than their parents. Allowing this seems to be counterproductive. So, the first scheme requires offspring to be more fit than both of their parents to be introduced into the next generation. The second scheme enhances the first scheme, but adds the requirement that the worst parent is to be removed from the population. The third scheme only requires an offspring to be more fit than one of its parents. Preselection is one of the earliest forms of NGAs.

Because of hardware limitations of the 1970s, tests on Cavicchio's algorithm were limited to very few individuals. In many cases population sizes were between 10 and 20 individuals (Cavicchio, 1970). Little research has been published with benchmarks on Cavicchio's NGA since the original research. It is difficult to determine how this NGA would perform with more modern benchmarks.

### *De Jong*

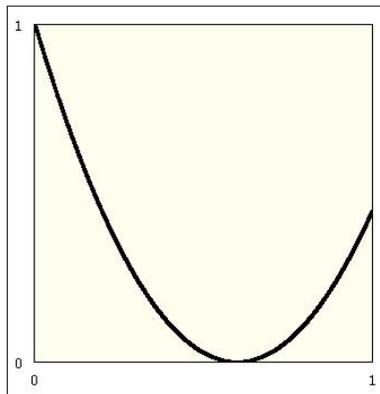
Some of the earliest works to address the problem of GAs converging globally even on multimodal domains were from De Jong (1975) and Holland (1975). De Jong's Elitist Model R2 introduces the strategy of including the best members of one generation in the next generation. After each generation is created, its least fit members are replaced by an equal number of the fittest members of the previous generation. The Elitist Model R2 replaces only one member from the previous population, but this idea can be expanded to

some predetermined fixed number of individuals. This influence can be seen in later NGAs (Li et al., 2002).



**Figure 2.** Graph of  $f(x) = 4(x - 0.5)^2$

It is easy to see why an elitist strategy would work for domains that have equally fit optima. Consider the equation  $f(x) = 4(x - 0.5)^2$ , where  $0 \leq x \leq 1$  as shown in Figure 2. Here there are two optima of equal fitness. A traditional GA will converge to one optimum or the other, but not both. An elitist strategy would preserve individuals of both optima. If considering an equation like  $f(x) = 2.8(x - 0.6)^2$ , where  $0 \leq x \leq 1$ , elitist strategies seem less useful. This equation is shown in Figure 3. Here there are two local optima,  $x = 0$  and  $x = 1$ . However,  $x = 1$  is a local, but not global optimum.



**Figure 3.** Graph of  $f(x) = 2.8(x - 0.6)^2$

An elitist strategy will probably not preserve individuals near the  $x = 1$  local optimum. These equations come from Goldberg and Richardson (1987).

Another NGA method described by De Jong (1975) is the Crowding Factor Model R5. The Crowding Factor Model R5 simulates an environment in which parents and offspring live together. To keep the population size stable, the system allows certain parents to die. This is done in crowded areas of the domain. In this method populations overlap one another. The Crowding Factor R5 method uses two parameters. *Generation gap* is the proportion of a population that is produced each generation. In De Jong's example it was 0.1, which means that the NGA produces enough individuals to increase the population size by 10% each generation. The second parameter is the *crowding factor*. For each new member of the population, an old member must be removed. The crowding factor is the number of old members considered for removal for each new member. In De Jong's example it was three. This NGA will randomly evaluate three old members for each new member. Of the three old members the one that is genetically similar to the new member is removed.

#### *Genetic Algorithm with Species*

Genetic Algorithm with Species (GAS) extends the crowding method concept by defining species (Jelasity & Dombi, 1998). Species are groups of individuals that are tracking a common optimum. Like other crowding methods GAS only allows crossover with individuals within the same species. This algorithm allows the population size to expand and contract for each generation. It also introduces the concept of individuals dying off. Traditional GAs have a generation die when the next generation is created.

GAS allows fit individuals to live longer than weak ones by allowing them to be members of multiple generations.

In GAS a species is defined as the triplet  $(o, l, S)$  (Jelasity & Dombi, 1998). The variable  $S$  is the population that makes up the species and  $o$  is the individual with the maximum fitness within the species. GAS uses a decreasing radius. The equation,  $R$ , defines the radius values as it decreases.  $R$  must always be greater than the maximum distance between two individuals and should approach 0. There is an index associated with the various radii, called the radius index. In the species  $l$  is the radius index when the species was defined. The radius for a given triplet  $(o, l, S)$  is  $R(l)$ .

The algorithm shown in Table 3 outlines the steps that GAS uses to create a new generation (Jelasity & Dombi, 1998).  $T$  is the current population and  $MP$  is a parameter that determines how large  $T$  can become.  $MP$  is not the size of  $T$ , rather the upper limit to the size of  $T$ . The algorithm for generation  $T$  will loop until the size of  $T$  is greater than  $MP$ . Within the loop two parents are selected,  $p1$  and  $p2$ . They produce two offspring,  $o1$  and  $o2$ . Parents and offspring are put back into the population.

Table 3. Genetic Algorithm with Species Algorithm

Line Number	Pseudocode
1	While (population size of $T < MP$ )
2	Select two parents, $p1$ and $p2$ within the same $(o, l, S)$
3	Create two offspring, $o1$ and $o2$
4	Put $p1, p2, o1$ and $o2$ back in population, $T$
5	End while loop
6	Dying off phase
7	Fusion

When the size of  $T$  reaches  $MP$ , the algorithm initiates a dying off phase. GAS uses a transformed fitness function  $f'$  to determine which individuals should die (Jelasity & Dombi, 1998). The function  $f'$  is defined as the following equation.

$$f'(e) = \frac{f(e) - f(\text{weakest\_individual\_in\_}S)}{\text{SizeOf}(S)}$$

For a given individual  $e$ ,  $f'(e)$  is calculated as the difference between  $e$ 's fitness and the fitness of the weakest individual in  $S$ , divided by the size of  $S$ . So, species with large population size will have a greater chance of having members die.

GAS has a process to decrease the number of species. This is called *Fusion*. After the dying off phase, GAS evaluates the existing species. If two are too close they are combined into a single species. A parameter is defined as *strict*, which is a radius index that determines how close two species must be in order to be merged into a single species. When two species are merged all of their members become members of the new species. If two species  $(o_1, l_1, S_1)$  and  $(o_2, l_2, S_2)$  are fused, the new  $o$  is the  $o_1$  or  $o_2$  that has the greatest fitness. The new  $l$  is the minimum of  $l_1$  and  $l_2$ . The new  $S$  is the union of  $S_1$  and  $S_2$ .

### *Species Conserving Genetic Algorithm*

Li et al. (2002) developed a NGA method called Species Conserving Genetic Algorithm (SCGA). It investigates how the concept of *elitism* can be applied to NGAs. This method differs with traditional GAs in two ways. Once a population is created, species are defined around individuals called *seeds*. A step to preserve species into the next generation is added to the usual selection, crossover and mutation found in GAs.

These two enhancements to the traditional GA algorithm allow SCGA to locate multiple optima.

Once a population is established using SCGA, species seeds are determined. This is done by evaluating each individual of the population from the fittest to the least fit. If no other species seed exists within a predefined distance, then the individual is added to the list of species seeds. Because it begins with the fittest individuals, it ensures that the seeds of the species are the most optimal members within the radius.

Table 4 shows the SCGA seed selection algorithm. In this algorithm  $X_s$  is the set of species seeds and  $\sigma_s$  is the distance that defines a species.

Table 4. Species Conserving Genetic Algorithm Seed Selection

Line Number	Pseudocode
1	Initialize algorithm by setting $X_s = \emptyset$
2	While (there are individuals in the population that have not been evaluated)
3	Find the best unevaluated individual, $x^*$
4	Set found = false
5	For every individual $x$ in $X_s$ do
6	Mark $x$ has having been evaluated
7	If distance $(x^*, x) \leq \sigma_s/2$ then
8	Set found = true
9	Break for loop
10	End If
11	End for loop
12	If found = false then
13	Add $x^*$ to $X_s$
14	End If
15	End while loop

After the next generation is created, SCGA conserves species. Each seed is compared to individuals in the next generation within the radius of the seed. If the seed is more fit than the weakest individual in this area, the seed replaces the individual. If there are no individuals in the species of the next generation, the seed replaces the least fit individual

of the next generation. By conserving these species, pressure is created that prevents global convergence and allows multiple optima to be generated. When the NGA finishes, the fittest species seeds are the optima.

Table 5 shows the SCGA species conserving algorithm. Like the algorithm in Table 3,  $X_s$  is the set of species seeds and  $\sigma_s$  is the distance that defines a species.

SCGA has one drawback. It requires a parameter that defines the distance from its seed that a species covers. The ideal radius value depends on the problem instance. That is often unknown before the NGA runs. Li et al. (2002) addressed this drawback in their research. It is their belief that it is better to have the parameter too large than too small. They recommend that the user informally compare the domain to one that is known. The input parameter should be set to double the distance between the optima of the known domain (Li et al., 2002).

Table 5. Species Conserving Genetic Algorithm Species Conservation

Line Number	Pseudocode
1	Mark all individuals as not being evaluated
2	For all $x$ in $X_s$ do
3	Select the least fit individual $y$ in the area of the domain that is $\sigma_s/2$ from $x$
4	If there is a $y$ that meets this condition then
5	If ( $f(y) < f(x$ ) then
6	Replace $y$ with $x$
7	End if
8	Else
9	Select the least fit individual $y$ in the new generation
10	Replace $y$ with $x$
11	Mark $x$ has having been evaluated
12	End for loop

### *Genetic Algorithm with Species and Sexual Selection*

Raghuwanshi and Kakde (2007) developed a method called Genetic Algorithm with Species and Sexual Selection (GAS3) that is a sexual GA, which means that it introduces the notion of gender into the NGA. GAS3 also uses the concept of species, which corresponds to niches. Species are formed around strong members. GAS3 also uses population overlapping, meaning that some members of a generation stay in the population pool with the next generation. The GAS3 algorithm has four steps that it performs on each species.

1. The first one is the *selection plan*, which determines the female member of the species. This is the member of the species with the highest fitness. All of the other members are males.
2. Then it performs the *generation plan*. The generation plan creates a set of offspring by randomly selecting males to reproduce with the female.
3. In the *replacement plan* the original group and the offspring group are merged together.
4. This new group goes through an *update plan*, which determines the female and males and removes the least fit members so the size of the population is constant.

Occasionally, GAS3 will reevaluate the species. If some species are not performing well, they will be merged with other species.

GAS3 has many interesting characteristics. Gender plays an important role in the algorithm. Having only one female individual, does not seem to model most biological species. Overlapping generations assists in the algorithm preventing premature convergence. GAS3 (Raghuwanshi & Kakde, 2007) was tested against a large set of 13

benchmark functions. Published results show that the algorithm performs very well against many commonly used multimodal functional optimization problems.

### *Crowding Clustering Genetic Algorithm*

Crowding Clustering Genetic Algorithm (CCGA) is a NGA developed to solve functional optimization problems for both local minimums and maximums (Ling et al., 2008). Similar to other methods, CCGA accomplishes this by promoting some members of one generation into the next generation to prevent genetic drift.

CCGA begins each iteration with typical selection, crossover and mutation operations. Each child is grouped with a parent who it is closest to, using some distance measurement. This leaves each parent associated with zero or more individuals in the child generation. Each of these sets is a *cluster*. The *cluster center* is the fittest individual in the cluster. This may be the individual with the smallest or largest objective value depending on if this is a minimization or maximization problem. The objective value of the fittest member is the *center value*. The largest distance between the cluster center and the other individuals in the cluster is the *cluster radius*. Clusters are sorted descending by the fitness of the cluster center. This ensures that members at the front of the list are the fittest. The sorted list of clusters is evaluated. Each one may be moved into a second list called *reserved clusters*. When a cluster is added to the reserved cluster list, the cluster radius is referred to as the *reserved cluster radius*. Either of two conditions can move a cluster into the reserved cluster list. The cluster is added to the reserved cluster list if its cluster center is outside all of the existing reserved cluster radii. The second condition for moving a cluster to the reserved cluster list is if it satisfies the

peak detection requirement, which is described later. When a cluster is added to the list of reserved clusters, its *reserved cluster radius* is set. This reserved cluster radius can be set to the minimum of the radius of the cluster being added or minimum distance from the cluster center to another reserved cluster center. A new generation is created by taking the cluster centers of the reserved cluster and generating enough uniformly distributed individuals to keep the population size stable. Each iteration of the algorithm creates a new reserved cluster list. These steps repeat until some predefined number of generations is reached.

Table 6. Crowding Clustering Genetic Algorithm

Line Number	Pseudocode
1	Create initial population uniformly distributed across solution space
2	Use traditional GA selection, crossover and mutation to create a new generation
3	For each parent, $P_j$ , construct $\{P_j, CS_j\}$
4	For each cluster $\{P_j, CS_j\}$
5	Set $CC_j$ to the fittest individual within each $\{P_j, CS_j\}$
6	Set $CR_j$ to the largest distance between individuals in $\{P_j, CS_j\}$ and $CR_j$
7	End for loop
8	Sort clusters descending according to their objective value
9	Set $RC = \emptyset$
10	For each cluster $\{P_j, CS_j\}$
11	If $(D(CC_j, RCC_i) > RCR_i$ for all $RCC_i$ in $RC$ ) or $(Peak(CC_j, RCC_i) = 1)$ then
12	Put $CC_j$ into $RC$
13	Set the $RC$ for $CC_j$ to $\min(CR_j, D(CC_j, RCC_i))$
14	End if
15	End for loop
16	Generate (population size – $RC$ size) of uniformly distributed individuals for the next generation
17	Repeat steps 2 through 16, until the termination condition is met

Table 6 shows the CCGA algorithm for a minimization functional optimization problem. In this algorithm the parent generation is  $P_j$ , where  $j = 1, 2, \dots$ , population size.

$CS_j$  is the set of children closest to the  $j$ th parent.  $CC_j$  is the cluster center and  $CR_j$  is the radius cluster associated with the  $j$ th parent.  $RC$  is the set of reserved clusters.  $RC_i$  is the  $i$ th reserved cluster radius.

Peak detection in CCGA is another way that a cluster can be added to the list of reserved clusters. It attempts to determine if individuals are tracking different peaks. A cluster satisfies the peak detection condition if the function  $Peak(CC, RCC_j)$  returns 1 for all  $j = 1, 2, \dots$  reserve cluster size. In these equations the cluster center is  $CC$ , the  $j$ th reserved cluster center is  $RCC_j$  and  $f$  is the objective function. The equation for the peak detection is defined by the following equation for minimization problems.

$$Peak(CC, RCC_j) = \begin{cases} 1 & \text{if } f\left(\frac{CC + RCC_j}{2}\right) > \frac{f(CC) + f(RCC_j)}{2} \\ 0 & \text{, otherwise} \end{cases}$$

For maximization problems peak detection is defined by the following equation.

$$Peak(CC, RCC_j) = \begin{cases} 1 & \text{if } f\left(\frac{CC + RCC_j}{2}\right) < \frac{f(CC) + f(RCC_j)}{2} \\ 0 & \text{, otherwise} \end{cases}$$

The CCGA algorithm is used to determine functional optimization. More specifically, it is used to search for functional minimums. Experiments performed by Ling et al. (2008) show that CCGA out perform other crowding methods.

### *Crowding Summary*

Crowding methods take very direct approaches to maintain genetic diversity. After locating individuals of interest, crowding methods put these individuals into the next generation. There is no chance that these individuals will not be represented in future generations. Fitness sharing methods take a very subtle approach to exploration. They only increase the chance the interesting individuals will be used for crossover. In contrast crowding methods take direct approaches to encouraging exploration.

This section highlights many crowding methods. Early crowding methods simply select interesting individuals and put them in the next generation. More recent crowding methods have complex algorithms to determine what individuals deserve to be preserved. Crowding methods closely resemble biological systems by combining parents and children in the same generation.

### **Other Niche Genetic Algorithm Methods**

While most NGAs fall into one of the two categories of fitness sharing or crowding schemes (Deb & Goldberg, 1989), there are some NGAs that do not exhibit either characteristic. Some of these NGAs are hybrid methods or are traditional GAs that were created for special purposes that happen to solve multimodal optimization. These NGAs provide unique looks at NGA research and introduce different approaches to multimodal optimization of continuous functions.

### *Fitness-based Neighbor Selection*

Ando and Kobayashi's (2005) method of Fitness-based Neighbor Selection (FNS) addresses many known limitations of NGAs. Many NGAs have parameters that need to be set in order to solve multimodal optimization problems. Often this is a radius that needs to be set to a value less than the distance between two species. This parameter is difficult to set prior to knowing where the optima are.

Ando and Kobayashi (2005) observed that integrals can be used to determine which peaks individuals are tracking. This observation is incorporated into the FNS algorithm. When trying to decide if an individual  $A$  is a neighbor of  $B1$  or  $B2$ , comparing the integral between  $A$  and  $B1$  to the integral between  $A$  and  $B2$  can be helpful. The one, assume  $B1$ , with the largest integral has a greater probability of being neighbors for maximization problems. This observation is true based on integrals measuring area beneath the objective function. The greater the area indicates that the objective function peaks and that  $A$  and  $B1$  are in the same neighborhood.

However, calculating integrals in higher dimensional space can be as challenging as locating local optima. FNS estimates these calculations using the Wilcoxon Rank-sum Test (Wilcoxon, 1945). FNS creates two sets of offspring. The first set of offspring is between  $A$  and  $B1$ . The second is between  $A$  and  $B2$ . The fitness of these two sets of offspring is used to define neighborhoods using the Wilcoxon Rank-sum Test.

### *Enhanced Evolutionary Tabu Search*

The Enhanced Evolutionary Tabu Search (EE-TS) is a metaheuristic technique that combines a Tabu Search with a GA (McLoughlin & Cedeno, 2005). This hybrid

technique which combines a Tabu Search and GA, is used in the research to solve the Quadratic Assignment Problem (QAP). The QAP is a problem that attempts to minimize cost when placing facilities into locations. Facilities accrue a cost based on how far they are from other facilities, but the costs may not be uniform. The QAP problem is somewhat different than traditional functional optimization problems. Most algorithms, including EE-TS, place facilities one at a time. As the algorithm runs more facilities are placed into different locations. At the beginning of the algorithm an individual represents one facility to location mapping. Then, as the algorithm runs, an individual represents more facility to location mappings. Finally the individual represents all facilities mapped to locations.

Tabu Searches are designed to prevent revisiting the same solution repeatedly. Repetition can occur when a series of optimal moves revisits a previous solution state (Glover, 1990b). If this happens the algorithm could enter an infinite loop or fail to explore promising regions of the domain. The Tabu Search uses a memory structure to record previous solution states and prevents them from being revisited (Glover, 1990b). EE-TS also evaluates for repetition to encourage exploration of the domain.

EE-TS begins with an initialization phase like other GAs (McLoughlin & Cedeno, 2005). As the algorithm runs it keeps track of a current candidate. As long as repetition is not occurring, the algorithm evaluates the neighborhood and selects a move that will increase the fitness the most. A move consists of swapping two facilities. After the move is identified, tournament selection picks an individual. Crossover is performed with this individual and the current candidate. If the child is fitter than the current candidate with the identified move applied to it, the child becomes the current candidate.

Otherwise, the current candidate with the identified move applied to it, is the new current candidate. This loop continues until the termination condition is met. After each loop the algorithm uses the tabu list to determine if repetition is occurring. If it is occurring, the algorithm identifies an individual through tournament selection. The new current candidate is the winner of the tournament and the old candidate. Finally, the tabu list is cleared. The loop repeats.

Table 7 shows the EE-TS algorithm (McLoughlin & Cedeno, 2005). In this algorithm  $i^*$  is the current candidate and  $i$  is a possible new current candidate. The variable *escape* is used to indicate if repetition is occurring. The variables *champion* and *move* are temporary variables to hold the winner of the tournament selection and a possible move.

Table 7. Enhanced Evolutionary Tabu Search Algorithm

Line Number	Pseudocode
1	Generate initial population, $P$
2	Set $i$ and $i^*$ to the fittest individual in $P$
3	Set <i>escape</i> to true if detection of repetition is discovered, otherwise set to false
4	If <i>escape</i> = false then
5	Set <i>move</i> to the best move
6	Set <i>champion</i> to winner of tournament selection
7	Set <i>child</i> to crossover of $i$ and <i>champion</i>
8	If $\text{fitness}(\text{child}) < \text{fitness}(i)$ with move <i>move</i> applied to it then
9	Set $i$ to <i>child</i>
10	Else
11	Set $i$ to $i$ with move <i>move</i> applied to it
12	Else
13	Set <i>champion</i> to winner of tournament selection
14	Set $i$ to crossover of $i$ and <i>champion</i>
15	Reset the tabu list and solution history
16	If $\text{fitness}(i) < \text{fitness}(i^*)$ then
17	Set $i^* = i$
18	Repeat steps 3 through 17 until termination condition

The EE-TS algorithm performs equally well as other QAP algorithms. Its value is that it locates the optima in fewer steps or iterations (McLoughlin & Cedeno, 2005). Because it is designed for the QAP problem, it is not suited for multimodal functional optimization. EE-TS represents a new class of hybrid GAs that incorporates other search techniques into them.

### *Hybrid Genetic Algorithm and Particle Swarm Optimization*

Recently hybrid algorithms have increased in popularity. Kao and Zahara (2008) created an algorithm that combines GAs and Particle Swarm Optimization (PSO). Hybrid approaches to multimodal optimization have shown promise by combining the best aspects of different types of algorithms.

PSO is another type of search algorithm. Unlike GAs that eliminate individuals after each generation, the individual in PSOs remain throughout the algorithm. Individuals move throughout the domain space to locate optima. Each individual tracks where in the domain space they have been and has the ability to communicate these locations to other individuals in the swarm. Individuals also have the ability to adjust their position in the domain based upon communication from other individuals in the domain. As a group, the swarm converges to the optima.

Kao and Zahara's (2008) algorithm uses both a GA and a PSO. It begins by randomly generating a population. Half of the population that has the greatest fitness is used in a standard GA. After the next generation is created, it is used to communicate with the second half of the population through PSO techniques. Ideally the offspring of the GA will have higher fitness than the second half. As a result the second half will adjust their

positions in the domain based upon their previous knowledge and the communication from the offspring of the first half. After they have adjusted their positions in the domain, the two halves are combined and reevaluated. This process continues until a termination condition is met.

Kao and Zahara's (2008) GA and PSO algorithm is shown in Table 8. Crossover is done by generating a uniform random number  $N$ . Then  $N$  proportion of the alleles are taken from one parent and  $1-N$  proportion from the other. The function  $Uniform(0, 1)$  is the function that generates the uniformly distributed random number between 0 and 1. In this algorithm the parameter  $P$  is the population size and the  $x$ 's are individuals.

Table 8. Genetic Algorithm and Particle Swarm Optimization Algorithm

Line Number	Pseudocode
1	Generate initial population of size $P$
2	While termination condition is not met do
3	Sort individuals by their fitness
4	Perform the following steps on the fittest $P/2$ individual
5	For all $j = 1$ to $P/2 - 1$ do
6	Create individual $x$ using $Uniform(0, 1)$ proportion of $x_j$ alleles and $(1 - Uniform(0, 1))$ proportion of $x_{j+1}$ alleles
7	Add $x$ to next generation
8	End for loop
9	Create individual $x$ using $Uniform(0, 1)$ proportion of $x_{P/2}$ alleles and $(1 - Uniform(0, 1))$ proportion of $x_1$ alleles
10	Add $x$ to next generation
11	End perform block
12	Apply 20% mutation on next generation
13	Adjust the $P/2$ least fit individual by PSO
14	Add these individual into the next generation
15	End while loop

This hybrid approach is novel and leverages the strengths of both GA and PSO methods. GAs are very effective at taking a set of fit individuals and creating a generation of more fit individuals. PSOs are effective at adjusting weak members of the

population to increase their fitness. The two algorithms work fluidly together to locate optima.

### *Cellular Genetic Algorithms*

Cellular Genetic Algorithms (cGA) were originally developed in the early 1990's to run GAs using parallel machines (Whitley, 1993). To take advantage of parallel processors the domain space was divided into squares. Individuals were only allowed to mate with individuals within its square or neighboring squares. By creating this grid across the domain space, crossover for each generation could be performed in parallel. This allowed GAs to converge much faster than traditional methods, which made them more practical for solving real world problems. This approach is based upon cellular automata (Whitley, 1993).

Table 9. Cellular Genetic Algorithm

<b>Line Number</b>	<b>Pseudocode</b>
1	While not termination condition
2	For $x = 1$ to $w$
3	For $y = 1$ to $h$
4	Get list of neighbors for individual $(x, y)$
5	Select parents $p1$ and $p2$ from list of neighbors
6	Create individual $i$ from $p1$ and $p2$
7	Mutate( $i$ )
8	If $\text{fitness}(i) > \text{fitness}(\text{individual}(x, y))$
9	Replace individual( $x, y$ ) with $i$
10	End for loop
11	End for loop
12	End while loop

The basic cGA algorithm is shown in Table 9 (Alba, Alfonso & Dorronsoro, 2005). This algorithm assumes that the domain space has been divided into a grid of width,  $w$ , and height,  $h$ . For each individual in the grid, the algorithm determines a list of

neighbors. In cGAs an individual is considered to be its own neighbor. A selection step identifies two individuals from this list,  $p1$  and  $p2$ . A new individual,  $i$ , is created using crossover. A mutation function will determine if mutation is needed based upon a mutation rate. If it is determined that mutation should occur, the function will perform the mutation. If the new individual  $i$  is more fit than the original individual, it will replace it. The algorithm evaluates every individual in the population. It will continue this process until a termination condition is met.

Because individuals are restricted to mating only with individuals close to them, cGAs prevents premature convergence and can be used for multimodal optimization problems (Nebro, Durillo, Luna, Dorronsoro, & Alba, 2006). This form of selection prevents individuals in one area of the domain from dominating other niches. Because of recent advancements in computational power, the parallel aspects of cGAs have been eclipsed by their ability to solve multimodal optimization problems. A variety of enhancements have been made to cGAs and multimodal optimization.

*Anisotropic Selection* in cGAs assigns probabilities of replacement to the squares around an offspring (Simoncini, Verel, Collard & Clergue, 2006). Individuals within a square perform a typical GA with selection, crossover and mutation. The offspring then replaces some of the old generation's individuals. Different probabilities are assigned to different geometric directions used in selection. There is a probability that selection will be made using a north or south square. There is a probability that selection will be made using an east or west square. The final probability is that the center square will be used for the selection. These probabilities guide the direction of the search in the local area of the domain. A control parameter,  $\alpha$ , is used to influence these three probabilities. The

following probabilities are used in determining the direction for selection (Simoncini et al., 2006).

Probability of center cell	$p_c = 0.2$
North or south cell	$\frac{(1-p_c)}{2}(1+\alpha)$
East or west cell	$\frac{(1-p_c)}{2}(1-\alpha)$

Once the direction is determined, tournament selection is used to select the individual for crossover. If the individual in the new generation is better than the individual selected for replacement, it will be replaced.

There are two cGAs that attempt to solve multiobjective optimization problems. These algorithms are Cellular Multiobjective Genetic Algorithm (cMOGA) and Multiobjective Optimization Cellular Genetic Algorithm (MOCeLL) (Alba & Dorronsoro, 2008). Both algorithms are very similar and use the same general approach.

MOCeLL is another type of cGA (Nebro et al., 2006). MOCeLL uses a *Pareto front*, which is an alternate population that contains optimal non-dominant individuals. The Pareto front has a maximum size and maintains genetic diversity. In MOCeLL selection, crossover and mutation take place according to normal cGA principles. The offspring are added into the next generation. Offspring may also be added to the Pareto front. When this Pareto front hits its maximum size, individuals are replaced using a crowding method, which increases genetic diversity (Nebro, Durillo, Luna, Dorronsoro & Alba, 2009). The final step in MOCeLL is to randomly replace members of the population with individuals from the Pareto front. This feedback ensures that dominant areas of the domain do not eclipse other optima. cMOGA is the same as MOCeLL, except it does not contain the feedback step (Alba et al., 2005).

cGAs research has shown that they are very effective in solving multimodal optimization problems. Although research in other NGA areas has been around twice as long as cGAs, results of cGA research are very impressive. Multiple methods have been developed based upon cGA principles.

#### *Novel Sexual Adaptive Genetic Algorithm*

GAS3 is not the only NGA that incorporates the concept of gender. Novel Sexual Adaptive Genetic Algorithm (NSAGA) has genders also (Zhang, Zhao & Wang, 2009). But it more evenly divides the number of males and females. Similar to biological organisms, individuals in NSAGA have gender based upon genetic characteristics. This gender selection more closely resembles genders in biological species.

NSAGA leverages an early evolutionary theory called the Baldwin effect (Baldwin, 1896). This theory proposes that an individual's fitness is not always limited to their biological characteristics. It is possible that through environmental influences an individual can increase its fitness. In an NGA however, environmental influences are not defined. NSAGA uses other individual's fitness to be this environmental influence. The Baldwin effect provides a new approach to NGAs.

NSAGA computes fitness as the weighted sum of three types of fitness: innate fitness, evaluation fitness and acquired fitness (Zhang et al., 2009). The first part of the fitness is the innate fitness,  $IF$ . This fitness excludes environmental influences. In NSAGA innate fitness is defined as the following, where  $f_{\min}^t$  is the minimum fitness and  $f_{\max}^t$  is the maximum fitness for generation  $t$ .

$$IF(x) = \frac{f(x) - f_{\min}^t}{f_{\max}^t - f_{\min}^t}$$

The second type of fitness is the evaluation fitness,  $EF$ . This fitness includes influences from the individual's parents. In the fitness function below  $w_1$  and  $w_2$  are parameters that weight each parent's influence and  $x_f$  and  $x_m$  are the individual's mother and father, respectively.  $IF$  is the innate fitness defined previously.

$$EF(x) = w_1 \frac{IF(x) - IF(x_f)}{IF(x_f)} + w_2 \frac{IF(x) - IF(x_m)}{IF(x_m)}$$

The final type of fitness is the acquired fitness,  $AF$ . This fitness derives from the Baldwin effect. Individuals within a niche may increase or decrease their fitness by some factor, between 0 and 1, of the average innate fitness of the members of the niche. This is shown in the following equation where  $c$  and  $P_b$  are parameters and  $rnd$  is a random number.

$$f_y^1(x) = \begin{cases} f_y(x) + c \cdot f_y^{ave}, & \text{if } rnd \leq P_b \\ f_y(x) - c \cdot f_y^{ave}, & \text{otherwise} \end{cases}$$

The equation that reflects the Baldwin effect is used in the third fitness function. The acquired fitness is given by the function below (Zhang et al., 2009).

$$AF(x) = \frac{f_y^1(x) - f_{y \min}^t}{f_{y \max}^t - f_{y \min}^t}$$

With the acquired fitness function individuals may have their fitness increase or decrease based upon the factors previously outlined. The final fitness of an individual is a weighted sum of the innate, evaluation and acquired fitness functions.

$$Fitness(x) = \beta_1 IF(x) + \beta_2 EF(x) + \beta_3 AF(x)$$

$\beta_1$ ,  $\beta_2$  and  $\beta_3$  are weights placed on each type of fitness.

Gender determination is an important part of NSAGA. It has great consequences because individuals may only mate with individuals of the opposite gender. Selecting an existing gene would divide genders to different areas of the domain. This is not desired. Gender is determined randomly thus giving all areas of the domain the ability to have both genders (Zhang et al., 2009). The parameter  $P_g$  is the probability that an individual is a male.  $P_g$  gives the algorithm more flexibility in controlling the proportion of males and females. Gender determination helps NSAGA preserve interesting areas of the domain for exploration.

The final important aspect of NSAGA is the selection process. The selection process goes beyond just limiting heterosexual selection. A parameter  $P_{elitism}$  is used to determine the fittest members of the population. NSAGA uses a different selection method for elite and non-elite individuals. For the elite individuals selection is done according to rank within the group. Of the remaining individuals selection is done through tournament selection.

NSAGA is a very unique NGA that resembles natural selection more than many other NGAs. Incorporating gender and the Baldwin effect make NSAGA a novel algorithm. Rather than extending existing fitness sharing or crowding methods, NSAGA takes a more accurate approach to modeling biological evolution.

#### *Other Niche Genetic Algorithm Summary*

This section briefly describes some NGAs that cannot be categorized as fitness sharing or crowding methods. Some NGAs have characteristics so different than standard

fitness sharing and crowding methods, that this third categorization has been defined. This other type of category introduces new concepts to NGAs and the multimodal optimization problem.

The algorithms described in this section introduce many new approaches to NGAs. Some of them are hybrid methods that combine the concept of GAs with other search techniques, like PSO. Others model biological theories, like the Baldwin effect. cGAs solve multimodal optimization problems through limiting selection to neighbors. These methods take a different approach to NGA research.

## **Summary**

This chapter provides an extensive literature review of NGAs. The first section describes relevant literature used in this research that is not an NGA. Three other sections describe different types of NGAs. A variety of methods from classical NGAs to modern methods have been presented. NGAs are normally organized into two groups. Sharing methods adjust fitness to encourage exploration. Crowding methods replace individuals with individuals of the previous generation based upon distance measurements. There are some NGAs that don't directly fall into either category. These algorithms are presented in a separate section.

This literature review provided the foundation for this research. Many of the methods here are compared to the new framework that is presented. Some of the concepts used in previous research are used in the creation of the new framework. These cases will be described in Chapter 3.

## **Chapter 3**

### **Methodology**

Chapter 3 describes the methodology for this research. The research method will be described, followed by the new framework that will be tested. The benchmark equations and performance criteria are defined. The next two sections describe the formatting for presenting of the results and the resources required to perform the research. The chapter concludes with a summary of the framework and methodology.

#### **Research Method Employed**

An experimental research methodology was used to conduct this research. A new framework was developed and compared against existing algorithms using benchmark equations and performance criteria. The experimental research methodology provided a basis to evaluate the performance of the new framework.

While the framework is new, many ideas and concepts are based upon existing algorithms. Chapter 2 describes existing NGA research. Conclusions derived from this body of knowledge were reflected in the new framework. The experiment was used to test the hypothesis used to develop the new framework.

## **Specific Procedures Employed**

A new NGA was needed to solve the types of equations described earlier and shown in Figure 1. This new NGA leveraged other NGAs and search methods. This new algorithm most closely resembles SCGA (Li et al., 2002). At a very high level the algorithm did not attempt to find all optima in a single pass. It defined some search areas to investigate. Once defined, it let the NGA run to find optima within these areas. These optima were placed on a tabu list, which prevented them from being revisited. The algorithm found optima in parallel each time it defined a set of searchable domain spaces.

### *DSGA Algorithm*

A traditional GA performs three general steps to create each generation. Selection, crossover and mutation allow the GA to converge to an optimum. The SCGA augments the traditional GA to include seed selection and seed conservation (Li et al., 2002). The new algorithm, Dynamic-radius Species-conserving Genetic Algorithm (DSGA), also uses this seed selection and seed conservation approach but differs from SCGA in three important ways. First, DSGA incorporates a tabu list to track optima and encourages exploration in other areas. Second, DSGA varies the value of the radius. Two strategies will be presented for varying the radius. Third, DSGA has two different strategies for seed selection. These two variations will encourage exploration. The variations of this framework will be presented later in this section.

Groups of individuals are formed around a fit member, called a seed. These groups cover a search area in the domain. A predefined radius is used around seeds to define which members are grouped with the seed. Every new generation redefines the seeds and

areas. By conserving these search areas, DSGA locates multiple optima and prevents a single dominant optimum from eclipsing the other ones.

Table 10. DSGA Parameters

<b>Abbreviation</b>	<b>Name</b>	<b>Description</b>
<i>N</i>	Population size	Number of individuals in each generation
<i>M</i>	Mutation Rate	Odds of gene mutating
<i>IS</i>	Initial Sigma	Initial value for sigma
<i>SD</i>	Sigma Delta	The amount that sigma will be changed each iteration
<i>RLC</i>	Reevaluate Loop Count	The number of times that the NGA will loop before it reevaluates the seed radius
<i>CL</i>	Convergence Limit	The number of individuals needed to determine that convergence has taken place

Table 10 shows the parameters of the algorithm. Like traditional GAs, there is a population size,  $N$ , and a mutation rate,  $M$ . Similar to SCGA (Li et al., 2002), DSGA uses a radius parameter. The radius is the minimum distance a strong individual must be away from all other clusters in order to create a new cluster. In DSGA sigma is the radius of the clusters. It is called the initial sigma, because the sigma varies as the algorithm runs. This radius is changed by the sigma delta,  $SD$ , to search for new optima. The Reevaluate Loop Count,  $RLC$ , is the number of times that the NGA loops before allowing the search areas and radius to be redefined. When determining if an optimum has been located the algorithm looks for identical individuals. If there are  $CL$  or more identical individuals, the algorithm concludes that an optimum has been located.

Table 11 shows the new algorithm, DSGA. Descriptions of the seed selection, seed conservation and radius altering approaches are described later in this section. Unlike other GAs, DSGA has two loops. The inner loop, lines 4 through 10, perform a typical GA with the enhancement of seed selection and seed conservation. Once this loop is finished the algorithm records any optima as an ordered pair of the optima and the radius

in the tabu list, line 11 and 12. Optima are any domain values which have  $CL$  or more identical individuals, line 11. Seeds that are not optimums are also added to the tabu list, but optima are marked as such. In line 18, the radius, associated with the seeds, is altered and the GA is run again. This will occur in the outer loop, lines 2 through 19, and ends when a termination condition is met. The algorithm varies the size of the radius to locate other optima.

Table 11. DSGA Algorithm

Line Number	Pseudocode
1	Initialization
2	While not termination condition
3	For (int $r = 1$ ; $r \leq RLC$ ; $r++$ )
4	Begin
5	Seed Selection
6	Selection
7	Crossover
8	Mutation
9	Seed Conservation
10	End For Loop
11	If there exists an individual $d$ with $CL$ or more identical individuals then
12	Add $(d, \sigma)$ pair to $best\_Xs$
13	Mark pair $(d, \sigma)$ as optimum
14	Replace $d$ and all of the identical individuals to $d$ with randomly generated individuals
15	End if
16	Add $(s, \sigma)$ pair to $best\_Xs$ for all $s$ that are seeds
17	Replace all individuals $s$ with randomly generated individuals
18	Alter radius $\sigma$
19	End

This algorithm differs from SCGA in many important ways. SCGA selects a single radius size and performs the algorithm in a single loop (Li et al., 2002). DSGA varies the radius to locate other optima. This new algorithm has also been augmented with a tabu list to prevent already located optima from being used as seeds in the future. These

changes should allow the algorithm to locate other optima that may have been missed in earlier iterations.

DSGA performs an initialization phase to create the first generation and initialize the tabu list. Table 12 shows the initialization steps. The variable *best\_Xs* is the tabu list. Originally, the seed radius,  $\sigma$ , is set to *IS*. This value will change by *SD* as the algorithm runs.

Table 12. DSGA Algorithm Initialization

Line Number	Pseudocode
1	Set list <i>best_Xs</i> to $\emptyset$
2	Set list <i>generation</i> to $\emptyset$
3	
4	For (int <i>i</i> = 1; <i>i</i> <= <i>N</i> ; <i>i</i> ++)
5	Begin
6	Create new string <i>y</i>
7	Randomly generate genes for <i>y</i>
8	Add <i>y</i> to <i>generation</i>
9	End For Loop
10	Set $\sigma$ to <i>IS</i>

#### *DSGA Seed Selection*

Seed selection is a critical part of locating optima. Each generation defines its own seeds. This algorithm for seed selection is shown in Table 13. In line 1, it begins with an empty list of seeds. It is beneficial to make the fittest individual in each niche a seed, but the algorithm also needs to explore other areas of the domain. So the algorithm evaluates individuals in an order defined by a *seed evaluation ordering* (*seo*) function. Possible implementations for *seo* will be presented later in this section. Each individual is evaluated as a candidate for seed selection based on this function. The algorithm determines if an individual is within  $\sigma$  distance to a currently established seed. If a seed exists within  $\sigma$  distance, the individual is not a seed, but a member of the seed's species.

If no established seeds exist within the  $\sigma$  distance of this individual, it will become a seed. The only exception to this rule, which is shown in lines 16 and 17, is if the individual is on the tabu list. Here it has already been determined that it has been investigated and conservation of it is not needed. It is prevented from being a seed.

Table 13. Seed Selection

Line Number	Pseudocode
1	Set list <i>seeds</i> to $\emptyset$
2	Sort <i>generation</i> descending by <i>seo</i> function
3	For (int $k = 0$ ; $k < \text{size of } generation$ ; $k++$ )
4	Begin
5	Set $K$ to the $k$ -th individual in <i>generation</i>
6	Set boolean <i>found</i> = true
7	For (int $m = 0$ ; $m < \text{size of } seeds$ ; $m++$ )
8	Begin
9	Set $M$ to the $m$ -th individual in <i>seeds</i>
10	If $\text{distance}(K, M) < \sigma$ then
11	Set <i>found</i> to true
12	Break
13	Else
14	Set <i>found</i> to false
15	End For Loop
16	If <i>found</i> = false and $K$ is not in best_Xs then
17	Add $K$ to <i>seeds</i>
18	End For Loop

This seed selection algorithm is identical to SCGA except in two areas. In line 16 DSGA prevents individuals on the tabu list from becoming seeds in the future. This is done because these areas of the domain have already been investigated. Using the tabu list encourages the algorithm to explore other areas of the domain. In line 2 individuals are sorted by a function called *seo*. DSGA uses two strategies to investigate unexplored areas of the domain. One strategy uses a standard fitness sharing approach; the other excludes individuals from becoming seeds if they are too close to individuals on the tabu

list. Each strategy is implemented by using different *seo* functions. In SCGA individuals are only sorted by fitness. These differences will allow the algorithm to find other optima.

### *DSGA Seed Conservation*

Once seeds are established there needs to be a means to preserve them into the next generation. This is called seed conservation and is shown in Table 14. After each generation is created, it goes through a seed conservation step. Each seed from the previous generation replaces a weak individual in the new generation. First the algorithm looks at individuals in the new generation, which are within  $\sigma$  distance of the seed. If there are individuals in the new generation, which meet this condition the seed replaces the weakest individual of this list. If there are no individuals within the seed's radius, then the seed replaces the weakest individual in the new generation. Every seed is promoted into the next generation, but this does not mean that this seed will be a seed in the next generation. It will have to be evaluated as any other individual.

Table 14. Seed Conservation for each Generation

Line Number	Pseudocode
1	For (int $p = 1$ ; $p \leq \text{size of seeds}$ ; $p++$ )
2	Begin
3	Set $P$ to the $p$ -th individual in <i>seeds</i>
4	Find $y$ such that it is the least fit individual with distance( $y, P$ ) $\leq \sigma$
5	If $y$ exists and $y$ is less fit than $P$
6	Replace $y$ with $P$
7	Else
8	Replace the least fit individual in new generation with $P$
9	End For Loop

DSGA uses many of the steps of SCGA to conserve seeds after each generation is created. The only exception is that Euclidean distance is used in this algorithm. The seed conservation phase of recording seeds does not exist in SCGA. This should allow the algorithm to investigate other areas of the domain to locate other optima.

Besides the algorithm parameters shown in Table 10, two components of the algorithm may vary. There are two strategies to alter the radius after the inner loop of the algorithm and two strategies to encourage further exploration of the domain. The different combinations of strategies allow the DSGA framework to create multiple algorithms. The two components of the DSGA framework allow it to be used in many domains.

#### *Varying Radius Strategies*

We will consider two strategies to vary the radius used in DSGA. This step is shown in Table 11, line 18. In this step of the algorithm, the radius will be changed by a constant value of sigma delta,  $SD$ . The two strategies differ in the way that the radius is changed by  $SD$ . One strategy consistently increases or decreases the radius. A second strategy may increase or decrease the radius as the algorithm runs.

In the first strategy the radius is increased or decreased by  $SD$ . It will either always increase or always decrease the radius. The method could start the radius very small and increase it incrementally after the inner loop completes. In this strategy the radius would start at  $IS$  and be increased by  $SD$  in every pass of the outer loop. Eventually the radius would increase to such a size that only one seed would be formed. This condition would be the termination condition shown in Table 11, line 2. Or the method could start with a

very large  $IS$  and decrease it by  $SD$  as the algorithm runs. Here the termination condition is not so obvious. A natural termination condition would be to terminate the first time the inner loop completes but does not find any additional optima.

The second strategy is to increase or decrease the radius by  $SD$  at the end of the inner loop. At the end of the inner loop the algorithm will decide if the radius should be increased or decreased by  $SD$ . There are many possible methods that could be used to determine how to vary the radius. One possibility for this approach is to base the radius change on the number of optima located. If no optima are located in the inner loop, then decreasing the radius will allow more seeds to form and should increase the chance of optima location. If optima are located, the approach would increase the radius. As in the previous approach the termination condition would be when the inner loop does not locate any additional optima. These two strategies allow DSGA flexibility in locating optima.

### *Exploration Approaches*

The second component of the DSGA framework encourages exploration in areas of the domain where optima have not been found. Two strategies will be presented. Both strategies fulfill this through the seed selection in the algorithm. Each approach accomplishes exploration by defining different *seo* functions. One *seo* function implements a fitness sharing algorithm. The other one excludes individuals from becoming seeds that are too close to individuals on the tabu list.

The first strategy eliminates individuals that are too close to existing seeds. It uses the ordered pair  $(o, r)$  on the tabu list. The optimum is  $o$  and the radius when the optimum

was located is  $r$ . In future iterations of the algorithm the radius will change. The variable  $r'$  is the current value of the radius when the strategy is executed. This approach excludes individuals from becoming seeds if their distance is within  $\min(r, r')$  of an element on the tabu list. In the equation below  $i$  is the individual being evaluated for seed selection,  $tl$  is the current tabu list,  $r'$  is the current radius and  $d$  is Euclidean distance. For all ordered pair  $(o, r)$  on the tabu list, the function will return the individual's fitness if the distance between  $o$  and  $i$  is greater than the  $\min(r, r')$ . If the distance is less than  $\min(r, r')$ , the function returns 0. This ensures that this individual will not be a seed.

This equation for  $seo$  is given below:

$$seo_{tl, r'}(i) = \begin{cases} fitness(i), & \text{if and only if there does not exist an } (o_l, r_l) \in tl, \text{ such} \\ & \text{that } d(o_l, i) \leq \min(r_l, r') \\ 0, & \text{otherwise} \end{cases}$$

The SCGA algorithm evaluates individuals in order of their fitness. This  $seo$  function performs the same functionality for DSGA, except it eliminates individuals within a minimum of  $r$  and  $r'$  distance of a seed on the tabu list.

The second strategy is very similar to Goldberg and Richardson's (1987) sharing function. A sharing function is an alternate way to determine fitness, called shared fitness. It weighs fitness based on the distance that the individual is to other individuals. In this approach fitness will be weighted based on the distance that the individual is to the individuals on the tabu list. The function can be defined many ways with its goal being to weigh individuals higher, the farther away they reside from the individuals on the tabu list. This approach does not necessarily have the strongest individuals as seeds; rather it selects individuals as seeds that are worth investigating.

The DSGA framework can support a variety of sharing functions for the second strategy. Goldberg and Richardson's (1987) fitness sharing function can easily be adapted to encourage exploration in DSGA. The niche count can be defined as the following:

$$m_i = \sum_{j=1}^{tabuSize} sh(d(i, o_j))$$

The parameters for this equation are described in Chapter 2. The only difference is that  $d(i, o_j)$  is the Euclidian distance between  $i$  and the  $j$ th individual on the tabu list. The function  $seo$  would be defined as the objective function divided by the niche count for individual  $i$ .

$$seo(i) = \frac{objective\_function(i)}{m_i}$$

This niche count will be smaller for individuals further away from the seeds on the tabu list. They will be more likely to be selected for crossover and be represented in future generations. This encourages exploration.

### *Set of Benchmark Optimization Problems*

After the algorithm was implemented, it was evaluated against a set of benchmarks. These benchmarks are examples of multimodal optimization problems. Prior literature shows a variety of test functions that can be used to solve multimodal optimization problems with NGAs. Some of the functions are Shubert (Ando & Kobayashi, 2005), Rosenbrock (Raghuwanshi & Kakde, 2007) and Ackley (Ling et al., 2008; Raghuwanshi & Kakde, 2007). However, one function is used most often. This function is given below:

$$f(x) = \sin^6(5.1\pi x + 0.5)$$

This function was first used in Goldberg and Richardson (1987) and this function, with minor modifications, has been used in many other research papers (Bernier, 1996; Lee, Cho & Jung, 1999; Miller & Shaw, 1996; Yin & Gerday, 1993).

Bernier (1996) generalized the Goldberg and Richardson (1987) equation. The new equation shown below can generate many different types of test cases.

$$f(x) = Rc^{-cx^2} \sin^6(k\pi x^p)$$

By defining different values for  $R$ ,  $c$ ,  $p$  and  $k$ , this equation can generate many interesting test cases similar to the one shown in Figure 1. The parameter  $c$  determines the rate of decay of each oscillation of the sine wave. In most cases,  $k$  determines the number of peaks.  $R$  controls the height of the highest peak and is set to 1.

Since the goal of this research was to develop a new NGA that can solve problems with arbitrarily close optima, while doing equally well with other optimization problems, test functions were needed to be selected in these two areas. Each function was associated with one of the goals. These equations are shown in Table 15.

Six test functions were based on Bernier's (1996) test functions. These are shown in Table 15 as F1 through F6. For the parameter  $(c, p, k)$ , the six groups of parameters were  $\{(0, 1, 5); (0, 3, 5); (0, 2, 10); (1, 3, 10); (2, 2, 5); (2, 1, 10)\}$ . The algorithm attempted to locate the local maximum of these six functions. These benchmarks have been used to test other NGAs that attempt to solve problems of this type.

The final two test functions completed the set. Function F7 in Table 15 is a general test case. Function F7 has a surface of high sides with a global and three local minimums in the center. This function was used to test Zhang, Shang, Gao and Dong's NGA

(2008). This test case tests general NGA functionality. Finally, function F8 is the function shown in Figure 1. As seen in Figure 1, this function has ever increasing optima that become arbitrarily close.

Table 15. Test Functions

	<b>Equation</b>	<b>Domain</b>	<b>Goal</b>
F1	$\max: f(x) = \sin^6(5\pi x)$	$0 \leq x \leq 1$	General
F2	$\max: f(x) = \sin^6(5\pi x^3)$	$0 \leq x \leq 1$	General
F3	$\max: f(x) = \sin^6(10\pi x^2)$	$0 \leq x \leq 1$	General
F4	$\max: f(x) = 1^{-x^2} \sin^6(10\pi x^3)$	$0 \leq x \leq 1$	Arbitrarily Close Optimum
F5	$\max: f(x) = 2^{-2x^2} \sin^6(5\pi x^2)$	$0 \leq x \leq 1$	Arbitrarily Close Optimum
F6	$\max: f(x) = 2^{-2x^2} \sin^6(10\pi x)$	$0 \leq x \leq 1$	General
F7	$\min: f(x, y) = 2x^2 - 1.05x^4 + \frac{x^6}{6} - xy + y^2$	$-3 \leq x \leq 3$ $-3 \leq y \leq 3$	General
F8	$\max: y = x \sin(x^2)$	$0 \leq y \leq 10$	Arbitrarily Close Optimum

These eight test functions cover a wide range of different multimodal functional optimization problems. Some are general test cases that can determine how a NGA handles typical functional optimization problems. Other test cases address functional optimization when optima are arbitrarily close. Use of these functions as benchmarks is supported by a wide variety of literature.

### *Performance Evaluation*

This section describes the performance goal of this research. Eight benchmark functions have been presented. The first seven benchmarks are used in other literature using a variety of performance criteria. DSGA was compared against one or more NGAs

that cite each benchmark. Most research uses benchmarks that highlight the algorithm's performance. Comparing DSGA against other algorithms by using each algorithm's benchmarks was an appropriate test.

The first six benchmarks were used in Bernier's (1996) research. This research had four performance criteria. The first criterion used was the  $X^2$ -like deviation. Ideally an NGA should have individuals distributed over the peaks relevant to the fitness of the peak. The  $X^2$ -like deviation is a measurement of how much a population deviates from this distribution. Because DSGA removes optima from the population once they are discovered and discourages them from being revisited, the criterion is not appropriate for DSGA. The next two criteria measure the proportion of the peaks that were located and the proportion of individuals outside the peaks. The proportion of peaks is the number of optima located divided by the number of optima. The proportion of individuals outside of the peaks is the number of individuals not tracking an optimum divided by the total number of individuals. The final criterion was the average fitness of the individuals in the last 50 generations.

The results of Bernier's (1996) research were the average of 10 runs for each of Bernier's algorithms: Biggest Difference Method and Biggest Proportion Method. Each run of the algorithm generated 200 generations of a population size of 100. The results for the criteria were the average of the last 50 generations for each benchmark.

Benchmark F7 has three local minimums and one global minimum, which is (0, 0). This benchmark was used in Zhang, Shang, Gao and Dong's (2008)  $hK_1$  Triangulation NGA. In this NGA there is a tuning parameter  $h$  that indicates the precision of the algorithm. The results of this research show the minimum points or most fit individuals

around the three local optima, excluding the (0, 0) minimum. The performance criterion was the objective value of the fittest individual around each niche. For this research DSGA used the same performance criterion and compare it results against the  $hK_1$  Triangulation NGA for  $h = 0.1$ .

Benchmark F8 is a function that has not been introduced in previous literature. Even though it is very close to Bernier's benchmarks, no present research has been conducted using it. This paper hopes to introduce this function as a future benchmark. The performance criteria used for benchmark F8 were the three criteria defined by Bernier (1996). These are the proportion of peaks located, proportion outside of the peaks and average fitness. These three performance criteria cover different characteristics of NGA behavior.

For comparison DSGA was compared against a number of other NGAs. When benchmarks have previously published work, results from the previous research were used in the comparison. Three additional NGAs were used in this research. The NGAs are Goldberg and Richardson's (1987) algorithm, Kao and Zahara's (2008) algorithm and SCGA (Li et al., 2002). These three algorithms were selected because they represent a variety of NGAs from a classic algorithm, like Goldberg and Richardson (1987) to a new algorithm, Kao and Zahara (2008). There are no published results for these three algorithms and the benchmark functions. As part of this research these algorithms were implemented and run against all eight benchmark functions. The published and newly obtained results were used to evaluate DSGA against the benchmarks.

Considering there are three approaches for varying the radius and two strategies for encouraging exploration, there are six distinct combinations of strategies for DSGA.

Each of the six strategies of DSGA were implemented and attempted to locate the optima of all eight benchmark functions. The performance criterion for each benchmark has been described in this section. This research used the corresponding population size, number of runs, number of generations and performance criteria as the algorithm that it is being compared against. The results of the six combinations of DSGA strategies were compared to the published results of the research cited in this section using the performance criteria shown in Table 16.

Table 16. Benchmark Algorithm Comparison

	<b>Algorithms Compared Against</b>	<b>Performance Criteria</b>
F1	Bernier Biggest Difference Method	Proportion of peaks
	Bernier Biggest Proportion Method	Proportion of points outside of peaks
	Goldberg and Richardson's Fitness Sharing	Average fitness
F2	Kao and Zahara Genetic Algorithm and Particle Swarm Optimization	
	Species Conserving Genetic Algorithm	
	Bernier Biggest Difference Method	Proportion of peaks
F3	Bernier Biggest Proportion Method	Proportion of points outside of peaks
	Goldberg and Richardson's Fitness Sharing	Average fitness
	Kao and Zahara Genetic Algorithm and Particle Swarm Optimization	
	Species Conserving Genetic Algorithm	
	Bernier Biggest Difference Method	Proportion of peaks
	Bernier Biggest Proportion Method	Proportion of points outside of peaks
	Goldberg and Richardson's Fitness Sharing	Average fitness
	Kao and Zahara Genetic Algorithm and Particle Swarm Optimization	
	Species Conserving Genetic Algorithm	

Table 16. Benchmark Algorithm Comparison Continued

	<b>Algorithms Compared Against</b>	<b>Performance Criteria</b>
	Bernier Biggest Difference Method	Proportion of peaks
	Bernier Biggest Proportion Method	Proportion of points outside of peaks
	Goldberg and Richardson's Fitness Sharing	Average fitness
F4	Kao and Zahara Genetic Algorithm and Particle Swarm Optimization Species Conserving Genetic Algorithm	
	Bernier Biggest Difference Method	Proportion of peaks
	Bernier Biggest Proportion Method	Proportion of points outside of peaks
	Goldberg and Richardson's Fitness Sharing	Average fitness
F5	Kao and Zahara Genetic Algorithm and Particle Swarm Optimization Species Conserving Genetic Algorithm	
	Bernier Biggest Difference Method	Proportion of peaks
	Bernier Biggest Proportion Method	Proportion of points outside of peaks
	Goldberg and Richardson's Fitness Sharing	Average fitness
F6	Kao and Zahara Genetic Algorithm and Particle Swarm Optimization Species Conserving Genetic Algorithm	
	Zhang, Shang, Gao, and Dong hK <sub>1</sub> Triangulation Algorithm	Fitness of best individual for each niche Proportion of peaks
F7	Goldberg and Richardson's Fitness Sharing Kao and Zahara Genetic Algorithm and Particle Swarm Optimization Species Conserving Genetic Algorithm	

Table 16. Benchmark Algorithm Comparison Continued

	<b>Algorithms Compared Against</b>	<b>Performance Criteria</b>
	Goldberg and Richardson's Fitness Sharing	Proportion of peaks
	Kao and Zahara Genetic Algorithm	Proportion of points outside of peaks
F8	and Particle Swarm Optimization	Average fitness
	Species Conserving Genetic Algorithm	

All of the algorithms shown in Table 16 have published results for the performance criteria with three exceptions. Goldberg and Richardson's Fitness Sharing method, Kao and Zahara Genetic Algorithm and Particle Swarm Optimization algorithm and SCGA did not have published results for these performance criteria. As part of this research these three algorithms were implemented. The implementations were run in an attempt to solve the benchmark functions.

### **Format for Presenting Results**

The results of this research were presented in the form of tables. There are eight performance benchmark optimization problems selected for this research. Each has between one and three performance criteria. The results contain one table for each benchmark optimization problem. The rows of the table are the selected algorithms chosen for comparison, along with the six different combinations of DSGA. The columns of the table are the performance criteria for the selected benchmark optimization problem. This method of presenting results allows for comparison between DSGA and other NGAs.

## **Resources Required**

There were few resources needed to conduct this research. The NGAs were developed in the Java programming language and ran on a desktop PC. This included implementations of DSGA, Goldberg and Richardson's Fitness Sharing method, SCGA and the Genetic Algorithm and Particle Swarm Optimization algorithm. Because this research was conducted through running trials of this new algorithm against other NGAs, no additional resources were needed. These resources were obtained to complete this research.

## **Summary**

While DSGA is not a tabu search, there are many parallels between the two techniques. Like the tabu search, DSGA investigates different areas of the domain space. New areas to investigate are selected based on previous areas. A tabu list is used to discourage redundant exploration of previously investigated areas of the domain. Unlike the tabu search, DSGA has no aspiration level. In a GA the only way to determine if a tabu move is superior is to create multiple generations based on the move. This makes aspiration levels difficult in GAs.

DSGA uses a tabu list, but not a complete tabu search to encourage exploration. As shown in Chapter 2 a tabu search contains a tabu list in addition to an aspiration condition. DSGA does not have an aspiration condition. The aspiration condition is not needed, because in DSGA moves are not completely eliminated for being on the tabu list. The seed selection algorithm encourages exploration in other areas of the domain, but does not prevent convergence to any specific area of the domain.

This new algorithm was designed to locate optima in functional optimization problems that have arbitrarily close optima. While it can locate multiple optima in a single pass, it uses multiple passes to locate all of them. After a set of optima are located, a tabu list is used to ensure that these optima are not revisited. This frees the algorithm to locate other optima. In problems that have arbitrarily close optima, it is important to prevent an optimum from eclipsing nearby optima. This algorithm attempts to overcome this problem by the exploration approaches described in this chapter.

## **Chapter 4**

### **Results**

Chapter 4 presents the results of this research. This first section describes the parameter settings and implementation methods. Where previous research did not publish parameters, values are selected. Parameters that are only specific to some algorithms are also covered. There is a section for each of the eight benchmark functions. Finally, there is a summary section.

Two of the criteria used in this research are measurements of recall and precision. Recall, defined as the number of optima identified divided by the total number of optima, is a measure of the algorithm's ability to discover optima. Bernier (1996) described this as the proportion of peaks found. Precision, defined as the total number of individuals tracking optima divided by the total number of individuals, is a measure of the algorithm's accuracy. Bernier (1996) described this as the proportion of individuals outside of the peak. Algorithms with a high proportion of individuals outside of the peak make it more difficult to determine what the optima are. These two measures provide insight into the usefulness of the algorithms.

#### **Parameter Settings and Implementation Methods**

NGAs have many parameters and implementation methods. Chromosome representation, population size and single or multiple point crossover decisions can

greatly affect the results of experiments in evolutionary algorithms (Burke, Gustafson & Kendall, 2004). NGA research often includes a section of the best parameter settings for a given algorithm. When comparing algorithms it is important to keep parameters consistent across experiments.

When previously published results were available for an algorithm, they were used instead of implementing the algorithm. This occurred with Biggest Difference Method, Biggest Proportion Method and  $hK_1$  Triangulation Algorithm. Results for Fitness Sharing; Genetic Algorithm and Particle Swarm Optimization; DSGA Increasing Radius, Seed Exclusion (DSGA (R+, S-)); DSGA Decreasing Radius, Seed Exclusion (DSGA (R-, S-)); DSGA Dynamic Radius, Seed Exclusion (DSGA (R $\Delta$ , S-)); DSGA Increasing Radius, Fitness Sharing (DSGA (R+, FS)); DSGA Decreasing Radius, Fitness Sharing (DSGA (R-, FS)); and DSGA Dynamic Radius, Fitness Sharing (DSGA (R $\Delta$ , FS)) were obtained from implementing these algorithms as part of this research.

Parameter settings and implementation methods for these results were determined by the following method. First, if results were shown from previously published research, then parameter settings and implementation methods from that research were used for the given benchmark. Second, in cases where the previous research did not state all parameters, ones were selected for the entire benchmark. Third, some algorithms have additional parameters that do not apply to other NGAs. In this case parameter values were selected and used consistently across the benchmark for all algorithms that have this parameter. This method of parameter selection should provide the most impartial comparison.

### *Previous Research Parameters and Implementations*

All results in this section came from the implementation of the algorithms with two exceptions. Results shown for Biggest Difference Method and Biggest Proportion Method came from Bernier (1996) research. Results shown for the hK<sub>1</sub> Triangulation Algorithm came from Zhang, et al. (2008). All published parameter values and implementation considerations for these algorithms were used in this research.

Functions F1 through F6 were used in Bernier (1996). The results shown in the following sections for Bernier's Biggest Different Method and Biggest Proportion Method came directly from Bernier (1996). Bernier (1996) used 30 chromosomes for each individual. The research used a population size of 100 and created 200 generations. The probability of a gene mutating was 0.001. Before determining if an individual is tracking an optimum a threshold must be defined. Bernier (1996) used 0.1, which is the threshold used in this research. Any individual that is within 0.1 of an optimum is considered tracking the optimum. These controlled parameters were used for all of the other algorithms used in F1 through F6.

Zhang, et al. (2008) did not publish parameter settings or implementation considerations that can be used in this research. As a result the parameter setting and implementation considerations will be describe below. Because there are no consistent parameters between the hK<sub>1</sub> Triangulation Algorithm and the other algorithms, it is difficult to compare the results. It is possible that other parameter values could change the results of the implemented algorithms.

### *Common Parameters and Implementations*

When previously published research did not provide values for some parameters or implementation considerations, they were selected and held consistent for all algorithms in the benchmark. In some cases they were held consistent across all of the benchmarks. This section describes the parameters selected for this research.

#### Chromosome Representation

All of the algorithms implemented for this research used binary chromosome representation, although DSGA can support binary and floating-point representation. This representation evenly divides the domain space providing greater precision as the number of chromosomes increases. Binary chromosome representation is a common method of representing individuals in a GA. This method of representation is often selected for its simplicity (Pang, 2006).

Binary chromosome representation allows for any number of chromosomes to cover an area of the domain. Assuming there are binary chromosomes  $b_{n-1}b_n \dots b_1b_0$ , an upper bound of  $UB$  and a lower bound of  $LB$ , the following equations shows the implementation of this representation (Janikow & Michalewicz, 1991).

$$x = LB + \frac{\sum_{i=0}^{n-1} b_i 2^i}{2^n - 1} (UB - LB)$$

The equation begins at the lower bound,  $LB$ . The factor  $(UB-LB)$  is the length of the domain that needs to be covered. Based upon what chromosomes are active a portion of the spanning area is added to the lower bound. The following factor of the equation produces a number between 0 and 1.

$$\frac{\sum_{i=0}^{n-1} b_i 2^i}{2^n - 1}$$

The precision of the equation can be represented by the following equation. The  $\Delta x$  term is the smallest value that  $x$  can change.

$$\Delta x = \frac{UB - LB}{2^n - 1}$$

As the number of bits,  $n$ , increases the domain is divided into smaller sections giving greater precision.

All of the algorithms in the research used crossover as a genetic operation. GAs can use single-point or multiple-point crossover. All of the implemented algorithms in this research used single-point crossover.

#### Fitness Function

Table 17 shows the fitness functions used in these trials for the research.

Table 17. Fitness Functions

<b>Benchmark Function</b>	<b>Fitness Function</b>
F1	F1
F2	F2
F3	F3
F4	F4
F5	F5
F6	F6
F7	$1 / (F7 + 1)$
F8	$F8 + 10$

In the case of F1 through F6 the fitness function was the benchmark function itself.

These are all maximization problems. F7 is a minimization problem. In this case the benchmark function of the fitness function should be inverse to each other. Benchmark function F8 has negative values, which can cause some issues for the selection process

(Beasley, Bull & Martin, 1993a). The fitness function for F8 was  $(F8 + 10)$ . For the range of  $x$  equal 0 to 10, this ensures all fitness values are positive.

#### Other Parameters

Function F7 was previously used in Zhang et al. (2008). This research did not provide parameters. As a result there were no control parameters for F7. The algorithms implemented by this research used a population size of 50, created 100 generations and had a gene mutation rate of 0.15625. Function F8 had no previously published results. All of the algorithm results for F8 used these same parameters.

All of the results from the algorithms implemented as part of this research were the average of 10 trials. Benchmark function F7 has a criterion of the best individual for each niche. The results shown in the research for the algorithms are the average of the best individual for each niche. Not all of the algorithms implemented were able to locate all of the optima in all trials. Although it is not specifically stated in Zhang et al. (2008), it is assumed that this algorithm located all of the optima.

#### *Algorithm Specific Parameters and Implementations*

Some parameters are specific to certain NGAs. In some cases they may span multiple NGAs used in this research, but not all of them. When this occurred values were selected, often from previous research, and held consistent across the benchmark. This section addresses algorithm specific parameters.

### Genetic Algorithm and Particle Swarm Optimization Parameters

Genetic Algorithm and Particle Swarm Optimization has three additional parameters (Kao & Zahra, 2008). When updating a weak member with the stronger member, two constants,  $C1$  and  $C2$ , are needed. These constants are weights to the factors when computing the new velocity. Kao and Zahra (2008) had them set to 2. This research kept the values at 2. The other parameter in this algorithm was the weight for the weak individual. This determines how much of the weak individual was maintained after the Particle Swarm Optimization step. Kao and Zahra (2008) calculated this as  $0.5 + Z / 2$  where  $Z$  is a uniform random number between 0 and 1. This research kept this calculation as well. These are the additional parameters for the Genetic Algorithm and Particle Swarm Optimization algorithm.

### Fitness Sharing Parameters

Four algorithms used a fitness sharing method: Goldberg and Richardson's Fitness Sharing, DSGA (R+, FS), DSGA (R-, FS) and DSGA (R $\Delta$ , FS). These algorithms all implement Goldberg and Richardson's (1987) algorithm. The implementation of these algorithms used the power law function, described in Chapter 2, as the fitness sharing function. The parameters  $\sigma_{\text{share}}$  and  $\alpha$  were set to 0.1 and 1 respectively, which were the same parameter values as in Goldberg and Richardson's (1987).

### Species Conserving Parameters

The SCGA algorithm had additional parameters. Since DSGA was based on SCGA, these parameters are also needed in DSGA. The parameter  $\sigma_S$  defines the diameter of the

neighborhood. In these trials  $\sigma_S$  was set to 0.2, which makes a radius of 0.1. In DSGA this parameter was *IS*. This radius was used for all benchmark functions, even F7 and F8 which have large domain areas.

### DSGA Specific Parameters

DSGA has additional parameters to the SCGA algorithm. In DSGA the radius changes as the algorithm runs. The Sigma Delta, *SD*, determines the change in the radius. The *SD* parameter was set to 0.015 in all trials. DSGA also has a parameter, *RLC*. *RLC* or Reevaluation Loop Count determines how many inner loops of the algorithm should be performed before the radius is reevaluated. In trials for F1 through F6 *RLC* was set to 50. This divides the total number of generations, 200, into four groups. In trials for F7 and F8 *RLC* was set to 25. This divided the total number of generations, 100, into four groups. After every *RLC* number of generations DSGA analyzes the last generation seeking optima. If there *CL* number of identical individuals, the individual is placed on the tabu list and considered an optima. *CL* was set to two in all trials. DSGA specific parameters used in the following trials are described above.

One of the methods to vary the radius of the DSGA framework is to increase or decrease the radius based upon information after each iteration of the inner loop completes. While there are many different ways that this can be implemented, one consistent method was used in this research. After each iteration of the inner loop completes the algorithm checks to see how many individuals were added to the tabu list through convergence. If two or more individuals were added to the tabu list, the radius was increased by *SD*. Otherwise, it was decreased by *SD*. This implementation was

selected to increase the chance of finding optima. If fewer than two areas of the domain converged, decreasing the radius would allow more seeds to be identified in future generations and should preserve more areas of the domain.

In most GAs the final generation contains the optima that the GA has located. That is not the case with DSGA. DSGA removes optima from the population through the use of the tabu list. Therefore the final generation will not contain the optima located. The tabu list contains the optima. The data provided for all benchmarks for the criterion of proportion of peaks located for DSGA came from analyzing the tabu list, not the last generation. The data provided for the F7 criteria of best individual in each niche also came from the DSGA tabu list. The data for all other criteria for DSGA came from the population.

This section describes all of the parameters used in this research. Controlled parameters are the parameter values used in prior research. When prior research provided parameter values, they were maintained throughout all trials. Some algorithms required additional parameter values. These values have been described. When a parameter existed in multiple algorithm, the parameter value was kept consistent across all trials for a given benchmark function.

## **Results of Algorithms on F1**

Benchmark function F1 is a sine wave with five evenly distributed local maximums all of equal magnitude. The results for the Biggest Difference Method and Biggest Proportion Method come from Bernier (1996) research. Table 18 shows the results for F1. Figure 4 is a chart of the precision and recall of the algorithms.

Table 18. Results for Equation F1

Algorithm	Recall	Precision	Average fitness
Bernier Biggest Difference Method	0.9800	0.7044	0.863015
Bernier Biggest Proportion Method	1.0000	0.6012	0.829098
Goldberg and Richardson's Fitness Sharing	0.9000	0.6570	0.8590
Kao and Zahara Genetic Algorithm and Particle Swarm	0.2000	0.9890	0.9855
SCGA	0.9800	0.9250	0.9630
DSGA (R+, S-)	0.9800	0.9376	0.9714
DSGA (R-, S-)	0.9800	0.9158	0.9754
DSGA (R $\Delta$ , S-)	0.9400	0.9537	0.9859
DSGA (R+, FS)	0.9800	0.9568	0.4839
DSGA (R-, FS)	0.9800	0.8749	0.9667
DSGA (R $\Delta$ , FS)	0.9800	0.9425	0.9754

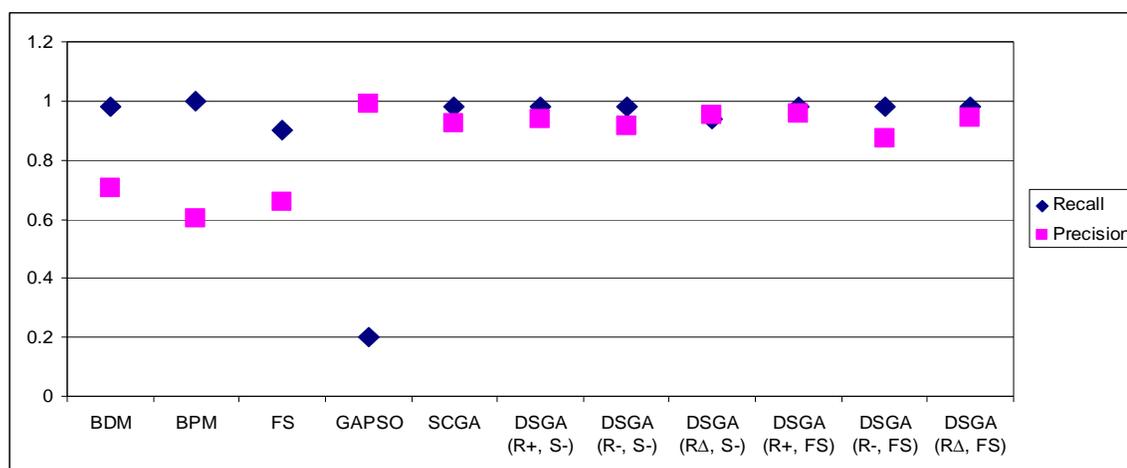


Figure 4. Chart of Recall and Precision for F1

Bernier's (1996) algorithms did not outperform all of the other algorithms in all criteria. No algorithm tested could locate as many peaks as Biggest Proportion Method, 100%. However, SCGA and five of the six DSGA algorithms located 0.9800 of them, which is the number that Biggest Difference Method found. The algorithm that had the fewest individuals outside of the peaks was the Genetic Algorithm and Particle Swarm Optimization algorithm. The algorithm with the highest average fitness was DSGA (R $\Delta$ , S-). However, DSGA (R+, FS) had the lowest average fitness.

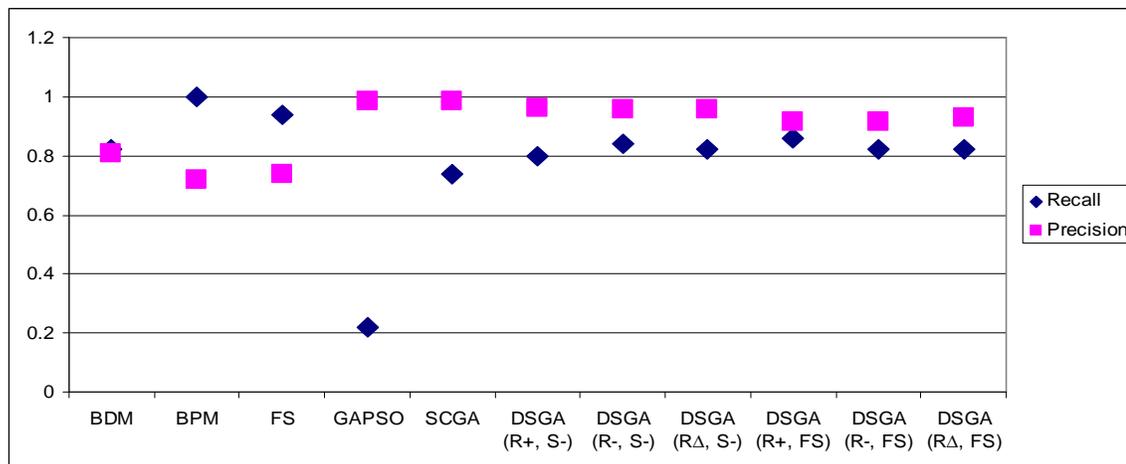
## Results of Algorithms on F2

The F2 function also is a sine wave that has five local maximums all of equal magnitude. However, in F2 the optima are increasingly closer together. Table 19 shows the results of the average of 10 trials for function F2. Data provided for Biggest Difference Method and Biggest Proportion Method comes from Bernier (1996). Figure 5 shows a chart of the precision and recall.

Table 19. Results for Equation F2

Algorithm	Recall	Precision	Average fitness
Bernier Biggest Difference Method	0.8220	0.8095	0.891414
Bernier Biggest Proportion Method	1.0000	0.7187	0.854766
Goldberg and Richardson's Fitness Sharing	0.9400	0.7400	0.8808
Kao and Zahara Genetic Algorithm and Particle Swarm	0.2200	0.9850	0.9831
SCGA	0.7400	0.9830	0.9801
DSGA (R+, S-)	0.8000	0.9626	0.9888
DSGA (R-, S-)	0.8400	0.9574	0.9862
DSGA (R $\Delta$ , S-)	0.8200	0.9577	0.9894
DSGA (R+, FS)	0.8600	0.9138	0.4897
DSGA (R-, FS)	0.8200	0.9156	0.9806
DSGA (R $\Delta$ , FS)	0.8200	0.9292	0.9834

As in F1 no algorithm could meet Biggest Proportion Method in locating 100% of the peaks. The closest algorithm for this criterion was the Fitness Sharing method with 0.9400 peaks located. The Genetic Algorithm and Particle Swarm algorithm had the fewest proportion of individuals outside of the peaks, but only found 22% of the optima. In the average fitness criterion all of the DSGA algorithms did well with the exception of DSGA (R+, FS).



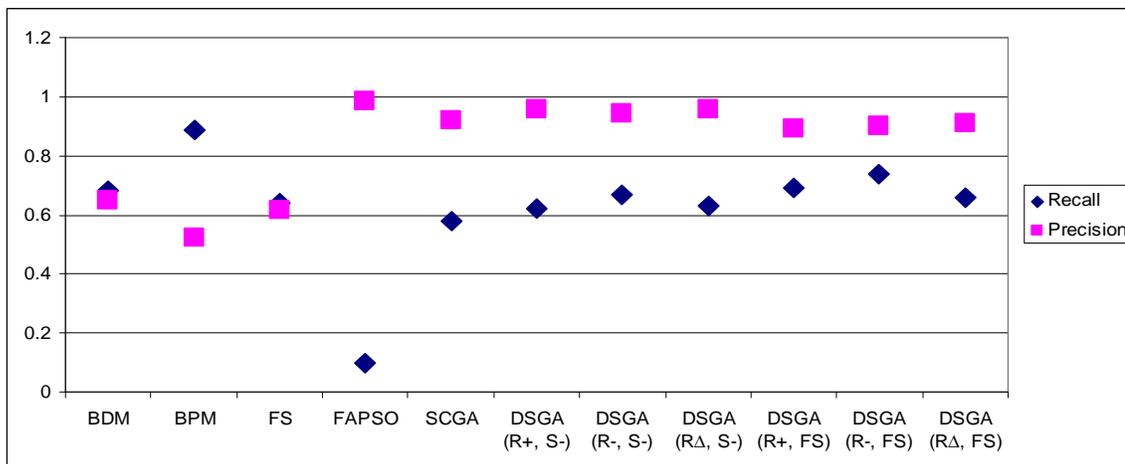
**Figure 5.** Chart of Recall and Precision for F2

### Results of Algorithms on F3

The function F3 is similar to F2 except that it has 10 optima instead of five. Table 20 shows the results for function F3 and Figure 6 is a chart of the results. As stated previously data for Biggest Difference Method and Biggest Proportion Method comes from Bernier (1996).

Table 20. Results for Equation F3

Algorithm	Recall	Precision	Average fitness
Bernier Biggest Difference Method	0.6822	0.6495	0.848098
Bernier Biggest Proportion Method	0.8880	0.5247	0.806407
Goldberg and Richardson's Fitness Sharing	0.6400	0.6170	0.8477
Kao and Zahara Genetic Algorithm and Particle Swarm	0.1000	0.9840	0.9868
SCGA	0.5800	0.9180	0.9658
DSGA (R+, S-)	0.6200	0.9570	0.8979
DSGA (R-, S-)	0.6700	0.9410	0.9786
DSGA (RΔ, S-)	0.6300	0.9561	0.9882
DSGA (R+, FS)	0.6900	0.8924	0.4889
DSGA (R-, FS)	0.7400	0.8995	0.9682
DSGA (RΔ, FS)	0.6600	0.9090	0.9770



**Figure 6.** Chart of Recall and Precision for F3

Biggest Proportion Method located all of the peaks in the function with DSGA (R-, FS) locating the second most peaks at 0.7400. As in the previous benchmarks the Genetic Algorithm and Particle Swarm Optimization algorithm had the least number of individuals outside of a peak. DSGA (RΔ, S-) had the greatest average fitness of the last 50 generations with a fitness of 0.9882.

#### Results of Algorithms on F4

The function F4 is similar to F3 except that the 10 optima are even closer together. Table 21 shows the results of the average of the 10 trials. Figure 7 is a chart of the precision and recall for the algorithms for F4.

For this function Biggest Proportion Method outperforms all of the other algorithms in proportion of peaks found by at least 0.3. As in all other functions the Genetic Algorithm and Particle Swarm Optimization algorithm produced the best results for the proportion of points outside of the peaks. The algorithm that had the greatest average fitness in the

last 50 generations was the DSGA (R+, S-) algorithm. The ranking of algorithms for the different criteria in F4 is very similar to that of F3.

Table 21. Results for Equation F4

Algorithm	Recall	Precision	Average fitness
Bernier Biggest Difference Method	0.5480	0.7327	0.742360
Bernier Biggest Proportion Method	0.89866	0.6738	0.611701
Goldberg and Richardson's Fitness Sharing	0.5800	0.7170	0.8546
Kao and Zahara Genetic Algorithm and Particle Swarm	0.1000	0.9800	0.9852
SCGA	0.4600	0.9440	0.9698
DSGA (R+, S-)	0.5200	0.9544	0.9881
DSGA (R-, S-)	0.5200	0.9438	0.9832
DSGA (R $\Delta$ , S-)	0.5100	0.9642	0.9834
DSGA (R+, FS)	0.6500	0.8814	0.4775
DSGA (R-, FS)	0.6400	0.9194	0.9750
DSGA (R $\Delta$ , FS)	0.6300	0.9041	0.9702

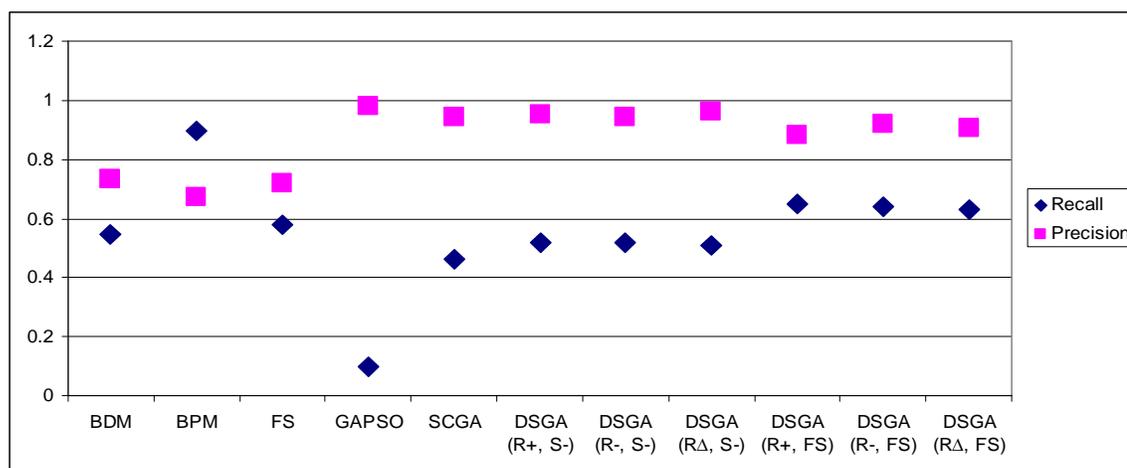


Figure 7. Chart of Recall and Precision for F4

### Results of Algorithms on F5

Function F5 is the first function that has optima of different magnitudes. It has five optima of decreasing fitness. Table 22 shows the results for function F5 for 10 trials of the algorithms implemented. Figure 8 is a chart of the recall and precision.

Table 22. Results for Equation F5

Algorithm	Recall	Precision	Average fitness
Bernier Biggest Difference Method	0.9336	0.8598	0.694831
Bernier Biggest Proportion Method	1.0000	0.8335	0.584304
Goldberg and Richardson's Fitness Sharing	0.9400	0.7580	0.5648
Kao and Zahara Genetic Algorithm and Particle Swarm	0.2000	0.9790	0.8165
SCGA	0.1000	0.9900	0.8615
DSGA (R+, S-)	0.2000	0.9541	0.8601
DSGA (R-, S-)	0.1000	0.9670	0.8588
DSGA (R $\Delta$ , S-)	0.1000	0.9521	0.8602
DSGA (R+, FS)	0.1000	0.9496	0.4302
DSGA (R-, FS)	0.1000	0.9511	0.8599
DSGA (R $\Delta$ , FS)	0.2000	0.9396	0.8587

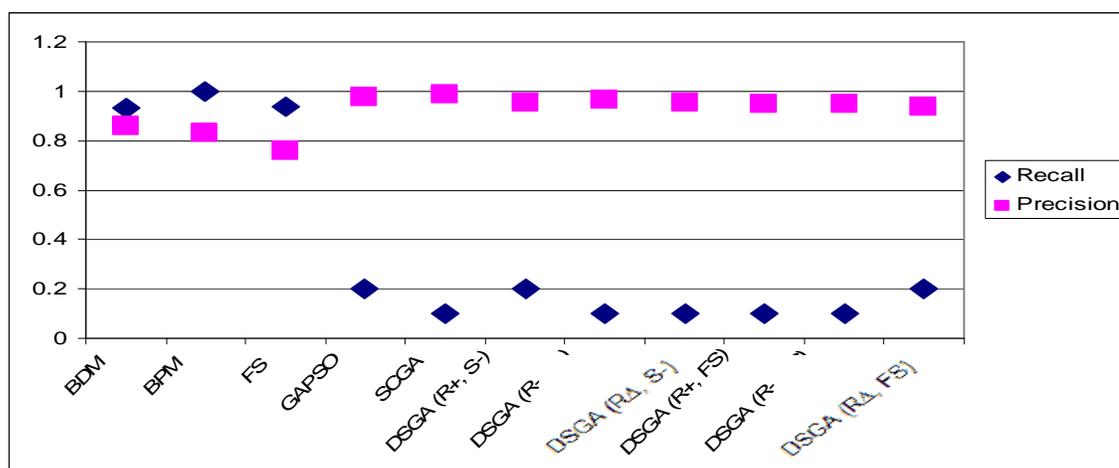


Figure 8. Chart of Recall and Precision for F5

Three algorithms did very well in locating peaks: Biggest Difference Method, Biggest Proportion Method and Fitness Sharing. Each located 0.9336 or more optima. All of the other algorithms did poorly finding no more than 0.2 optima. SCGA had the most number of individuals tracking a peak with 0.99. The algorithm with the best average fitness was DSGA (R $\Delta$ , S-).

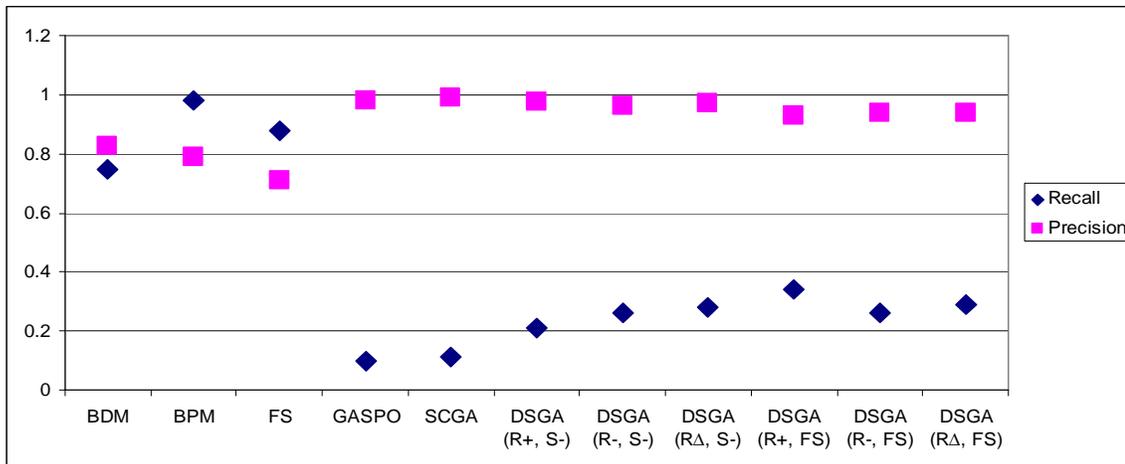
## Results of Algorithms on F6

Function F6 has 10 optima of decreasing fitness. The results for function F6 are shown in Table 23. The recall and precision are shown in Figure 9.

Table 23. Results for Equation F6

<b>Algorithm</b>	<b>Recall</b>	<b>Precision</b>	<b>Average fitness</b>
Bernier Biggest Difference Method	0.7480	0.8247	0.682606
Bernier Biggest Proportion Method	0.9788	0.7891	0.614887
Goldberg and Richardson's Fitness Sharing	0.8800	0.7110	0.6799
Kao and Zahara Genetic Algorithm and Particle Swarm	0.1000	0.9810	0.9345
SCGA	0.1100	0.9900	0.9860
DSGA (R+, S-)	0.2100	0.9738	0.9759
DSGA (R-, S-)	0.2600	0.9608	0.9816
DSGA (R $\Delta$ , S-)	0.2800	0.9692	0.9739
DSGA (R+, FS)	0.3400	0.9281	0.4885
DSGA (R-, FS)	0.2600	0.9370	0.9822
DSGA (R $\Delta$ , FS)	0.2900	0.9390	0.9708

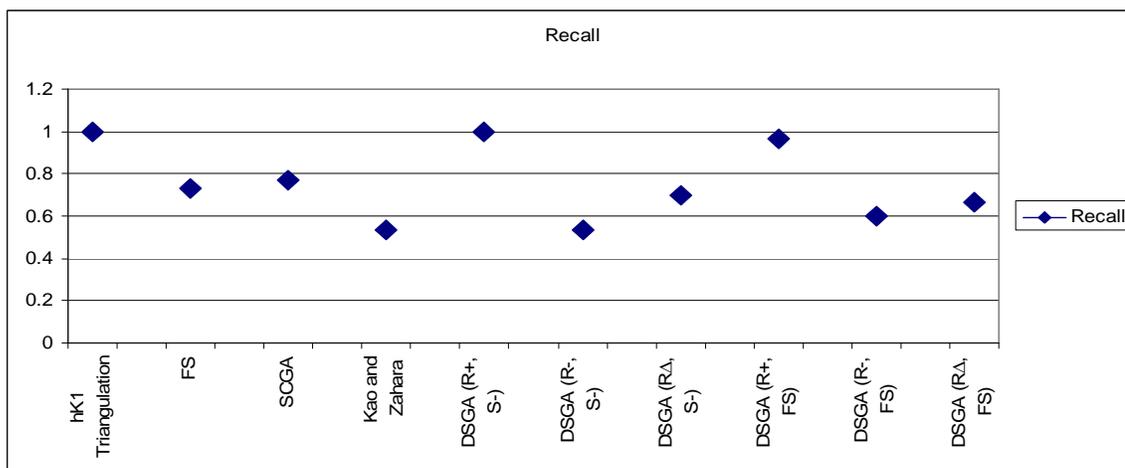
Once again Biggest Difference Method, Biggest Proportion Method and Fitness Sharing did very well at locating peaks and the other algorithms did not. SCGA did the best at having the least number of individuals outside of the peaks and also had the highest average fitness. The DSGA algorithms did poorly at locating peaks, finding no more than 0.3400 of them. However, they did very well at having very high average fitness of the last 50 generations.



**Figure 9.** Chart of Recall and Precision for F6

### Results of Algorithms on F7

The benchmark function F7 has three local optima. The global optimum is at (0, 0). One criterion for this function is the  $F(x, y)$  value of the best individual for each of the three optima. Since this is a minimization problem, smaller values are advantageous. The data for the  $hK_1$  Triangulation Algorithm came from Zhang, et al. (2008). The data for the other algorithms came from the implementation of the algorithms for this research. The results of this test can be seen in Table 24. Figure 10 shows the recall.



**Figure 10.** Chart of Recall for F7

Table 24. Results for Equation F7

<b>Algorithm</b>	<b>F(x, y) of best individual for each niche</b>	<b>Recall</b>
Zhang, Shang, Gao and Dong hK <sub>1</sub> Triangulation Algorithm	0.000015 0.003706 0.003706	1.0000
Goldberg and Richardson's Fitness Sharing	0.527280 0.077355 0.94090	0.7333
SCGA	0.415233 0.000919 0.439676	0.7667
Kao and Zahara Genetic Algorithm and Particle Swarm	1.082841 0.000003 1.001006	0.5333
DSGA (R+, S-)	0.318560 0.000545 0.320355	1.0000
DSGA (R-, S-)	0.326982 0.003262 0.313548	0.5333
DSGA (RΔ, S-)	0.372950 0.001551 0.355833	0.7000
DSGA (R+, FS)	0.324833 0.004004 0.324392	0.9667
DSGA (R-, FS)	0.325207 0.002342 0.317107	0.6000
DSGA (RΔ, FS)	0.306499 0.002572 0.319992	0.6667

The algorithms that were implemented did not find all of the optima in all of the trials. This was the reason that the proportion of peaks criterion is included in this results section. Zhang, et al. (2008) did not specifically state how many optima their algorithm located. It is assumed that all trials located all three optima.

After averaging the sum of 10 trials for each algorithm implemented, the  $hK_1$  Triangulation Algorithm did the best for optima 1 and 3. Optimum 2 was the global minimum of (0, 0). For this optimum the Genetic Algorithm and Particle Swarm Optimization algorithm performed best. Of the algorithms implemented only DSGA (R+, S-) found all of the peaks in all 10 trials. This is impressive since it is only an assumption that the  $hK_1$  Triangulation Algorithm located all of them.

### **Results of Algorithms on F8**

Function F8 has been discussed in Chapter 1 and is the best example of the types of functions that DSGA was developed to solve. This function has arbitrarily close optima. Between the  $x$  values of 0 and 10, there are 16 optima. Most of the optima are within the radius value of other optima. All of the algorithms for this benchmark function were implemented as part of this research. The results of the average of 10 trials can be seen in Table 25.

DSGA overwhelmingly outperformed the other algorithms in many of the criteria. All six of the DSGA algorithms found more peaks than any of the other algorithms. The Genetic Algorithm and Particle Swarm Optimization algorithm had the least number of individuals outside of a peak and had the highest average fitness for the last 50

generations. However, for proportion of points outside of peaks and average fitness, all DSGA algorithms did better than the Fitness Sharing algorithm.

Table 25. Results for Equation F8

Algorithm	Recall	Precision	Average fitness
Goldberg and Richardson's Fitness Sharing	0.4375	0.2800	1.6982
Kao and Zahara Genetic Algorithm and Particle Swarm	0.0875	0.8040	7.7072
SCGA	0.8625	0.7640	6.7078
DSGA (R+, S-)	0.9625	0.5930	6.4542
DSGA (R-, S-)	0.9688	0.5050	5.9423
DSGA (RA, S-)	0.9313	0.5479	6.6220
DSGA (R+, FS)	0.9500	0.3736	5.4074
DSGA (R-, FS)	0.9500	0.4225	5.7020
DSGA (RA, FS)	0.9563	0.3691	5.4805

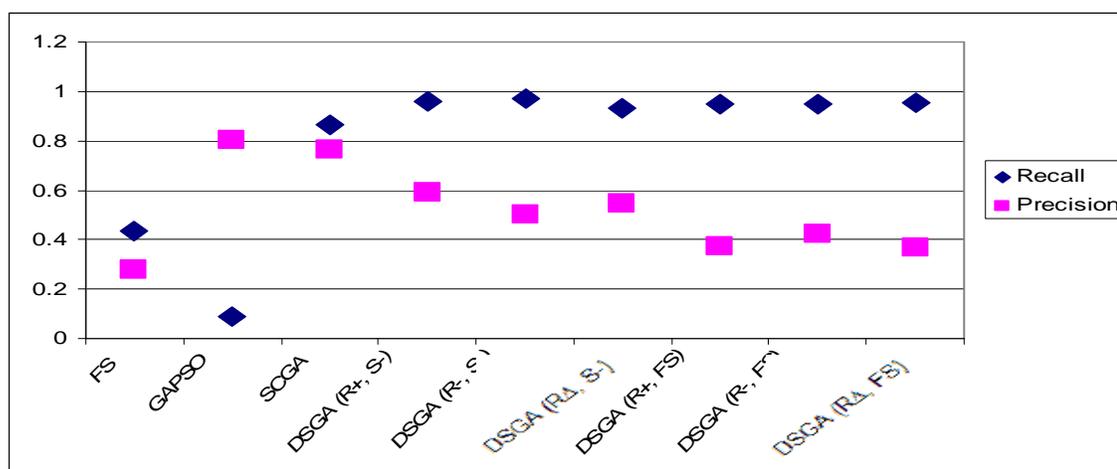


Figure 11. Chart of Recall and Precision for F8

## Summary of Results

This chapter provides the results of six algorithms derived from the DSGA framework compared to six other NGAs. Research results came from eight benchmark functional optimization problems, seven of which had been used in prior research. The benchmark functions covered many different functional optimization problems, including

minimization and maximization problems and two and three dimensional problems. Each optimization problem used two or three criteria.

Results from the Genetic Algorithm and Particle Swarm Optimization algorithm did poorly for most criteria. These results do not correspond to other results for this algorithm (Kao & Zahra, 2008). Two factors could explain this. First the controlled parameters used in Bernier (1996) may not be the best parameter settings for this algorithm. Perhaps with a different mutation rate or population size this algorithm would have located more optima. Kao and Zahra (2008) noted that higher mutations rates increase the algorithm's ability to locate optima. Second, results published in Kao and Zahra (2008) represented individuals as a vector of real numbers instead of a binary implementation. This research kept the chromosome representation as binary since the other algorithms were coded using binary representations. Other research indicates that chromosome representation can affect results in GAs (Golub, 1996). This could explain the poor performance of the Genetic Algorithm and Particle Swarm Optimization algorithm.

Each algorithm performed differently against this set of benchmark functions. Some performed consistently well, others performed poorly. Appendix A Table 26 shows how the algorithms ranked for each benchmark and criteria. A ranking of one is the best performing algorithm. Higher ranking algorithms did not perform as well as lower ranking ones for the given criteria. Results shown in this table for function F7 with criteria of  $F(x, y)$  of best individuals for each niche, shows the results of the sum of the three best  $F(x, y)$  values. While algorithm performance varied widely, no single algorithm proved superior.

## Chapter 5

### Conclusions, Implications, Recommendations and Summary

This chapter discusses the conclusions, implications, recommendations and summary of this research. There is a section in this chapter for each of these four topics. The conclusion section analyzes the results against the hypothesis. The implications section discusses the impact of this research and the contribution to the field. The recommendations section presents future research ideas. The summary section summarizes this research.

#### Conclusions

The DSGA framework was developed to solve functional optimization problems for continuous functions when the optima are arbitrarily close. The framework allows for the creation of multiple algorithms. There are two categories of strategies. The first category is how to change the radius as the algorithm runs. The second category addresses how to encourage exploration in the domain. The DSGA framework provides a foundation for the building of a variety of algorithms to solve for arbitrarily close optima.

The first goal of this research was to develop an algorithm to solve for arbitrarily close optima. The benchmark function F8 which is shown in Figure 1 is an example of one such function. In this example a majority of the optima are within the radius value of each other. All six DSGA algorithms did well in locating optima for F8. They located

93.13% to 95% of the optima. The three other algorithms tested only located 86.25%, 43.75% and 8.75% of the optima. The DSGA framework is remarkably good at locating arbitrarily close optima. Results for F8 indicate that this goal was met by the DSGA framework.

The second goal of this research was to develop an algorithm that will work equally well for other types of problems. Excluding F8, there were 20 combinations of benchmark functions and criteria. Six benchmark functions had three criteria and one benchmark function had two criteria. Of these 20 combinations there were six combinations in which a DSGA algorithm was ranked one. So, in 30% of the cases a DSGA algorithm outperformed all other algorithms. Of the 14 combinations in which DSGA was not ranked one, it was ranked two in seven combinations. While DSGA algorithms did not always rank number one, results seem to indicate that it does equally well against other types of problems.

The first hypothesis of this research was that finding optima in phases is a better strategy for locating arbitrarily close optima. All results showed that this is a good strategy for these types of problems. Consider the DSGA (R+, S-) and the DSGA (R+, FS) algorithms. The beginning radius value was 0.1 and it increased by 0.015 each of the four phases that the algorithm performed. This means that the four values of the radius were 0.1, 0.115, 0.130 and 0.145. Of the 16 optima for F8 all but two had other optima within these four radii. Multiple optima within a radius will cause problems for NGAs (Ando & Kobayashi, 2005). But the two DSGA algorithms located 96.25% and 95.0% of the optima. This occurred because each phase located some optima and removed them from the search through the two exploration strategies to allow the algorithm to locate the

other optima with the remaining phases. The SCGA algorithm, which performed only one phase, only located 86.25% of the optima. The approach to solving arbitrarily close optima problems in phases is supported by these results.

The second hypothesis was that traditional NGAs do poorly against arbitrarily close optima because of their use of a static radius. Changing the radius as the algorithm runs compensates for the difficulty in solving these types of problems. Results from this research confirmed this hypothesis. DSGA and SCGA are very similar. SCGA has a static radius and DSGA has a dynamic radius. With the exception of benchmark function F1, a DSGA algorithm found as many or more optima as SCGA. In F1 the distance between optima was greater than the radius. As more optima exist within the radius, the ability for SCGA to locate optima decreased to about half of what DSGA located. Varying the radius as the algorithm runs helped in adjusting for poorly chosen radius values.

Determining which DSGA strategies were the best is difficult. All six of the DSGA algorithms performed against the benchmarks equally well. One exception to this observation is the DSGA (R+, FS) algorithm. This algorithm consistently had an average fitness about half of what the other DSGA algorithms had. The average fitness criterion was the average fitness of the last 50 generations. This is a difficult criterion for DSGA, because DSGA removes optima from the population when they are placed on the tabu list. The DSGA (R+, FS) algorithm did perform well at locating the optima. One explanation for this low average fitness could be in the order that DSGA (R+, FS) located the optima. It could have located the fittest optima first and be left with the least fit ones

in the final generations. With a few exceptions all six DSGA algorithms performed equally well across the benchmark criteria.

This research has demonstrated that the DSGA framework is very effective at solving problems with arbitrarily close optima. Its ability to solve other types of problems is comparable to other NGAs. Many factors attribute to DSGA's ability to solve such problems. Finding optima in phases and then removing them from the search space allows the algorithm to decompose the problem and find answers iteratively. The use of a tabu list to store areas of the domain that have been investigated and found to be optimal allows DSGA to encourage exploration into other areas of the domain. Changing the radius as the algorithm executes compensates for poor radius choices that limit other NGAs. DSGA even proved successful when all of the radius values had multiple optima within them. Results for DSGA showed that it was successful at solving many types of functional optimization problems.

### **Implications**

The results of this research can be useful in a variety of areas. The DSGA framework has been shown to be successful in locating optima for problems with arbitrarily close optima. When it is known or suspected that a function has arbitrarily close optima a DSGA algorithm would be appropriate in locating maximums and minimums. Results of this research show that it locates more optima than other NGAs for these types of problems.

Another area that the DSGA framework has implications in is when there is little or no knowledge of where the optima are located. Without proper parameter settings many

NGAs have difficulty locating optima. DSGA algorithms do very well locating optima even when the radius parameter is set incorrectly. Other NGAs have difficulties locating optima when a poor radius parameter is selected.

This research also introduced a new benchmark function, F8. Results against other NGAs showed that this function is difficult to solve for many NGAs. This function could be used in future research to test other NGAs.

### **Recommendations**

While the results of this research support the hypotheses, there are still unanswered questions about this approach. More research could be done to provide a better understanding of the value that DSGA has. The following are some areas where more research is recommended.

DSGA has been tested against eight benchmark functions. While seven of the eight functions have been used in other NGA research, DSGA has not been applied to real-world problems. Future research could be done to test DSGA against real-world problems like those outlined in Chapter 1: handwriting matching, electromagnetic system design and data mining classification.

DSGA has a variety of parameters and implication considerations. In addition to traditional GA parameters like population size, number of chromosomes and number of generations, there are specific parameters like radius and radius delta. The fitness sharing strategy in this researched used the power law function, but many other functions could be used to implement this strategy. More research with other parameters and implementation considerations could be conducted.

The DSGA framework enhances the SCGA algorithm. It enhances it based upon a few principles: locating optima in phases and then excluding optima in future generations, use of a tabu list to store optimal candidates and changing the radius as the algorithm runs. All of these have shown to be very useful enhancements to the SCGA algorithm, but they could be applied to other NGAs. Research that applies these principles to another NGA would provide additional evidence that this approach is correct.

These recommendations highlight some additional areas of research that could be undertaken. The DSGA framework has generated six algorithms that prove to be very useful for some types of problems. However, they have only been tested against eight functional optimization problems. Additional research can better define the usefulness of DSGA.

## **Summary**

GAs can be useful tools for searching large, complex domain spaces. GAs do very well when searching for a single optimum. But when they attempt to locate multiple optima, they often fail. GAs have two competing forces that act upon the population. Mutation expands the area of the domain that is being searched. This exploration increases the area of the search space. Selection and crossover eliminate areas of the domain and focus the search on ever shrinking areas of the domain. This exploitation reduces the area of the search space. In every GA selection and crossover eventually win out and the population converges.

### *Classify NGAs*

NGAs are a specific type of GA that employ novel methods to prevent the exploitation force from removing optima in the domain space. Currently there are many NGAs, which can be classified as fitness sharing methods, crowding methods and other methods. This research provides multiple examples of all three categories.

In fitness sharing methods special fitness functions are used. These functions alter the fitness of individuals based upon how far they are from other individuals in the population. More isolated individuals are given preference to increase their chances of being selected for crossover. This provides preservation for individuals that are in low populated areas of the domain.

A second category of NGAs is crowding methods. In crowding methods individuals from one generation are promoted into the next generation. These individuals are often the fittest individuals in a specific area of the domain. Crowding methods prevent the exploitation forces of fit optimum from eclipsing weaker optima by directly maintaining interesting individuals.

There are some NGAs that do not easily fit into the fitness sharing or crowding categories. The other category groups these methods. Some of these methods are other GAs that solve multiple optima problems, like Cellular Genetic Algorithms. Many methods in this category are hybrid methods. These methods combine GAs with other search algorithms, like the Particle Swarm Optimization and the Tabu Search. This other category classifies NGAs that solve multiple optima problems but do not use fitness sharing or crowding approaches.

### *DSGA Framework*

One problem that many NGAs have is that when optima become arbitrarily close they have difficulty locating all of the optima. Most NGAs have some radius parameter. When searching the domain the parameter is used to determine how large an area of the domain should be to make it worth preserving. The algorithm assumes that if two individuals are within the radius, they are tracking the same optima. But this may not be the case. An issue arises when no matter how small the radius is set to; there is some area of the domain that has multiple optima within the radius (Ando & Kobayashi, 2005). When this happens one optimum is often preserved and the others are lost. This makes functional optimization problems of continuous functions that have arbitrarily close optima difficult for NGAs to solve.

DSGA is a new NGA framework developed to solve functional optimization problems of continuous functions that have arbitrarily close optima. The DSGA framework is based upon the SCGA algorithm. The SCGA algorithm is a crowding NGA, but was not developed to specifically address problems of arbitrarily close optima. The enhancements made to SCGA are supported by other research.

SCGA is a crowding NGA. It identifies interesting individuals within a population. These individuals are called seeds. Seeds get promoted into the next generation. Seed selection begins by sorting a population by the fitness of each individual. Individuals are evaluated from the fittest to the least fit. A radius parameter is used to define the area around a seed. As individuals are evaluated, if they are not within the radius of an existing seed, the individual is added to the list of seeds. SCGA uses normal selection, crossover and mutation. When the next generation is created SCGA replaces members of

this new generation with individuals on the list of seeds. Each seed replaces the weakest individual in the new generation that is within the radius of the seed in the domain space. After all of the seeds are promoted into the next generation the list of seeds is emptied and individuals must compete again to be a seed. This allows SCGA to preserve these individuals into the next generation.

DSGA enhances SCGA in a number of ways. DSGA does not attempt to locate optima in a single loop. It runs a series of generations in an attempt to locate some optima. Optima and seeds are placed on a short term memory structure, called a tabu list. Then it encourages exploration into other areas of the domain to locate undiscovered optima. DSGA has a radius parameter, which often is a limitation for most NGAs. DSGA overcomes the problem of having multiple optima within the radius, by varying the radius as the algorithm runs. DSGA uses two strategies to vary the radius and two methods to encourage exploration.

DSGA has two strategies for varying the radius. It has two parameters concerning the radius. DSGA has a radius parameter and a radius delta parameter. After a series of generations are created, DSGA changes the radius by the radius delta parameter. The two strategies for varying the radius are to always increase or decrease the radius and vary the radius based upon some condition. The condition to vary the radius is arbitrary, but the strategy was developed to use run-time information to determine if the radius should be increased or decreased.

There are two strategies for encouraging exploration in DSGA. One is based upon the fitness sharing method. A fitness function is defined in such a way that it decreases an individual's fitness the closer that the individual is to members of the tabu list. This

differs from other fitness sharing algorithms that vary the fitness based upon how close individuals are to other individuals in the population. This encourages exploration into other areas of the domain. The second strategy for encouraging exploration prevents individuals from being seeds. If an individual is within the radius of an individual on the tabu list, it is excluded from being a seed. These two strategies encourage exploration in DSGA.

### *Research Results*

The research had two goals and two hypotheses. The first goal was to develop an NGA that could solve problems with arbitrarily close optima. The second goal was that this NGA would perform as well as other NGAs for other types of problems. The first hypothesis was that finding optima in phases, increases a NGAs chances of finding arbitrarily close optima. The second hypothesis was that NGAs often miss optima in problems with arbitrarily close optima because of static radius. Eight functional optimization problems for continuous functions were used to test these goals and hypotheses.

DSGA was compared to six other NGAs with eight benchmark functional optimization problems. Each benchmark function had two or three criteria to be judged against. One specific function had ever increasing arbitrarily close optima. In one area of this domain the function had multiple optima within the radius.

The results of this research support the two hypotheses and show that the two goals were met. Each of the six combinations of DSGA strategies located more optima than any of the other algorithms tested for the benchmark function with arbitrarily close

optima. It even located more optima than SCGA, which shows that the ability to locate arbitrarily close optima was not inherent in SCGA. Rather this ability came from the enhancement that this research made in DSGA. For the other seven benchmark functions DSGA performed equally well as other algorithms. This research indicates that locating optima in phases works better for arbitrarily close optima and that static radius often prevent other NGAs from locating such optima.

### *Conclusions*

DSGA is a new NGA framework that was designed specifically to locate optima in problems that have arbitrarily close optima. For problems in which multiple optima existed within the radius, all DSGA algorithms located more optima than any of the other algorithms used. DSGA does a respectable job against other functional optimization problems. The results of this research show that the DSGA framework does very well against functional optimization problems.

The DSGA performance comes from two factors. Locating optima in phases and then encouraging exploration away from the located optima, simplifies the problem. This makes locating optima easier. Varying the radius as the algorithm runs compensates for poor radius choices. These two characteristics of DSGA make it a useful search technique.

DSGA is a new NGA framework. It was developed to solve for functional optimization of continuous functions when the optima are arbitrarily close. However, DSGA results for problems that do not have arbitrarily close optima were comparable to

other NGAs. The DSGA framework provides a new NGA approach that leverages existing NGA research.

## Appendix A

## Ranking of Algorithms

Table 26. Ranking of Algorithms

Benchmark Criteria	Rank	Ranked Algorithm
F1 Proportion of Peaks	1	Bernier Biggest Proportion Method
	2	Bernier Biggest Difference Method
	2	SCGA
	2	DSGA (R+, S-)
	2	DSGA (R-, S-)
	2	DSGA (R+, FS)
	2	DSGA (R-, FS)
	2	DSGA (R $\Delta$ , FS)
	3	DSGA (R $\Delta$ , S-)
F1 Proportion of points outside of peaks	4	Goldberg and Richardson's Fitness Sharing
	5	Kao and Zahara Genetic Algorithm and Particle Swarm
	1	Kao and Zahara Genetic Algorithm and Particle Swarm
	2	DSGA (R+, FS)
	3	DSGA (R $\Delta$ , S-)
	4	DSGA (R $\Delta$ , FS)
	5	DSGA (R+, S-)
	6	SCGA
	7	DSGA (R-, S-)
	8	DSGA (R-, FS)
	9	Goldberg and Richardson's Fitness Sharing
10	Bernier Biggest Difference Method	
11	Bernier Biggest Proportion Method	
F1 Average fitness	1	DSGA (R $\Delta$ , S-)
	2	Kao and Zahara Genetic Algorithm and Particle Swarm
	3	DSGA (R-, S-)
	3	DSGA (R $\Delta$ , FS)
	4	DSGA (R+, S-)
	5	DSGA (R-, FS)
	6	SCGA
	7	Goldberg and Richardson's Fitness Sharing
	8	Bernier Biggest Difference Method
	9	Bernier Biggest Proportion Method
10	DSGA (R+, FS)	

Table 26. Ranking of Algorithms Continued

<b>Benchmark Criteria</b>	<b>Rank</b>	<b>Ranked Algorithm</b>
F2 Proportion of Peaks	1	Bernier Biggest Proportion Method
	2	Goldberg and Richardson's Fitness Sharing
	3	DSGA (R+, FS)
	4	DSGA (R-, S-)
	5	Bernier Biggest Difference Method
	6	DSGA (R $\Delta$ , S-)
	6	DSGA (R-, FS)
	6	DSGA (R $\Delta$ , FS)
	7	DSGA (R+, S-)
8	SCGA	
9	Kao and Zahara Genetic Algorithm and Particle Swarm	
F2 Proportion of points outside of peaks	1	Kao and Zahara Genetic Algorithm and Particle Swarm
	2	SCGA
	3	DSGA (R+, S-)
	4	DSGA (R $\Delta$ , S-)
	5	DSGA (R-, S-)
	6	DSGA (R $\Delta$ , FS)
	7	DSGA (R-, FS)
	8	DSGA (R+, FS)
	9	Bernier Biggest Difference Method
	10	Goldberg and Richardson's Fitness Sharing
	11	Bernier Biggest Proportion Method
F2 Average fitness	1	DSGA (R $\Delta$ , S-)
	2	DSGA (R+, S-)
	3	DSGA (R-, S-)
	4	DSGA (R $\Delta$ , FS)
	5	Kao and Zahara Genetic Algorithm and Particle Swarm
	6	DSGA (R-, FS)
	7	SCGA
	8	Bernier Biggest Difference Method
	9	Goldberg and Richardson's Fitness Sharing
	10	Bernier Biggest Proportion Method
	11	DSGA (R+, FS)

Table 26. Ranking of Algorithms Continued

<b>Benchmark Criteria</b>	<b>Rank</b>	<b>Ranked Algorithm</b>
F3 Proportion of Peaks	1	Bernier Biggest Difference Method
	2	DSGA (R-, FS)
	3	DSGA (R+, FS)
	4	Bernier Biggest Proportion Method
	5	DSGA (R-, S-)
	6	DSGA (RΔ, FS)
	7	Goldberg and Richardson's Fitness Sharing
	8	DSGA (RΔ, S-)
	9	DSGA (R+, S-)
	10	SCGA
	11	Kao and Zahara Genetic Algorithm and Particle Swarm
F3 Proportion of points outside of peaks	1	Kao and Zahara Genetic Algorithm and Particle Swarm
	2	DSGA (R+, S-)
	3	DSGA(RΔ, S-)
	4	DSGA (R-, S-)
	5	SCGA
	6	DSGA (RΔ, FS)
	7	DSGA (R-, FS)
	8	DSGA (R+, FS)
	9	Bernier Biggest Difference Method
	10	Goldberg and Richardson's Fitness Sharing
	11	Bernier Biggest Proportion Method
F3 Average fitness	1	DSGA (RΔ, S-)
	2	Kao and Zahara Genetic Algorithm and Particle Swarm
	3	DSGA (R-, S-)
	4	DSGA (RΔ, FS)
	5	DSGA (R-, FS)
	6	SCGA
	7	DSGA (R+, S-)
	8	Bernier Biggest Difference Method
	9	Goldberg and Richardson's Fitness Sharing
	10	Bernier Biggest Proportion Method
	11	DSGA (R+, FS)

Table 26. Ranking of Algorithms Continued

<b>Benchmark Criteria</b>	<b>Rank</b>	<b>Ranked Algorithm</b>
F4 Proportion of Peaks	1	Bernier Biggest Proportion Method
	2	DSGA (R+, FS)
	3	DSGA (R-, FS)
	4	DSGA (R $\Delta$ , FS)
	5	Goldberg and Richardson's Fitness Sharing
	6	Bernier Biggest Difference Method
	7	DSGA (R+, S-)
	7	DSGA (R-, S-)
	8	DSGA (R $\Delta$ , S-)
	9	SCGA
	10	Kao and Zahara Genetic Algorithm and Particle Swarm
F4 Proportion of points outside of peaks	1	Kao and Zahara Genetic Algorithm and Particle Swarm
	2	DSGA (R $\Delta$ , S-)
	3	DSGA (R+, S-)
	4	SCGA
	5	DSGA (R-, S-)
	6	DSGA (R-, FS)
	7	DSGA (R $\Delta$ , FS)
	8	DSGA (R+, FS)
	9	Bernier Biggest Difference Method
	10	Goldberg and Richardson's Fitness Sharing
	11	Bernier Biggest Proportion Method
F4 Average fitness	1	DSGA (R+, S-)
	2	Kao and Zahara Genetic Algorithm and Particle Swarm
	3	DSGA (R $\Delta$ , S-)
	4	DSGA (R-, S-)
	5	DSGA (R-, FS)
	6	DSGA – Dynamic Radius; Fitness Sharing
	7	SCGA
	8	Goldberg and Richardson's Fitness Sharing
	9	Bernier Biggest Difference Method
	10	Bernier Biggest Proportion Method
	11	DSGA (R+, FS)

Table 26. Ranking of Algorithms Continued

<b>Benchmark Criteria</b>	<b>Rank</b>	<b>Ranked Algorithm</b>
F5 Proportion of Peaks	1	Bernier Biggest Proportion Method
	2	Goldberg and Richardson's Fitness Sharing
	3	Bernier Biggest Difference Method
	4	Kao and Zahara Genetic Algorithm and Particle Swarm
	4	DSGA (R+, S-)
	4	DSGA (RΔ, FS)
	5	SCGA
	5	DSGA (R-, S-)
	5	DSGA (RΔ, S-)
	5	DSGA (R+, FS)
F5 Proportion of points outside of peaks	5	DSGA (R-, FS)
	1	SCGA
	2	Kao and Zahara Genetic Algorithm and Particle Swarm
	3	DSGA (R-, S-)
	4	DSGA (R+, S-)
	5	DSGA (RΔ, S-)
	6	DSGA (R-, FS)
	7	DSGA (R+, FS)
	8	DSGA (RΔ, FS)
	9	Bernier Biggest Difference Method
	10	Bernier Biggest Proportion Method
11	Goldberg and Richardson's Fitness Sharing	
F5 Average fitness	1	SCGA
	2	DSGA (RΔ, S-)
	3	DSGA (R+, S-)
	4	DSGA (R-, FS)
	5	DSGA (R-, S-)
	6	DSGA (RΔ, FS)
	7	Kao and Zahara Genetic Algorithm and Particle Swarm
	8	Bernier Biggest Difference Method
	9	Bernier Biggest Proportion Method
	10	Goldberg and Richardson's Fitness Sharing
	11	DSGA (R+, FS)

Table 26. Ranking of Algorithms Continued

<b>Benchmark Criteria</b>	<b>Rank</b>	<b>Ranked Algorithm</b>
F6 Proportion of Peaks	1	Bernier Biggest Proportion Method
	2	Goldberg and Richardson's Fitness Sharing
	3	Bernier Biggest Difference Method
	4	DSGA (R+, FS)
	5	DSGA (RΔ, FS)
	5	DSGA (RΔ, S-)
	6	DSGA (R-, S-)
	7	DSGA (R-, FS)
	8	DSGA (R+, S-)
	9	SCGA
	10	Kao and Zahara Genetic Algorithm and Particle Swarm
F6 Proportion of points outside of peaks	1	SCGA
	2	Kao and Zahara Genetic Algorithm and Particle Swarm
	3	DSGA (R+, S-)
	4	DSGA (RΔ, S-)
	5	DSGA (R-, S-)
	6	DSGA (RΔ, FS)
	7	DSGA (R-, FS)
	8	DSGA (R+, FS)
	9	Bernier Biggest Difference Method
	10	Bernier Biggest Proportion Method
	11	Goldberg and Richardson's Fitness Sharing
F6 Average fitness	1	DSGA (R-, FS)
	2	DSGA (R-, S-)
	3	DSGA (R+, S-)
	4	DSGA (RΔ, S-)
	5	DSGA (RΔ, FS)
	6	Kao and Zahara Genetic Algorithm and Particle Swarm
	7	Bernier Biggest Difference Method
	8	Goldberg and Richardson's Fitness Sharing
	9	Bernier Biggest Proportion Method
	10	DSGA (R+, FS)
	11	SCGA

Table 26. Ranking of Algorithms Continued

<b>Benchmark Criteria</b>	<b>Rank</b>	<b>Ranked Algorithm</b>
F7 F(x, y) of Best Individual for Each Niche	1	Zhang, Shang, Gao and Dong hK1 Triangulation Algorithm
	2	DSGA (R $\Delta$ , FS)
	3	DSGA (R+, S-)
	4	DSGA (R-, S-)
	5	DSGA (R-, FS)
	6	DSGA (R+, FS)
	7	DSGA (R $\Delta$ , S-)
	8	SCGA
	9	Goldberg and Richardson's Fitness Sharing
F7 Proportion of Peaks	1	Zhang, Shang, Gao and Dong hK1 Triangulation Algorithm
	1	DSGA (R+, S-)
	2	DSGA (R+, FS)
	3	SCGA
	4	Goldberg and Richardson's Fitness Sharing
	5	DSGA (R $\Delta$ , S-)
	6	DSGA (R $\Delta$ , FS)
	7	DSGA (R-, FS)
	8	Kao and Zahara Genetic Algorithm and Particle Swarm
F8 Proportion of Peaks	8	DSGA (R-, S-)
	1	DSGA (R-, S-)
	2	DSGA (R+, S-)
	3	DSGA (R $\Delta$ , FS)
	4	DSGA (R+, FS)
	4	DSGA (R-, FS)
	5	DSGA (R $\Delta$ , S-)
	6	SCGA
	7	Goldberg and Richardson's Fitness Sharing
8	Kao and Zahara Genetic Algorithm and Particle Swarm	
F8 Proportion of points outside of peaks	1	Kao and Zahara Genetic Algorithm and Particle Swarm
	2	SCGA
	3	DSGA (R+, S-)
	4	DSGA (R $\Delta$ , S-)
	5	DSGA (R-, S-)
	6	DSGA (R-, FS)
	7	DSGA (R+, FS)
	8	DSGA (R $\Delta$ , FS)
	9	Goldberg and Richardson's Fitness Sharing

Table 26. Ranking of Algorithms Continued

<b>Benchmark Criteria</b>	<b>Rank</b>	<b>Ranked Algorithm</b>
F8 Average fitness	1	Kao and Zahara Genetic Algorithm and Particle Swarm
	2	SCGA
	3	DSGA (R $\Delta$ , S-)
	4	DSGA (R+, S-)
	5	DSGA (R $\Delta$ , S-)
	6	DSGA (R-, FS)
	7	DSGA (R $\Delta$ , FS)
	8	DSGA (R+, FS)
	9	Goldberg and Richardson's Fitness Sharing

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