Automated Design of Hyper-Heuristics Components Applied to the PSP Problem

Vidal D. Fontoura Federal University of Parana (DInf-UFPR), Curitiba - PR, Brazil Email: vdfontoura@inf.ufpr.br Aurora T. R. Pozo Federal University of Parana (DInf-UFPR), Curitiba - PR, Brazil Email: aurora@inf.ufpr.br Roberto Santana Hermida University of the Basque Country Barrio Sarriena s/n 48940 Leioa Bizkaia Email: roberto.santana@ehu.es

Abstract—The Protein Structure Prediction (PSP) problem is one the modern most challenging problems from science. Hence, many heuristics strategies have been applied in order to find protein structures that minimize its free energy. However, these strategies have difficulties on finding the optimal solutions to the longer sequences of amino-acids, due to the complexity of the problem and the huge amount of local optima. The hyperheuristics framework are usually useful in this kind of context since they try to combine different heuristics strengths into a single framework. However, there is lack of works whose aim the automated design of hyper-heuristics components. This paper proposes the GEHyPSP which aims to the generation, through grammatical evolution, of selection mechanisms and acceptance criteria for a hyper-heuristic framework applied to PSP problem.

I. INTRODUCTION

Proteins plays a fundamental role in nature, they are involved in many important functions for the living cells. Proteins are structures, composed by amino-acids that guarantee the correct operation of wide range of biological entities. These structures are the result of the so-called protein folding process, in which an unfolded chain of amino-acids will be transformed into its final/native structure.

The protein structure prediction (PSP) has a wide range of biotechnology and medical applications. For instance: synthesis of new proteins and folds [29], structure based synthesis of new drugs [8], refinement of theoretical models obtained by comparative modeling [21], [16], and obtaining experimental structures from incomplete nuclear magnetic resonance data [25].

The determination of the native structure of a protein is a hard task even for the modern super computers. The difficulty is due to the huge search space to test all possible configurations that a given protein can adopt. Many models to represent the proteins structures have been proposed and can be used to simulate the folding process. Although there exist extremely detailed models, these representations are computational very expensive. Hence, many authors [6], [12], [18], [28], [24], [7], [11] have used simplified models to represent the protein structures. A useful model for this purpose is the Hydrophobic-Polar (HP) model presented by Lau and Dill [17]. This model reduces the amino-acids that composes the proteins into just two types hydrophobic (H) or polar (P). Two-dimensional (2D) or three-dimensional (3D) grids can be used to represent the

conformations that a protein can adopt. Each conformation has an associated energy [28].

Many heuristics strategies were developed to find conformations of minimum energy of the HP model. Among the approaches that have been applied are genetic algorithms [28], ant colony optimization [26], [27], estimation distribution algorithms [?], and others [9].

Although there are different strategies already proposed for the PSP, all of them face difficulties to reach the optimal configurations when the length of amino-acids sequences increases. This fact motivates the study here presented. In this kind of context that hyper-heuristics are usually a good option in order to better explore a complex search space. In general the hyper-heuristics frameworks propose a strategy to select, from a pool, the heuristic which is more appropriate for a given moment. Combining the strengths and weakness of a set of heuristics. Also the hyper-heuristics frameworks define a move acceptance criterion, which is responsible to decide on whether accept or reject when worst solutions are generated by the heuristics. Theses strategies are high level heuristics from a hyper-heuristic framework.

Usually the high level heuristics are human-designed approaches although a study [23] propose the automate design of those components. This study presented promising results and also inspired the idea of automating the design of high level heuristics to a hyper-heuristic framework to the PSP problem. Since the PSP with the HP-2D model presents a very complex landscape automating the design of hyper-heuristic framework could save time, money and human effort when design a hyper-heuristic strategy to the PSP. Another motivation of the present study is that the hyper-heuristics frameworks have been applied to other domains problems and achieving good results in the bin packing, personnel scheduling, flowshop, TSP, MAXSAT and VRP problems [23].

Burke et al. [5] recently defined hyper-heuristics as "an automated methodology for selecting or generating heuristics to solve hard computational search problems". Over the years, these methodologies have demonstrated success in solving a wide range of real world problems. However, there is not an in-depth investigation of the suitability of this approach to solve the PSP problem.

We introduce the GEHyPSP: an automated mechanism for generation of high level heuristics to a hyper-heuristic framework to solve the PSP problem. The high level heuristic includes two main components: a selection mechanism and an acceptance criterion. Both components are generated using grammatical evolution (GE) [22], which is a kind of genetic programming (GP) [13]. A set of experiments was designed to evaluate the approach using eleven PSP instances. The results are compared with the previous works of the literature.

This paper is organized as follows: in the next section we briefly introduce the Protein Structure Prediction problem and a review of the related works is also presented. Section III presents a background of hyper-heuristics, GP and the GE algorithm. Thereafter, in Section IV, the GEHyPSP is introduced. In Section V, the experimental benchmark and numerical results of the conducted experiments are presented. Finally, in Section VI, the conclusions of the research are given, and further work is discussed.

II. PROTEIN STRUCTURE PREDICTION

Proteins are macromolecules composed by an alphabet of twenty different amino acids, also referred to as residues. An amino acid is formed by a peptide backbone and a distinctive side chain group. The peptide bond is defined by an amino group and a carboxyl group connected to an alpha carbon to which a hydrogen and side chain group are attached.

Under specific conditions, the protein sequence folds into a unique native 3-D structure. Each possible protein fold has an associated energy value. The thermodynamic hypothesis states that the native structure of a protein is the one for which the free energy achieves the global minimum. Based on this hypothesis, many methods [6], [12], [14], [18], [28] that search for the protein native structure define an approximation of the protein energy and use optimization algorithms that look for the protein fold that minimize the energy. These approaches mainly differ in the type of energy approximation employed and in the characteristics of the protein modeling.

A. The HP Model

The protein structures can be very complex. Detailed representations of proteins exist and can be used to model the protein folding, these representations are computationally very costly. Having this in mind, Lau and Dill [17] created a model called *Hydrophobic-Hydrophilic* Model (HP Model), to represent the proteins using simplifications. The model can be used either to represent proteins in a 2D space or 3D space.

The HP model considers two types of residues: hydrophobic (H) residues and hydrophilic or polar (P) residues. A protein is considered a sequence of these two types of residues, which are located in regular lattice models forming self-avoided paths. Given a pair of residues, they are considered neighbors if they are adjacent either in the chain (connected neighbors) or in the lattice but not connected in the chain (topological neighbors).

For the HP model, an energy function that measures the interaction between topological neighbor residues is defined as $\epsilon_{HH}=-1$ and $\epsilon_{HP}=\epsilon_{PP}=0$. The HP problem consists of finding the solution that minimizes the total energy. In the

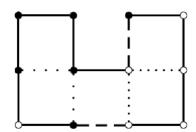


Fig. 1. One possible configuration of sequence PPPHPHHHHHPH in the HP model. There is two HH (represented by a dotted line with wide spaces), one HP (represented by a dashed line) and two PP (represented by dotted lines) contacts.

linear representation of the sequence, hydrophobic residues are represented with the letter H and polar ones, with P. In the graphical representation, hydrophobic proteins are represented by black beads and polar proteins, by white beads. Figure 1 shows the graphical representation of a possible configuration for the sequence PPPHPHHHHPH in a 2D space. The energy that the HP model associates with this configuration is -2.

Among many works that explores the PSP problem, here are some examples of the approaches that have been used to solve it.

A multimeme algorithm (MMA) presented in [14] is an EA combined with a group of local search methods. For each individual in the population, the MMA, selects a local search method that is more suitable with the individual. Used first to find solutions for the functional model protein, the strategy was later improved with fuzzy-logic-based local searches, leading the algorithm to achieve improved results in the PSP problem.

In [12], the author uses pruned-enriched Rosenbluth method (PERM), also known as chain growth algorithm, that is based on growing the sequence conformation by adding particle by particle, aiming to increase good configurations and eliminating bad ones.

The ant colony optimization (ACO) was applied to the PSP problem using the HP-2D model in [26], [27]. This strategy, uses artificial ants to build conformations for a given HP instance. A local search method is then applied to further improve the solutions and also maintain the quality of the solutions.

The work of [24] applies EDAs as an efficient evolutionary algorithm that learns and exploits the search space in the form of probabilistic dependencies. New ideas were introduced in [24] for the application of EDAs to 2D and 3D simplified protein folding problem. The obtained results showed that EDAs can achieve superior solutions compared with other well-known population based optimization algorithms.

The present paper proposes the use of a grammatical evolution GE to generate high level heuristics for a hyper heuristic framework that will be applied to the PSP problem.

III. BACKGROUND

This section will present the background context in order to provide the readers with the necessary information, for a good comprehension of the concepts and techniques used in this paper.

A. Hyper-Heuristics

Mainly, a generic hyper-heuristic framework is composed of two main components known as high-level and low-level heuristics. Figure III-A shows a general scheme of hyper-heuristic frameworks. The high and low levels are separated by a domain barrier which means that the low-level component is problem dependent while the high-level component does not require any knowledge of the problem domain. Ideally, to apply a hyper-heuristic framework to a different problem domain: it is possible to do it only by replacing the low-level component, no changes should be required at the high-level. It is responsibility from the high-level component to manage the selection or generation of which heuristic should be applied at each decision point. The low-level component corresponds to a pool of heuristics or heuristic components [23].

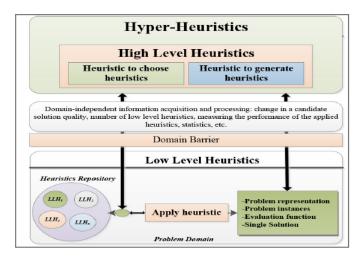


Fig. 2. General scheme of hyper-heuristic framework.

Based on the nature of the search space, hyper-heuristics can be classified [5] as either selecting or generating heuristics for the underlying problem. Next will be analyzed both kinds of high-level heuristics.

• Selecting Heuristics: The majority of hyper-heuristic frameworks define high level heuristics to select low level heuristics. In general, these frameworks define a selection mechanism and also an acceptance criterion. These strategies use information from the statistical history of low level heuristics applications. For instance, a greedy selection considers only the number of improvements that a low level heuristic obtained [3]. In contrast, the choice function [3] is a reward-based strategy, which considers both the number of improvements and the elapsed time since the last application.

 Generating Heuristics: In this case, the hyper-heuristic framework starts with a set subcomponents from low level heuristics and has the goal of building new low level heuristics with it. GP is reported as a good strategy to combine and generate new heuristics for the SAT, scheduling, and bin-packing problems [23].

A study presented by Sabar et al. [23], proposes an on-line approach to generate selection mechanisms and acceptance criteria (high level heuristics) to a hyper-heuristic framework, using GEP [9] (gene expression programming a variant of GP). The experiments presented by Sabar et al. [23], using the 6 problem domains provided by the HyFlex hyper-heuristic framework [20], showed impressive results in comparison with other human-designed hyper-heuristic strategies from the state-of-art. This encouraging result inspired the present study which has the objective of off-line generate selection mechanisms and acceptance criteria, through a grammatical evolution process, for a hyper-heuristic framework to solve PSP problem.

B. Genetic Programming (GP)

GP is a sub-field from the program synthesis which uses ideas from the evolution theory to produce programs [3].

Grammatical Evolution (GE) is a relatively new technique from the evolutionary computing, introduced by Ryan el al. [22] and it is a type of GP where programs are evolved using a genetic algorithm. The chromossome encode production rules of the grammar. GE uses a mapping mechanism between the genotype (coded individuals by integer vectors) and phenotype (generated programs to solve a problem). The BNF notation is used to represent a grammar from a language in form of production rules. A BNF grammar consists in a set of terminals, which are items that are allowed to the language, for instance: +, -, *, /, etc and non-terminals, which can be expanded into one or more terminals. The Grammar can be expressed as a tuple N, T, P, where N is a set of non-terminals presented in the Grammar 1 as $\langle expr \rangle$, $\langle op \rangle$, $\langle pre - op \rangle$ and $\langle var \rangle$, T a set of terminals and is presented in the Grammar 1 as +, -, /, *, Sin, Cos and Tan. Finally, P is the set of production rules that maps the elements N to T and is presented in Table I.

 $\label{thm:condition} \textbf{TABLE} \ \textbf{I}$ Production rules and the number of choices allowed

Production Rules	Number of choices
$\langle expr \rangle$	4
$\langle op \rangle$	4
$\langle pre - op \rangle$	3
$\langle var \rangle$	1

Ryan et al. [22] proposed the use of genetic algorithm (GA) to control which choices should be made, in this sense allowing the GA to select which production rules should be utilized. An individual (chromosome) consists in variable-length integer vector and it is called genotype. Algorithm 1 presents the pseudocode of a grammatical evolution program. Lines 3 to 6 represents the initialization of the algorithm: a randomly generated population is created and then mapped

$$\langle expr \rangle \qquad ::= \langle expr \rangle \langle op \rangle \langle expr \rangle \qquad (0)$$

$$| (\langle expr \rangle \langle op \rangle \langle expr \rangle) \qquad (1)$$

$$| \langle pre-op \rangle (\langle expr \rangle) \qquad (2)$$

$$| \langle var \rangle \qquad (3)$$

$$\langle op \rangle$$
 ::= + (0) (1) (2) (3)

$$\langle pre\text{-}op \rangle$$
 ::= Sin (0)
 | Cos (1)
 | Tan (2)

$$\langle var \rangle$$
 ::= X (0)

Grammar 1. Sample grammar to demonstrate how to decode integer vectors in computer programs

to programs using the GF (Grammar File). Those programs are evaluated and a fitness value is assigned for each individual in the population. Line 7 represents the main loop of the algorithm and the evolution process occurs within this loop. Line 8,9,10,11 and 12 occurs: the parent selection; the crossover between the parents, the application of the prune and duplication operators and finally, the mutation operator is applied. Line 13 represents the mapping between the offspring (integer vectors) to programs using the GF. Line 14 and 15 represents the fitness assignment to the individuals from the population. Line 16 represents the population replacement with the fittest individuals. Finally, line 18 return the program with the best fitness found.

Algorithm 1: Pseudo code from the Grammatical Evolution

```
1 Input: GF - Grammar File
2 begin
        population \leftarrow Create population;
3
        programs \leftarrow \text{Maps the } population \text{ to programs using } GF;
4
        Execute the program;
        Assign fitness value to the solutions of population according with
        the output obtained by the respective decoded program;
7
        while Stop condition not reached do
            parents \leftarrow Select individuals for crossover;
             offspring \leftarrow Crossover(parents);
             Apply the Prune operator to the offspring;
10
             Apply the Duplicate operator to the offspring;
11
12
             Apply the Mutation operator to the offspring;
13
            programs \leftarrow \text{Maps } of fspring \text{ to programs using } GF;
             Execute programs;
14
15
             Assign fitness value to solutions of fspring according with
             the output obtained by the respective decoded program;
             population \leftarrow Replacement;
16
17
        end
        return Best program from the population;
18
19 end
```

For the sake of comprehension the chromosome mapping process will be illustrated using the Grammar 1. The Algorithm 2 presents the general template of the generated programs. The expression $\langle expr \rangle$ presented on line 2 is replaced by mathematical expressions coded by the chromosomes (integer vectors).

```
Algorithm 2: General template for the generated algorithms
```

```
1 float symb(float x)
2 a = \langle expr \rangle;
3 return a;
```

Now suppose the following integer vector:

This vector will be used to decode the chromosome (genotype) into a piece of code (phenotype) using the Grammar 1. Table I summarizes the number of choices associated within each production rule from Grammar 1.

There are 4 options of production rules that can be selected for the expression $\langle expr \rangle$. In order to select which option, the first value from the vector should be used. The value is 220 and its remainder divided by 4 (four options that can be selected for the expression $\langle expr \rangle$) results in 0, which means that the first option should be selected $\langle expr \rangle \langle op \rangle \langle expr \rangle$. Note that the first expression is also $\langle expr \rangle$ and following the same logic we should expand using the next value from the integer vector and applies its remainder by the division of the number of options. The remainder of 203 % 4 = 3, which indicates that we should select the fourth option: $\langle var \rangle$ which is a terminal. The $\langle var \rangle$ has just one option associated with it and is the value X. Placing this selection in the original expression we have $X\langle op \rangle \langle expr \rangle$.

Next it is necessary to decode the non-terminal expression $\langle op \rangle$. The next value from the integer vector is 17 and again we have 4 options (+|-|/|*). The result is equal to 1, and indicates to select the: -. Re-writing the expression we got: $X - \langle expr \rangle$. This process should continue until all the non-terminals be expanded to terminals. In this example the resulting expression (phenotype) is: X - Sin(X). Note that not all of the genes were necessary to obtain the phenotype. In these cases, the genes that were not used are discarded. Moreover, the opposite case can occur: if a chromosome does not contain the necessary number of genes to map to a program. In this scenario the strategy is re-utilize the genes starting from the first one.

IV. GEHYPSP

In this section will be presented the GEHyPSP: an offline grammatical evolution application for generating high level heuristics of hyper-heuristics framework to the Protein Structure Prediction problem. The approach presented next is based on the study by Sabar et al. [23]

The high level heuristics are composed by a selection mechanism and an acceptance criterion. These high level heuristics use information related with the history of low level heuristics applications. Data about the improvements obtained by the low level heuristics, number of times since the last application of a low level heuristic and the fitness difference between the current solution and the generated one are examples of information used by selection mechanisms and acceptance criteria. Furthermore, two terminal sets (one for the selection mechanisms and another for the acceptance criteria) were defined according with the information that can be extracted during the search progress. The selection terminals are presented next:

- RC (Reward Credit): The reward that a given heuristic should receive based on its performance. The improvement is calculated, for the i_{th} heuristic, using M(i) = (|f1-f2|/f1)*100 if f2 < f1, where f1 is the current fitness and f2 is the fitness of the solution generated by the i_{th} heuristic.
- C_{best}: Number of times that the i_{th} heuristic updated the best known solution. This terminal is useful to systematically improve the current local minimum.
- $C_{current}$: Number of times that the i_{th} heuristic updated the current solution. This terminal is useful to keep the search near to the current solution.
- C_{accept}: Number of times that the generated solution by the i_{th} heuristic was accepted by the acceptance criterion. This terminal favors heuristics that can escape from local minimum.
- C_{ava} : The average of previous improvements made by the i_{th} heuristic during the search progress. This terminal favors heuristics that made big improvements in average.
- C_r : Number of times that the i_{th} heuristic was selected to be applied.

A specific terminal set for generating acceptance criteria was also defined.

- Delta: The difference between the quality of the current solution and the generated solution.
- PF: The quality of the previous solution.
- CF: The quality of the current solution.
- CI: Current iteration.
- TI: Total number of iterations.

Using these statistics a as terminals and a function set containing the following arithmetic operations: addition, subtraction, multiplication and division, a grammar was designed to support the generation of the high level heuristics. The designed grammar to generate selection mechanisms and acceptance criteria is presented in the Grammar 2.

For the initialization of the terminal set data: all heuristics were executed once and the data was calculated for each terminal. Every consecutive iteration will update the terminal set data and this information will be used during the search progress.

A. Fitness Function

In order to evaluate the generated individuals during the search progress, a fitness function was designed. The fitness

```
\langle hh\text{-selection}\rangle ::= \langle selection\text{-mechanism}\rangle
                          (acceptance-criterion)
\langle selection-mechanism \rangle :== \langle selection-terminal \rangle
                          ⟨selection-mechanism⟩
                                                              ⟨math-function⟩
                          (selection-mechanism)
                          (\langle selection-mechanism \rangle
                                                              ⟨math-function⟩
                          ⟨selection-mechanism⟩)
\( \selection-terminal \rangle :== RC \ | Cbest \ | Ccurrent \ | Caccept \ |
                          Cava | Cr
\langle math\text{-}function \rangle :== + | - | * | \%
\langle acceptance\text{-}criterion \rangle ::== \langle acceptance\text{-}terminal \rangle
                          (acceptance-criterion)
                                                              (math-function)
                          (acceptance-criterion)
                          (\langle acceptance-criterion \rangle
                                                              ⟨math-function⟩
                          ⟨acceptance-criterion⟩)
\langle acceptance\text{-}terminal \rangle :== PF \mid CF \mid CI \mid TI
```

Grammar 2. Designed grammar to generate high level heuristics

function consisted on running the generated high level heuristics, within a hyper-heuristic framework, against three random selected instances from a set of eleven instances. Each run will be executed for one minute and will return the best HP solution found. The fitness value associated with the returned solution is then normalized between 0 and 1. The fitness of an individual, of the GE, it is the sum of the three outcomes from each execution of the three randomly selected instances. Hence, the best possible fitness value is 3 and the worst is 0. The motivation behind executing the high level heuristic (individual) against three HP instances is that executing with only one might not be sufficient to train a high level heuristic to obtain good results for various instances of the HP model.

B. Stopping Criterion

To stop the GE process a maximum number of evaluations was setup to 60000. This value was defined based on previous work [22] where the grammatical evolution general process was introduced for the first time.

C. Low Level Heuristics

1) Representation of the problem: There are different ways of representing a protein configuration (or protein structure) within the HP-2d model. According to Krasnogor et al. [15] the relative representation has a better potential to achieve superior results. The relative representation stats that: the movements in the grid for each amino-acid are represented always based on the previous, that is why it is named relative. There are 3 possible movements within the HP-2d model: forward (F), left (L) and right (R). Hence, the following integer codification was used $F \to 0$, $L \to 1$ and $R \to 2$. Thereby, the allowed alphabet can be represented as $\{0,1,2\}$.

- 2) Low Level Heuristics Set: The low level heuristics set was selected from previous studies [1], [7], [6], [10] that explore the PSP. It consists of the following operators:
 - Two Points Crossover (2X): This operator selects, randomly, two crossing points splitting the individuals in 3 pieces. The genes between the selected positions are exchanged between the parents in order to generate the offspring. [1].
 - Multiple Points Crossover (MPX): Similar with 2X although with c points of crossing. We use c = int(n*0.1), ..., where n is the sequence length. The MPX is useful to promote a wide structural diversity [23].
 - Segment Mutation (SMUT): It changes a random number (5 to 7) of consecutive genes to distinct values.
 This heuristic introduces large changes in the protein conformation, and it has a great probability of creating collisions.
 - Exhaustive Search Mutation (EMUT): This heuristic selects a random gene and changes to all possible directions and it will keep the change that achieves the best improvement. This operator has a trade-off because it demands four fitness evaluations instead of just one. However, this heuristic has great potential of improving the input solution.
 - Local Move Operator (LM): This heuristic exchanges directions between two consecutive random selected genes.
 There are some conditions in order to execute this heuristic, for instance, the new directions can not create redundant movements.
 - Loop Move Operator (LPM): Similar with LM, this heuristic exchanges directions between two genes that are five genes of distance between each other.
 - Opposite Mutation (OM): This heuristic exchanges directions, for the opposite, within a sequence of genes between two genes (i, j) randomly selected.
- 3) Backtrack Repair: Since the low level heuristics have a great probability of generating solutions with collisions [1]. Whenever it is possible, to repair a solution using a backtrack strategy the solution is maintained otherwise it is penalized.
- 4) Memory Mechanism: Sabar et al. [23] suggested that the use of a memory of solutions to the problem would be more effective than relying on a single solution and may restrict the ability of dealing with large and heavily constrained search spaces, as it is widely known that single solution based methods are not well suited to cope with large search spaces and heavily constrained problems [2].

V. EXPERIMENTS

In this section, we will present and discuss the conducted experiments in order to evaluate the GEHyPSP proposed by this paper. Three groups of experiments were designed/executed and will be described next. The first group was only concerned on generating selection mechanisms. The second group was designed only to generate acceptance criteria. The third group was developed to generate both selection and acceptance mechanisms. All experiments were executed 30 times

TABLE II
INSTANCES AND THE RESPECTIVE SIZE OF EACH ONE

Inst	1	2	3	4	5	6	7	8	9	10	11
Size	20	24	25	36	48	50	60	64	85	100	100

because of the stochastic behavior of the GEHyPSP. The results obtained by the groups were compared with each other. Also a comparison with a state-of-art hyper-heuristic will be presented. The results were compared with the Generic Intelligent Hyper-heuristic (GIHH) presented by Misir el al [19].

In the first group of experiments (GEHyPSP-1) only selection mechanisms are generated. The acceptance criterion was fixed with a "better or equal" acceptance [4]. The goal of this group is to evaluate the ability of our approachx to generate selection mechanisms using a fixed acceptance criterion.

The second group of experiments, GEHyPSP-2, consisted on generating only acceptance criteria. The selection mechanism was fixed using the best selection mechanism found in the first group. Consequently, this group of experiments depends on the output from the first group. The goal of this experiment was to evaluate the generation of acceptance criteria separately from the selection mechanism using a fixed one

The third group of experiments GEHyPSP-3 was designed to generate both selection mechanisms and acceptance criteria. The goal of this group was evaluate the ability of generating selecting mechanisms along with acceptance criteria. Differently from GEHyPSP-2, this group of experiment does not depend on any output from previous experiments since it generates both mechanisms without fixing any component.

For each group of experiments eleven instances (amino-acid sequences from the HP model) selected from the previous studies with the PSP problem [6], [12], [18], [28], [24], [7], [11]. The sizes (number of residues) of the HP instances

used in the experiments are presented in table II. For the sake of space, the sequences are not reproduced here. They can be obtained from [24]. Also, for each group of experiments the training phase consisted on executing the GE process with three ramdonly selected instances from the eleven available instances. In the validation phase the best generated selection mechanisms and acceptance criteria were executed together against all eleven instances.

A. Results from GEHyPSP-1

The best individual found in the GEHyPSP-1 was the following selection mechanism: RC*Ccurrent*Cava-Cr and it was executed 30 times with a time limit of 10 minutes against the eleven instances. Table III presents the average, standard deviation, minimum and maximum of 30 executions of the best individual found in the GEHyPSP-1 experiment. The last row denoted with $O(x^*)$ is the best known value for each sequence. It is possible to note analyzing the last two lines that only for the smaller instances the generated selection mechanism achieved good results. For the sequences 8, 9, 10

 $\label{thm:conditional} TABLE~III \\ Results~from~the~best~individual~found~in~GEHyPSP-1 \\$

Inst	1	2	3	4	5	6	7	8	9	10	11
Avg	8.1	7.6	6.7	11.9	17.4	16	30	28.3	40.1	35.6	35.9
St Dv	0.3	0.5	0.5	0.7	1	1.4	1.7	2	2.7	2.1	2.9
Min	8	7	5	11	15	13	25	23	34	32	27
Max	9	9	7	13	19	20	33	32	46	40	41
$O(x^*)$	9	9	8	14	23	2.1	36	42.	53	48	50

 $\label{thm:conditional} TABLE\ IV$ Results from the best individual found in GEHyPSP-2

	Inst	1	2	3	4	5	6	7	8	9	10	11
_	Avg	8	7.6	6.7	10.1	14.8	14.7	27	25.7	38.3	32.8	30.
_	St Dv	0.6	0.4	0.6	0.7	1.5	1.4	2.0	2.5	3.3	3.7	3.4
_	Min	7	7	5	8	12	12	23	22	31	26	24
	Max	9	8	8	11	17	18	31	31	44	40	37
	O(x*)	9	9	8	14	23	21	36	42	53	48	50

and 11 the results obtained are very far from the best known results.

B. Results from GEHyPSP-2

The best individual found in the GEHyPSP-2 was the following acceptance criterion: ((TI/Delta)/((Delta *((TI/Delta)/CI)*Delta/Delta*TI)-CI) and combined with the best individual from the GEHyPSP-1 the hyperheuristic framework was executed 30 times for each one of the eleven instances with a time limit of 10 minutes. Table IV presents the results for the best generated acceptance criterion in GEHyPSP-2 and also using the best selection mechanism generated in GEHyPSP-1. Again, looking to the two last lines it is possible to notice that the hyper-heuristic framework, using the best generated selection mechanisms and acceptance criterion, was not able to reach the best known results for the larger sequences. And comparing Tables III and IV it is possible to visualize that there is a slight difference between the results, favoring the generated selection mechanism using the "better or equal" acceptance criterion.

C. Results from GEHyPSP-3

The best individual generated by the experiment GEHyPSP-3 was both a selection mechanism and an acceptance criterion. And they are presented below:

Selection Mechanism:
$$(((((Caccept/RC) * Cr/Caccept)/RC * Cr)/Caccept/RC) * Cr)/Caccept$$

Acceptance Criterion: $((((CI/PF) * Delta/CI)/PF * Delta)/CI/PF) * Delta)/CI$

Table V presents the results, of the best individual generated by the experiment GEHyPSP-3, of 30 executions with a time limit of 10 minutes. Once again the values presented did not achieve good results when analyzing the most difficult instances. Furthermore, when comparing Table V with Tables III and IV it is possible to see that generating both selection mechanisms and acceptance criteria produces worse results than generating them separately.

D. Comparison with a state-of-art hyper-heuristic

In order to compare the generated high level heuristics with a already proposed hyper-heuristic strategy a state-of-art

 $TABLE\ V \\ Results\ from\ the\ best\ individual\ found\ in\ GEH YPSP-3$

Inst	1	2	3	4	5	6	7	8	9	10	11
Avg	7.6	7.0	5.7	9.7	13.8	12.7	24	24.2	31.6	27	26.4
St Dv	0.6	0.7	0.8	0.9	1.3	0.9	1.1	1.6	1.7	1.7	2
Min	7	6	4	7	12	11	21	21	29	24	24
Max	9	8	8	11	18	15	26	28	37	31	31
O(x*)	9	9	8	14	23	21	36	42	53	48	50

TABLE VI
BEST RESULTS FOUND BY GEHYPSP, BEST RESULTS FOUND BY GIHH
AND THE BEST KNOW RESULTS $O(x^*)$

	Inst	1	2	3	4	5	6	7	8	9	10	11
_	GEHyPSP	9	9	7	13	19	20	33	32	46	40	41
	GIHH (max)	9	9	8	13	22	21	35	37	49	43	45
_	O(x*)	9	9	8	14	23	21	36	42	53	48	50

human-designed hyper-heuristic framework [19] was selected. This framework produced the best results applied to the 6 domain problems from HyFlex framework [20]. Misir et al. [19] provided us the source code from the GIHH and then it was executed against the PSP problem.

Table VI presents the best results, for each instance, found by the best generated high level heuristic with GEHyPSP, the best results found by the GIHH [19] and finally the best know results (O_x(*)). The GEHyPSP and the GIHH were executed 10 minutes. It is possible to note that the GEHyPSP found worse results than the GIHH as the sequences length increase. However, the GIHH is a human-designed hyperheuristic which demanded several years of research in order to be completed and achieved good results in the six domains problems provided by HyFlex. Despite the good performance obtained by the GIHH in the domains from HyFlex, it is possible to notice that the GIHH was also unable to achieve the best known results in many instances.

E. Discussion

In order to better investigate the relationship between the generated selection mechanisms and acceptance criteria another experiment was designed. From the 30 executions of the GEHyPSP-2 ten random individuals were selected to be further analyzed. These individuals were re-executed in debugging mode in order to check its behavior. Note that only the acceptance criterion was different between them because the selection mechanism was fixed, using the best generated in the previous experiment GEHyPSP-1. From 10 individuals 7, were accepting only better or equal solutions just like the fixed acceptance criterion that was used in the GEHyPSP-1. The difference between the individuals and a fixed acceptance criterion was that individuals were slower than the fixed, because it is required to execute arithmetical functions and in the other hand only a simple if was evaluated. But thinking in the behavior they were exactly the same. It was also noticed that 2 individuals, the worst of the group, were always accepting worst solutions just like the "all moves" described by Burke et al. [3]. Finally, one individual was never accepting any solution.

These experiments showed that the GE managed, several times, to find different acceptance criteria with the same behavior of human-designed strategies.

VI. CONCLUSION

In this work the GEHyPSP, an automatic way of generating high level heuristics to a hyper-heuristic framework for the PSP problem, was presented and evaluated. The PSP is a very challenging problem with a high number of local optima and a very complex landscape. Many authors explored the PSP problem with heuristic methods. However, very often the proposed heuristic approaches are unable to find the best known results when executing against longer sequences. Usually, the hyper-heuristic frameworks fits well in this kind of complex scenario. Hence, the goal of this paper was to generate, using a grammatical evolution strategy, selection mechanisms and acceptance criteria to a hyper-heuristic framework and evaluate its performance and behavior with a set of eleven HP instances. Three groups of experiments were executed, using three randomly selected HP instances, in order to generate the high level heuristics and later the best individuals found in the experiments were executed using all the eleven HP instances.

Three groups of experiments were executed: first generating only selection mechanisms within a fixed acceptance criterion; second generating only acceptance criteria using the best selection mechanism found in the first; finally both high level heuristics were generated in parallel. The results showed that better high level heuristics were found when generating them separately. Unfortunately, when analyzing the behavior of the generated high level heuristics against the eleven instances it was possible to see that they were not able to achieve the best known results for the longer sequences. However, when comparing with a good state-of-art human-designed hyperheuristic framework (GIHH) [19] the results are slightly close. This fact shows that it is possible to automate the creation of high level heuristics and obtain results close to the state-of-art hyper-heuristics frameworks.

Another finding of this work was the behavior of the best generated acceptance criteria. It was noticed that it behaves just like "a better or equal" human-designed move acceptance strategy. Also some of the generated acceptance criteria were always accepting worst solution and this fact impacted in the individual fitness. This fact demonstrates that the GEHyPSP was able to generate acceptance criteria with the same behavior of simple human-designed move acceptance strategies. However, in order to obtain better results it might be necessary to improve the GEHyPSP to generate more complex selection mechanisms and acceptance criteria to couple with the landscape complexity of the PSP problem.

REFERENCES

- César Manuel Vargas Benítez. Um algoritmo genético paralelo para o problema de dobramento de proteínas utilizando o modelo 3dhp com cadeia lateral. 2010.
- [2] Christian Blum, Jakob Puchinger, Günther R Raidl, and Andrea Roli. Hybrid metaheuristics in combinatorial optimization: A survey. Applied Soft Computing, 11(6):4135–4151, 2011.

- [3] Edmund K Burke, Michel Gendreau, Matthew Hyde, Graham Kendall, Gabriela Ochoa, Ender Özcan, and Rong Qu. Hyper-heuristics: A survey of the state of the art. *Journal of the Operational Research Society*, 64(12):1695–1724, 2013.
- [4] Edmund K Burke, Mathew R Hyde, Graham Kendall, Gabriela Ochoa, Ender Ozcan, and John R Woodward. Exploring hyper-heuristic methodologies with genetic programming. In *Computational intelligence*, pages 177–201. Springer, 2009.
- [5] Edmund K Burke, Matthew Hyde, Graham Kendall, Gabriela Ochoa, Ender Özcan, and John R Woodward. A classification of hyper-heuristic approaches. In *Handbook of metaheuristics*, pages 449–468. Springer, 2010.
- [6] Fábio L Custódio, Hélio JC Barbosa, and Laurent E Dardenne. Investigation of the three-dimensional lattice HP protein folding model using a genetic algorithm. Genetics and Molecular Biology, 27(4):611–615, 2004.
- [7] Fábio Lima Custódio, Helio JC Barbosa, and Laurent Emmanuel Dardenne. A multiple minima genetic algorithm for protein structure prediction. Applied Soft Computing, 15:88–99, 2014.
- [8] Ian W Davis and David Baker. Rosettaligand docking with full ligand and receptor flexibility. *Journal of molecular biology*, 385(2):381–392, 2009
- [9] Candida Ferreira and U Gepsoft. What is gene expression programming, 2008.
- [10] Mario Garza-Fabre, Eduardo Rodriguez-Tello, and Gregorio Toscano-Pulido. Multiobjectivizing the hp model for protein structure prediction. In European Conference on Evolutionary Computation in Combinatorial Optimization, pages 182–193. Springer, 2012.
- [11] Mario Garza-Fabre, Gregorio Toscano-Pulido, and Eduardo Rodriguez-Tello. Locality-based multiobjectivization for the hp model of protein structure prediction. In *Proceedings of the 14th annual conference on Genetic and evolutionary computation*, pages 473–480. ACM, 2012.
- [12] Hsiao-Ping Hsu, Vishal Mehra, Walter Nadler, and Peter Grassberger. Growth algorithms for lattice heteropolymers at low temperatures. *The Journal of chemical physics*, 118(1):444–451, 2003.
- [13] John R Koza. Genetic programming: on the programming of computers by means of natural selection, volume 1. MIT press, 1992.
- [14] Natalio Krasnogor, BP Blackburne, Edmund K Burke, and Jonathan D Hirst. Multimeme algorithms for protein structure prediction. In *Parallel Problem Solving from NaturePPSN VII*, pages 769–778. Springer, 2002.
- [15] Natalio Krasnogor, W Hart, Jim Smith, and D Pelta. Protein structure prediction with evolutionary algorithms. 1999.
- [16] Elmar Krieger, Keehyoung Joo, Jinwoo Lee, Jooyoung Lee, Srivatsan Raman, James Thompson, Mike Tyka, David Baker, and Kevin Karplus. Improving physical realism, stereochemistry, and side-chain accuracy in homology modeling: four approaches that performed well in casp8. Proteins: Structure, Function, and Bioinformatics, 77(S9):114–122, 2009.
- [17] Kit Fun Lau and Ken A Dill. A lattice statistical mechanics model of the conformational and sequence spaces of proteins. *Macromolecules*, 22(10):3986–3997, 1989.
- [18] Cheng-Jian Lin and Shih-Chieh Su. Protein 3 d hp model folding simulation using a hybrid of genetic algorithm and particle swarm optimization. *International Journal of Fuzzy Systems*, 13(2):140–147, 2011.
- [19] Mustafa Misir. Intelligent hyper-heuristics: a tool for solving generic optimisation problems. 2012.
- [20] Gabriela Ochoa, Matthew Hyde, Tim Curtois, Jose A Vazquez-Rodriguez, James Walker, Michel Gendreau, Graham Kendall, Barry McCollum, Andrew J Parkes, Sanja Petrovic, et al. Hyflex: A benchmark framework for cross-domain heuristic search. In *Evolutionary computation in combinatorial optimization*, pages 136–147. Springer, 2012.
- [21] Bin Qian, Angel R Ortiz, and David Baker. Improvement of comparative model accuracy by free-energy optimization along principal components of natural structural variation. *Proceedings of the National Academy of Sciences of the United States of America*, 101(43):15346–15351, 2004.
- [22] Conor Ryan, JJ Collins, and Michael O Neill. Grammatical evolution: Evolving programs for an arbitrary language. In *Genetic Programming*, pages 83–96. Springer, 1998.
- [23] Nasser R Sabar, Masri Ayob, Graham Kendall, and Rong Qu. Automatic design of a hyper-heuristic framework with gene expression programming for combinatorial optimization problems. *IEEE Transactions on Evolutionary Computation*, 19(3):309–325, 2015.
- [24] Roberto Santana, Pedro Larrañaga, and Jose A Lozano. Component weighting functions for adaptive search with edas. In 2008 IEEE

- Congress on Evolutionary Computation (IEEE World Congress on Computational Intelligence), pages 4066–4073. IEEE, 2008.
- [25] Yang Shen, Robert Vernon, David Baker, and Ad Bax. De novo protein structure generation from incomplete chemical shift assignments. *Journal of biomolecular NMR*, 43(2):63–78, 2009.
- [26] Alena Shmygelska, Rosalia Aguirre-Hernandez, and Holger H Hoos. An ant colony optimization algorithm for the 2d hp protein folding problem. In *Ant Algorithms*, pages 40–52. Springer, 2002.
- [27] Alena Shmygelska and Holger H Hoos. An improved ant colony optimisation algorithm for the 2d hp protein folding problem. In Advances in Artificial Intelligence, pages 400–417. Springer, 2003.
- [28] Ron Unger and John Moult. Genetic algorithms for protein folding simulations. *Journal of molecular biology*, 231(1):75–81, 1993.
 [29] Ling Wang, Eric A Althoff, Jill Bolduc, Lin Jiang, James Moody,
- [29] Ling Wang, Eric A Althoff, Jill Bolduc, Lin Jiang, James Moody, Jonathan K Lassila, Lars Giger, Donald Hilvert, Barry Stoddard, and David Baker. Structural analyses of covalent enzyme–substrate analog complexes reveal strengths and limitations of de novo enzyme design. *Journal of molecular biology*, 415(3):615–625, 2012.