POSL: A Parallel-Oriented Solver Language

THESIS FOR THE DEGREE OF DOCTOR OF COMPUTER SCIENCE

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Part I

PRESENTATION

PRIOR WORKS LEADING TO POSL

In this chapter are presented Prior works leading to POSL. In Section 1.1 we present a brief work where we applied the problem subdivision approach to solve the k-medoids problem in parallel, as a first attempt aiming for the right direction in order to find the proper approach. Finally we present in Section 1.2 a study applying the PARAMILS tool in order to find the optimum parameter configuration to Adaptive Search solver.

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1.1 Domain Split

Usually, to solve some problem using parallelism, our first thought is either partitioning the problem into a set of sub-problems, or dividing its search space. In both cases, the idea is to solve a set of problems, all of them smaller and easer than the original one, and combining all the solution to obtain the solution of the original problem. In [83] a study of the impact of space-partitioning techniques on the performance of parallel local search algorithms to tackle the k-medoids clustering problem is presented. The authors use a parallel local search, in order to improve the scalability of the sequential algorithm, which is measured, in terms of the quality of the solution, with respect to the sequential algorithm. In this work two main techniques for domain partitioning are presented: space-filling curves, used to reduce any n-dimensional representation into a one-dimension space; and k-Means algorithm. We found that the used methods for domain partitioning do not take into account the number of clients associated to each new sub-domain. This results in an unbalanced distributions of workload phenomenon. For that reason the goal of this study was designing some ideas to tackle the same problem. i

The k-medoids problem aims to select a subset M of k points (the medoids) from a set of points S, such that the average distance from any point to its closest medoid is minimized. It finds a lot of applications in the industry, like in resource allocation, data mining, among others. It is quite similar to k-means problem, except that the set of medoids M is forced to be a subset of S.

We propose Algorithm 1, which represents the backbone of our idea. This algorithm takes a set of \mathbb{R}^2 points (representing the locations of the clients) and returns a partition of size K. Such a set of points is called a *domain* and the partition a *sub-domain*. At each intermediate step i, we have a set (list) of sub-domains. The algorithm takes the most populated one, splits it into two (or four, depending on the strategy) new sub-domains and includes them in

ⁱThis work falls within the framework of the Ulysses project between France and Ireland.

the list. The stop condition will depend on the fallowed approach (see below).

```
Algorithm 1: Domain_Split

input : U: Set of client locations

output: Q = \{Q_i\}_{i=1...K}: K subsets of U

1 A \leftarrow U

2 Q.Insert(A)

3 repeat

4 A \leftarrow Q.GetNext() /* It also removes the returned element

*/

5 <math>[a_1, a_2] \leftarrow Split(A)

6 Q.Insert(a_1, a_2)

7 until < some \ condition >
```

First of all, we make clear some details of the algorithm exposed above:

- Insert(...) Inserts a set (or two) into the data structure.
- GetNext() Returns the next sub-set to be divided, tacking into account the *split strategy* (see below).
- Split(...) Returns two sub-domains as a subdivision of the given domain (parameter).

In the next sub-sections we answer two main question arising at this point:

- a) How to split the each sub-domain? (<Split> function on line 5) It refers to, given a set of points (locations), how to decide which of them will be included into one sub-domain and which of them into the other.
- b) How much to split each sub-domain? (<some condition> on line 3) It refers to decide when to stop splitting the domain.

1.1.1 Domain Splitting. General point of view

In order to split the domain, we can think in some approaches, taking into account the number of available cores and the number of metro-nodes we want to place in the system. In the article of A. Arbelaez and L. Quesada a domain split taking into account the number of available cores for parallel calculus is proposed. In our approach we intend to extend this idea keeping in mind also the number of metro-nodes to allocate. Following, we propose three variants to face the problem:

a) one metro-node per core: In this variant we can assign one metro-node to each core, and in this case, splitting the domain in K sub-domains (K is the number of

cores). It means that the algorithm will compute the best position for a metro-node in a current sub-domain.

- In this case we only have to replace the line 3 by something like: for $i \leftarrow 1$ to N do ..., where N is the number of metro-nodes.
- The ideal scenario here is when N = K which is not probable at all. So we only should study the case when they are different
- In that case, we need to distribute efficiently the metro-nodes into the domain subdivisions, but here, one possible scenario arise: it can be happen that, depending on the followed domain-split strategy, we were trying to allocate a metro-node into an area with a few clients. This produce a very local point of view of the problem. That is the reason why we propose the following *second variant*.
- b) Incomplete partition: To split a sub-domain if it can generate sub-sub-domains containing at least C clients. It means, for example, if there exist in list a sub-domain with 8 clients, and the number C is fixed in C=5, this sub-domain can not be divided, because it will generate for sure a sub-sub-domain with less that 5 clients. In this case more than one metro-node would be allocated in some of those specific sub-domains, because we will have more metro-nodes than sub-domains in our model.
 - Of course, using this variant we can find situations described in the first variant. In that case we should proceed consequently.
- c) **Combination**: This variant is just a combination of the two previous variants. Then, we will be working with two parameters:
 - $C \to \text{Lower}$ bound of clients for new sub-domains. It means that a sub-domain can be divided iff the new produced sub-domains will contains more than C clients.
 - M → Lower bound of metro-nodes to be allocated in a sub-domain. In this case
 we will split the domain while it can be ensure that at least M metro-nodes will
 be assigned to each sub-domain.

1.1.2 | Split strategies

In this sub-section we first discuss three ideas to split the domain. They take a sub-domain and produce other two, dividing the space vertically or horizontally, depending on the shape of the current sub-domain. In other section we expose another strategy to follow, in which the subdivision of a sub-domain produces four sub-domains.

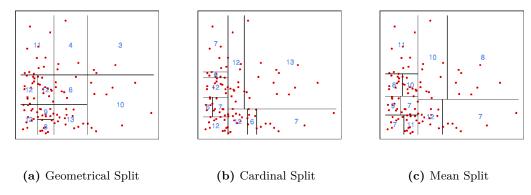


Figure 1.1: Split domain of point normally distributed $\mathcal{N}(0, 0.35)$

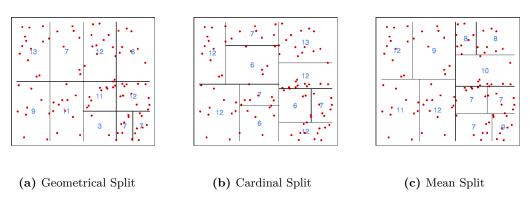


Figure 1.2: Split domain of point uniformly distributed

In all of them the number of clients in each sub-domain is taken into account, but in different ways. A common feature between them is that, at the moment of splitting a given sub-domain, it will be done in a perpendicular way (either to the x-axis or to the y-axis, depending on the characteristic of the sub-domain to be divided). The reference point to divide the sub-domain is called the *cut point* of the sub-domain.

a) Geometrical Split:

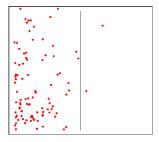


Figure 1.3: If we assume C=3 then this set can't be divided

- <u>Split criterion</u>: Geometrical → Dividing the region in tow parts with the same area.
- Cut point: The middle point of the x-axis (or y-axis, depending on the shape)
- <u>Result</u>: Two equal regions, but with different amount of clients (see Figure 1.1a). In the next step, the next sub-domain to divide will be, from those already divided, the most populated one (the sub-domain with more clients).
- <u>Problem</u>: Maybe we need to divide a sub-domain because it has a lot of client, but in the other hand it generate a sub-domain with a few clients, as you can see in the Figure 1.3.
- <u>Benefit</u>: Very fast split. If we attack scenarios with point uniformly distributed, the behavior is pretty much desirable, as we can see in the Figure 1.2a.

b) Cardinal Split

- <u>Split criterion</u>: The number of clients, i.e., a sub-domain is divided in such a way that in the resulting sub-sub-domains there will be the same number of clients.
- <u>Cut point</u>: The current sub-domain is ordered and the x-axis (or y-axis, depending on the shape) of the element (location of the client) right in the center is selected to be the *cut point*
- Result: Two regions with the same (± 1) amount of clients (see Figure 1.1a). In the next step, the next sub-domain to divide will be, from those already divided, the most populated one (the sub-domain with more clients).
- <u>Problem</u>: The subdivision process is a bit more costly: we need to group the clients on both sides of a *perfect pivot*ⁱⁱ.
- Benefit: It guarantees the same cardinality in both new sub-sub-domains.

c) Mean Split

- Split criterion: This is a mid-point strategy between the two previous. The goal is to find a *cut point* to group the elements of the current sub-domain, but in a easy way (fast), for that reason we do not compute the exact middle point to produce two sub-sub-domains with the same cardinality, as in the previous approach, instead of that we work with his expected value: the arithmetic mean.
- <u>Cut point</u>: We compute the mean of the x-axis (or y-axis) of the elements (locations) of the current sub-domain, and it will be the $cut\ point$

[&]quot;Element of a set with the same number of elements lower and greater than him

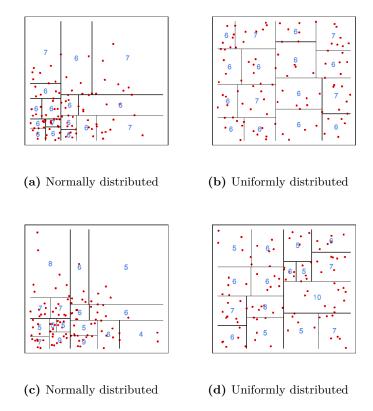


Figure 1.4: Domain divided in $2^4 = 16$ sub-domains: (a) - (b) Cardinal; (c) - (d) Mean.

- <u>Result</u>: Two regions with not exact information about their sizes or their cardinality.
- Problem: *Idem*.
- Benefit: The *cut point* can be obtained by a O(n) number of operations. As we can see in the preliminary resultsⁱⁱⁱ (Figures 1.4c and 1.4d), the behavior of this technique is near to the *cardinal split* technique, at least for normally and uniformly distributed sets of point.

1.1.3 Conclusion

In this section we present a theoretical and not validated work where we applied the search space subdivision approach to solve k-medoids problem in parallel. We have proposed some different strategies and we have showed graphically some characteristics of them.

iii The experiments are coded in R[119].

1.2 Tunning methods for local search algorithms

In this section we present our results applying Paramiles (version 2.3)^{iv}. Paramiles (first introduced by Hutter, Hoos and St utzle in 2007), is a stochastic local search approach for automated algorithm configuration. The source is available on internet and includes some examples that you can run and see how the tool works. In addition, it brings a complete User Guide with a compact explanation about how to use it with a specific solver [109, 108]. In this study we used it to tune *Adaptive Search* solver^v.

The first step was building a *wrapper* in C++ language, in order to tune more than one problem with the same code. The goal of doing this is using the tool to tune the solver, i.e., finding the best parameter configuration for a specific problem, but also the best parameter configuration to solve any kind of benchmark (a general parameter configuration).

Following we present in Table 1.1 the parameter list that we worked with:

Parameter Type Description -P PERCENT probability to select a local min (instead of staying on a plateau) -f NUMBER freeze variables of local min for NUMBER swaps -F NUMBER freeze variables swapped for NUMBER swaps LIMIT reset some variables when LIMIT variable are frozen -1 PERCENT reset PERCENT of variables -p

Table 1.1: Adaptive Search parameters

In this section, we explain in details the implementation and the experimentation process.

1.2.1 Using ParamILS

To use the tool PARAMILS, we have installed Ruby! 1.8.7 in our computer. We used a laptop Dell XPS 15(Intel Core i7-4702HQ 2.2 GHz, 16384 MB, Dual-channel DDR3L 1600 MHz) with UBUNTU 14.4. To run the tool, we needed to use the following command line:

```
>> ruby param_ils_2_3_run.rb -numRun 0 -scenariofile /\dots/<scenario_file > -validN 100
```

Where <scenario_file> is the name of the file where we have to put all the information that PARAMILS needs to tune the solver (the *tuning scenario file*). We explain its content in the next section.

ivOpen source program (project) in Ruby, available at http://cs.ubc.ca/labs/beta/Projects/ParamILS

^vAn implementation from Daniel Díaz available at https://sourceforge.net/projects/adaptivesearch/

1.2.2 Tuning scenario files

The tuning scenario file is a text file with all needed information to tune the solver using PARAMILS. It includes where to find the solver binary file, the parameters domains, etc. In our case, the tuning scenario file looks like the following:

```
algo = ./QtWrapper_wrapper
execdir = /.../src
deterministic = 0
run_obj = runtime
overall_obj = mean
cutoff_time = 50.0
cutoff_length = max
tunerTimeout = 3600
paramfile = instances/all_intervals-params.txt
outdir = instances/all_intervals-paramils-out
instance_file = instances/.../all_intervals-lower-instances.txt
test_instance_file = instances/.../all_intervals-upper-instances.txt
```

We explain in details each line in this file:

- algo \rightarrow An algorithm executable or a call to a wrapper script around an algorithm that aims the input/output format of *ParamILS* (the wrapper).
- execdir → Directory to execute algo from: "cd <execdir>; <algo>"
- run_obj → A scalar quantifying how "good" a single algorithm execution is, such as its required runtime.
- overall_obj → While run_obj defines the objective function for a single algorithm run, overall_obj defines how those single objectives are combined to reach a single scalar value to compare two parameter configurations. Implemented examples include mean, median, q90 (the 90% quantile), adj_mean (a version of the mean accounting for unsuccessful runs: total runtime divided by number of successful runs), mean1000 (another version of the mean accounting for unsuccessful runs: (total runtime of successful runs + 1000 x runtime of unsuccessful runs) divided by number of runs—this effectively maximizes the number of successful runs, breaking ties by the runtime of successful runs; it is the criterion used in most of Frank experiments), and geomean (geometric mean, primarily used in combination with run_obj = speedup. The

empirical statistic of the cost distribution (across multiple instances and seeds) to be minimized, such as the mean (of the single run objectives). vi

- cutoff_time → The time after which a single algorithm execution will be terminated unsuccessfully. This is an important parameter: if chosen too high, lots of time will be wasted with unsuccessful runs. If chosen too low the optimization is biased to perform well on easy instances only.
- tunerTimeout → The timeout of the tuner. Validation of the final best found parameter configuration starts after the timeout.
- ullet paramfile o Specifies the file with the parameters of the algorithms.
- outdir \rightarrow Specifies the directory ParamILS should write its results to.
- instance_file \rightarrow Specifies the file with a list of training instances.
- test_instance_file \rightarrow Specifies the file with a list of test instances.

Another important file that we have to compose properly is the algorithm parameter file, just following the instruction from [109] –[...] each line lists one parameter, in curly parentheses the possible values considered, and in square parentheses the default value [...]. Our algorithm parameter file looks like follows:

```
P {20, 25, 30, 35, 40, 45, 50, 55, 60} [50]

f {0, 1, 2, 3} [1]

F {0, 1, 2, 3} [0]

1 {0, 1, 2, 3} [1]

p {1, 2, 3, 5, 10, 20} [5]
```

In the current $Adaptive\ Search$ implementation, the solver binary file and the problem instance are the same thing. It means that we only have to use the following command to solve the All-intervals problem of size K, for example:

```
>> ./all-intervals K
```

So, to use PARAMILS we modified a little the code: now our solver takes the size parameter from a text file. That way, the instance file is a text file only containing a number.

The solver we want to tune using Paramiles (Adaptive Search in this case), must aims specific input/output rules. For that reason instead of modifying the current code of Adaptive Search implementation, we preferred to build a wrapper.

viWe use mean but maybe we can experiment with other values

1.2.3 Building the wrapper

The algorithm executable must follow the input/output criteria presented below:

Launch command:

```
>> <algo_exectuable> <instance_name> <instance-specific_information> ...  
<cutoff_time> <cutoff_length> <seed> <params>
```

- <algo_exectuable> Solver
- <instance_name> In our case, a text file containing only the problem size
- <instance-specific_information> We don't use it
- <cutoff_time> Cut off time for each run of the solver (see above)
- <cutoff_length> We don't use it
- <seed> Random seed
- params> Parameters and its values

Exmaple:

```
>> ./QtWrapper_320.txt "" 50.0 214483647 524453158 -p 5 -l 1 -f 1 -P 50 -F 0
```

Output:

```
>> <solved >, <runtime >, <runlength >, <best_sol >, <seed>
```

- <solved> SAT if the algorithm terminates successfully. TIMESOUT if the algorithm times out.
- <runtime> Runtime
- <runlength> -1 (as Frank Hutter recommended)
- <best_sol> -1 (as Frank Hutter recommended)
- < cutoff_length > We don't use it
- <**seed**> Used random seed

Exmaple:

```
>> SAT, 2.03435, -1, -1, 524453158
```

To build the wrapper we have followed a simple algorithm: launch two concurrent process. In the parent process we translate the input of the wrapper to the input of Adaptive Search solver. The solver is executed, and the runtime is measured. After that we post the output properly. In the child process a sleep command is executed for <runtime> seconds and after that, if the parent process has not finished yet, it is killed, posting a time-out message. See Algorithm 2 for more details.

```
Algorithm 2: Costas Wrapper
```

```
input : Pth_{\pi}: problem instance path, k: cut off time, s: random seed, \theta: parameters configuration
   \mathbf{output}: PiLS_{out}:  Output in a PARAMILS way
 1 fork() /* Divide the execution in two processes
                                                                                                          */
 2 if <in child process> then
       t_0 \leftarrow \texttt{clock\_TIC()}
       strCall \leftarrow build\_str("./AS\_Wrapper %1 -s %2 %3", Pth_{\pi}, s, \theta)
 4
       systemCall(strCall)
 5
       t_e \leftarrow \texttt{clock\_TOC()}
 6
       killProcess(<parent process>)
 7
       t \leftarrow t_e- t_0
 8
       return paramilsOutput(SAT, t, s)
 9
10 else
       sleep(k)
11
12
       killProcess(<child process>)
       return paramilsOutput(TIMESOUT, k, s)
13
14 end
```

1.2.4 Using the wrapper

In this section we explain how to use our wrapper to be able to tune easily instances of All-Interval Series and Costas Array problems. The All-Interval Series Problem^{vii} is the problem of finding a vector $s = (s_1, \ldots, s_n)$, given $n \in \mathbb{N}$, such that s is a permutation of the vector $(0, 1, \ldots, n-1)$ and the interval vector $v = (|S_2 - s_1|, |S_3 - s_2|, \ldots, |S_n - S_{n-1}|)$ (called an all-interval series of size n) is a permutation of the vector $(1, 2, \ldots, n-1)$. The Costas Array Problem consists in finding a Costas array, which is an $n \times n$ grid containing n marks such that there is exactly one mark per row and per column and the n(n-1)/2 vectors joining each couple of marks are all different (see below for more details about this problems).

viiCSPlib:007 (http://www.csplib.org/Problems/prob007/)

1.2.4.1 Factory call

The first step is to implement the class ICALLFACTORY. Here, the string-binary-name for the command call is statically obtained. We present, as example, the class All_IntervalCallFactory:

```
// all_interval_call_factory.h
class All_IntervalCallFactory: public ICallFactory
{
    public:
        std::string BuildCall();
        std::string BuildDefaultCall();
};
```

```
// all_interval_call_factory.cpp
#define ALGO_EXECUTABLE "./all-interval"
#define DEFAULT_CALL "./all-interval _100.txt"

std::string All_IntervalCallFactory::BuildCall()
{
    return ALGO_EXECUTABLE;
}
std::string All_IntervalCallFactory::BuildDefaultCall()
{
    return DEFAULT_CALL;
}
```

All we have to do is to define our new macro ALGO_EXECUTABLE (DE-FAULT_CALL is not being used)

1.2.4.2 | Main method

Let's suppose now that we want to run an algorithm called *mySolver* that receives a file as parameter, called *my_instance_size.txt* (this is mandatory). We have to create (as we've explained before) the class My_SolverCallFactory and defining the macro as follows:

```
#define ALGO_EXECUTABLE "./mySolver"
```

Now, the main method would be exactly like this:

```
int main(int argc, char* argv[])
{
    shared_ptr<ICallFactory> problem =
        make_shared<My_SolverCallFactory>();
    shared_ptr<TuningData> td =
            (make_shared<TuningData> (argc, argv, problem));

    shared_ptr<ADWrapper> w (make_shared<ADWrapper>());
    string output = w->tune(td);

    cout << output << endl;
    return 0;
}</pre>
```

1.2.5 Results

In this section we present the results of applying PARAMILS to the resolution of *All-Interval Series* and *Costas Array* problems through *Adaptive Search*. In both cases, we need to chose a set of *training instances*, to train the tuner, and a set of *test instances*, used to obtain the parameter setting.

1.2.5.1 Tuning Adaptive Search for All-Intervals Series Problem

Study cases:

a) The training instances set is composed by instances of All–Intervals problems of order N with

```
N \in \{100, 110, 120, 130, 140, 150, 160, 170, 180\}
```

b) The test instances set is composed by instances of All–Intervals problems of order N with

```
N \in \{190, 200, 210, 220, 230, 240, 250, 260, 265\}
```

- c) The time-out for each run is 50.0 seconds
- d) The test quality is based on 100 runs

In a **First Experiment** we use the following *parameters domains*:

- P {41, 46, 51, 56, 60, 66, 71, 76, 80}
- F, f, l {0, 1, 2, 3}
- p {5, 10, 15, 20, 25, 30, 35}

Table 1.2 shows results using a time-out of 5.5 hours (20,000 seconds), and Table 1.3 shows results using a time-out of 1 hour. In the second case we where able to perform more runs, due to the available time, but in both cases the training qualities are not so different. However, we can se the difference int the test qualities, and conclude that results using 5 hours of time-out are more reliables.

Ini	itial o	config	gurat	ion	Fin	Final best configuration				Training	Test	
F	P	f	1	p	F	Р	f	1	p	quality	of runs	quality
0	66	1	1	25	0	80	2	1	35	0.79666	1780	8.274
2	56	2	2	20	1	80	1	1	10	0.795	1637	5.508
0	41	0	0	5	1	80	3	0	15	0.789	1547	5.8478
3	80	3	3	35	1	80	2	0	10	0.880686	1258	6.15398

Table 1.2: All-Intervals Series: tunerTimeout = 20,000 seconds

Ini	itial c	onfig	gurat	ion	Fin	al be	st co	nfigu	ration	Training	Number	Test
F	Р	f	1	p	F	Р	f	1	p	quality	of runs	quality
0	66	1	1	25	0	80	0	1	25	0.815	384	5.8191
0	66	1	1	25	1	80	1	1	35	0.737	452	6.267
0	66	1	1	25	1	56	0	1	35	1.03	371	9.056
0	66	1	1	25	0	76	0	1	20	0.814	385	4.915
0	66	1	1	25	0	80	3	1	20	0.76	469	5.417
2	56	2	2	20	0	41	0	1	10	0.919	239	18.364
2	56	2	2	20	0	56	1	1	20	0.819	407	5.409
2	56	2	2	20	1	80	1	1	35	0.772	457	5.43
2	56	2	2	20	1	80	0	1	10	0.858	504	5.566
2	56	2	2	20	0	80	1	1	10	0.7845	562	18.944
0	41	0	0	5	0	41	1	0	10	0.9749	367	5.97813
0	41	0	0	5	0	41	1	0	10	0.885	450	5.706
0	41	0	0	5	0	41	1	0	10	0.906	335	18.707
0	41	0	0	5	0	41	1	0	10	0.995	335	19.558
0	41	0	0	5	0	41	0	0	5	0.855	404	5.686
3	80	3	3	35	0	66	3	1	25	0.9118	230	26.585
3	80	3	3	35	0	80	1	1	10	0.732	310	7.875
3	80	3	3	35	0	80	0	1	20	0.816	303	7.2896
3	80	3	3	35	1	80	3	1	35	0.821	327	6.812
3	80	3	3	35	0	80	0	1	30	0.9203	443	5.401

Table 1.3: All-Intervals Series: tunerTimeout = 3,600 seconds

In a **Second Experiment** we decide to enlarge a bit more the parameters domains and use a time-out of 5 hours. The **Parameters domains** are the following:

- P {10, 20, 30, 40, 50, 60, 70, 80, 90}
- F, f, l {0, 1, 2, 3, 4, 5, 6, 7, 8}
- p {10, 20, 30, 40, 50, 60, 70}

In	itial d	config	gurat	ion	Fin	al be	st co	nfigu	ration	Training	Number	Test
F	Р	f	1	p	F	Р	f	1	p	quality	of runs	quality
0	10	0	0	10	0	40	7	0	50	0.883188	936	6.3191
0	10	0	0	10	0	80	2	1	40	0.774659	1584	5.45674
0	10	0	0	10	0	40	2	0	10	0.96885	1104	6.82643
4	60	4	4	40	0	60	8	1	40	0.90358	1566	5.48127
4	50	4	4	40	0	80	5	1	20	0.78536	1662	11.5649
3	50	4	2	30	0	90	6	1	70	0.79440	1395	5.08108
0	90	0	0	10	1	90	6	1	10	0.859569	1379	5.4286
0	90	0	0	10	1	90	6	1	30	0.80738	1117	5.47126
8	90	8	8	60	0	80	5	1	10	0.834934	1384	5.5377
5	30	2	3	60	0	90	1	0	20	0.862013	1707	5.21837
3	20	2	4	60	0	80	6	1	10	0.805604	1630	5.4467
6	70	1	3	50	0	80	5	1	10	0.792600	1344	5.46558
6	40	1	3	30	1	80	7	0	20	0.822703	1977	5.41185

Table 1.4: All-Intervals Series: tunerTimeout = 18,000 seconds

The results presented in Table 1.4 show better results in terms of test quality with respect to Table 1.2. For that reason, in the **FINAL Experiment**, only the results obtained in those tables were took into account (also because they were obtained by using longer times-out). As it can be observed in those tables, *Adaptive Search* seems to show a good behavior if the parameters \mathbf{F} , \mathbf{P} and \mathbf{l} are in the following sets: $\mathbf{F} \in \{0, 1\}$, $\mathbf{P} \in \{80, 90\}$ and $\mathbf{l} \in \{0, 1\}$.

In that sense, a specific configuration was extracted from the results above, and 60 runs of Adaptive Search were performed solving All–Intervals (N = 600) benchmark:

- 30 using the default parameter configuration (-F 0 -P 66 -f 1 -l 1 -p 25)
- 30 with an optimal parameter configuration extracted from the Tables 1.2, 1.4 (-F 0 -P 80 -f 6 -l 1 -p 10)

Table 1.5 shows results by using the default parameter settings, and Table 1.6 shows the results by using the parameter configuration found by PARAMILS, and it is clear that the default configuration shows better results than ParamILS's one, in terms both of runtime mean and standard deviation Using the default parameter settings, $Adaptive\ Search\ can$ obtains best results int terms of $mean\ and\ slowest\ run$. However, using the PARAMILS found

parameter settings, it reached a *fastest* run two times faster than the one using the default parameter settings.

37.210	411.300	112.510	171.000	73.770				
327.880	214.910	124.910	482.740	530.440				
212.660	99.370	287.400	533.540	18.410				
197.290	1016.950	110.230	566.480	1362.010				
94.860	819.700	434.460	620.600	95.920				
80.580	333.370	121.590	489.700	248.370				
mean: 341.005333								
	sprea	ad: 310.4	44635					

Table 1.5: All-Intervals Series: Default configuration runtimes (secs)

154.460	264.530	169.840	26.990	108.790				
550.210	104.900	31.100	9.870	1242.900				
678.760	475.570	201.200	622.410	297.960				
526.930	375.620	293.380	598.850	350.270				
540.290	252.940	673.630	423.030	589.210				
32.080	254.640	2034.020	571.100	207.090				
mean: 422.085667								
	spre	ead: 404.61	8226					

Table 1.6: All-Intervals Series: Paramiles configuration runtimes (secs)

1.2.5.2 Tuning Adaptive Search for Costas Array Problem

Study cases:

- a) The training instances set is composed by instances of Costas Array problems of order N with $9 \le N \le 15$
- b) The test instances set is composed by instances of Costas Array problems of order N with $14 \le N \le 19$
- c) The cutoff for each run was 60.0 seconds
- d) The test quality is based on 100 runs

The **First Experiments** with this benchmark was using the following parameter domains:

- P {10, 20, 30, 40, 50, 60, 70, 80, 90}
- F, f, l {0, 1, 2, 3, 4, 5, 6, 7, 8}

• p {5, 10, 20, 30, 40, 50, 60, 70}

Table 1.7 shows results selecting directly a time-out of 5 hours (18,000 seconds). In this case the training quality of the solutions is better, but do not observe any improvement in the test quality. We can see also how *Adaptive Search* seems to be not sensitive to parameters \mathbf{F} and \mathbf{p} , i.e. they don't change during the tuning process. On the other hand, the tuner seems to find some optimum values for the other parameters: $\mathbf{P} \in \{$ 80, 90 $\}$, $\mathbf{f} \in \{$ 4, 5 $\}$ and $\mathbf{l} = 2$.

In that case also, an specific configuration was extracted from the results showed in Table 1.7, and 60 runs of Adaptive Search were performed solving Costas Array (N = 20) benchmark:

- 30 using the default parameter configuration (-F 0 -P 50 -f 1 -l 0 -p 5)
- 30 with an optimal parameter configuration extracted from the Table 1.7 (-F 3 -P 90 -f 5 -l 2 -p 30)

In	itial o	onfig	gurat	ion	Fin	al be	st co	nfigu	ration	Training	Number	Test
F	P	f	1	p	F	Р	f	1	p	quality	of runs	quality
0	10	0	0	5	2	90	2	2	5	0.0493699	957	5.8461
0	10	0	0	5	0	90	5	2	5	0.0509388	1783	6.52742
0	10	0	0	5	0	90	5	2	5	0.049901	1759	5.21828
3	40	4	4	30	3	90	5	2	30	0.053974	856	6.3539
4	50	3	5	20	4	90	5	2	20	0.0500355	2000	5.4047
4	60	5	3	50	4	60	5	3	50	0.0520575	2000	6.09106
8	90	8	8	70	8	80	4	2	70	0.052685	550	3.85682
8	90	8	8	70	8	80	4	2	70	0.054104	536	4.17855
8	90	8	8	70	8	80	4	2	70	0.0497819	1284	3.90945
3	10	1	6	60	3	90	5	2	60	0.054934	2000	6.81675
5	70	6	1	10	5	90	4	2	10	0.0499895	2000	4.07365
1	30	5	7	5	1	90	4	2	5	0.0525747	1237	2.70091
7	80	2	0	70	7	90	5	2	70	0.0502264	212	5.2637

Table 1.7: Costas Array: tunerTimeout = 18,000 seconds

Table 1.8 shows the results by using the default parameter configuration, and Table 1.9 shows the results by using the parameter configuration found by *ParamILS*. One more time, "in the mean", the default configuration outperforms PARAMILS's.

1.2.6 Conclusion

The conclusion of this study is that the tunning process by hand in this case was more effective than using PARAMILS. Results show that default parameters used in the current

452.980	91.420	31.510	827.860	96.670					
635.030	295.830	272.360	151.040	170.660					
183.550	161.340	91.240	426.470	62.020					
138.090	236.030	2.850	187.240	21.510					
165.370	90.440	195.580	15.390	229.720					
170.840	174.210	30.520	6.570	115.880					
mean: 191.007									
	spread: 185.362								

Table 1.8: Default configuration runtimes (secs)

546.260	263.230	17.200	29.220	495.940				
237.340	187.760	7.810	43.120	94.370				
59.930	128.690	247.810	265.010	231.260				
209.640	465.340	21.840	8.740	1264.610				
57.700	122.890	450.610	229.580	174.540				
414.080	402.250	91.150	677.190	58.640				
mean: 250.125								
	\mathbf{sp}	read 263	.539					

Table 1.9: ParamILS configuration runtimes (secs)

Adaptive Search implementation are more effective and consistent than those found by PARAMILS for both benchmarks (All-Interval Series and Costas Array problems).

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