Heuristic Crossover Operator for Evolutionary Induced Decision Trees

Sašo Karakatič and Vili Podgorelec
Institute of Informatics FERI, University of Maribor, Smetanova 17, Maribor, Slovenia

Keywords: Genetic Algorithm, Decision Tree, Crossover.

Abstract: In this paper we propose an innovative and improved variation of genetic operator crossover for the classification decision tree models. Our improved crossover operator uses heuristic to choose the tree node that is exchanged to construct the children solutions. The algorithm selects a single node based on the classification accuracy and the usage of that particular node. We evaluate this method by comparing it with the results of the standard crossover method where nodes for exchange are chosen at random.

1 INTRODUCTION

The decision tree classification method is one of the classifiers used in the machine learning field that is notorious for its ease of interpretation by human users and the ability to generalize the problem solutions (Cantu-Paz & Kamath 2003). Many methods for the construction of decision trees were proposed in the past, but our paper is focused on the evolutionary method of genetic programming. Genetic programming is a variation of the genetic algorithm where individuals in the generation are presented with a tree structure. Each individual presents one solution and the process of evolution consists of several genetic operators which create new individuals (crossover) or try to modify the existing ones (mutation) (Koza 1992). Our paper focuses on the crossover operator which is a method that generates new individuals from the parent trees in such a way, that some characteristics from both parents are represented in the child tree – therefore mimicking the natural process of mating, where DNA of humans is constructed from the genetic material from both parents.

Crossover is heavily dependent on the type of representation used for individuals. Due to the standard representation of decision trees being done with a tree structure instead of an array, our focus was on the crossover for tree genotype – genetic programming. The crossovers on the tree are done by exchanging the subtrees from one individual to another, with the method of choosing these subtrees being the main question. We propose an innovative crossover method that tries to eliminate the weak parts of the tree individual, based on the partial accuracy of the subtrees and usage of that subtree.

Rest of the paper is organized as follows. We start with the background overview of existing research done on the crossover operators of the decision tree models and other similar tree based problems. We continue with the description of the proposed improvement in the crossover operator, where we present the idea of the innovative crossover method. Following that, the experiment on the standard benchmark datasets, to evaluate the quality of our proposed improvement is presented. The results are backed up by the statistical methods and are supplemented with the interpretation of the results and the discussion. In the conclusion we summarize the paper and present the final verdict. Possibilities for further research are also discussed.

2 RELATED WORK

Koza (Koza 1992) describes the random subtree exchange crossover as a blind operator, due to no context being included in the crossover location choosing. He also introduced constrained crossover that exchanges only two matching subtrees (Koza & Noyes 1994). The positive effect of this is the maintenance of the context, which allows the subtree to evolve faster to its optimal form. However, this has its own limitations. Namely, there is no new genetic material entering the subtrees which eventually leads to the local optimum.
D'haeseleer takes a different approach, where he assigns a location to each node. He tries to extend the initial context preserving crossover from Koza by limiting the crossover so that only subtrees on same location can be exchanged with its strong context preserving crossover. As he states himself, the greatest weakness of this crossover is the resulting limited diversity, so he expands it with the weak context preserving crossover where the nodes exchanged have to have the same parent (D’haeseleer 1994). The main disadvantage of this type of location based crossover is that most of the times the nodes on the same location are not in the same context.

Iba and Garis (Iba & de Garis 1996) used a heuristic search for subtrees, where the evaluation of each node was calculated with the mean square error value. By this criterion, the worst subtrees are replaced with the best subtrees. This method is similar to the one proposed in the paper. However, the search for the worst subtrees can select nodes in high depth levels, thus constructing large trees which was never evaluated on classification problem. In addition to that, we believe that substitution for best subtrees eventually leads to less diversified population of solution meaning that this kind of method has a potential to get stuck in local optimum early on in the evolution process.

One very similar method to ours was made by Hengpraprohm and Chongstitvatana where they exchanged the worst subtrees with the best ones. They evaluated the subtrees, followed by the pruning and comparison of the fitness pre and post (Hengpraprohm & Chongstitvatana 2001). Same problems may arise here, as stated previously, replacing the worst subtrees with the best ones can prematurely result in local optima.

Another context aware crossover was done by Majeed and Ryan in (Majeed & Ryan 2006a) and (Majeed & Ryan 2006b), where they make multiple children from two parent trees and select only one of them as the real successor that advances to the next generation. The crossover point in the second parent is chosen at random, while in the first parent every legal point is used. Surveys of evolutionary algorithms for decision trees by Barros et. al. (Barros et al. 2012) and Espejo at. al. (Espejo et al. 2010) do not mention any other heuristic crossover operators.

Expansion of good building block ideas came from Zhang (Zhang et al. 2007) with his looseness controlled operator, where he used local hill climbing search to construct and evaluate subtrees. He introduces the evaluation function called “sticky” that evaluates how good do particular nodes perform together in parent-child relationship and perform crossover based on the stickiness values.

3 CROSSOVER IMPROVEMENT

Genetic operator crossover is one of the most impacting operators on the final solution of the whole process. Process of mating begins with two parent individuals that are selected (with the selection operator) for the creation of the child individual. The standard procedure is to select a random subtree in the first parent tree and exchange it for a random subtree from the second parent tree; the resulting tree is considered as the child individual. Our goal was to improve this process with heuristics in hopes of improving the results.

![Figure 1: Evaluation of nodes in the whole tree.](image)
The nature of the decision tree is such, that we can calculate the basic performance metrics of each individual subtree, which is based on the instances that were processed by that particular subtree. Our hypothesis was that we can select the worst performing subtree from the first parent and replace it, so that the worst part of the tree will always be selected for the crossover and eliminated. This will produce the solutions with at least an equal and hopefully even better quality based on the classification metrics.

For the evaluation of subtrees we initially used the overall classification accuracy of the subtree on its instances. The results showed that this produced somewhat better results based on the mean accuracy of the 100 experiments, but the results were not statistically significant. This crossover chose the nodes with worst accuracy, which were terminal in most instances and thus failed to make an impact by ignoring the nodes that were used the most but were in the higher levels of the tree. Mistakes made by the nodes in the tree on the lower levels impact the results much more than the nodes on the higher levels, so a change which would force the algorithm to choose more nodes on lower levels of decision tree was necessary.

\[ e = \text{usage} + \text{error rate} \]  

(1)

Based on this experience we introduced another metric to the evaluation of the subtrees – the usage of the root nodes on the subtrees. We reasoned that this process should replace the subtrees with the worst accuracy while at the same time giving more emphasis on more used subtrees (near the root of the tree) and essentially ignoring the nodes near the end of the tree which are the unfortunate consequences of randomness and processes just a few classification instances. Combined evaluation of nodes is calculated as the sum of the error rate \((1 - \text{accuracy})\) and the usage of the node as is show in the equation 1. Usage of the nodes is a percentage of classification instances processed in the node during the training phase of the genetic algorithm. This process ignores the root node. The evaluation of the nodes in the tree is shown in the Figure 1, where node with the black background is the node with the highest crossover evaluation and is chosen for the exchange with the random subtree with the second parent.

### 4 EXPERIMENT

<table>
<thead>
<tr>
<th>Dataset</th>
<th>Instances</th>
<th>Avg standard cx</th>
<th>Avg improved cx</th>
<th>p</th>
<th>Max standard</th>
<th>Max improved</th>
</tr>
</thead>
<tbody>
<tr>
<td>autos</td>
<td>205</td>
<td>.4705</td>
<td>.5044</td>
<td>&lt;.001</td>
<td>.62</td>
<td>.67</td>
</tr>
<tr>
<td>glass</td>
<td>214</td>
<td>.5744</td>
<td>.5814</td>
<td>.136</td>
<td>.73</td>
<td>.77</td>
</tr>
<tr>
<td>diabetes</td>
<td>768</td>
<td>.6382</td>
<td>.6444</td>
<td>.789</td>
<td>.70</td>
<td>.70</td>
</tr>
<tr>
<td>iris</td>
<td>150</td>
<td>.9447</td>
<td>.9480</td>
<td>.629</td>
<td>1.00</td>
<td>1.00</td>
</tr>
<tr>
<td>primary-tumor</td>
<td>339</td>
<td>.2601</td>
<td>.2741</td>
<td>.007</td>
<td>.38</td>
<td>.38</td>
</tr>
<tr>
<td>sonar</td>
<td>208</td>
<td>.6067</td>
<td>.6740</td>
<td>&lt;.001</td>
<td>.80</td>
<td>.80</td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>Dataset</th>
<th>No. of classes</th>
<th>Avg standard cx</th>
<th>Avg improved cx</th>
<th>p</th>
<th>Max standard</th>
<th>Max improved</th>
</tr>
</thead>
<tbody>
<tr>
<td>autos</td>
<td>7</td>
<td>.2069</td>
<td>.2460</td>
<td>&lt;.001</td>
<td>.35</td>
<td>.41</td>
</tr>
<tr>
<td>glass</td>
<td>7</td>
<td>.2623</td>
<td>.2913</td>
<td>&lt;.001</td>
<td>.43</td>
<td>.50</td>
</tr>
<tr>
<td>diabetes</td>
<td>2</td>
<td>.4899</td>
<td>.5121</td>
<td>.015</td>
<td>.62</td>
<td>.62</td>
</tr>
<tr>
<td>iris</td>
<td>3</td>
<td>.9443</td>
<td>.9478</td>
<td>.597</td>
<td>1.00</td>
<td>1.00</td>
</tr>
<tr>
<td>primary-tumor</td>
<td>22</td>
<td>.0321</td>
<td>.0419</td>
<td>&lt;.001</td>
<td>.08</td>
<td>.09</td>
</tr>
<tr>
<td>sonar</td>
<td>2</td>
<td>.5501</td>
<td>.6510</td>
<td>&lt;.001</td>
<td>.79</td>
<td>.79</td>
</tr>
</tbody>
</table>
implemented and tested on six standard classification benchmark datasets. Characteristics of the used dataset and the results of the experiments are shown in the Table 1 and Table 2. Algorithm ran 10 times on 10 folds on each dataset (100 times in total) in order to produce reliable results and minimize the chance of randomness. Accuracies in the Table 1 are means of accuracies on all 100 runs. In addition to mean of all accuracies, we calculated the average Fscore of all classes on the dataset for each run and included the mean of those in the Table 2, for comparison. Independent samples Mann-Whitney U test was conducted to ensure the validity of the results. We added the p-values for each test. Both, the results that are statically significant (p < 0.05) and the instances where our improved algorithm achieved better best resulting classification model are bolded. Crossovers are labelled as CX.

Basic genetic algorithm settings were the same for all of the datasets and were set as follows: evolution process ran for 2000 generations, with 100% chance of crossover and 10% chance of basic random mutation. Each generation consisted of 150 individual solution decision trees, where the best individual automatically advanced to next generation (one elite individual). Selection operator was chosen to be binary tournament, and the fitness evaluation method was based on the accuracy of the individuals.

Table 1 and Table 2 show, that our proposed crossover operator not only matched, but improved the classification results on both metrics, the accuracy and the average Fscore, on used datasets. In datasets glass (p = 0.136), diabetes (p = 0.789) and iris (p = 0.597) the difference was not as dramatic as in other datasets, but average accuracy achieved is still better. Based on these results we can conclude that the proposed improvements to the crossover methods made significant improvements, which is unlikely due to chance, on chosen datasets. Out of six datasets, our algorithm matched 4 of them in the best resulting model and was better in other 2 datasets in the overall accuracy. In the Table 2 and Figure 2, the same comparison is made but on the Fscore classification metric, where our algorithm found better average Fscore in all 6 datasets, where 5 of them are statistically significant improvements. In 3 out of 6 dataset, our proposed algorithm achieved better maximal Fscore and matched the original crossover on the maximal score on 3 other datasets. The dataset iris was an outlier here, because both crossover methods achieved in perfect accuracy and Fscore. As can be seen from Figure 2 no further improvement is possible on iris dataset. This also explains why no statistically significant improvement could be made with the iris dataset.

Let us look in details why our improvement results in better solutions in accuracy and Fscore. Figure 3 compares two samples (from dataset sonar), one from each crossover, through evolution. Vertical axis shows accuracy of the best solution in the population. From results shown in Figure 3 we can
see that algorithm with the standard crossover reaches its optimum much faster than the algorithm with our improved crossover. Figure 3 confirms that indeed, the algorithm with standard crossover reached its optimum already around 1400th generation, while algorithm with our improved crossover continues to run for another 600 generations and reaching its optimum in generation 1860. The difference between accuracy is ~0.5 in favour of our proposed crossover.

5 CONCLUSIONS

We proposed the innovative crossover operator that does not choose the place for the crossover process based at random but rather on the accuracy and usage of the subtree. As mentioned in the overview in the second chapter, no similar method was found in the tree classification models. This innovative method was tested on several datasets in our experiment, with the results showing significant increases in the metrics of overall classification accuracy and average Fscore of final tree models. Based on this we can assume that these result can be reproduced on other classification problem datasets. To corroborate this statement further studies are planned on more diverse, unbalanced, and otherwise unusual datasets. Additional research option would be to try and replace the overall accuracy in the evaluation of nodes with another metric such as Fscore or similar classification or decision tree evaluation metric.

REFERENCES


