

An in-depth study of the syntactic analysis of graphs using  
search matching algorithms, to the degree of sub graphs

Student number:

email:

Supervisor: Dr Linda Marshall

Dept Computer Science, University of Pretoria

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# 1 Graph Basics

## 1.1 Graph Overview

A Graph in Mathematics and Computer Science is defined as a pair  $G = (V, E)$ , where  $V$  is the set of vertices and  $E$  is the set of edges, formed by pairs of vertices with each other. Figure 1 demonstrates the structural attributes of a simple graph.

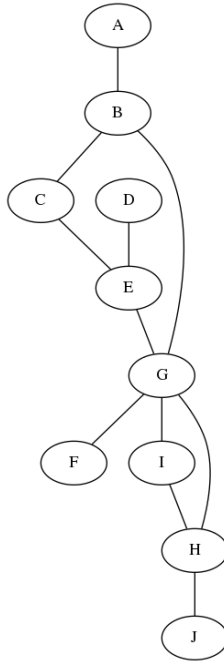


Figure 1: Representation of a graph

Graphs can either be directed or they can be undirected. This means that the edges in the graph could have an absence of direction as in the 1 above, or they could have a direction showing from which vertex the edge is coming from, and to which vertex the edge is going to. Figure 2 demonstrates a directed graph, commonly known as a digraph. This characteristic is demonstrated by the edge 3, that goes from node  $B$  to node  $G$ , and also by edge 2 that goes from node  $B$  to node  $C$ .

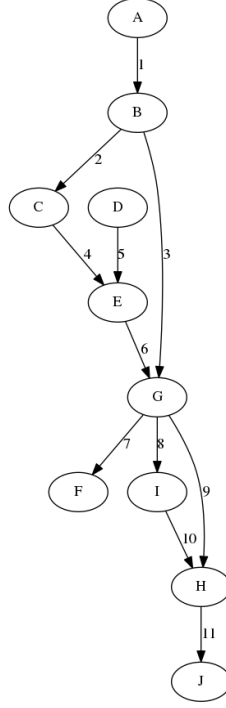


Figure 2: Representation of a digraph

Note that the undirected graph mentioned above is commonly referred to as a bigraph because the direction of its edges could be perceived as to going in both direction as it is not specified.

In this paper, the focus is primarily on directed graphs, and they shall be referred to as digraphs from here onward.

## 1.2 Digraph

A Digraph in mathematics is defined as is a pair of disjoint vertices and edges  $(V, E)$ , and their respective mappings that comprises of two components, namely the initial vertex and terminal vertex of each edge i.e. each edge has a initial vertex:

$$Ei \rightarrow Vi \quad (1)$$

and a terminal vertex:

$$Vj \rightarrow Ej \quad (2)$$

for some vertices  $Vi, Vj$  in  $V$  and edges  $Ei, Ej$  in  $E$  [7] refer to the figure 2.

### 1.3 Graph representation

Graphs are represented in a variety of ways, from adjacency lists, incident matrices and adjacency matrices. The algorithms that are studied in this paper make use of adjacency matrices and adjacency list representations of graphs.

#### 1.3.1 Adjacency matrices

An adjacency matrix is a  $n \times n$  matrix  $A$ , with  $A(i, j) = 1$  if  $(i, j) \in E$  [9]. This means that wherever there is an edge in the graph, it is denoted by a 1 in the matrix, places in the matrix where there is an absence of an edge, are denoted by 0.

Figure 3 depicts the association between the graph in figure 1 and its adjacency matrix.

	A	B	C	D	E	G	F	I	H	J
A	0	1	0	0	0	0	0	0	0	0
B	1	0	1	0	0	1	0	0	0	0
C	0	1	0	0	1	0	0	0	0	0
D	0	0	0	0	1	0	0	0	0	0
E	0	0	1	1	0	1	0	0	0	0
G	0	1	0	0	1	0	1	1	1	0
F	0	0	0	0	0	1	0	0	0	0
I	0	0	0	0	0	1	0	0	1	0
H	0	0	0	0	0	1	0	1	0	1
J	0	0	0	0	0	0	0	0	1	0

Figure 3: Representation of a graph and its associated adjacency matrix

#### 1.3.2 Adjacency list

An Adjacency list is vertices of a graph, of which each vertex is connected to the list. The vertices in an adjacency list point to their own list of edges that they are connected to (i.e. the list contains the edges that connect them to other vertices).

Figure 4 depicts the association between the graph in figure 1 and its adjacency list.

A	B				
B	A	C	G		
C	B	E			
D	E				
E	C	D	G		
G	B	E	F	I	H
F	G				
I	G	H			
H	G	I	J		
J	H				

Figure 4: Representation of a graph and its associated adjacency list

#### 1.4 Supergraphs and subgraphs

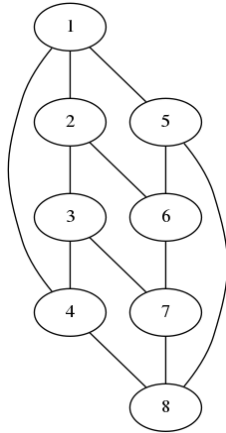
Let  $G_A$  be a graph defined as follows  $G_A = (V_A, E_A)$  and let  $G_B$  be another graph that is defined as follows  $G_B = (V_B, E_B)$  where  $V_A, V_B$  are sets of vertices and  $E_A, E_B$  are sets of edges. In graph theory, a graph  $G_A$  is said to be a subgraph of graph  $G_B$ , and graph  $G_B$  is said to be a supergraph of graph  $G_A$  if all the vertices and edges that are in graph  $G_A$  are also in graph  $G_B$ , that is [3]:

- (1)  $V_A \subseteq V_B$ , and
- (2) Every edge of  $G_A$  is also an edge in  $G_B$ .

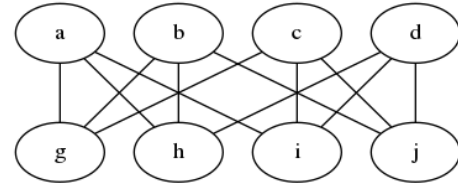
In figure 1, the graph constructed by vertices  $D, E, G, F$  and  $I$  is the sub-graph of the entire graph. And thus the graph is a super-graph of the sub-graph constructed by the vertices, namely  $D, E, G, F$  and  $I$ .

#### 1.5 Graph Isomorphism

Two graphs are said to be isomorphic if they are syntactically similar to each other *iff* there is a bijection between their respective nodes which make each edge of  $G_A$  correspond to exactly one edge of  $G_B$ , and vice versa [12], i.e. the graphs are structurally the same to each other. This property is demonstrated in figures 5b.



(a) Isomorphic graph on the left



(b) Isomorphic graph on the right

Figure 5: A demonstration of two isomorphic graphs

The two graphs look very different, but when they are further inspected, it is evident that the two are a representation of the same structural scheme or maybe even the same graph that has been rearranged.

Consider the vertex 1 from the left-most graph, it has three edges going to and from vertices 2, 4 and 5, and now consider the vertex  $a$  from the right-most graph, it also has three vertices going to and from it, namely  $g, h$  and  $i$ . These two vertices have the same structural, but from two different graphs, the same goes for all the other vertices in both graphs .i.e each vertex in the left-most graph can be associated with one in the right-most graph as we did for vertices 1 and  $a$ .

## 2 Search Matching Algorithm

### 2.1 Introduction

Subgraph isomorphism can be determined using a brute-force approach on the tree representation of a graph  $G_A$ . Though this technique is effective, it is however not efficient, this is because all the possible permutation subgraphs of a graph  $G_A$  are tested against the graph  $G_B$  to determine if there are subgraphs in graph  $G_A$  that are isomorphic to graph  $G_B$ . The number of subgraphs of a graph  $G_A$  increase at an exponential rate with every addition of a vertex  $V_n$  into the graph, thus the total number of subgraphs that can be evaluated are

$$ST = 2^{n/2} \quad (3)$$

where  $ST$  is the total number of subgraphs and  $n$  the number of vertices in the graph  $G_A$ .

The matching process is computationally expensive due to this very fact, that is the more vertices there are in the graph  $G_A$ , the more expensive it becomes to detect the subgraph isomorphisms because of the amount of subgraphs it has and thus must be evaluated.

This paper explores two graphs that are very effective with regards to graph isomorphism and subgraph isomorphism detection. The algorithms that are investigated are the Ullman Algorithm and the VF2 algorithm.

## 2.2 Ullman Algorithm

### 2.2.1 Algorithm

The Ullman algorithm was developed by J.R.Ullman and was published in his paper titled "An Algorithm for Subgraph Isomorphism" [1]. The algorithm performs graph matching on an adjacency matrix representation of both the graphs, and uses the depth search first (DSF) recursive approach to traverse through the graphs and perform the graph matching process. The Ullman algorithm improves the efficiency of the brute-force approach at detecting subgraph isomorphisms by deductively eliminating nodes in the tree that are in graph  $G_A$ , but are not in graph  $G_B$ , thus reducing the number of subgraphs that are matched against graph  $G_B$  to determine isomorphism.

The algorithm starts by building a starting adjacency matrix  $M_0$  using the two adjacency matrix representations of graphs  $G_A$  and  $G_B$  using the following procedure.

- (1) Construct a  $n * m$  matrix where  $n$  is the number of rows of graph  $G_B$  and  $m$  is the number of columns of graph  $G_A$ .
- (2) Set all the entries in the matrix to the value of 1.
- (3) Apply the following rule: Set the values in  $M_0$  to 0 for all  $M_{0ij}$  where the degree of a vertex in graph  $G_A$  at  $j$  is greater than the degree of the same vertex in graph  $G_B$  .i.e.

$$deg(Ai) < deg(Bj). \quad (4)$$

A more formal representation of this rule is as follows

$$f(x) = \begin{cases} 1, & \text{if } deg(Ai) \geq deg(Bj) \\ 0, & \text{otherwise } \forall i,j \end{cases}$$

When the starting matrix has been constructed, we systematically permute matrices  $M^d$  from the starting matrix  $M_0$  where  $d$  represents the depth of the generated matrix. The procedure of generating the permuted matrices follows a depth search first (DSF) recursive approach where the stopping condition (leaf matrices) conform to the following form:

- (1)  $M$  contains only 0's and 1's.
- (2) There is exactly one 1 in each row.
- (3) Not more than one 1 in each column.

An demonstration of how the permutation matrices are generated is demonstrated in figure 6.



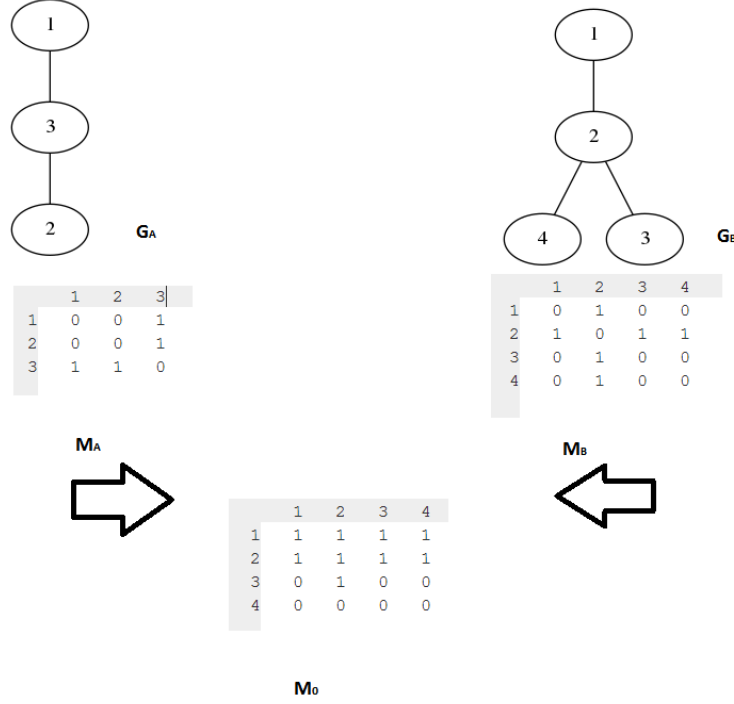


Figure 6: Demonstration of how a permutation matrix is generated from two graphs

Once all the permutation matrices have been generated, each one of the matrices is matched with a graph  $C$ , that is obtained from the dot product of the permuted matrix and the graph  $G_A$ . The formula for calculating graph  $C$  is follows:  $C = M_n(M_n \cdot G_A)^T$ , where  $G_A$  = input graph and  $M_n$  = permuted matrix  $M_n$  in  $M^d$ , obtained from the starting matrix  $M_0$   $(M_n \cdot G_A)^T$  = the transpose of the dot product of the permuted matrix  $M_n$  and the graph  $G_A$ . If there is a single instance of the matrix  $C$ , that is calculated using some permuted matrix  $M_n$  obtained from the starting matrix  $M_0$ , that is equal to matrix  $G_B$ , then  $G_B$  is isomorphic to  $G_A$ . Thus  $G_B$  is isomorphic to  $G_A$  iff  $G_B$

$$ij = 1 \rightarrow Cij = 1 \forall i, j \quad (5)$$

If non of the generated permuted matrices can satisfy this condition, then  $G_B$  is not isomorphic to  $G_A$ .

### 2.2.2 Ullman Pseudo code

```

Step 1   $M = M^0$ ,  $d = 1$ ,  $H_1 = 0$ ,
        for all  $i = 1, \dots, p_\alpha$  set  $F_i = 0$ ;
        refine  $M$ , if exit FAIL then terminate algorithm;
Step 2. If there is no value of  $j$  such that  $m_d = 1$  and  $f_j = 0$  then go to step 7,
         $M_d = M$ ;
        if  $d = 1$  then  $k = H_1$  else  $k = 0$ ;
Step 3   $k = k + 1$ ,
        if  $m_{dk} = 0$  or  $f_k = 1$  then go to step 3;
        for all  $j \neq k$  set  $m_{dj} = 0$ ;
        refine  $M$ ; if exit FAIL then go to step 5;
Step 4  If  $d < p_\alpha$  then go to step 6 else give output to indicate that an isomorphism has been found;
Step 5  If there is no  $j > k$  such that  $m_{dj} = 1$  and  $f_j = 0$  then go to step 7,
         $M = M_d$ ;
        go to step 3,
Step 6   $H_d = k$ ,  $F_k = 1$ ;  $d = d + 1$ ;
        go to step 2,
Step 7. If  $d = 1$  then terminate algorithm,
         $F_k = 0$ ;  $d = d - 1$ ;  $M = M_d$ ,  $k := H_d$ ;
        go to step 5;

```

Figure 7: Pseudo code of the Ullman algorithm [1]

## 2.3 VF2 Algorithm

### 2.3.1 Algorithm

The VF2 algorithm was introduced by L.P.Cordella, P.Foggiaa, C.Sansone and M.Vento [11]. The algorithm is suitable for graph matching and isomorphic determination, including subgraph isomorphic determination on large graphs, this is attributed to the Data structures that the algorithm uses and the manner in which they are used [11], this feature is discussed later in the paper.

The algorithm performs the matching process by attempting to find a mapping  $M$ , of vertices in graph  $G_A$  which correspond to vertices in graph  $G_B$ . The mapping is then used to determine if the two graphs are completely syntactically similar(isomorphic), partially syntactically similar or have no structural similarities at all.

#### 2.3.1.1 Matching and Mapping definition

A mapping  $M$  is defined as a set of pairs  $(n, m)$ , where  $n$  is a vertex from  $G_A$  and  $m$  a vertex from  $G_B$ , thus  $n \subseteq G_A$  and  $m \subseteq G_B$ .

The isomorphism determining properties of the mapping are defined as follows, a mapping  $M \subset N_A * N_B$  is isomorphic *iff*  $M$  is a bijection, that preserves the branching structure of  $G_A$  and  $G_B$ , where  $N_A$  is a set of vertices from  $G_A$  and  $N_B$  a set of vertices from  $G_B$ .

The mapping  $M \subset N_A * N_B$  is subgraph isomorphic *iff*  $M$  is isomorphic to  $G_A$  and a subgraph of  $G_B$ .

#### 2.3.1.2 Mapping Procedure

The mapping  $M$  comprises of state based partial solution morphisms  $M(s)$  for each state  $s$ . The process of finding the mapping  $M$  that is described above uses State Space Representations (SSR) [12]. The partial solution morphisms  $M(s)$  selects two subgraphs from  $G_A$  and  $G_B$ , namely  $G_A(s)$  and  $G_B(s)$  respectively. The subgraphs comprises of only vertices that are present in the partial solution  $M(s)$  for the state  $s$  as well as the edges joining them together.

The algorithm starts with an initial state  $s_0$  that has no mapping between the two graphs, thus  $M(s_0) = NULL$ . The algorithm then computes a set of candidate pairs  $P(s)$ . Each candidate  $p$  in the set is checked against the feasibility function that is discussed in the following chapter, if  $p$  is successful then it is added to the state  $s$ . And the successor  $s'$  is computed using a combination of the predecessor state and the candidate  $p$ , thus:

$$s' = s \cup p \tag{6}$$

The process of generating successor states is a recursive procedure that makes use of the depth first traversal for graphs. When a path has been exhausted and a solution has not yet been found, the algorithm uses backtracking to explore the alternative paths [11,13].

### 2.3.1.3 Definition of the set $P(s)$ and of the feasibility function $F(s, n, m)$

The VF2 algorithm generates the states with close consideration that only some of the states are consistant with the desired morphisms [12]. The algorithm avoids inconsistant states by making use of a set of rules in it's state generation procedure, thus ensuring that only consistant state are generated, these rules are refered to as feasibility rules.

The algorithm uses a function called a feasibility function to test that an additon of a pair  $(n, m)$  to a state will be consistant. If the addition of the pair passes all the feasibility rules, the algorithm will return a true value, if not, a false value indicating that the procedure results in an inconsistant successor state  $s'$ , and thus that state  $s'$  will not be explored by the algorithm.

A further filter can be applied in the consistent states to rule out those states whose successor states will be inconsistant, this apporoach is employed by adding a additional rules called *k-look-ahead* rules [12]. They check to see if the current state  $s$  will have a consistant successor state after  $k$  steps, i.e. they check to see if the states from  $s$  to  $s^k$  are consistant with the desired morphisms.

### 2.3.1.4 Condidate Pairs

The candidate pairs are obtained by considering the vertices that are connect to  $G_A(s)$  and  $G_B(s)$ , the sub-graphs of  $G_A$  and  $G_B$  in the state  $s$ . The vertexs are used to form the pairs  $(n, m)$  as defined above. In order explain how the pairs are formed, we must first introduce the following definitions: Let:

- (1)  $T_{Ain}(s)$  be the set of vertexs that are not yet in the partial mapping  $M(s)$  and are the origin of the edges from graph  $G_A$
- (2)  $T_{Bin}(s)$  be the set of vertexs that are not yet in the partial mapping  $M(s)$  and are the origin of the edges from graph  $G_B$
- (3)  $T_{Aout}(s)$  be the set of vertexs that are not yet in the partial mapping  $M(s)$  and are the destination of the edges from graph  $G_A$
- (4)  $T_{Bout}(s)$  be the set of vertexs that are not yet in the partial mapping  $M(s)$  and are the destination of the edges from graph  $G_B$

The pair  $(n, m)$  is made by vertex  $n$  from  $T_{Aout}(s)$  and  $m$  from  $T_{Bout}(s)$ . If the any of the sets is empty, then we consider the vertex  $n$  from  $T_{Ain}(s)$  and  $m$  from  $T_{Bin}(s)$ . In the case where that graphs are not connected, the pairs will be made by all the vertex not yet contained in either  $G_A(s)$  and  $G_B(s)$ . These pairs form the entries in the set  $P(s)$  for that respective state  $s$ .

### 2.3.1.5 The feasibility rules

The feasibility rules that are used to ensure that the states that are evaluated play a role in improving the performance, by preventing inconsistent states from being explored and thus optimizing the execution of the algorithm. There are five general feasibility rules defined as  $R_{pred}, R_{succ}, R_{in}, R_{out}$  and  $R_{new}$  respectively.

The feasibility functions check for two main things, firstly they check the consistency of the partial solution in the successor state  $s'$ , namely  $M(s')$ . Rules  $R_{\text{pred}}$  and  $R_{\text{succ}}$  are the rules used for this checking.

The remaining rules are used for pruning the search space for different levels of look ahead. The  $R_{\text{in}}$  and  $R_{\text{out}}$  are used to look ahead one level and determine which of those successor states are consistent, and  $R_{\text{new}}$  is used for the same purpose, but for a look ahead level of two. The definition for each rule is as follows:

$$\begin{aligned} R_{\text{pred}}(s, n, m) \iff & \\ & (\forall n' \in M1(s) \cap \text{Pred}(GB, n) \exists m' \in \text{Pred}(GA, m) | (n', m') \exists M(s) \wedge \\ & \forall m' \in M2(s) \cap \text{Pred}(GA, m) \exists n' \in \text{Pred}(GB, n) | (n', m') \exists M(s)), \quad (7) \end{aligned}$$

$$\begin{aligned} R_{\text{succ}}(s, n, m) \iff & \\ & (\forall n' \in M1(s) \cap \text{Succ}(GB, n) \exists m' \in \text{Succ}(GA, m) | (n', m') \exists M(s) \wedge \\ & \forall m' \in M2(s) \cap \text{Succ}(GA, m) \exists n' \in \text{Succ}(GB, n) | (n', m') \exists M(s)), \quad (8) \end{aligned}$$

$$\begin{aligned} R_{\text{in}}(s, n, m) \iff & \\ & (\text{Card}(\text{Succ}(GB, n) \cap \text{TBin}(s)) \geq \text{Card}(\text{Succ}(GA, m) \cap \text{TAin}(s))) \cap \\ & (\text{Card}(\text{Pred}(GB, n) \cap \text{TBin}(s)) \geq \text{Card}(\text{Pred}(GA, m) \cap \text{TAin}(s))) \quad (9) \end{aligned}$$

$$\begin{aligned} R_{\text{out}}(s, n, m) \iff & \\ & (\text{Card}(\text{Succ}(GB, n) \cap \text{TBout}(s)) \geq \text{Card}(\text{Succ}(GA, m) \cap \text{TAout}(s))) \cap \\ & (\text{Card}(\text{Pred}(GB, n) \cap \text{TBout}(s)) \geq \text{Card}(\text{Pred}(GA, m) \cap \text{TAout}(s))) \quad (10) \end{aligned}$$

$$\begin{aligned} R_{\text{new}}(s, n, m) \iff & \\ & \text{Card}(\tilde{N}A(s) \cap \text{Pred}(GB, n)) \geq \text{Card}(\tilde{N}B(s) \cap \text{Pred}(GA, n)) \wedge \\ & \text{Card}(\tilde{N}A(s) \cap \text{Succ}(GB, n)) \geq \text{Card}(\tilde{N}B(s) \cap \text{Succ}(GA, n)) \quad (11) \end{aligned}$$

In order for a state to be considered consistent, it must pass a combination of all of the five rules, namely:

$$F_{syn}(s, n, m) = R_{pred} \wedge R_{succ} \wedge R_{in} \wedge R_{out} \wedge R_{new} \quad (12)$$

where  $F_{syn}(s, n, m)$  the feasibility function that is invoked upon the state  $s$ .

### 2.3.2 VF2 Pseudo code

```

PROCEDURE Match( $s$ )
  INPUT:  an intermediate state  $s$ ; the initial state  $s_0$  has  $M(s_0)=\emptyset$ 
  OUTPUT: the mappings between the two graphs

  IF  $M(s)$  covers all the nodes of  $G_2$  THEN
    OUTPUT  $M(s)$ 
  ELSE
    Compute the set  $P(s)$  of the pairs candidate for inclusion in  $M(s)$ 
    FOREACH  $p$  in  $P(s)$ 
      IF the feasibility rules succeed for the inclusion of  $p$  in  $M(s)$  THEN
        Compute the state  $s'$  obtained by adding  $p$  to  $M(s)$ 
        CALL Match( $s'$ )
      END IF
    END FOREACH
    Restore data structures
  END IF
END PROCEDURE Match

```

Figure 8: Pseudo code of the VF2 algorithm [11]

## 2.4 Conclusion

The Ullman and VF2 algorithm that are discussed above both follow a similar approach in attempting to perform graph matching. They both construct a tree from the adjacency representation of their input graphs and use the depth first tree traversal techniques to evaluate the graphs, this is done very effectively by both the algorithms.

Though both algorithms are effective in their own respective regards, they are optimized rather differently and thus differ in their degree of complexity. The VF2 algorithm optimizes its execution by performing a look-ahead operation of two states from its current states in an attempt to ignore paths that will result in inconsistent states.

The Ullman algorithm on the other hand optimizes its execution by not computing all possible sub-graphs of some graph  $G$ , but reduces the computed matrices by initially computing a matrix  $M_0$  from the input graphs and then ensuring that all the matrices that are computed from  $M_0$  are tested so as to prevent the algorithm from exploring a branch that will not result in a graph or sub-graph isomorphism

## **3 VFLibGraph Library**

### **3.1 Introduction**

The VFLibGraph library is a graph matching library that is written in C++, it was developed at the Intelligent Systems and Artificial Vision Lab (SIVALab) which is situated in the University of Naples [13]. The library was originally developed to test the VF algorithm [13], the predecessor of the VF2 algorithm, but the library has since evolved to include some of the latest graph matching algorithms such as the Schmidt and Druffel algorithm.

### **3.2 Using the library**

The library provides interfaces to the implemented algorithms that it has in its employ, and it also has a comprehensive documentation of how each algorithms interface can be constructed, as well as the matching process between two input graphs.

We have constructed interfaces for the Ullman algorithm and the VF2 algorithm to test against our data set, which comprises on graphs of various number of vertices and edges.

## 4 Experimentation

### 4.1 Data Set

The graph data that is used in the experimentation was created by David S. Johnson, Cecilia R. Aragon, Lyle A. McGeoch and Catherine Schevon [14]. It was created in 1991 to study optimization by simulated annealing. One of the graphs that used was chosen to be experimented on to study the subject algorithms, namely the Ullman and VF2 algorithms. The graph used can be found here [14].

The graph is a large direct graph that comprises of 250 vertices and 3218 edges [14], and it is used as the super-graph in the experiments, and will be referred to as the subject-graph from here on out. Figure 9 depicts a diagram of how the graph looks like.

The algorithms both need two graphs in order to determine if they are isomorphically related to each other or not, thus subgraphs from our object graph are required for this comparison. These subgraphs are then generated from the object graph and they all meet the condition that:

- They are either partially or completely isomorphic in relation to one or other digraph inside the set.

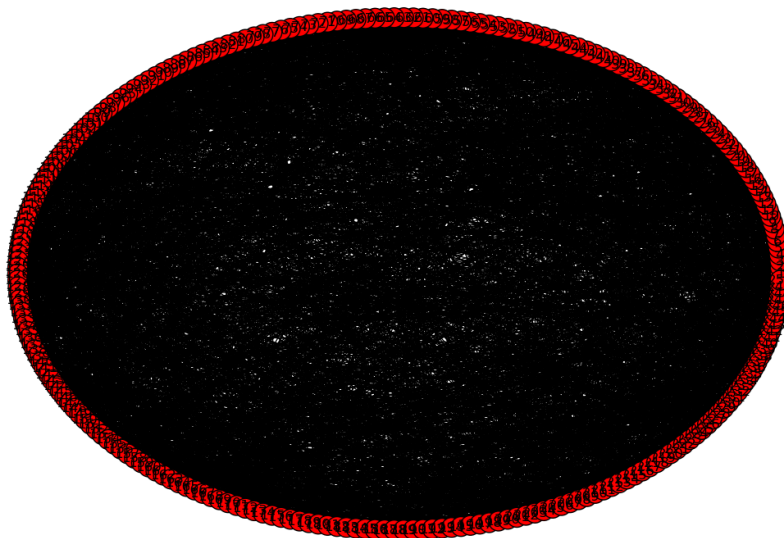


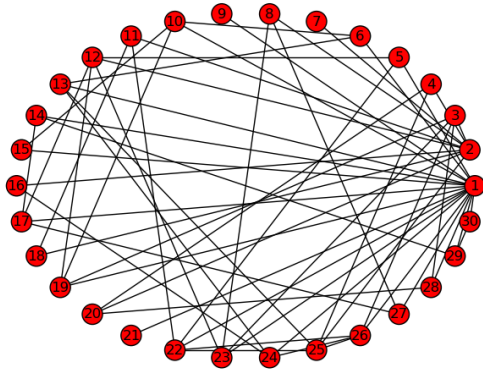
Figure 9: The super-graph used in the experiment comprising of 250 vertices



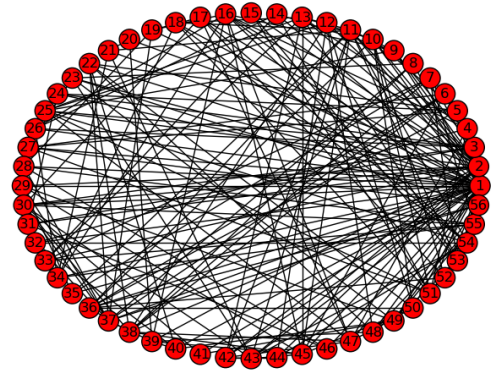
The subgraphs will be referred to as subject-graphs from here on out.

The generated subgraphs have the following number of vertices: 30,56,75,92,109,121,148,166,181,197,211 and 222. Each one of the graphs is sub-graph isomorphic to the object graph, and will be compared with it to determine how long it takes to find the isomorphic relationship, and how much virtual memory [15] is used to find the relationship.

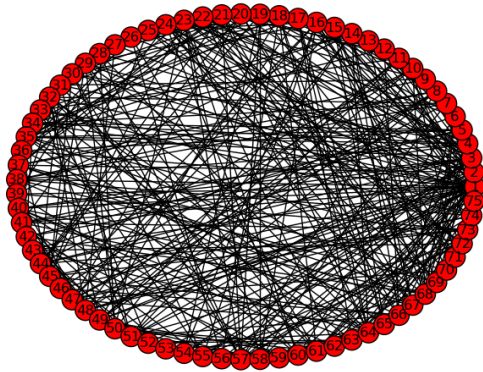
Figure 10 depicts how some of the generated subgraphs look relative to figure Figure 9.



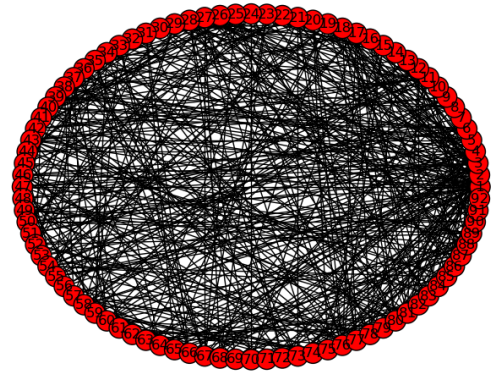
(a) Generated subgraph with 30 vertices



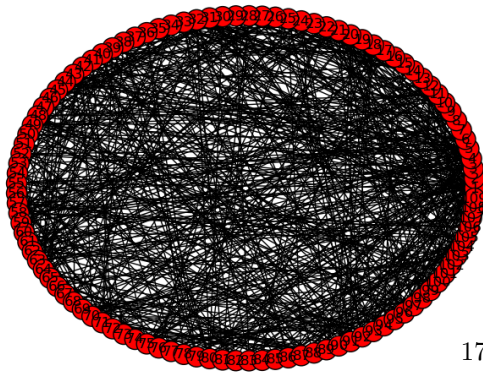
(b) Generated subgraph with 56 vertices



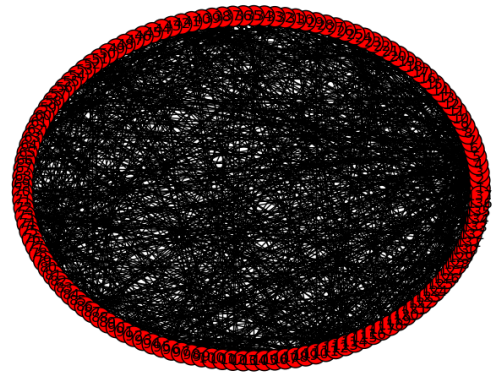
(c) Generated subgraph with 75 vertices



(d) Generated subgraph with 92 vertices



(e) Generated subgraph with 109 vertices



(f) Generated subgraph with 121 vertices

Figure 10: A demonstration of two isomorphic graphs

## 4.2 Experiment

The experiment that is conducted on the Ullman [1] and the VF2 [11] algorithm uses a quantitative approach that evaluates the algorithms against each other. The search graph matching algorithms are implemented and experimented on, using two digraph or disjoint objects at a time. The experiment for that is employed, measured and records the algorithms ability and performance of the algorithms in cases where Graph  $G_A$  and Graph  $G_B$  are either partially or completely isomorphic.

Several experiments are done on the two algorithms so that we can understand their behaviour in different conditions, and these experiments that are conducted are done on both joint and disjoint graphs. The experiments that are performed are explained below.

### 4.2.1 The comparison criteria

The evaluation performed on the algorithm results is based on two criteria's. These are use to weigh the algorithms against each other to determine which amongst them is better. The criteria's are listed below.

**Space efficiency:** This factor measures the amount of virtual memory (RAM) that is used by the algorithm in the graph matching process of its execution. The unit used to measure the amount of memory by the respective algorithms is measured in *bytes (b)*.

**Time efficiency:** This factor measures the time taken by the algorithm to start and *successfully* complete the graph matching process of its execution. The unit used to measure the amount of time taken by the respective algorithms is measured in *millisecond (ms)*.

### 4.2.2 Description Conducted Experiments

This section lists each of the experiments that where conducted for both algorithms as well as provides an explains of each experiment.

#### 4.2.2.1 Performance evaluation for different vertice numbers

This experiment uses data that is mentioned in chapter 4.1. The experiment is conducted on joint and disjoint graphs. This experiment is intended on evaluating the behaviour of the two algorithms performance for graphs of different vertice numbers. The number of the vertices in the graphs range from 55 to 250 vertices in the graphs.

#### 4.2.2.2 Performance evaluation for different edge numbers

This experiment uses generated graph data comprising of 1000 vertices and 499500 edges. The generated graph data is defined using the following rules.

- (1) Define a set of vertices  $V$  for some graph  $G$ .
- (2) Each vertice  $v$  in  $V$  is connect to every other vertice in  $V$
- (3) Non of the edges in the graph  $G$  are reflexive.

This experiment is intended on evaluating the behaviour of the algorithms for when they are searching for subgraphs of various sizes proportional to some supergraph  $G_{SUP}$ .

The subgraphs are generated from the supergraph  $G_{SUP}$  for different percentages e.g. a generated subgraph that is 10% of the supergraph  $G_{SUP}$  graph.

### 4.2.3 Experiment implementation

The experimentation process is accomplished in phases and each phase is explained below. We will refer to Graphs  $G_A$  and  $G_B$  in this context.

#### 4.2.3.1 Generate two graphs, $G_A$ and $G_B$

In the first phase of the experimentation process, a supergraph  $G_A$  and a subgraph  $G_B$  are generated. The two graphs are built together with their associate vertices and edges.

#### 4.2.3.2 Syntactic comparison

The comparison that is done, is dependent on the relationship between Graph  $G_A$  and  $G_B$ .

- (1)  $G_A$  is compared with the whole of  $G_B$  for syntactical similarity.
  - (I) A new subgraph of Graph B is generated.
  - (II) The subgraph is compared with Graph  $G_A$  for syntactical similarity.
    - (A) The time taken to perform the comparison is recorded.
    - (B) The amount of memory used by the algorithm to perform the comparison is recorded.
- (2) If Graph  $G_A$  and  $G_B$  are not completely syntactically similar, Graph  $G_A$  is compared with all the possible subgraphs of Graph B.

### 4.2.4 Result presevation

The amount of time and memory required by both alogrithm to complete the comparison procedure of each respective experiment is stored and graphically represented. The results from the algorithms are weighed against a comparison criteria, and it is this criteria that is used to evaluate the algorithms against each other.

Once the results have been obtained and throughly evaluated, then the algorithms with the best perform per criteria field are recorded.

And from that set, the best algorithm is chosen overall relative to the others based on the criteria.

## 5 Experiment Results

This section presents the results of the experiments described in Chapter 4.2.2. An evaluation of the results is performed before the algorithms can be compared against each other based on the criteria specified in chapter 4.2.1, namely the time and space efficiency of the algorithms.

### 5.1 Joined graph results for the evaluation of different vertex numbers experiment

This section presents the time and memory results for experiment described in section 4.2.2.1. The experiment is conducted on joined graphs.

#### 5.1.1 Memory Results

This section compares the efficiency of the two algorithms in terms of their memory. The amount of virtual memory used by the two algorithms are graphically represented, analysed and a conclusion of as to which of the two algorithm is more efficient in terms of memory is reached based on figure 11.

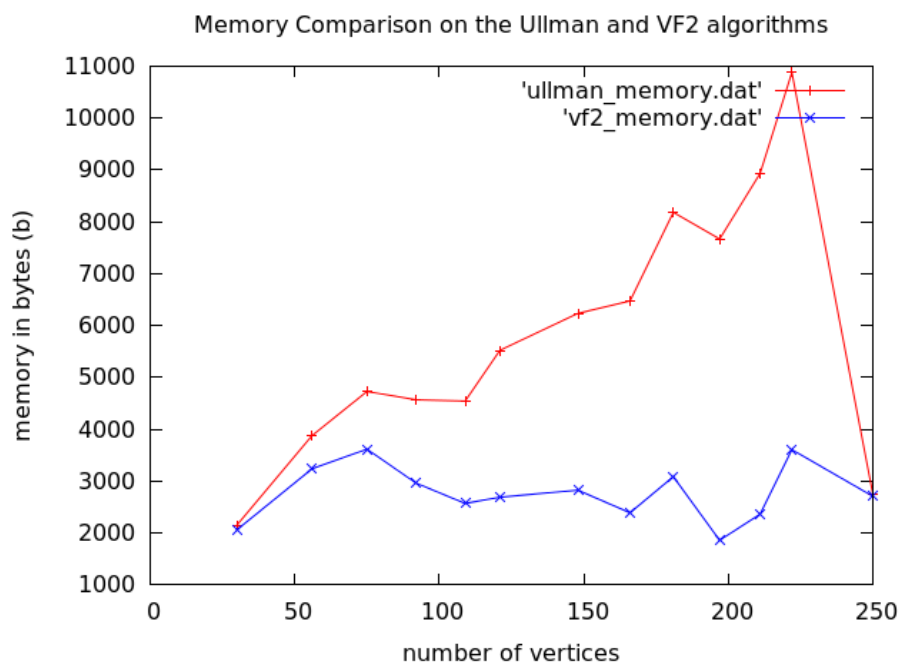


Figure 11: Graph depicting the results for the memory comparison between the algorithms for joined graphs

#### 5.1.1.1 Analysis of results

Figure 11 depicts the comparison of the Ullman and the VF2 algorithms on the criteria of memory used during the execution of the graph matching processes of the respective algorithms.

The virtual memory used by the the Ullman algorithm is depicted by the *red* line in figure 11 and the *blue* line represents the virtual memory used by the VF2 algorithm.

Figure 11 depicts that the memory usage of the two algorithms is approximately the same for smaller graphs of 30 - 40 vertices. But from approximately 40 and above, it is clear that the performance of the VF2 algorithm is exceptionally superior to that of the Ullman algorithm. This deduction is derived from the observation on figure 11, we can see the more vertices there are in the graph, the more virtual memory resources is required by both graphs, but the Ullman algorithm requires a greater amount of resources than those used by the VF2 algorithm.

The superiority of the VF2 algorithm for this data set is limited, this is because when the number of vertices in the graph is approximately 220, the required memory resources for both algorithms start dropping rapidly, and they finally end up requiring the same amount of memory resources.

#### 5.1.1.2 Conclusion

Figure 11 has provided with a graphical representation of the virtual memory required by both algorithms, and thus a way to make deductions about the efficiency of the two algorithms in terms of memory.

From 11, it is clear that the VF2 is algorithm is more memory efficient than the Ullman algorithm as it requires the least amount of virtual memory overall for its execution.

#### 5.1.2 Time Results

This section compares the efficiency of the two algorithms in terms of their time. The amount of taken by the two algorithms are graphically represented, analysed and a conclusion of as to which of the two algorithm is more efficient in terms of taken is reached based on figure 14.

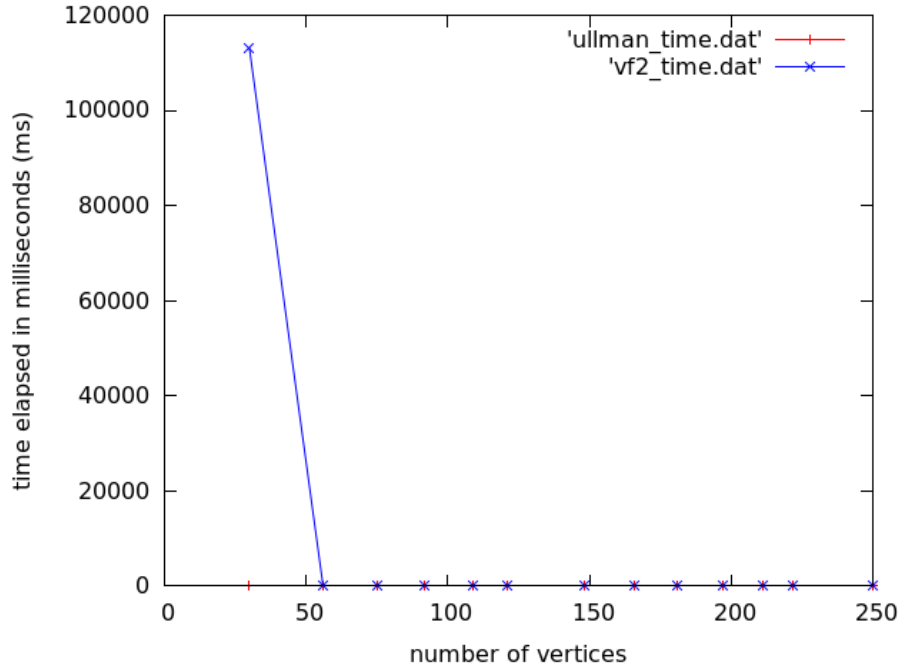


Figure 12: Graph depicting the results for the time comparison between the algorithms for a set of joined graphs

#### 5.1.2.1 Analysis of results

Figure 14 depicts the comparison of the Ullman and the VF2 algorithms on the criteria of the time taken during the execution of the graph matching processes of the respective algorithms.

The amount of time taken by the Ullman algorithm is depicted by the *blue* line in figure 14, and the time taken by VF2 algorithm is depicted by the *red* line in figure 14.

Figure 14 depicts that the amount of time taken by the Ullman algorithm to successfully complete its graph matching processes is very small and consistent for the data set used in the experimentation.

The time taken by the VF2 algorithm however is very high for small sets of vertices, approximately 40 - 60 vertices. But as the number of vertices increase, the amount of time declines rapidly and it equal to that of the Ullman algorithm.

#### 5.1.2.2 Conclusion

Based on figure 14, it can be deduced that the time efficiency of the Ullman and VF2 algorithms are approximately equal to each other overall, with the exception that for small graphs, the Ullman

algorithm is more efficient than the VF2 algorithm in terms of the time taken to complete the graph matching procedures.



## 5.2 Disjoined graph results for the evaluation of different vertex numbers experiment

This section presents the time and memory results for experiment described in section 4.2.2.1. The experiment is conducted on disjoined graphs.

### 5.2.1 Memory Results

This section depicts the memory results for the two algorithms on a set disjoined graphs of various vertex numbers.

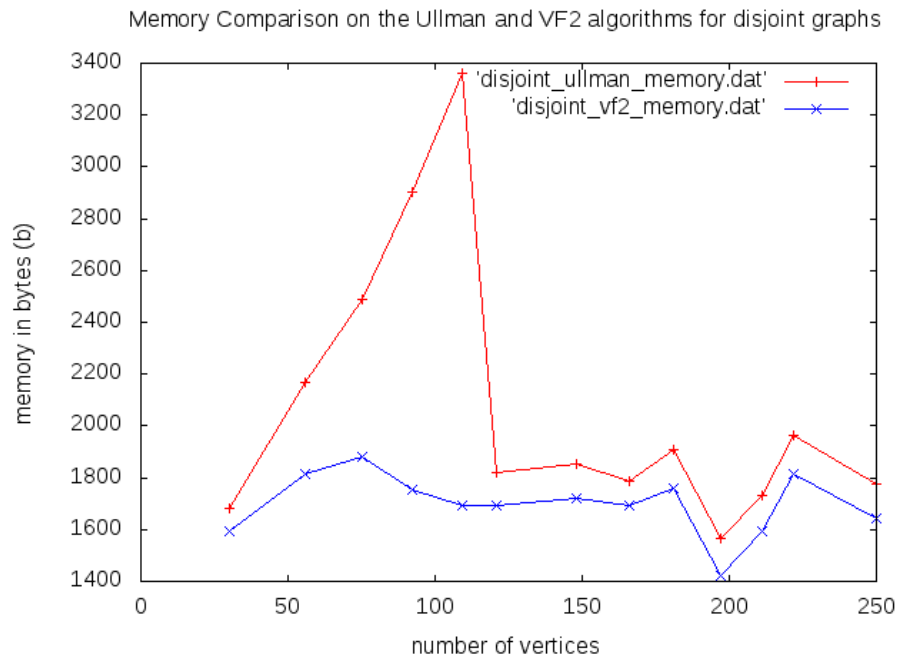


Figure 13: Graph depicting the results for the memory comparison between the algorithms for a set of disjoint graphs

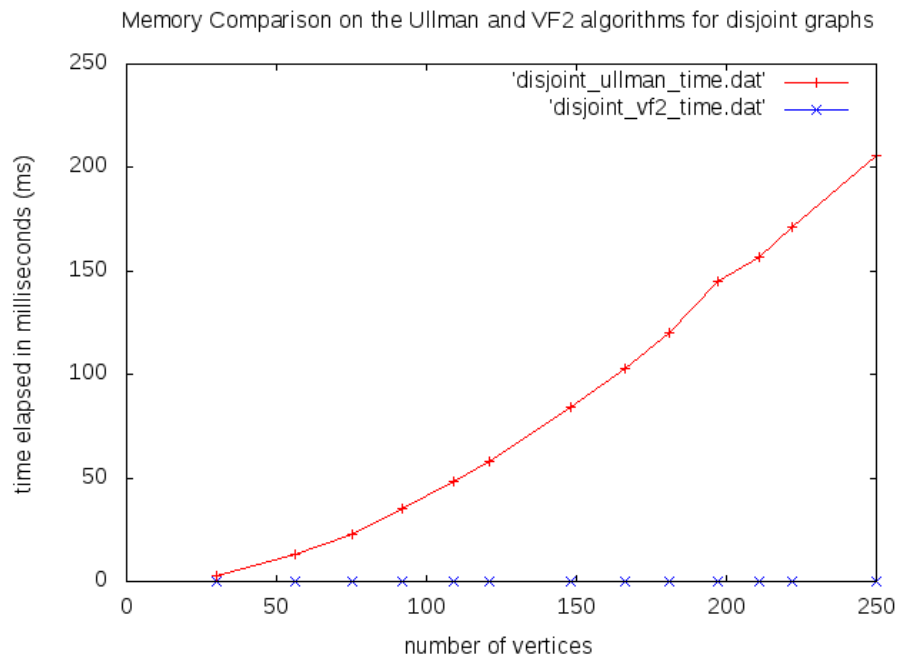


Figure 14: Graph depicting the results for the time comparison between the algorithms for a set of disjoint graphs