A New Method for Subgraph Detection - SubGraD

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ABSTRACT

A new method of subgraph detection is introduced in this paper, this new approach is SubGraD. Firstly consider the model graph (query graph) and make the possible sets called model sets starting from the chosen initial node or starter. Similarly, for the source graph (reference graph), all the possible sets called reference sets could be made. Our aim is to make the reference set on the basis of the model set. If it is possible to make the reference set, then it is said that query graph has been detected in the source graph.

KEYWORDS

model graph, reference graph, starter, model set, reference set

1. Introduction

Graphs are data structures which are proved to be effective way of representing objects [1]. One of the problems of interest, with graphs, is matching a model graph (query graph) against a reference graph (source graph). Graph matching is the process of finding a correspondence between the nodes and the edges of two graphs that satisfies some (more or less stringent) constraints ensuring that similar substructures in one graph are mapped to similar substructures in the other [2]. Many graph databases are growing rapidly in size. The growth is both in the number of graphs and the sizes of graphs (the number of nodes and the number of edges). There is a critical need for efficient and effective graph querying tools for querying and mining these growing graph databases.

We are introducing a new method called SubGraD, which is helpful in graph (directed graph) mining. Here we take the model graph first and make the model sets. Suppose that there are two nodes in the model graph (query graph), and then corresponding model set will have only two elements. Then we break the reference graph in such a way that each set of reference set will have sets of two elements. We are assuming that the smallest model set is the set which has two DOI: 10.5121/ijccsa.2012.2403

elements corresponding to two nodes of the query graph. We are not considering the case where the self loop is present. In this paper, we are trying to make sets for reference set by using the SubGraD algorithm. If reference set could be made, then query graph is present in reference graph otherwise not. Related works are discussed in section 2. In section 3, SubGraD algorithm is given. Model sets have been made in section 4. And in section 5, it is tried to make Reference sets so that we could find out whether the query graph is present on the source graph or not.

2. RELATED WORK

Graph matching is the concept that has been intensively used in various applications such as in the field of social networks, road networks, network topology, chemical structures, graph grammars and semantic networks or pattern recognition. Some of the related works associated with different fields for graph mining is discussed in this section.

In Fuzzy set theory, there are so many works proposed for graph matching, for example Perchant and I. Bloch [3-6] gave methods for graph matching using fuzzy set theory, and Hwan gave an idea for sub-graph matching [7]. Fuzzy attributed graph models are suggested for very different image type representations, such as fingerprint verification [8].

Fernandez and Valiente proposed a method to represent attributed relational graphs, the maximum common subgraph and the minimum common super graph of two graphs by means of simple structures [9]. Bunke [10] suggested the relationship between graph edit distance and the maximum common subgraph, graph edit distance computation is equivalent to solving the maximum common subgraph problem.

Cross [11] described an outline for performing relational graph matching using genetic algorithm. He showed that Bayesian consistency measure could be proficiently optimized using a hybrid genetic search process that includes a local search strategy using a hill-climbing step. This hill-climbing step accelerates convergence considerably. Cross extended this idea in [12] that is also a convergence analysis for the problem of attributed graph matching using genetic search.

Myers and Hancock gave an idea that in attributed graph matching problems usually there are more than one valid and satisfactory solution, and they proposed a method to obtain different solutions at the same time during the genetic search, using an appropriately modified genetic algorithm [13-14].

Hancock and Kittler [15-16] used an iterative approach called probabilistic relaxation, and considering binary relations and assuming a Gaussian error. Later a Bayesian perspective was considered for both unary and binary attributes by Christmas [17], Gold and Rangarajan [18], Wilson and Hancock [19-20]. Williams [21] presented a comparative study of various deterministic discrete search-strategies for graph-matching, which was based on the previous Bayesian consistency measure in [19-20] and Tabu search was proposed as a graph matching algorithm.

3. SUBGRAD ALGORITHM

Graph Matching techniques are important and very general form of pattern matching that finds

realistic use in areas such as image processing, pattern recognition and computer vision, graph grammars, graph transformation, bio computing, search operation in chemical structural formulae database, etc. SubGraD Algorithm is one such graph matching method. SubGraD works like a graph decomposition algorithm. In which source graph is decomposed. The main idea of decomposing the directed graph is that, the smaller the graph structure, the easier the matching process and this will result in lower complexity than matching the original graph.

In SubGraD method, an adjacency matrix [A] is drawn for a query graph (model graph). Element A_{ij} of [A] is 1 if an edge is present from node i to node j else 0. Using [A], model set M is prepared. Model set M is a set, consisting all sets (i, j) as its element where A_{ij} is 1. On the basis of model set we will try to make the sets for reference set.

SubGraD

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Make the adjacency matrix [A] of n \times n
Make Model Set M as,

For i: 1 to n
For j: 1 to n
M = (i, j), where A_{ij} = 1
END FOR j
END FOR i
Make the adjacency matrix [S] of m \times m
Try to make the reference set R corresponding to Model Set M
Try to break adjacency matrix [S] into k sets R_1, R_2, \ldots, R_k, where S_{ij} = 1 (row i and column j has entry 1) and each R_k has the same number of elements as M has R = (R_1, R_2, \ldots, R_k)
If reference R set can be made, then query graph is detected in source graph Otherwise query graph doesn't exist
END SubGraD
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4. MODEL SETS

Suppose we have the query graphs as shown in Figure 1, and our purpose is to detect these query graphs in a given source graph. First step will be to draw the adjacency matrices for query graphs. In first query graph (Figure 1(a)), there are two nodes a and b, hence Adjacency matrix [A] will be 2×2 . A single edge is present from a to b so the adjacency matrix [A] will have A_{ab} entry as 1 and rest of the elements (A_{aa} , A_{ba} , A_{bb}) as 0. Likewise, the adjacency matrices Figure 2(b), 2(c) and 2(d) are drawn for query graphs Figure 1(b), 1(c) and 1(d) respectively.

Next step is to prepare Model sets using these adjacency matrices. The adjacency matrix of Figure 1(a) has four elements A_{aa} , A_{ab} , A_{ba} , A_{bb} (Figure 2(a)). The model set corresponding to this adjacency matrix is (a, b), since only A_{ab} entry is 1. Similarly for query graph (Figure 1(b)), the model set is (a, b) and (b, c), since there are two entries which are 1 in adjacency matrix (i.e. Figure 2(b)). Model sets for query graphs of Figure 1, is shown in TABLE 1.

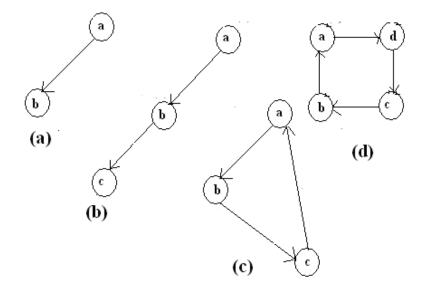


Figure 1. Query graphs

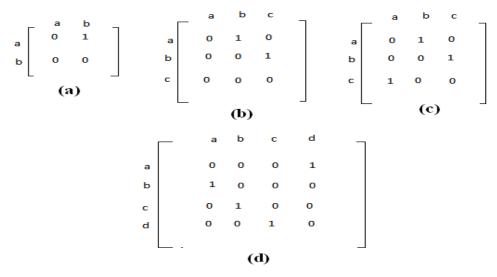


Figure 2. Adjacency matrices for query graphs

Table 1. Model Sets for Query Graphs.

Model set for Query Graph Figure 1(a)	Model set for Query Graph Figure1(b)	Model set for Query Graph Figure 1(c)	Model set for Query Graph Figure 1(d)
(a, b)	(a, b) (b, c)	(a, b) (b, c)(c, a)	(a, d) (b, a) (c, b) (d, c) Or
			(a, d) (d, c) (c, b), (b, a)

MODEL SET I MODEL SET II MODEL SET III MODEL SET IV

5. REFERENCE SETS

Let us consider the source graph present in Figure 3, now we will try to make reference set for source graph. Firstly select the Model set, we are taking MODEL SET I (from Table 1), there is only an edge that is present from node a to node b, so we are applying SubGraD algorithm, which selects the set of row i and column j, where the entry is 1 in adjacency matrix [S], of source graph (i.e. Figure 4). That means we are trying to make reference set R which consists of set $R_1,R_2,...,R_n$. In source graph for row 1, and for column 2, 3, 4 and 6 the entries are 1, so the reference set for row 1 is (1,2),(1,3),(1,4) (1,6). Similarly for row 2, the reference set is (2,4),(2,6), for row 3, the reference set is (3,5), for row 4, the reference set is (4,6), for row 5, the reference set is (5,1) and for row 6, the reference set is (6,5). Corresponding to MODEL SET I, the Reference set R I (Table 2) have been found. Hence the query graph is detected in source graph.

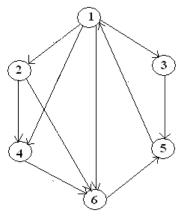


Figure 3. Source graph

	1	2	3	4	5	6
1 2 3 4 5 6	0 0 0 0 1	1 0 0 0 0	1 0 0 0 0	1 1 0 0 0	0 0 1 0 0	1 1 0 1 0 0
0	0	U	U	U	1	U

Figure 4. Adjacency matrix of Source graph

Now suppose we want to find out whether query graph (Figure 1 (b)) is present or not, take the MODEL SET II from TABLE 1, i.e. (a, b), (b, c). In MODEL SET II we can see that element b is presented twice in the set. This type of element is called middle element. Actually, it is the node which has an incoming edge as well as an outgoing edge (in our case the node b has an incoming edge from node a and an outgoing edge to node c from Figure 1(b)). Now from source graph we have to make such a reference set $R = (R_1, R_2)$, where R_1 should have the set (i, k) and R_2 should have set (k, j). Here element k is the middle element. Firstly we start from row 1 in source graph

(Figure 4), S_{12} , S_{13} , S_{14} and S_{16} are equal to 1, that means entries from row 1 to column 2, 3, 4 and 6 is 1. So next we will take element 2, 3, 4 and 6 as middle element, and will now consider row 2, row 3, row 4 and row 6. From adjacency matrix [S] (Figure 4), we will see that which Sij entry is 1, here i is equal to 2, 3, 4 and 6. For row 2, S_{24} and S_{26} elements are 1. So in consequence there are two elements, i.e. S_{12} and S_{24} , which are 1. So the reference set R is made which has an element (1, 2), (2, 4), where 2 is the middle element. Similarly other elements of the reference set can be made. So reference set for row 1 is (1, 2)(2, 4), (1, 2)(2, 6), (1, 3)(3, 5), (1, 4)(4, 6), (1, 6)(6, 5). For row 2, the reference set is (2, 4)(4, 6), (2, 6)(6, 5). For row 3, the reference set is (3, 5)(5, 1). For row 4, the reference set is (4, 6)(6, 5). For row 5, the reference set is (5, 1)(1, 2), (5, 1)(1, 3), (5, 1)(1, 4), (5, 1)(1, 6). For row 6, the reference set is (6, 5)(5, 1). For row 1, 2, 3, 4, 5 and 6 we can make the reference set for MODEL SET II, hence the query graph Figure 1(b) exists in the source graph at all nodes. Reference set II for MODEL SET II can be seen by Table 2.

Now suppose we want to find out whether query graph (Figure 1 (c)) is present or not, take the MODEL SET III from Table 1, i.e. (a, b) (b, c)(c, a). In MODEL SET III, we can see that there are middle elements b and c, and a cycle is formed. Cycle formation means that the starting node is also the ending node. It should be noted that the order of the set is not important, it means that, (a, b) (b, c) (c, a) is same as (a, b) (c, a) (b, c). But the former one satisfies the definition of cycle formation, so we should try to write in the order as the former one has. From adjacency matrix S_{ij} , we could see that, S_{13} , S_{35} , S_{51} , S_{16} , S_{65} , S_{51} entries are 1. Reference set for row 1, is (1, 3)(3, 5)(5, 1) and (1, 6)(6, 5)(5, 1). Hence for node 1, the query graph has been detected twice. Elements 1, 3 and 5 corresponds to the elements a, b, c respectively, which forms a cycle. The null entry in Table 2 denotes that for that particular node, the query graph could not be detected. It could be seen from Table 2, for node 2 there is no reference set for MODEL SET III. Hence the query graph (Figure 1 (c)), is not present at node 2 of the source graph. Similarly for node 4 of source graph, this query graph cannot be detected.

For the query graph (Figure 1 (d)), again the cycle is formed. But here the model set M is consists of four sets. So now we will try to make reference set R which should have R_1 , R_2 , R_3 and R_4 in such a way that they should form a cycle. For example from adjacency matrix [S], it could be seen that S_{46} , S_{65} , S_{51} and S_{14} are 1, and they form a cycle. Hence query graph is detected at node 4. But at node 2 and node 3 it is not present. Hence if we could not able to make the reference set for particular query graph, then query graph could not be detected in the source graph.

Nodes	Reference Set I for MODEL SET I	Reference Set II for MODEL SET II	Reference Set III for MODEL SET III	Reference Set IV for MODEL SET IV
1	i.(1, 2), ii.(1, 3), iii.(1,4), iv.(1,6)	i.(1, 2)(2, 4) (1, 2)(2, 6) ii.(1, 3)(3, 5) iii.(1, 4)(4, 6) iv.(1, 6)(6, 5)	i.(1,3)(3,5)(5,1) ii.(1,6)(6,5)(5,1)	i.(1,2)(2,6)(6,5)(5,1) ii.(1,4)(4,6)(6,5)(5,1)
2	i.(2, 4),	i.(2,4)(4,6)	-	-

Table 2. Reference sets for source graph.

	ii.(2, 6)	ii.(2,6)(6,5)		
3	(3,5)	(3,5)(5,1)	(3,5)(5,1)(1,3)	-
4	(4,6)	(4,6)(6,5)	-	(4,6)(6,5)(5,1)(1,4)
5	(5,1)	i.(5,1)(1,2) ii.(5,1)(1,3) iii.(5,1)(1,4) iv.(5,1)(1,6)	i.(5,1)(1,3)(3,5) ii.(5,1)(1,6)(6,5)	(5,1)(1,2)(2,6)(6,5) same as (1,2)(2,6)(6,5)(5,1)
6	(6,5)	(6,5)(5,1)	(6,5)(5,1)(1,6) same as (1,6)(6,5)(5,1)	(6,5)(5,1)(1,2)(2,6) same as (1,2)(2,6)(6,5)(5,1)

5. CONCLUSION

In this paper we have introduced an efficient approach for graph matching SubGraD. We first make the model set M for query graphs with the help of adjacency matrix and then corresponding to that model set, we try to make reference set R for source graph. If reference set could be made, then query graph is detected in source graph otherwise it does not exist in the source graph. Actually in this paper, we are simply decomposing the source graph. Making the reference set means is that we are decomposing the source graph, in the sets, which has the same number of elements as in the model set.

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