

Vertex Cover Optimization Using a Novel Graph Decomposition Approach

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Received: 10 January 2022; Accepted: 30 March 2022

Abstract: The minimum vertex cover problem (MVCP) is a well-known combinatorial optimization problem of graph theory. The MVCP is an NP (nondeterministic polynomial) complete problem and it has an exponential growing complexity with respect to the size of a graph. No algorithm exists till date that can exactly solve the problem in a deterministic polynomial time scale. However, several algorithms are proposed that solve the problem approximately in a short polynomial time scale. Such algorithms are useful for large size graphs, for which exact solution of MVCP is impossible with current computational resources. The MVCP has a wide range of applications in the fields like bioinformatics, biochemistry, circuit design, electrical engineering, data aggregation, networking, internet traffic monitoring, pattern recognition, marketing and franchising etc. This work aims to solve the MVCP approximately by a novel graph decomposition approach. The decomposition of the graph yields a subgraph that contains edges shared by triangular edge structures. A subgraph is covered to yield a subgraph that forms one or more Hamiltonian cycles or paths. In order to reduce complexity of the algorithm a new strategy is also proposed. The reduction strategy can be used for any algorithm solving MVCP. Based on the graph decomposition and the reduction strategy, two algorithms are formulated to approximately solve the MVCP. These algorithms are tested using well known standard benchmark graphs. The key feature of the results is a good approximate error ratio and improvement in optimum vertex cover values for few graphs.

Keywords: Combinatorial optimization; graph theory; minimum vertex cover problem; maximum independent set; maximum degree greedy approach; approximation algorithms; benchmark instances

1 Introduction

The Minimum Vertex Cover Problem (MVCP) is a subset of NP complete problems. Solution of NP class of problems is one of the seven outstanding millennium problems stated by the Clay Mathematics institute. The solution of these problems can be verified in polynomial time scale, but time complexity for solving these problems grow exponentially with size of the problems [1]. The MVCP



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involves finding a set U such that $U \subset V$. Here the set U has the smallest possible cardinality in a graph $G = G(V, E)$ such that V is a set of vertices and E is a set of edges of the graph. For the set U to be a cover of graph, every edge of the graph is connected to at least one element of U . The set U is called a minimum vertex cover of G [2]. The problem has an exponentially growing complexity since the number of combinations which are required to be verified grows as $n_v!$ where n_v represents the number of vertices in the graph. Due to exponential growth in complexity of the problem, it is almost impossible to exactly solve the problem in a realistic time scale. Therefore, solving these problems via Brute force method i.e., checking all the possible combinations is not feasible. However, one may opt for an approximate solution of these problems in a reasonably quick time.

The MVCP has a wide range of applications; for example, cyber security, setting up or dismantling of a network, circuit design, biochemistry, bioinformatics, electrical engineering, data aggregation, immunization strategies in network, network security, internet traffic monitoring, wireless network design, network source location problem, marketing and franchising, pattern recognition and cellular phone networking [3–9].

Due to its wide range of applications, the MVCP has received special attention in the scientific community. Several approximate algorithms for solving the problem have been proposed, e.g., the depth first search algorithm, the maximum degree greedy algorithm, the edge weighting algorithm, the deterministic distributed algorithm, the genetic algorithm, the edge deletion algorithm, the support ratio algorithm, the list left algorithm, the list right algorithm and iterated local search algorithm etc [10]. Since all these algorithms provide approximate results with certain accuracy, there is a certain space to improve accuracy and to reduce complexity by introducing faster and more accurate algorithms. The scientific community around the globe has proposed approximate solutions of the problem with polynomial complexity. Some of the efforts by scientific community are described in following paragraph.

Jiaki Gu et al. proposed an algorithm that uses a general three stage strategy to solve the minimum vertex cover problem. Their method includes graph reduction, finding minimum vertex cover of bipartite graph components and finally finding the vertex cover of actual graph [11]. Changsheng Quan and coworkers proposed an edge waiting algorithm to solve MVCP. They claim that their algorithm has a fast-searching performance for solving large-scale real-world problem [12]. Shaowei Cai et al. in their work proposed a heuristic algorithm that make use of a preprocessing algorithm, construction algorithms and search algorithms to solve the MVCP. They claim that their algorithm is fast and accurate as compared to other existing heuristic algorithms [13]. Chuan Luo et al. proposed an algorithm that uses a highly parametric framework and incorporates many effective local search techniques to solve the MVCP. According to their claim their algorithm performs better for medium size graph and is competitive for large sized graphs [14].

Jinkun Chen and coworkers proposed an approximate algorithm based on rough sets. They use a Boolean function with conjunction and disjunction logics [15]. Cai. S et al. in their work use an edge weighting local search technique for finding an approximate MVC [16]. Khan, I and coworkers proposed an algorithm that works by removal of nodes to find a maximum independent set yielding an approximate MVC [17]. Arstrand, M et al. formulated a deterministic distributed algorithm to solve the MVCP. The authors solved the problem for two approximate solutions of the MVC during $(\Delta + 1)^2$ synchronous communication rounds where Δ represents an upper bound of maximum degree [18]. Bar-Yehuda et al. used Dijkstra algorithm in their work in order to solve the problem [19]. Genetic algorithm has been used for the solution of the problem by Bhasin et al. Their algorithm demonstrated advantage of handling graphs when compared to the reported literature algorithms. The

authors mentioned that the algorithm is unable to tackle some problems due to which they proposed the usage of Diploid Genetic Algorithms as an extension [20]. Support Ratio Algorithm (SRA) used a heuristic approach to solve the MVCP in which Balaji and coworkers used an adjacency binary matrix to represent a graph. The complexity of the algorithm has been $O(n_e n_v^2)$, where n_e is number of edges and n_v is the number of vertices. The authors claim that the support ratio algorithm has been found better for large scale problems compared to the reported algorithms [21]. Kettani and co-workers introduced a novel heuristic algorithm to find MVC. The author suggested to use their algorithm for other graph optimization problems including maximum clique problem [22]. Xu and Kumar proposed a solver for the minimum weighted vertex cover problem (MWVC). Their algorithm reformulated a series of SAT (satisfiability) instances using a primal-dual approximation algorithm as a starting point [23]. Ruizhi Li, and coworkers proposed a local search algorithm with tabu strategy and perturbation mechanism for generalized vertex cover problem [24]. For hypergraphs and bounded degree graphs Halperin and co-workers proposed an algorithm to find minimum vertex. They used semi-definite programming and introduced a new rounding technique for this purpose [25]. Cai, S. et al. have reported an algorithm based on local search (NuMVC) that has been found efficient in finding MVC. They introduced two new processes that involve a two-way exchange and an edge weighting mechanism [26].

The literature survey indicates variety of results as far as complexity and accuracy of algorithms are concerned. Onak and Rubinfeld developed a randomized algorithm for maintaining an approximate maximum cardinality matching with a time complexity of $O(\log^2 n_v)$ [27].

In particular, the present work is more suitable for two dimensional graphs with triangular grid structures. The two-dimensional triangular grid graphs are very common in telecommunications, in molecular biology, in configurational statistics of polymers and in various other fields [28–30]. We are proposing here a new way to find edges shared by such triangular grid structures and use these subgraphs to simplify the MVC for such graphs.

2 Definitions

A graph can be represented as a matrix M_e (Edge Matrix or Adjacency Matrix) such that;

$$M_e = [e_{ij}], \text{ where } e_{ij} = \begin{cases} 1 & \text{if } (i,j) \in E \\ 0 & \text{if } (i,j) \notin E \end{cases}$$

Here i and j are numbered vertices such that $i, j \in V$.

The set $N(k)$ is the set of neighbors of the k^{th} vertex in the graph. The set of edges connected by the k^{th} vertex is represented by k^{th} row or k^{th} column of the edge matrix. The removal of the k^{th} row and k^{th} column from the edge matrix M_e is equivalent to the removal of the k^{th} vertex and all of its incident edges from the graph.

From here onward let H_o denote a Hamiltonian cycle with odd number of edges (subgraphs of the form triangles, pentagons and heptagons etc.), H_T denote a Hamiltonian cycle with three edges (triangles only), a common or shared edge here is defined as an edge that is shared by more than one Hamiltonian cycles of the form H_o . A Hamiltonian cycle with three edges i.e., H_T can be represented as $e_{ij}e_{jk}e_{ki} = 1$, where i, j, k are the three vertices of that Hamiltonian cycle. Similarly, for any odd number of vertices i, j, k, l, \dots, z one can represent the Hamiltonian cycle H_o as $e_{ij}e_{jk}e_{kl} \dots e_{zi} = 1$.

Any vertex $w \in V_c$ is said to be covered, V_c denotes here the vertex cover of a graph. $A \setminus B$ implies all those elements of a set A which are not elements of set B and a shared vertex is defined here as a vertex that has a degree greater than two.

3 Proposed Work

A graph can be divided into a number of subgraphs such that;

$$G(V, E) = \bigcup_{i=1}^r G_i(V_i, E_i) \quad (1)$$

One may construct these subgraphs G_i 's such that $E_i \cap E_j = \emptyset$, that is these subgraphs do not share any edge. Yet, there are two possibilities, $V_i \cap V_j = \emptyset$ or $V_i \cap V_j \neq \emptyset$ for $i, j \in \{1, 2, 3, \dots, r\}$. If $V_i \cap V_j = \emptyset$, ($i, j = 1, 2, 3, \dots, r$), $C = \bigcup_{i=1}^r C_i$ is minimum vertex cover of G , where C_i is the minimum vertex cover of G_i . But if $V_i \cap V_j \neq \emptyset$, a union of the intersections of each such pair will yield a set of vertices that may not be covered by individual subgraphs. In that case C does not represent the minimum vertex cover of the graph G . However, the union does not necessarily require to be union of intersection of all the pairs, rather union of intersection of fewer pairs may yield an optimum solution. Let us denote the optimum union set as U . Covering all vertices of U and removing from the graph G will leave $U_i \cap G_j = \emptyset$, i.e., all these subgraphs become disjoint and vertex cover becomes $C = U \bigcup_{i=1}^r C_i$. Let $U \subseteq V$ be a set with a minimum cardinality among the other sets, exclusion of which assures $G_i \cap G_j = \emptyset$ (where subgraphs G_i and G_j become either isolated paths or Hamiltonian cycles). However, to conveniently decompose a graph into a number of subgraphs and to find the set U is difficult.

A solution for graph decomposition is proposed that is based on Lemma and Theorem proved below.

Lemma: An edge that has both of its vertices covered must belong to a Hamiltonian cycle or a path with odd number of edges.

Proof: A graph can be decomposed into a number of subgraphs that may be a Hamiltonian cycle or a path with odd or even number of edges. For even number of edges a Hamiltonian cycle or path does not require both vertices of any edge to be covered. For a Hamiltonian cycle with odd number of sides only one of the edges must have both vertices covered. For paths with odd number of sides both vertices of one of those edge may or may not require to be covered. Hence, if both vertices of an edge are covered that edge may either be an edge of a Hamiltonian cycle or a path with odd number of edges.

Theorem: The exact minimum vertex cover of a graph G can be found by constructing a subgraph from edges $(x, y) \in H_o$ and finding minimum vertex cover of the subgraph, where $x, y, (x, y) \in G$

Proof: For a given set of vertex cover there may exist at most two types of edges depending on the vertices being covered i.e., an edge with both of its vertices covered and an edge with a single vertex covered. The edge with both vertices covered essentially belongs to a subgraph of the form of a path or a Hamiltonian cycle with odd number of edges as proved in the Lemma. In case of two different set of vertex covers i.e., an optimum vertex cover and an approximate one, there can be at most five cases, which are listed below;

Case 1: An edge (x_1, y_1) covered by both vertices x_1 and y_1 in the set of optimum vertex cover and covered by a single vertex (either x_1 or y_1) in approximate set of vertex cover.

Case 2: An edge (x_2, y_2) covered by a single vertex (either x_2 or y_2) in the optimum set and covered by both vertices (x_2 and y_2) in the approximate set.

Case 3: An edge (x_3, y_3) covered by both vertices x_3 and y_3 in both sets.

Case 4: An edge (x_4, y_4) covered by a single but different vertex in both sets.

Case 5: An edge (x_5, y_5) covered by a single and same vertex in both sets.

For both of the sets the number of cases like case 3, 4 and 5 are the same i.e., such cases do not cause any difference on the cardinality of both sets. In cases 1, 2 and 3 at least one of the sets have its edges covered by both vertices. Since case 3 is the same for both the sets, the only two cases that can cause difference on the cardinality of both sets are case 1 and 2. For the approximate set the number of cases like case 2 is greater than or equal to the number of cases like case 1. This leads to the fact that a large number of cases like case 2 can increase cardinality of an approximate set compared to the optimal set. All edges that belong to subgraphs that are either paths or Hamiltonian cycles with odd number of edges are candidates of having both vertices in the vertex cover. By simply separating all such edges the optimization problem simplifies.

This leads to a conclusion that minimum vertex cover optimization depends only on optimization of subgraphs formed by Hamiltonian cycles or paths with odd number of sides.

Based on the Theorem a new approach is proposed to find $u \in U$, approximately, and C_i exactly, and hence an approximate minimum vertex cover of the form $C = U \bigcup_{i=1}^r C_i$. In principle, any subgraph of the form H_o must have both vertices of one of its edges in the vertex cover. This work is based on finding edges that satisfies $e_{ij} \in H_T$. Removal of these edges from the graph assures $E_i \cap E_j = \emptyset$ which is followed by removal of common or shared vertices to result in disjoint graphs satisfying $G_i \cap G_j = \emptyset$.

Our aim here is to find edges, which are common in more than one triangular Hamiltonian cycle. We refer such edges as shared edges. To find such shared edges in a large graph we are proposing a new approach. The new approach is based on decomposition of a graph $G(V, E)$ into two subgraphs G_U and G_S in such a way that an edge $e_{ij} \in G_S$ satisfies $e_{ij} \in H_o$ for at least two subgraphs of the form H_o (referred here as shared edges) in G , whereas all edges other than shared edges forms G_U . A greedy approach can then be used for selecting a vertex from G_S . The selected vertex is then covered. After covering such vertices some of the edges (e_{mn}) may no longer satisfy the condition $e_{mn} \in H_o$ and hence are moved from G_S to G_U , the removal of vertices from of G_S will eventually result in $G_S = \emptyset$. At this stage the uncovered graph G_U will contain shared vertices, isolated polygons and/or isolated paths. This subgraph can further be divided into two subgraphs G_I and G_{SV} , where G_{SV} consists of all the shared vertices and their adjacent edges, and G_I consists of one or more than one isolated Hamiltonian cycles or paths. Both the subgraphs G_{SV} and G_I are covered using maximum degree greedy approach. The greedy approach exactly covers the subgraph G_I . This work is limited to find subgraph of the form of H_T . i.e., the set of all shared edges of all triangles in a graph. This can be done by finding common neighbors of all edges of the graph. For vertices i and k that form edge e_{ik} , the common neighbors can simply be found by an intersection of a subgraph $N(i)$ with subgraph $N(k)$. The set $N(i) \cap N(k)$, is found by carrying out AND operation of i^{th} row or column with k^{th} row or column of the edge matrix M_e . One can write; $N(i) \cap N(k) = e_{ij} \wedge e_{jk}$, where $j = 1, 2, 3, \dots, n_v$ and n_v is the total number of vertices in the graph. The intersection yields the common neighbors of i^{th} and k^{th} vertices. One can evaluate square of the edge matrix as;

$$S_e = M_e^2 = \left[\sum_{j=1}^n e_{ij}e_{jk} \right]_{n_v \times n_v}$$

An element of matrix S_e , i.e., $s_{ik} = \sum_{j=1}^n e_{ij}e_{jk}$ has the bounds $0 \leq s_{ik} \leq n_v - 2$, where $i \neq k$. In matrix S_e the diagonal elements s_{jj} ($j = 1, 2, 3, \dots, n_v$) represent the degree of vertex j , whereas off diagonal elements s_{ik} are the number of common neighbors of the vertices i and k , or the number of triangles sharing the edge e_{ik} if the edge exists. In other words, each term in a given element s_{ik} corresponds to a

Hamiltonian cycle H_T i.e., $e_{ij}e_{jk}e_{ki} = 1$. One can also sort Hamiltonian cycles H_O with odd number of edges greater than H_T by using $e_{ij}e_{jk}e_{kl} \dots e_{zi} = 1$ condition, where $\{i, j, k, l, \dots, z\}$ is a set of odd number of vertices. However, the current work is limited to find all Hamiltonian cycles of the form H_T . There are $n_v(n_v - 1)$ off-diagonal elements of S_e . Since S_e is symmetric i.e., $s_{ik} = s_{ki}$, only $n_v(n_v - 1)/2$ elements of S_e are calculated. A maximum of $n_v(n_v - 1)/2$ subgraphs can be constructed for each possible pair of vertices i and k . Each of these subgraphs consists of set of vertices $\{i, k, N(i) \cap N(k)\}$ and contains $2s_{ik} + e_{ik}$ number of edges. However, construction of such subgraphs is beyond the scope of this work, therefore we restrict ourselves to the evaluation of weighted matrix S_e only. The subgraphs are completely covered by the vertices i and k . However, depending on the value of s_{ik} the vertices i and k may or may not be the minimum vertex cover of that subgraph. The matrix S_e does not contain any information about common neighbors other than the total number of common neighbors two vertices can have.

An element-by-element multiplication (just corresponding element multiplication of two matrices) of matrix S_e with that of the matrix M_e results in a matrix T_e such that each element $t_{ik} = e_{ik}s_{ik}$ of T_e represents number of triangles that share the edge e_{ik} . The elements of matrix T_e are classified as;

$$t_{ik} = \begin{cases} 0 & N(i) \cap N(k) = \emptyset \\ 1 & |N(i) \cap N(k)| = 1 \\ \geq 2 & |N(i) \cap N(k)| \geq 2 \end{cases}$$

For example, $t_{ik} \geq 2$ corresponds to a subgraph which forms two or more than two triangles with a shared edge e_{ik} . For all the cases with $t_{ik} \geq 2$, the vertices i and k are the minimum vertex cover of the subgraph. For $t_{ik} = 0$, implies either the edge is not shared or $e_{ik} \notin G$. For $t_{ik} = 1$, the subgraph forms a single triangle. The value $t_{ik} = 0$, implies that $e_{ik} \notin H_T$, but $e_{ik} \in H_O$ may still be possible. A subgraph G_s of all edges satisfying $e_{ik} \in H_T$ can be generated using the weight matrix T_e for $t_{ik} \geq 2$. However, if all the edges of a graph are shared edges, the condition $t_{ik} \geq 2$ will produce a graph of shared edges same as the original graph i.e., $G_s = G$. For such graphs one can modify the condition as $t_{ik} \geq t_{min} + n$, where t_{min} is minimum number of triangles sharing a single edge in that graph and n is a small number. This modification will generate a reduced graph of shared edges. A vertex cover of the reduced subgraph G_s removes all shared edges of the form $e_{ik} \in H_T$ from G . However, the condition $e_{ik} \in H_O$ may still not be satisfied.

Removal of any subset of vertices from a graph requires removal of the corresponding row or column from the edge matrix M_e . This leads to calculation of new square matrix S_e . However, instead of calculating S_e from scratch, a low-cost solution is proposed here to reduce the simulation time.

Proposition: S_e^* being the square of matrix M_e^* can be calculated from S_e , where M_e^* is the reduced graph after removal of l^{th} row and l^{th} column from M_e .

Proof: Since

$$S_e = \left[\sum_{j=1}^n e_{ij}e_{jk} \right]_{n_v \times n_v} \tag{1}$$

S_e can be decomposed as

$$S_e = \left[\sum_{\lambda=1}^{l-1} e_{i\lambda}e_{\lambda k} \right]_{n_v \times n_v} + [e_{il}e_{lk}]_{n \times n} + \left[\sum_{\mu=l+1}^n e_{i\mu}e_{\mu k} \right]_{n_v \times n_v}$$

where $\lambda = 1, 2, 3, \dots, l - 1$ and $\mu = l + 1, l + 2, \dots, n_v$

Let $F = [e_{il}e_{lk}]_{n \times n}$ is an $n_v \times n_v$ matrix that can be obtained by multiplying l^{th} column and l^{th} row of the edge matrix M_e . One can write;

$$S_e = \left[\sum_{\lambda=1}^{l-1} e_{i\lambda}e_{\lambda k} \right]_{n_v \times n_v} + \left[\sum_{\mu=l+1}^n e_{i\mu}e_{\mu k} \right]_{n_v \times n_v} + F \tag{2}$$

Let

$$\hat{S}_e = \left[\sum_j e_{ij}e_{jk} \right]_{n_v \times n_v} \tag{3}$$

where $j = 1, 2, 3, \dots, l-1, l+1, \dots, n_v$

Eqs. (2) and (3) yields;

$$S_e = \hat{S}_e + F \tag{4}$$

All elements in l^{th} row and l^{th} column from \hat{S}_e are zeros, therefore, removing l^{th} row and l^{th} column from \hat{S}_e yields S_e^* , which is the required matrix of order $(n_v - 1) \times (n_v - 1)$.

$$S_e^* = \left[\sum_j e_{ij}e_{jk} \right]_{(n_v-1) \times (n_v-1)} \tag{5}$$

Reduced matrix S_e^* is multiplied element by element with M_e^* to evaluate T_e^* . As previously discussed, the square matrix elements s_{ik} is the number of common neighbors of the vertices i and k , the value s_{ik} can be used as a weight to select vertices to be covered.

Algorithms

The decomposition of a graph G into G_U and G_S (The subgraph containing all shared edges of triangular structures) is accomplished by transforming the matrix $T_e \rightarrow W_e$ such that $[W_e]_{ij} = \min(1, [T_e]_{ij})$ and $t_{ik} \geq 2$. The algorithm is divided into three stages. In first stage a vertex with highest degree in the subgraph G_S is found and covered. The first stage is terminated when $G_S = \emptyset$. The subgraph G_U does not contain any shared edge that belongs to triangular structures but it may still have edges shared by subgraphs of the form H_o . In second stage, the vertex with highest degree is found from G_U and covered. After covering all the shared vertices of the graph, the graph is left with isolated paths or polygons. In third stage, a vertex with degree 2 is found and covered. The removal of a vertex from G in all three stages may lead to leaves in the residual graph. Therefore, these leaves are removed by covering their adjacent vertex.

The proposed algorithms are described in the following section. Algorithm 1 and Algorithm 2 are abbreviated as ASE and ASER such that ‘A’ stands for Algorithm, ‘SE’ stands for ‘Shared Edges’ and ‘R’ stands for ‘Reduction Strategy’.

Algorithm 1: (ASE)

- Input: Graph $G(V, E)$.
 - Output: Minimum Vertex Cover of graph $G(V, E)$.
- 1: Read data and construct edge matrix M_e and W_e
 - 2: while size $(M_e) > 1$
-

(Continued)

Algorithm 1: Continued

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3:   while (size ( $W_e$ ) > 1)
4:       Perform operation (A0), (A1), (A2), (A3) and (A4)
5:   end
6:   while (deg ( $v \in V$ )  $\geq$  2)
7:       Perform operation (A5), (A6), (A3) and (A4)
8:   end
9: end

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(A0) Evaluate W_e using matrices M_e , S_e and T_e for $t_{ik} \geq 2$ or $t_{ik} \geq t_{\min} + 3$

(A1) Find a vertex with highest degree from W_e and cover the vertex

(A2) Evaluate matrix F for the vertex found in step A1. Using Eq. (4) evaluate reduced matrix S_e using matrix F and the matrices M_e and S_e from step A1. Reduce matrix M_e .

(A3) Remove all leaves from the graph

(A4) Remove all isolated vertices from the graph

(A5) Find deg ($u \in U$), i.e., degree of vertex u in present set of vertices U

(A6) Find the vertex u that has the maximum degree in present graph, cover the selected vertex u and remove from the graph.

Complexity

Total complexity of the algorithm is $(34n_v^3 - 99n_v^2 + 170n_v - 120) / 24 \cong 1.42n_v^3$.

To reduce complexity of ASE a reduction strategy is proposed. The reduction strategy consists of splitting graph into two subgraphs and finding independent set of 1st subgraph and taking union of that independent set with 2nd subgraph and finding independent set of the union set. However, this graph splitting is not random. For splitting one can list the set of edges $E = \{(\lambda, \mu) : \mu \in N(\lambda), \forall \lambda, \mu \in E\}$. Construct two sets of vertices $L = \{\lambda : \forall (\lambda, \mu) \in E\}$ and $R = \{\mu : \forall (\lambda, \mu) \in E\}$. Find $L \cap R$. One can see that the set, $I_L = L \setminus L \cap R$ is an independent set, because there are no two vertices in I_L , that are neighbors of each other. Similarly, $I_R = R \setminus L \cap R$ is also an independent set.

Under normal circumstances the list of edges E may not provide reasonably large I_L and I_R , therefore, one may need to prepare the list the edges E so that I_L and I_R are sufficiently large. One may opt for an alternative strategy to find sufficiently large I_L , I_R and a small $V_{res} = L \cap R$ by using a low-cost algorithm that can find an approximate minimum vertex cover. The low-cost algorithm outputs an independent set and residual set of vertices. Therefore, one can use the low-cost algorithm twice to find I_L and $(V \setminus I_L)$ and again to find I_R and $(V \setminus I_L) \setminus I_R$ and remaining set of vertices $V_{res} = L \cap R$. Now one can construct a subgraph from V_{res} , that is $G_{res} = G_{res}(V_{res}, E_{res})$. An independent set I_{res} can be found using Algorithm 1 (ASE). A second subgraph can be constructed from $I_s = I_L \cup I_R \cup I_{res}$, that is $G_s = G_s(I_s, E_s)$. Using Algorithm ASE again one can construct the final independent set. Complexity of this algorithm depends on the cardinality n_L of I_L and n_R of I_R . The cardinality of set of vertices of the 1st subgraph is $n_1 = n_v - n_L - n_R$ and cardinality of set of vertices of 2nd subgraph is $n_2 = n_{res} + n_L + n_R$, where n_{res} is the cardinality of the set I_{res} . The complexity of algorithm ASER is $[34(n_1^3 + n_2^3) - 76(n_1^2 + n_2^2) + 170(n_1 + n_2) - 240] / 24$.

Algorithm 2: (ASER)

· Input: Graph $G(V, E)$.
 · Output: Minimum Vertex Cover of graph $G(V, E)$.

- 1: Perform operation (B0) to find I_{s1} and perform operation (B1) to find residual graph G_{R0} ,
- 2: Perform operation (B0) using G_{R0} as input to find I_{s2} and perform operation (B1) to find residual graph G_R
- 3: Perform operation (B3) and find an independent set I_{s3} using G_R as input
- 4: Perform operation (B2) to find I_s using I_{s1} , I_{s2} and I_{s3} and construct subgraph G_s from I_s
- 5: Using G_s as input perform step (B3) and find the independent set I_{ind} and vertex cover V_c
- 6: Using I_{ind} and V_c perform operation (B4) and find the final maximum independent set I_{max}
- 7: end

(B0) Find an independent set I_s of a graph using a low-cost algorithm

(B1) Construct residual subgraph $G_R(V_R, E_R)$ from $V_R = V - I_s$.

(B2) Construct a single set using three sets as; $I_s = I_{s1} \cup I_{s2} \cup I_{s3}$

(B3) Find maximum independent set of given subgraphs using algorithm 1

(B4) Find a subset I_c of the vertex cover V_c , such that I_c contains vertices that individually can go to independent set I_{ind} . Find independent set of I_c using Algorithm1. Move all vertices of that independent set to I_{ind} to construct the final independent set.

The low-cost algorithm simply selects and adds a vertex to an independent set followed by searching vertices that can be moved to the independent set. This algorithm has a complexity of $n_v^2/2$. However, one is free to choose the low-cost algorithm in accordance with the suitability.

4 Results and Discussion

In this section both the algorithms are tested and analyzed for their accuracy and complexity for benchmark graphs taken from [31,32] and [33]. The simulations are performed on a computer with 1.61 GHz processor and 8.00 GB RAM using sequential programming.

The results from the 72 benchmark graphs referred above are organized in the form of three tables. [Tabs. 1](#) and [2](#) contain optimum vertex cover and accuracy in the form of error ratios for the algorithms of respective graphs. [Tab. 3](#) contains a comparison of results of the algorithm ASE with three well know algorithms. The error ratio is defined here as the value of minimum vertex cover for a given graph obtained from each algorithm divided by the optimum vertex cover (n_o) of that graph. The condition $t_{ik} \geq 2$ has been used to generate shared edges graph (G_s) for 64 benchmark graphs. For some of the benchmark graph all the edges are shared edges, therefore, the condition $t_{ik} \geq 2$ will yield $G_s = G$. For such graphs the condition is modified to $t_{ik} \geq t_{min} + 3$, where t_{min} is minimum number of triangles sharing a single edge in that graph. This modification results in reduced shared edges graph. The condition $t_{ik} \geq t_{min} + 3$ is used to generate G_s for eight such graphs and their error ratios are given in [Tab. 2](#).

Table 1: The calculated MVC and error ratio for the proposed algorithms for $t_{ik} \geq 2$ (c stands for clq and cc for clq_compliment)

Sr. #	Benchmark Graph	n_v	n_o	ASE (ϵ_r)	ASER (ϵ_r)	Sr.#	Benchmark graph	n_v	n_o	ASE (ϵ_r)	ASER (ϵ_r)
1	graph50_01	50	30	1	1	33	p_hat700_2	700	656	0.995	0.998
2	graph50_02	50	30	1	1	34	Jhonson8-2-4c	28	24	1	1
3	graph50_03	50	30	1	1	35	Jhonson8-4-4c	70	56	1	1
4	graph50_05	50	27	1	1	36	Jhonson16-2-4c	120	112	1	1
5	graph50_06	50	38	1	1	37	Jhonson32-2-4c	496	480	1	1
6	graph50_07	50	35	1	1	38	sanr200_0_7	200	182	1.005	1.016
7	graph50_08	50	29	1	1	39	sanr200_0_9	200	158	1.025	1.038
8	graph50_09	50	40	1	1	40	sanr400_0_5	400	387	1.005	1.01
9	graph50_10	50	35	1	1	41	sanr400-0.7c	400	379	1.013	1.013
10	graph100_01	100	60	1	1	42	frb35_17_2	595	560	1.016	1.02
11	graph100_02	100	65	1	1	43	c125	125	91	1.011	1.022
12	graph100_03	100	75	1	1	44	c250	250	206	1.024	1.019
13	graph100_04	100	60	1	1	45	c500	500	443	1.02	1.025
14	graph100_05	100	60	1	1	46	brock200_2	200	188	1.011	1.016
15	graph100_06	100	80	1	1	47	hamming6-2_cc	64	32	1	1
16	graph100_07	100	65	1	1	48	hamming6-4_cc	64	60	1	1
17	graph100_08	100	75	1	1	49	hamming8-2_cc	256	128	1	1
18	graph100_09	100	85	1	1	50	hamming8-4_cc	256	240	1	1
19	graph100_10	100	70	1	1	51	hamming10-2_cc	1024	512	1	1
20	graph200_01	200	150	1	1	52	c_fat200_1	200	188	1	1
21	graph200_02	200	125	1	1	53	c_fat200_2	200	176	1	1
22	graph200_03	200	175	1	1	54	c_fat200_5	200	142	1	1
23	graph200_04	200	140	1	1	55	c_fat500_1	500	486	1	1
24	graph200_05	200	150	1	1	56	c_fat500_2	500	474	1	1
25	graph500_01	500	350	1	1	57	c_fat500_5	500	436	1	1
26	graph500_02	500	400	1	1	58	MANN_a27	378	252	1.003	1.003

(Continued)

Table 1: Continued

Sr. #	Benchmark Graph	n_v	n_o	ASE (ϵ_r)	ASER (ϵ_r)	Sr.#	Benchmark graph	n_v	n_o	ASE (ϵ_r)	ASER (ϵ_r)
27	graph500_03	500	375	1	1	59	MANN_a45	1035	1032	1	1
28	graph500_04	500	300	1	1	60	C2000_9	2000	1920	1.013	1.012
29	graph500_05	500	290	1	1	61	C4000_5	4000	3982	1.001	1.001
30	p_hat300_1	300	292	0.997	1.007	62	MAAN-a81	3321	2221	1.002	1.002
31	p_hat300_2	300	275	0.996	1.004	63	Ca_GrQc	4158	2208	1	1.005
32	p_hat300_3	300	264	0.992	1.004	64	Bio-dmela-mtx	7393	2630	1.000	1.009

Table 2: The calculated MVC and error ratio for the proposed algorithms for $t_{ik} \geq t_{min} + 3$

Sr. #	Benchmark Graph	n_v	n_o	ASE (ϵ_r)	ASER (ϵ_r)	Sr.#	Benchmark Graph	n_v	n_o	ASE (ϵ_r)	ASER (ϵ_r)
1	graph50_04	50	40	1	1	5	DSJC500_5	500	487	1.004	1.004
2	p_hat700_1	700	689	1.004	1.003	6	DSJC1000_5	1000	986	1.002	1.002
3	p_hat700_3	700	638	1	1	7	keller4	171	160	0.981	0.981
4	frb30_15_5	450	420	1.012	1.014	8	keller5	776	749	0.995	0.995

Table 3: Comparison of ASE with MDG, MVSA and MtM

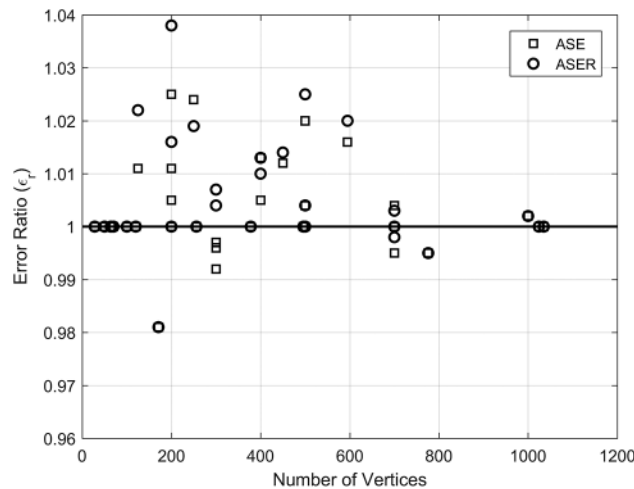
Sr. #	Benchmark Graph	n_v	n_o	ASE (ϵ_r)	MDG (ϵ_r)	MVSA (ϵ_r)	MtM (ϵ_r)
1	p_hat300_1	300	292	0.997	1.003	1.006	1.005
2	p_hat300_2	300	275	0.996	1.010	1.014	1
3	p_hat300_3	300	264	0.992	1.01	1.03	1.008
4	p_hat700_2	700	656	0.995	1.003	1.006	1.003
5	sanr200_0_7	200	182	1.005	1.005	1.021	1.005
6	sanr200_0_9	200	158	1.025	1.005	1.031	1.058
7	sanr400_0_5	400	387	1.005	1.003	1.005	1.003
8	sanr400-0.7c	400	379	1.013	1.007	1.005	1.008
9	frb35_17_2	595	560	1.016	1.014	1.008	1.009
10	c125	125	91	1.011	1.022	1.043	1.033
11	c250	250	206	1.024	1.0145	1.0145	1.0242
12	c500	500	443	1.02	1.07	1.042	1.011
13	brock200_2	200	188	1.011	1.0213	————	————
14	C2000_9	2000	1996	0.999	————	————	————

(Continued)

Table 3: Continued

Sr. #	Benchmark Graph	n_v	n_o	ASE (ϵ_r)	MDG (ϵ_r)	MVSA (ϵ_r)	MtM (ϵ_r)
15	p_hat700_1	700	689	1.004	1.004	1.004	1.004
16	frb30_15_5	450	420	1.012	1.009	1.009	1.014
17	DSJC500_5	500	487	1.004	1.008	1.004	1.004
18	DSJC1000_5	1000	986	1.002	—	—	—
19	keller4	171	160	0.981	1.025	1	1
20	keller5	776	749	0.995	1.02	1.007	1.007
	Average (ϵ_r)			1.005	1.014	1.015	1.012

Fig. 1 shows the error ratio (ϵ_r) obtained from both the algorithms plotted against the number of vertices for 67 benchmark graphs (excluding the graph with number of vertices greater than 2000 for better visibility of the figure). The solid line in Fig. 1 corresponds to the error ratio for optimum vertex cover. It can be seen that the error ratio for graphs with fewer number of vertices is less accurate compared with that of the higher number of vertices. This suggests that the algorithms get better and better with increased number of vertices. It can also be noted that the worst-case error ratio (ϵ_r) for algorithms ASE and ASER are 1.025 and 1.038 respectively. An interesting finding of this study is the reported optimum error ratio for few benchmark graphs in literature is found slightly less accurate compared to the value calculated in this work. This can be seen in Fig. 1 as ASE and ASER lies below the optimum error ratio line on six and three occasions, respectively.

**Figure 1:** Accuracy of the suggested algorithms, solid line represents reported optimum values

As can be seen from the Tabs. 1 and 2 that out of 72 instances ASE and ASER both give an error ratio ($\epsilon_r = 1$) or better on 55 and 50 instances, respectively. The average error ratios (ϵ_r) for ASE and ASER are 1.0017 and 1.0030 respectively.

In ASE, since only edges $e_{mn} \in H_T$ are covered, all edges satisfying $e_{mn} \in H_O$ are not covered with certainty. This may lead to erroneous results. After covering all edges from subgraphs of the form H_T , shared vertices are selected with the priority of highest degree of the residual graph, and the vertices are

moved to vertex cover one by one. The selection may not be accurate since it uses the greedy approach. However, the approach will produce exact minimum vertex cover for the residual graph that is left with isolated Hamiltonian cycles or paths.

Supplementary data describes the complexity of the proposed algorithms and contains two parametric numbers or reduction parameters p_1 and p_2 . Supplementary data also contains time taken by each of the algorithm to find the minimum vertex cover. The reduction parameter is the ratio of approximate complexity of the form $(\cong 1.41n_{1,2}^3)$ of an algorithm with that of $(\cong 1.41n_v^3)$, and can be calculated as $p_{1,2} = n_{1,2}^3/n_v^3$, where n_1 and n_2 are defined in the previous section. These reduction parameters represent a time reduction of an algorithm with reduction strategy (ASER) to the same without reduction strategy (ASE).

The comparison between ASE and ASER is given in Fig. 2, which shows that if either $p_1 \rightarrow 1$ or $p_2 \rightarrow 1$, a difference between their simulation times approaches to zero. For graphs hamming6_2_clq_compliment, hamming8_2_clq_compliment and hamming10_2_clq_compliment, one can see that $p_1 \rightarrow 1$ and $p_2 \rightarrow 0$, hence no significant time difference is observed. Similarly, for graphs p_hat700_1, MANN_a27, MANN_a45 and C2000_9 reduction parameters $p_1 \rightarrow 0$ and $p_2 \rightarrow 1$, forcing the time difference to approach to zero. A noticeable difference in simulation time can be observed for the cases where neither $p_1 \rightarrow 1$ nor $p_2 \rightarrow 1$. The reduction parameters p_1 and p_2 for each of the benchmark graphs are plotted in Fig. 2.

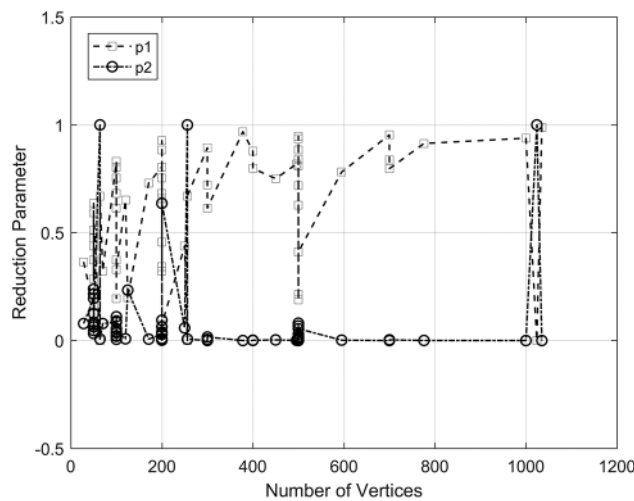


Figure 2: Reduction parameters p_1 and p_2 plotted against number of vertices

The simulation time for the proposed algorithms is plotted against the number of vertices in Fig. 3. A cubic fit function of the form $f(n_v) = Cn_v^3$ is also plotted to show the complexity trend of both the algorithms. It can be seen that the algorithms (ASE and ASER) follow the cubic fit trend. Since computer takes a small amount of preprocessing time before each simulation, one can see that for low values of n_v the simulation time is large compared to $f(n_v)$, whereas for large values of n_v the simulation time matches with $f(n_v)$. One can also see that simulation time for ASER is occasionally smaller than $f(n_v)$, reflecting a success in reduction strategy.

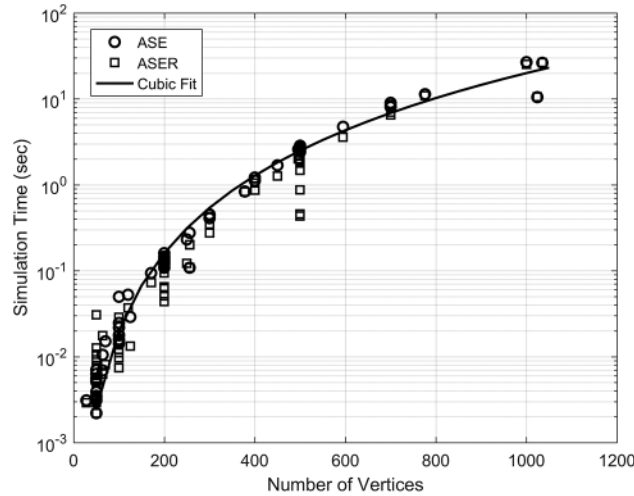


Figure 3: Simulation time vs. the number of vertices for the proposed algorithms

In [Tabs. 1](#) and [2](#) the performance of the algorithm ASE is evaluated using 72 benchmark graphs. On 48 instances our algorithms yield optimal values. In [Tab. 3](#) we have compared the results for 20 benchmark graphs for which ASE does not yield optimal values with three well known algorithms MDG (maximum degree greedy) [34], MVSA (modified vertex support algorithm) [35] and MtM (Min-to-Min) [36]. The average ratio error for these 20 benchmark graphs presented in the [Tab. 3](#) for ASE, MDG, MVSA and MtM are, 1.005, 1.014, 1.015 and 1.012, respectively. This shows that the algorithm ASE clearly outperforms the three algorithms in comparison.

5 Practical Implementation

For step-by-step analysis of the algorithm a simple real-life example is selected. Crypto or digital currency market currently has a market capital of billions of US dollars. There are hundreds of crypto currencies with billions of USD daily volume. Market data analysis of these currencies is becoming harder and harder with growth in data. However, almost all of these currencies are paired with each other for trading. Therefore, any market fluctuation is coupled within these crypto currencies. In principle, one can represent these currencies and their trading pairs in the form of a graph and can find minimum vertex cover of the graph to select only few currencies for crypto market data analysis. We have selected ten crypto currencies and their trading pairs in order to simplify the problem. These crypto currencies and their trading pair are presented in the form of a graph is shown in [Fig. 4a](#).

The graph is decomposed in subgraphs with bold dark lines showing shared edges and dashed lines as the rest of the graph. The shared edge graph is simply a triangle in this case and each vertex has a degree 2 in the subgraph. A single vertex (in this case vertex 1) is covered, i.e., $V_c = \{1\}$ and we are left with only a single edge (e_{23}) in the subgraph. One vertex (vertex 2) of the remaining edge (e_{23}) is covered, yielding $V_c = \{1, 2\}$. Since there are no more shared edges in the graph, therefore we are left with the subgraph of the form shown in [Fig. 4b](#). From here on ward the vertices are simply covered on the basis of their degree. Since the vertex 3 has the highest degree therefore, it is covered, yielding $V_c = \{1, 2, 3\}$ and hence the entire graph is covered. A conclusion of this process is that, only crypto currencies numbered 1, 2 and 3 represent the complete variation of the market for only these ten crypto currencies. However, in order to completely analyses the entire crypto market a minimum vertex cover of a graph representing entire market has to be found.

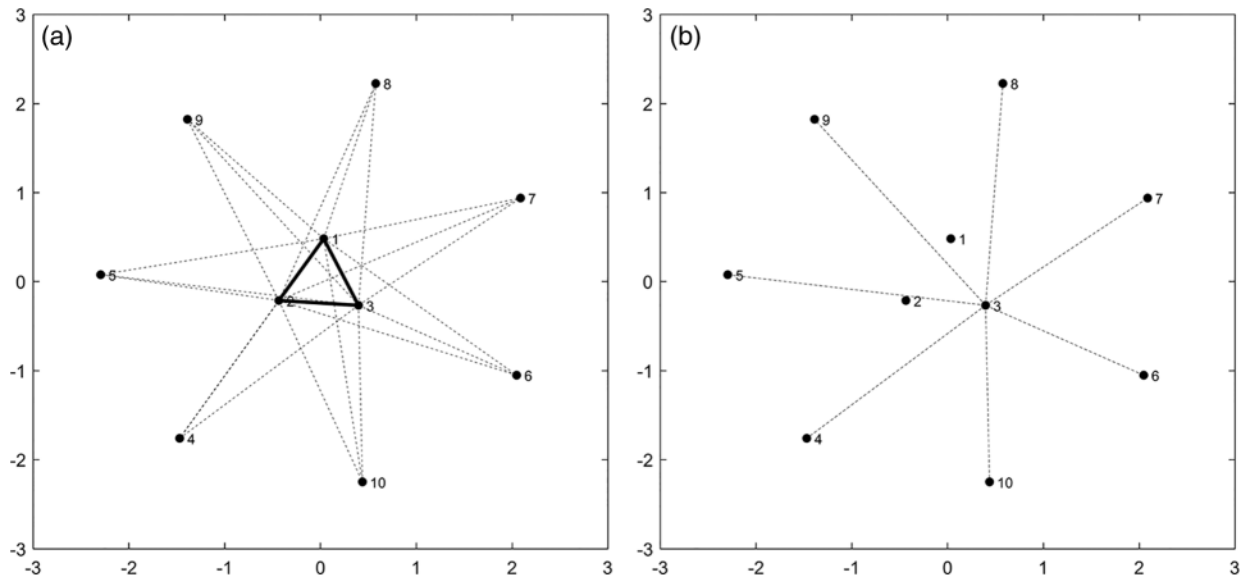


Figure 4: (a) Crypto currencies and their trading pair (b) Subgraph with no shared edge

6 Summary

The proposed approach simplifies a graph to isolated paths or polygons (Hamiltonian cycles) by moving shared edges followed by shared vertices to the vertex cover. This process forms three subgraphs i.e., a subgraph containing shared edges, a subgraph with shared vertices and finally a simplified subgraph. The final simplified subgraph is either a single Hamilton cycle or path or a number of isolated Hamiltonian cycles or paths. The vertex cover of the simplified subgraph can be exactly found in a short time using maximum degree greedy approach. The accuracy of finding the first two subgraphs depends on the sequence of covering the vertices. The proposed strategies are capable to search for the sequence of selection of vertices to an approximate extent only. However, using this approach the problem can be broken successfully into three smaller problems, which are relatively easy to handle. The worst-case error ratio (ϵ_r) for algorithms ASE and ASER are 1.025 and 1.038, respectively. The average error ratios (ϵ_r) for ASE and ASER are 1.0017 and 1.0030 respectively. The algorithms have a maximum complexity of approximately $1.42n_v^3$. Both the algorithms (ASE and ASER) improve optimum error ratio for few graphs compared to the values reported in literature [17,34–36].

7 Future Work

The proposed approach may work even better if all the edges shared by subgraphs of the form H_o can be found and removed with certainty (i.e., in correct sequence) and if one may find a better strategy (or sequence) to remove shared vertices other than the greedy approach. A future work is suggested to find shared edges among all subgraphs of the form H_o .

Funding Statement: The authors received no specific funding for this study.

Conflicts of Interest: The authors declare that they have no conflicts of interest to report regarding the present study.

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