An Edge-Based Approach for Virtual Network Embedding Based on the Graph Edit Distance

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RESEARCH

An Edge-based Approach for Virtual Network Embedding Based on the Graph Edit Distance

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Abstract

Virtualization and resource isolation techniques have enabled the efficient sharing of networked resources. How to control network resource allocation accurately and flexibly has gradually become a research hotspot due to the growth in user demands. Therefore, this paper presents a new edge-based virtual network embedding approach to studying this problem that employs a graph edit distance method to accurately control resource usage. In particular, to manage network resources efficiently, we restrict the use conditions of network resources and restrict the structure based on common substructure isomorphism and an improved spider monkey optimization algorithm is employed to prune redundant information from the substrate network. Experimental results showed that the proposed method achieves better performance than existing algorithms in terms of resource management capacity, including energy savings and the revenue-cost ratio.

Keywords: virtual network embedding; graph theory; resource management; graph edit distance

1 Introduction

Network virtualization [1] is an integral component of current and future networks and promotes the development of networked technologies such as cloud computing [2]. Due to virtualization, network operators have opportunities to consolidate their equipment into standardized high-volume components. This is reflected by efficiently utilizing substrate network (SN) resources through sharing among several virtual networks (VNs) [3]. Virtualization introduces great flexibility in terms of “where” virtual network requests (VNRs) can be embedded. However, exploiting this resource allocation problem poses a fundamental algorithmic challenge. In general, the resource allocation problem is known as the virtual network embedding (VNE) problem [4], which can be posed as a graph-theoretic problem: both the virtual network, describing a workload, and the substrate network, describing the physical infrastructure, can be modeled as graphs. Fig.1 shows a physical network of numerous nodes and edges, receiving a VNR that arrives dynamically.

The virtual network (left), the so-called request graph, describes the resource requirements both on the nodes and on the links; the substrate network graph (right), the so-called substrate graph, describes the physical infrastructure and its resources. To realize the embedding for VNR demands, the VNE algorithm searches for candidate physical nodes that can satisfy these demands while maintaining similar connectivity as in the virtual requests. More importantly, the more similar the

hosted physical graph is to the request graph, the fewer redundant resources are used. However, this has been largely overlooked in previous studies [8]–[10].

The classic approach to handling the VNE problem was usually conducted by focusing on the availability of resources [11]–[15], which embeds request nodes onto physical nodes with the most available resources and embeds each request edge via a path in the substrate network connecting the respective endpoints. As the internet grew, it became unprofitable to search directly for the most promising computing resources in the entire physical network. Therefore, another strategy, performing resource allocation according to predefined constraints by building a substrate resource management region [16]–[18], was developed. By using a feedback loop, spectral clustering and other methods, several nodes and links in the substrate graph are formulated into a collection, which is used to allocate all or part of the networked resource requirements in the VNR. Applications range from mapping testbeds over embedding batch-processing tasks to embedding service function chains and have different mapping restrictions for nodes and edges, not only in capacity [5] but also latency [6] and power consumption [7]. Therefore, a flexible method that can provide predictable performance guarantees on all resources, i.e., nodes and edges, is needed to deal with various requirements. Nevertheless, the VNE problem is a wellknown NP-hard problem [21] and cannot be solved in polynomial time [22]. Traversing the substrate network to find the highest potential substrate region to allocate the VNR is the most effective method for solving this problem [23]–[28]. However, the information of the substrate graph is not always useful, and some information is redundant. To quickly find the most accurate VNE solution possible, a fast and efficient optimization method is needed.

Consequently, the main motivation of this paper is to introduce a new edge-based VNE algorithm (VNE_MR) that accurately and flexibly controls the networked resources usage. Different from previous algorithms that focus on the nodes, we abstract the virtual network as a request graph, denoted by G_v , and employ an edge-

based graph edit distance method CSLGED [29] following the links to edit G_v to the physical resource management region with a similar size to G_v , denoted by $R(G_v)$. In particular, to control the hosted physical network structure efficiently and support various requirements flexibly, we restrict the use conditions of network resources and restrict the structure of $R(G_v)$ based on common substructure isomorphism. In addition, to find the highest potential $R(G_v)$ to host the G_v as the VNE solution, a modified spider monkey optimization (SMO) [30] algorithm is designed to prune redundant information from the substrate network and to find the highest potential $R(G_v)$ in a parallel manner. To our knowledge, our work is the first attempt to model the problem of finding physical resources for VNRs using graph theory and transform it into a problem of finding the minimum graph editing distance based on edges.

The remainder of the paper is organized as follows. Section 2 reviews related studies. Embedding preliminaries are defined in Section 3. Section 4 introduces the proposed VNE_MR algorithm in detail. An explanation of the proposed algorithm and a discussion of its results are presented in Section 5. Finally, Section 6 concludes the paper and highlights certain future work.

2 Related work

In existing VNE research, a common method focuses on resource availability when addressing the resource allocation problem and managing substrate resources.

Francesco et al. [11] used Markov reward processes based on greedy heuristics to present a two-stage VNE algorithm. The key element of this approach is a node-ranking algorithm, which associates a metric that considers the number of local resources available in the vicinity of a given node. Similarly, Gong et al. [12] devised a metric called global resource capacity to quantify the embedding potential of each substrate node and proposed an efficient heuristic VNE algorithm. With regard to resource availability, the most referenced approach is that suggested by Chowdhury et al. [13], which embeds virtual nodes onto SN nodes based on their residual capacities and coordinates edge embedding using the multicommodity flow algorithm. With the increase in delay-sensitive services [31]–[32] on the network, a smaller and guaranteed transmission delay is required. Therefore, researchers are starting to consider delay as an important factor when considering the VNE problem. Based on the Google PageRank algorithm and considering link propagation delay as a node-link constraint in the VNE problem, Cao et al. [14] designed a VNE algorithm, denoted as VNE-TAGRD, to manage the underlying network resources. Using Markov reward theory, Francesco et al. [15] defined a notion of similarity between nodes and proposed a VNE algorithm with delay and placement constraints to map virtual nodes to their most similar physical nodes.

Another common approach is to build a resource management unit to allocate resources according to predefined constraints. For example, to manage resources, Khaled et al. [16] proposed a resource set format, called segments, to manage and match resources in segments. They modeled the VNE problem as a two-stage mapping problem and introduced a VNE algorithm called OPaCoVNE to solve the resource management problem while considering the end-to-end delay as the embedded constraint. Chen et al. [17] developed a feedback system to identify a stable

active subset of substrate resources and then mapped the virtual network in active areas to increase the number of dormant nodes and links, thereby reducing energy consumption. He et al. [18] used spectral clustering based on field theory to extract substrate network features and manage physical resources. Then, they developed dynamic regions of interest to find embedding areas with energy-saving potential for virtual networks.

It is well known that not all of the information in the substrate graph is positive to the VNE problem, and some of the information is redundant, which is not conducive to quickly finding the highest potential substrate resources [9]. Therefore, whether focusing on resource availability during VNE or focusing on building a VNE resource management unit, metaheuristic algorithms have played an important role. For example, to reduce VNE energy consumption and ensure VNE solution validity, Xu et al. [25] devised a topology-aware mapping strategy to manage resources by mimicking information interactions between bacteria during foraging and built a candidate set of VNE solutions. Then, they enhanced the bacterial foraging optimization bacteria to search for the optimal solution. To manage substrate resources, Jahani et al. [26] designed a D-reachability algorithm to assign reachability rank to resources, modeled VNE as a multiobjective optimization problem, and pursued a nondominated sorting heuristic genetic algorithm to solve it. After continually calculating the fitness, nondominated sorting of the population, calculating crowding distance, crossover, and mutation sorting and replacing, the optimal solution is finally output. Su et al. [27] devised a topology-aware measure to rank nodes based on random walks and then designed a greedy algorithm that matches virtual nodes to physical nodes according to this rank. To find the best embedding, they devised a fast-convergent discrete particle swarm optimization.

3 Preliminaries

In this section, VN and SN are modeled as the request graph and the substrate graph, respectively. Some propositions about resources, i.e., nodes and links in the VNE problem, are reviewed (Propositions 1-7). In addition, Proposition 8, presenting the common substructure isomorphism, is specifically proposed for our VNE_MR algorithm.

3.1 Substrate network model

We model the substrate network as an undirected graph $G_s = (N_s, L_s)$. This paper also refers to it as a substrate graph, where N_s is the set of all substrate nodes. Each substrate node $m \in N_s$ is characterized by its node capacity-available CPU, denoted as $C(m)$. With respect to substrate links, we consider any link as a pair of nodes, and each substrate link l_{mn} has a finite bandwidth $B(l_{mn})$ and a substrate link propagation delay $D(l_{mn})$.

3.2 VN request model

In the VNE research area, each VN can also be modeled as a weighted graph $G_V = (N_V, L_V)$. This paper also refers to it as a request graph, where N_V is the set of all virtual nodes and L_V is the set of all virtual links. Each virtual node $M \in N_V$ is characterized by the required CPU, denoted as $C(M)$. With respect to virtual links,

each virtual link L_{MN} has a required bandwidth $B(L_{MN})$ and required virtual link propagation delay $D(L_{MN})$. By adding the time attributes (e.g., maximum waiting time, arrival time, duration time, leaving time), the VN is extended to a VNR.

3.3 VNE propositions

•Proposition1: node mapping

$$\forall M \in G_V, m \in G_s : x_M^m = \begin{cases} 1, & \text{if } M \text{ is mapped on } m \\ 0, & \text{otherwise} \end{cases} \quad (1)$$

In (1), if $x_m^M = 1$, virtual node M is mapped to physical node m .

$$\forall M \in G_V : \sum_{m \in G_s} x_m^M = 1 \quad (2)$$

$$\forall m \in G_s : \sum_{M \in G_V} x_m^M \leq 1 \quad (3)$$

Equations (2) and (3) ensure that a virtual node must correlate with just one substrate node.

•Proposition 2: node capacity

$$\forall M \in G_V, \forall m \in G_s : x_m^M \cdot C(M) \leq C(m) \quad (4)$$

As shown in (4), to successfully embed a VN, all the virtual nodes must be embedded on substrate nodes with adequate capacity.

•Proposition 3: link mapping

$$\forall LMN \in G_v, \forall lmn \in G_s : x_{LMN}^{lmn} = \begin{cases} 1, & \text{if } LMN \text{ is mapped on } lmn \\ 0, & \text{otherwise} \end{cases} \quad (5)$$

In (5), if $c(f) = g(f) + h(f)$, virtual link L_{MN} is mapped on physical link l_{mn} .

$$\begin{aligned} &\forall L_{MN} \in G_v, \quad \forall l_{mn} \in G_s : \\ &\sum_{l_{mn}} x_{LMN}^{l_{mn}} - \sum_{LMN} x_{LMN}^{l_{mn}} = \begin{cases} 1, & \text{if } x_M^m = 1 \\ n, & \text{if } x_N^n = 1 \\ 0, & \text{otherwise} \end{cases} \end{aligned} \quad (6)$$

Equation (6) specifies that the physical path can be split.

•Proposition 4: link bandwidth and delay

$$\forall L_{MN} \in G_v, l_{mn} \in G_s : x_{LMN}^{l_{mn}} \cdot B(L_{MN}) \leq B(l_{mn}) \quad (7)$$

$$\forall L_{MN} \in G_v, l_{mn} \in G_s : x_{LMN}^{l_{mn}} \cdot D(L_{MN}) \geq D(l_{mn}) \quad (8)$$

As shown in (7) and (8), to successfully embed G_v , all the virtual links must be embedded on substrate links with available bandwidth and delay.

•**Proposition 5: node energy consumption**

$$\forall m \in G_v : E_m = \begin{cases} m_{\cdot base} + pl \cdot \tau, & \text{if the selected physical node } n \text{ is open} \\ 0, & \text{otherwise} \end{cases} \quad (9)$$

where $m_{\cdot base}$ is the baseline power without any central processing unit (CPU) load, pl represents the energy proportion factor, $m_{\cdot max}$ denotes the total power at maximum capacity ($pl = m_{\cdot max} - m_{\cdot base}$), and τ denotes the CPU utilization of node m ($\tau = C(M)/C(m)$). When the node is powered off or in the hibernation state, the energy consumption of the node is 0.

•**Proposition 6: link energy consumption**

$$\forall l_{mn} \in G_s, \\ E_m = \begin{cases} l_{mn \cdot base}, & \text{if the selected physical link } n \text{ is open} \\ 0, & \text{otherwise} \end{cases} \quad (10)$$

$l_{mn \cdot base}$ indicates the link energy consumption, which is generally constant. When the link is powered off or in the hibernation state, the link energy consumption is 0.

•**Proposition 7: the status of substrate resources**

The variable s^t is a binary variable that demonstrates that the substrate node or link is turned on/off at time t and is described in (11). It should be noted that the turned-on (not-off) servers should have a greater chance of mapping because they need less electrical energy to host new nodes.

$$s^t = \begin{cases} 1, & \text{if the physical resource is turned on at time } t \\ 0, & \text{if the physical resource is turned off at time } t \end{cases} \quad (11)$$

As shown in (11), if $s^t = 1$, the status of the substrate node or link is turned on.

•**Proposition 8: common substructure isomorphism**

Given two graphs $G_1 = (N_1, L_1)$ and $G_2 = (n_2, l_2)$, if $\exists f: N_1 \rightarrow n_2$. Then, we must be able to find link $L_1 \rightarrow l'$ among them induced by f , where $L_1 = U_{(M,N) \in N} \{L_{MN}\}$, $l' = U_{(m,n) \in N'} \{L_{mn}\}$, where $l' \subseteq l_2$. The two graphs $G_1 = (N_1, L_1)$ and $G' = (n_2, l')$ have the same structure; thus, these nodes and links are a common substructure isomorphism of G_1 and G_2 induced by f . As a rule of thumb, the common substructure is typically a minimally connected graph and is not unique in the sense that it can be determined by different maps.

Fig.2 shows the common substructure isomorphism of G_1 and G_2 with four nodes colored green, where nodes A and B are mapped to nodes a and b, and edge A-B is mapped to edge a-b. Then, nodes B, C and D need to map to nodes b, c, and d, and edges B-C and B-D need to map to edges b-c and b-d, subsequently following a graph edit function. For this graph edit problem, we set this process beginning with the source node, N_{source} (the A in G_1) and n_{source} (the a in G_2), which are always

selected by a given method. Node N_a (B in G_1), which has been mapped to n_a (b in G_2), is denoted as an active node, and its task is to explore new nodes. Nodes that are surrounded by active nodes are called *passive nodes*, which are denoted as N_p (the C in G_1). Its task is to act as the "next" node to be mapped. The nodes in G_2 that have not been exploited are called *free nodes*, denoted as n_f . The edge in G_1 between N_a and N_p is marked as L_{ap} , and the edge in G_2 between n_a and n_p is marked as l_{ap} .

4 Proposed VNE_MR algorithm

In VNE_MR, instead of focusing on research availability and searching the whole substrate network, we narrow the solution space to a certain range by building resource management regions for solving the VNE problem efficiently; however, considering various constraints, the problem is still very complex.

Proposition. Building a resource management region with resource constraints is NP-hard.

Proof We use a reduction from the well-known NP-hard problem MAX-SAT [33]. An instance of MAX-SAT is defined by (CF, w) , where CF is a set of Boolean clauses such that each clause $C \in CF$ is a disjunction of literals with a positive weight $w(CF)$. Let $X = \{x_1, \dots, x_{2n}\}$ be the set of Boolean variables in the clauses of CF . A literal is a variable $x \in X$ or its negation $\bar{x}$. For each $x_i \in X$, let $x_i = 1$ ($x_i = 0$ resp.), $x_{n+i} = \bar{x}_i = 1 - x_i$, a clause $C_j = x_{j_1} \vee x_{j_2} \vee \dots \vee x_{j_{k_j}} \in CF$ can be considered a function on x as follows:

$$C_j = C_j(x) = 1 - \prod_{i=1}^{k_j} (1 - x_{j_k}) \quad (12)$$

Thus, its goal, as shown in (13), is to find an assignment of these variables that maximizes the weight of the clauses. The value of a truth assignment CF is defined

as:

$$\text{Max} \left(\sum_{C_j \in CF} w(C_j) C_j(x) \right) \quad (13)$$

If we solve the resource management region, denoted as $R(G_v)$, and build the problem by considering the resource constraints (Propositions 1 to 7) as variables, taking add one physical resource to $R(G_v)$ as a clause and taking the energy consumption inverse value as the clause weight, then we can obtain a VNE solution of MAX-SAT: $x_1 = 1$ (or 0) if the capacity of n_a is more than N_a ; $x_2 = 1$ (or 0) if the capacity of n_p satisfies the requirement of N_p ; $x_3 = 1$ (or 0) if n_p is turned on; $x_4 = 1$ (or 0) if the bandwidth of l_{ap} satisfies the constraint of L_{ap} ; $x_5 = 1$ (or 0) if the delay of l_{ap} satisfies the constraint of L_{ap} ; $x_6 = 1$ (or 0) if l_{ap} is turned on. The cost of the solution is equal to the sum of the costs of the clauses, which describes the energy consumption of this $R(G_v)$. This ends the **proof**. $\square$

Therefore, given that building $R(G_v)$ with resource constraints is an NP-hard problem, we turn our attention to finding a feasible solution that is near optimal by delimiting the highest potential physical region in the substrate network. However, when delimiting $R(G_v)$, it is not hard to conclude that restrictions in a resource management region that are too tight (in extreme cases, treating each node and link as a unit) will most likely result in rejecting some virtual demands; or restrictions in a resource management region that are too loose (in extreme cases, treating the whole substrate graph as a unit) will occupy many unnecessary physical resources and identify additional hidden nodes, accordingly increasing embedding costs and consuming more power. Therefore, when we delimit $R(G_v)$, an edge-based graph edit distance method CSLGED is used to ensure that the size of $R(G_v)$ is nearly the same as the size of G_v . The main advantage in delimiting $R(G_v)$ relies on its ability to "copy" the information of G_v into $R(G_v)$ step by step so that $R(G_v)$ can be controlled at a predicted size. Among this process, we use this approach to edit each element, i.e., virtual nodes and links, in G_v to their counterparts in $R(G_v)$ and calculate the edit distance from G_v to $R(G_v)$, considering the following constraints, namely, CPU, bandwidth, and end-to-end delay to find a physical region to host G_v .

More details about the steps in our VNE algorithm (VNE_MR) are discussed in the following subsections. The method of delimiting the $R(G_v)$ in the substrate graph is introduced in Section 4.1, and Section 4.2 introduces an improved SMO algorithm for pruning the substrate graph rapidly to find the highest potential $R(G_v)$ that is the final near-optimal VNE solution.

4.1 Delimiting the resource management region

In the previous VNE algorithm, the mapping process typically begins with node mapping and then embeds the links by connecting the respective endpoints, which ignores the network topology constraints on mapping location and may scatter the position of the mapped nodes, leading to a high probability that a substrate link may cross several substrate nodes and yield more resources to be occupied. Therefore,

when solving the VNE problem, we propose a method for delimiting $R(G_v)$ by using an edge-based graph edit distance CSLGED to edit G_v to $R(G_v)$, which focuses on links and taking their end nodes follow directly as a byproduct to strictly control the node locations and ensure the coordination between nodes and links.

The motivation behind using the link-based approach is to facilitate allocating all virtual resources belonging to a specific VNR on the corresponding substrate resources that have enough resources to host the demands of the virtual nodes and edges without using any additional hidden substrate resources, which guarantees using networked resources accurately. Before discussing the implementation of this method, we provide a brief introduction to CSLGED's basics. CSLGED is an edge-based mapping method that is used for computing graph edit distances via common substructure isomorphism enumeration and uses a depth-first backtracking search combined with heuristics to reduce memory requirements and quickly prune a large portion of the mapping search space. The goal is to transform graph G_1 into graph G_2 ($G_1 \neq G_2$) with minimal cost $c(f)$, which is defined as:

$$c(f) = g(f) + h(f) \quad (14)$$

where f is a function of finding the common substructure isomorphism between G_1 and G_2 ; $c(f)$ denotes the total edit cost on the mapped vertices and edges; $g(f)$ denotes the induced edit cost on the mapped vertices and edges in the common substructure isomorphism; and $h(f)$ is the edit cost of the remaining part (i.e., remaining nodes and edges in the graph that are not in the common substructure isomorphism). It is not difficult to find that in this typical graph edit distance computing problem, there are no weights on the nodes or links. However, in the VNE problem, not only are structural constraints embedded in the connection relationship between nodes and links but also resource constraints for them, including capacity, bandwidth and delay, and they are typically nonuniform. Therefore, when editing G_v into $R(G_v)$, we aim to minimize energy consumption (as shown in Formula 15, where energy consumption is posed as the edit distance), which can best illustrate the cost of this process because every time a node or link is turned on, there's a corresponding energy consumption, and formulate the bounds for $R(G_v)$ by presenting the restrictions from two aspects: resource restrictions and structure restrictions. The formulation is given below.

Objective

$$C = \sum_{(m,n,lmm) \in R(G)} E_m + El_{mn} \quad (15)$$

Resources restrictions: Capacity constraints:

$$\forall M \in G, \quad \forall m \in R(G_v) : x_M^m \cdot C(M) \leq C(m) \quad (16)$$

Bandwidth constraints:

$$\forall L_{MN} \in G_v, \forall l_{mn} \in R(G_v) : x_{L_{MN}}^{l_{mn}} \cdot B(L_{MN}) \leq B(l_{mn}) \text{ and } m \neq n \quad (17)$$

Delay constraints:

$$\begin{aligned} \forall L_{MN} \in G_v, \forall l_{mn} \in R(G_v) : \\ x_{L_{MN}}^{l_{mn}} \cdot D(L_{MN}) \geq D(l_{mn}) \text{ and } m \neq n \end{aligned} \quad (18)$$

Mapping constraints:

$$\begin{aligned} \forall M \in G_v, m \in G_s : \\ x_M^m = \begin{cases} 1, & \text{if } M \text{ is mapped on } m \\ 0, & \text{otherwise} \end{cases} \end{aligned} \quad (19)$$

$$\begin{aligned} \forall l_{mn} \in G_s, \\ E_m = \begin{cases} l_{mn} \cdot \text{base} & \text{if the selected physical link } n \text{ is open} \\ 0, & \text{otherwise} \end{cases} \end{aligned} \quad (20)$$

Structure restrictions: Inspired by CSLGED, we designed the structural restriction of $R(G_v)$, which is called the common substructure isomorphism constraint:

$$\begin{aligned} \forall (M, N, L_{MN}) \in G_v, \forall (m, n, l_{mn}) \in R(G_v) : \\ \text{if } x_{L_{MN}}^{l_{mn}} = 1; x_M^m = 1 \end{aligned} \quad (21)$$

Constraints (16)-(20) ensure the legal use of resources in $R(G_v)$; please refer to Section 3.3 for the specific meaning of the symbols. Constraint (21) is inspired by Proposition 8 in Section 3.3, and it ensures that links are allowed to match only if their composing nodes are consistent with the previously matched nodes and the nodes at both ends of the link satisfy the resource constraints. Under the resources and structure restrictions, even though the link map space seems to be relatively large, resource and structure restrictions prune considerable redundant information and sharply reduce the search space.

Algorithm 1 delimiting the resource management region

```

Input:
 $G_s$  and  $G_v$ 
while  $t \neq 0$  do
  for each VNR arriving at the SN randomly at time  $t$  do
    Formulate  $Set_L$ 
    for each  $L_{vw}$  in  $Set_L$  do
      if  $B(l_{vw}) > B(L_{VW})$  and  $D(l_{vw}) < D(L_{VW})$  then
        if  $C(n_p) > C(N_p)$  then
          match  $L_{VW}$  to  $l_{vw}, w$  to  $W$ 
        else
          extend to the neighbor node of  $w$ 
        end if
      else
        split the other link of  $V$  with sufficient bandwidth, and select the first node whose capacity
        meets the constraint as  $W$  from the neighbor node of the node at the other end of the
        link
      end if
      calculate  $g_{vne}(f)$  using eq.15
      update the physical resource capacity
    end for
    use the shortest path algorithm to match the remaining links calculate  $h_{vne}(f)$  using eq.15
    record  $c_{vne}(f)$  and  $c_{vne}(f) = g_{vne}(f) + h_{vne}(f)$ 
  end for
end while

```

Since our method is link-based, editing the virtual link set in G_v step by step to a physical link set is our task, and during this process, the parameters of both links must be unanimous (including the end node following directly); additionally, the values also satisfy the demands. Next, we provide more information about the steps for delimiting $R(G_v)$ by editing G_v into a similar-sized substrate region under the resource and structure restrictions. The pseudocode is shown in Algorithm 1.

At each time t , if a VNR arrives, first, we need to structure the common substructure isomorphism of G_v and $R(G_v)$ according to the CSLGED. For G_v , in accordance with the structure restrictions, a minimally connected graph without loops is obtained by using the Prim algorithm (an edge-based minimum spanning tree method), which takes the link bandwidth reciprocal as the weight. Thus, we can obtain a subset of $G_v = (N, L)$, denoted as $G'_v = (N, L')$, where $L' \subseteq L$. Then, we start from N_{source} and output the link set $Set_L = \{L'_1, L'_2 \dots L'_n\}$ (where $n < |L|$) of G'_v in a breadth-first manner with the link's bandwidth as the weight (Fig. 3). Next, starting from the adjacent link of n_{source} that is matched to the N_{source} (introduced in Section 4.2), we search the physical resources that meet the requirements of each link in Set_L in the same breadth-first manner and output the physical resources set $Set_l = \{l'_1, l'_2 \dots l'_n\}$ corresponding to Set_L . In this process, we check links one-to-one directly: if each element in Set_l has enough resources to satisfy the demands of its counterpart in Set_L , the first parameter in G_v is compared to the first parameter in $R(G_v)$, the second parameter to the second, and so on for all the remaining parameters. If the resource constraint check results are true, that is, each virtual link found a matching substrate link, as the nodes follow directly as the byproduct, embedding the virtual nodes and edges is realized together in full coordination onto the corresponding physical substructure isomorphism. Therefore, the resource matching in the common substructure isomorphism can be regarded as a one-stage mapping that highlights the link status. Note that because each virtual node can map to only one physical node, while the links can be split, an energy-saving region must have a compact structure that identifies as few links as possible

between a fixed number of physical nodes. However, CSLGED backtracks the edge mapping space in a depth-first manner, which violates our requirement. Therefore, we modify the search method to proceed in a breadth-first manner.

To facilitate understanding, an example of matching one virtual link in G_v to its counterpart in $R(G_v)$ is illustrated in Fig. 4-Fig. 6. The request graph is composed of three links and tree nodes, and the substrate graph is composed of six links and five nodes, in which a virtual node N_a is mapped to the physical node n_a . Next, when matching the adjacency links L_{ap1} and L_{ap2} of N_a (Fig. 4), if l_{ap1} satisfies all constraints of L_{ap1} , i.e., both the bandwidth and delay of l_{ap1} meet the L_{ap1} requirements, and n_{p1} at the other end of l_{ap1} also meets the capacity requirements of virtual node N_{p1} , then we match L_{ap1} to l_{ap1} . Similarly, if L_{ap2} meets all the requirements of l_{ap2} , we match L_{ap2} to l_{ap2} . However, if the bandwidth constraint cannot be met, then we match L_{ap1} to another adjacent link of N_a with sufficient bandwidth (Fig. 5), l_{ap2} is split to host L_{ap1} and L_{ap2} , and N_{p1} is matched to an n_{p2} -adjacent node that satisfies the N_{p1} capacity constraint. If the link's resource constraint is satisfied, but the capacity of the node at the other end of this link is not satisfied (Fig. 6), then we match N_{p1} to an adjacent node of n_{p1} that meets the capacity constraint.

Next, Equation (15) is used to calculate the editing distance $g_{vne}(f)$ during the editing process of the substructure isomorphism of G_v into that of $R(G_v)$, that is, the energy consumption of the physical resources turned on in $R(G_v)$ to hold this virtual request.

Finally, we edit the remaining part of G_v excluding the common substructure isomorphism to $R(G_v)$ by finding their shortest paths in the common substructure isomorphism and calculate the energy consumption of the physical resources that turn on in this region, denoted as $h_{vne}(f)$. Given that the remainder is only partial connection information between the nodes, to ensure the VNE service quality, we relax the restriction of embedding when editing this part: only the resource constraints are retained.

In general, using $R(G_v)$ to host VNRs provides an absolute advantage in controlling the resources usage by limiting $R(G_v)$ to a scale that is almost as small as G_v so that only a few of the necessary resources are opened. Moreover, it can also provide

flexible and predictable services for users. Since the bounds of $R(G_v)$ are set by the resource constraints and the structure constraints, we can make a new bound for $R(G_v)$ by changing the constraints in the restrictions and find a new $R(G_v)$ to host G_v while fitting the new requirement.

4.2 Searching the highest potential resource management region

Different from the previous metaheuristic algorithm that traverses the solution space with a fixed population [34]-[35]. SMO is an algorithm that simulates the intelligent foraging behavior of spider monkeys with a fission-fusion social structure (FFSS). In FFSS, a swarm is a social grouping pattern in which individuals form temporary subgroups whose members belong to a larger unit group. There can be fluid movement between subgroups and unit groups such that group composition and size change frequently and vice versa based on the real-time situation of the search to control the optimization speed. Therefore, when we search for the highest potential $R(G_v)$ in a parallel way in the SN, we take advantage of this feature of SMO to converge to a potential physical region quickly by searching with a small-size group in the SN at the beginning. When the search is stagnant, we expand the population size by splitting the group into a larger group to explore and repeat the previous

search phases until we find the near-optimal solution. Therefore, during this optimization process, we can balance the exploration and exploitation capability of the optimization algorithm while maintaining the convergence speed.

There are six stages in the algorithm: the local leader phase, global leader phase, local leader learning phase, global leader learning phase, local leader decision phase and global leader decision phase. The first and second phases are used to explore the search region while generating the new positions, i.e., the new solutions to this problem, for all the group members by using self experience, local/global leader (the spider monkey with the greatest fitness in its subgroup/unit group) experience and group member experience to promote the exploitation. The third and fourth phases, are used to determine whether the local best and global best solutions are updating in a predefined number of trials. If not, then the solution is considered stagnated. The fifth and sixth phases are used to avoid the stagnation or premature convergence of the local solutions/global best solution and control the population size. If the local best solution is not updated in a predefined number of trials (*LocalLeaderLimit*) then all the members of that group are reinitialized. If the global best solution is not updated within a predefined number of trials (*GlobalLeaderLimit*) then the group is divided into smaller subgroups.

To clarify how we use and enhance SMO to find the highest potential $R(G_v)$, a step-by-step scenario graph describing the overall process is shown in Fig. 8 and the implementation process is explained below:

Initialization:

Initially, we generate a population of P spider monkeys (SM s) as the VNE solutions, i.e., delimit P $R(G_v)$ in the substrate graph. SM_i represents the i th SM in the population, i.e., i th $R(G_v)$. The local/global leader represents $R(G_v)$ with the greatest fitness in its subgroup/swarm and MG represents the maximum number of groups in the swarm. However, as this optimization process starts with a single group having all the SM s, in the beginning, the local leader and global leader are both the same SM . Moreover, because the goal of this study is to find the most energy-saving $R(G_v)$, each SM_i corresponds to the i th $R_i(G_v)$. Therefore, the fitness of each SM_i is equal to the energy consumption of physical resources opened in $R_i(G_v)$:

$$\text{Fitness} = E(R(G_v)) = \sum_{n \in G_v} E(n) + \sum_{l \in R(G_v)} E(l) \quad (22)$$

In addition, according to the previous section, the following conclusion can be easily drawn: to build an energy-saving $R(G_v)$ with a compact structure in the SN, a proper *source node* is essential. Therefore, to ensure $R(G_v)$ has a compact structure, the region must have short links, and a high utilization rate is high; thus, Equation (23) is presented to calculate the virtual node N_{source} of G_v , which is the node with the shortest distance sum from all other nodes in the graph; the formula for calculating N_{source} is:

$$N_{source} = \text{Min}(\text{sum}(\text{hops from other nodes to the } N_{source})) \quad (23)$$

Then, P physical nodes, which are the candidate mapping nodes of N_{source} , are randomly selected in the SN and each physical node is marked as $n_{sourcei} = 1 \dots P$ to build the $R(G_v)$ (Fig. 7 (a): ③ - ①). A detailed description is provided in Lines 1-5 of Algorithm 3.

Local leader phase (LLP) and global leader phase (GLP):

To accelerate the convergence speed, in the LLP, every SM generates its current position based on the experience information of the local leader as well as local group members. The fitness value of the obtained new position is calculated. The

Algorithm 2 Local leader decision and global leader decision phase

```

for  $k$  th group in  $MG$  do
  if  $LLC_k >$  local leader limit then
     $LLC_k = 0$ 
    for  $SM_i$  in the  $k$  th group do
      if  $U(0,1) \geq pr$  then
        random select a  $n_{sourcei}$  from the substrate network and initialize  $SM_i$ 
      else
        initialize  $SM_i$  using (19)
      end if
    end for
  end if
end for
if global limit count  $>$  global leader then
  global limit count = 0
  if number of groups  $< MG$  then
    divide the population into subgroups
  else
    combine all the groups to make a single group
  end if
  Update Local Leader's position
end if

```

position update equation for the i th SM (which is a member of the k th subgroup) is:

$$SM_{\text{new}i} = SM_i \oplus [r_1 \cdot (LL_k \ominus SM_i) + r_2 \cdot (SM_r \ominus SM_i)] \quad (24)$$

where LL_k represents the k th subgroup leader position. SM_r is the r th SM which is chosen randomly within the k th group such that $r \neq i$. r_1 and r_2 represent the two random variables that are uniformly distributed in $[0, 1]$ and $r_1 + r_2 = 1$. As VNE is a discrete problem, numeric computing is divided into taking one step (in one-hop units) toward LL_k or SM_r . As shown in Fig. 7(a), after we calculate the updating direction of each SM_i ($i=1,2,3$) by using the contents in the brackets of (17), we move SM_i from the original position (black dotted circle in Fig. 7(a) - ①) to this direction by moving the center node n_{source} of $R_i(G_v)$ (yellow node in Fig. 7(a) - ②) to the calculated direction by one hop (green node in Fig. 7 (a) - ③). $\ominus$ denotes the AND operation. $\oplus$ denotes the calculation of the shortest path between two n_{source} of $R(G_v)$.

In the global leader phase, all SMs update their position using the experiences of the global leader and local group members. The position update equation for this phase is as follows:

$$SM_{\text{new}i} = SM_i \oplus [r_1 \cdot (GL \ominus SM_i) + r_2 \cdot (SM_r \ominus SM_i)] \quad (25)$$

Global leader learning (GLL) phase and local leader learning (LLL) phase:

The local leaders and global leader are updated in this phase by applying greedy selection in its subgroup and the population, respectively. Furthermore, in the LLL phase, we check whether the position of the local leader in its group is updating, and if not, then the *local limit count* is incremented by 1. In the GLL phase, we check whether the position of the global leader in the whole population is updating, and if not, then the *global limit count* is incremented by 1.

Local leader decision (LLD) phase and Global Leader Decision (GLD) phase:

If any local leader position is not updated to the threshold *local leaderlimit*, then the members in that subgroup update their positions either by random initialization or by collective information from the global leader and local leader through (19), based on the *pr*. The pseudocode of this phase for the *k*th group is shown in Lines 1-8 of Algorithm 2.

$$SM_{\text{newi}} = SM_i \oplus [r_2 \cdot (GL \ominus SM_i) + r_3 \cdot (SM_i \ominus LL_k)] \quad (26)$$

where *LLCk* is the trial counter for the local best solution of the *k*th group.

In the global leader decision phase, if the position is not updated to the threshold *global leader limit*, then the population is divided into smaller groups (Fig. 7 (b)). At first, the population is divided into two groups, and then three groups, etc.; each time, the local leaders in the newly formed groups are elected until *MG*'s limit is reached. In the case in which the maximum number of groups is formed, then the global leader combines all the groups to form a single unit group and further selects the position with the highest fitness as the solution to the problem. The operation in this phase is shown in lines 9-15 of Algorithm 2. The complete pseudocode of the proposed strategy is given in Algorithm 3.

Algorithm 3 VNE_MR

```

1: Input:
    $G_s$  and  $G_v$ ,  $P$ ,  $MG$ , localleaderlimit, globalleaderlimit, pr.
2: calculate the  $N_{\text{source}}$  of  $G_v$  use eq.23
3: randomly select  $Pn_{\text{source}i}$  ( $i = 1 \dots P$ )
4: for each  $n_{\text{source}i}$  do
5:   call Algorithm 11 to delimit  $R_i(G_v)$  corresponding to the  $SM_i$ 
6:   use (22) to calculate  $SM_i$ 's fitness
7: end for
8: Select global leader and local leaders by applying greedy selection
9: while (termination criteria is not satisfied) do
10:  for  $k$  in  $MG$  do
11:    generate a new position for each  $SM_i \in k$  th subgroup using eq.18
12:    produce new position for each  $SM_i \in k$  th subgroup using eq.19
13:    update global leader and local leaders; record the position and the fitness
14:    call Algorithm 2 to update all the groups
15:  end for
16: end while

```

Finally, for SMOs to play a better role in VNE problems, several points must be noted.

First, traversing the substrate network to find the optimal region quickly is a discrete problem that needs to be solved. To speed up convergence, we set every *SM* to use its self-experience, leader experience and group members experience to update its position in the LLP and GLP instead of having a certain probability to update.

Second, during the optimization process, if the capacity of a selected n_{source} does not meet the requirements of its corresponding virtual node N_{source} , this subgroup to which it belongs is pruned (Fig. 7 (a)). Regardless of which direction we select on the update, we must obey resource and structure restrictions as the proposed bounds to achieve rapid pruning from the surrounding physical topology. After all, if one constraint is not satisfied, then this $R(G_v)$ cannot be used, and the next endeavor will be meaningless.

Third, the search speed will be faster with a large population; however, gaining this advantage may sacrifice some computational time. If there are N SM s in the population, then the time complexity of VNE_MR is $O(|N_v| + |L_v| + |L_{R(G_v)}| + |N_{R(G_v)}|^2) \cdot |P|$, where $|N|$ represents the number of nodes, $|P|$ represents the number of links, and $|L|$ represents the total number of SM s used in this algorithm, which has a direct effect on the computational time. Therefore, initially, there is a small group in our algorithm so every newly generated position is attracted toward the best position; the population is only divided into smaller subgroups to expand exploitation when regeneration stagnates.

5 Performance evaluation

In this study, we designed two sets of experiments to demonstrate the resource management ability of the proposed VNE_MR algorithm. In the first group, to verify the overall performance of VNE_MR, VNE_MR is compared to VNE_TAGRD [14] and OPaCoVNE [16], all of which have the same goal: to solve the virtual network embedding problem using fewer resources and consuming less power. Table 1 provides a high-level comparison of VNE_MR and those methods, listing their strategies and resource constraints.

In the second group, VNE_MR is compared to EE_CTA [26], which also uses a metaheuristic algorithm to find the near-optimal solution in the substrate network, to verify that our optimization method design based on SMO decreases the computational time while ensuring the quality of service. (Of course, VNE_MR still requires considerable computational time compared to other methods without optimization, but it's worth to make the sacrifice for controlling networked resources accurately, we will not discuss this consensus issue in detail.) VNE_MRWD is a VNE algorithm without delay requirements, which is designed as VNE_MR's comparison experiment. It is presented to redelimit a new $R(G_v)$ by changing the resource constraint model introduced in Section 4.1 and then solving the VNE problem by using the improved SMO to find the highest potential $R(G_v)$ to host G_v . Therefore, it is designed to verify the flexibility of VNE_MR. Table 2 provides a high-level comparison between VNE_MR and those methods, listing their resource constraints and other parameters.

Table 1 TABLE 1 COMPARING VNE_MR TO VNE_TAGRD AND PaCoVNE

Item	VNE_MR	VNE_TARGET	OPaCoVNE
Embedding strategy	building a resource management unit by delimiting an $R(G_v)$ similar in size to G_v	focusing on the availability of resources	building a resource management unit by treating pairs of nodes and their edge as a segment
Power consumption	Yes	Yes	Yes
End-to-end delay	Yes	Yes	Yes

Table 2 TABLE 2 COMPARING VNE_MR TO VNE_MRWD AND EE_CTA

Item	VNE_MR	VNE_MRWD	EE_CTA
Optimization strategy	Based on the SMO	Based on the SMO	Based on the Non-dominated Sorting Genetic algorithm
Power consumption	Yes	Yes	Yes
End-to-end delay	Yes	No	No

5.1 Simulation Settings

Substrate network and virtual network topologies were randomly generated using the GT-ITM tool, and the mean probability of a pair of two nodes being connected

was set equal to 0.5. The substrate network included 100 nodes, *CPU* and *BW* resources were given as real numbers uniformly distributed between 50 and 100, and delay in each substrate edge was randomly selected between 1 and 25. The number of virtual nodes per VNR was randomly determined by a uniform distribution between 2 and 10. The virtual *CPU* and *BW* were real numbers uniformly distributed between 0-20 and 0-50, respectively, while delay in each virtual edge was randomly selected between 20 and 100. The VNR arrival process was simulated as a Poisson process, and its average arrival rate was set to 5 VNRs per 100 time units. Its residence time obeyed an exponential distribution with an average value of 500 time units. To map a sequence of 2,000 VNRs, our simulation lasted for approximately 40,000 time units. A time window was equal to 2,000 time units. To obtain final representative and stable experimental results, we ran all of our simulations in 20 time windows and collected the experimental data after 5 time windows. To ensure acceptable convergence of the SMO used in VNE_MR, we investigated the values of best and average fitness in different generations. The parameters used in the SMO were $P = 5$, $MG = 3$, $localleaderlimit = 2$, $globalleaderlimit = 3$, and $pr = 0.9$.

5.2 VNE Performance Metrics

The comprehensive ability to manage the resources of the VNE problem can be judged in terms of the following metrics.

1. Average number of open nodes (N_{ON}):

$$N_{ON} = \frac{\sum_{i=1}^{n_T} n_{oi}}{n_T}$$

where n_T represents the number of all valid time periods from 0 to T and n_{oi} represents the number of physical nodes that are active in effective period i .

2. The average number of open links (N_{OL}) is

$$N_{OL} = \frac{\sum_{i=1}^{n_T} l_{oi}}{n_T}$$

where l_{oi} represents the number of open substrate links at time period i .

3. The average utilization of CPUs (U_{CPU}) is

$$U_{CPU} = \frac{\sum_{i=1}^{n_T} n_{CPUi}}{n_T}$$

where n_{CPUi} represents the CPU utilization of the node resources at time slice i .

4. The average utilization of bandwidth (U_{BW}) is

$$U_{BW} = \frac{\sum_{i=1}^{n_T} l_{BW_i}}{n_T}$$

where l_{BW_i} represents the bandwidth utilization of the link resources at time slice i .

5. The average energy consumption (E) is

$$E = \frac{\sum_{i=1}^{n_T} \left(\sum_{N \in N_v, n \in n_s} E(n) + \sum_{L \in L_v, l \in l_s} E(l) \right)}{n_T}$$

where $(\sum_{N \in N_v, n \in n_s} E(n) + \sum_{L \in L_v, l \in l_s} E(l))$ is the substrate resource consumption at time period i .

6. The average ratio of revenue and cost (R_{RC}) is

$$R_{RC} = \frac{\sum_{i=1}^{A_T} \text{Revenue}_{Ri}}{\sum_{i=1}^{A_T} \text{Cost}_{Ri}}$$

where A_T represents the number of virtual networks that are received within the time period 0 to T and Revenue_{Ri} and Cost_{Ri} represent the revenue and cost, respectively, of the mapped virtual network Ri .

$$\text{Revenue}_{Ri} = \sum_{N \in N_v} C_{\text{accept}}(N) + \sum_{L \in L_v} B_{\text{accept}}(L)$$

$$\text{Cost}_{Ri} = \sum_{n \in N_v} x_M^m \cdot C(n) + \sum_{l \in E_v} x_{LMN}^{l_{mn}} \cdot B(l)$$

where $C_{\text{accept}}(N)$ represents the number of requested CPU resources for virtual node N that has been accepted and $B_{\text{accept}}(L)$ represents the amount of requested bandwidth for virtual link L that has been accepted.

5.3 Discussion of VNE_MR's overall performance

In the first group, compared to other algorithms, VNE_MR has the smallest number of open nodes (≈ 50), indicating that the proposed algorithm achieves good results in node opening control (Fig.9). We calculate node utilization while collecting the number of open nodes (Fig.10). The VNE_MR node utilization rate (73%)

exceeds those of the other algorithms (VNE_TAGRD $\approx 62\%$, OPaCoVNE $\approx 70\%$). As shown in Fig.11, after 5 time windows, the number of open links with VNE_MR is only 100 units, while that with VNE_TAGRD is 230 units, which is more than twice as high. The link resource utilization rate (40%) of VNE_MR is much higher than those of the other algorithms (VNE_TAGRD $\approx 28\%$, OPaCoVNE $\approx 32\%$) (Fig.12). According to Propositions 5 and 6, we can conclude that these performances directly affect the final average energy consumption (Fig. 13). This result likely occurs because VNE_TAGRD, a two-stage algorithm, only focuses on resource availability and ignores the coordination between nodes and links. Nodes selected in this manner may be scattered, which results in the use of unnecessary substrate resources. OPaCoVNE manages resources in several resource management units, named segments so that the original topology structure of G_v is disturbed when requirements are allocated in these segments. As the nodes and links are firmly bound in the segment, flexibility and fault tolerance are markedly reduced. Thus, these segments can only be separated to find the appropriate physical resource, which leads to opening hidden resources. Fig. 13 shows the energy consumption results. The energy consumption of VNE_MR exhibits an apparent advantage over other algorithms. In addition to the energy savings shown, the ratio of revenue and cost is another advantage of VNE_MR (Fig.14). The ratio is maintained at 86%, and only 50 nodes and 100 links are turned on. The energy consumption is also the lowest among all of the tested algorithms because we use an edge-based graph edit distance method to edit G_v to the physical $R(G_v)$ in the substrate graph, keeping the size of $R(G_v)$ similar to that of G_v , so that when mapping a VNR, only the minimum network resources need to be turned on, and no hidden resources are used. In addition, if this direct approach fails to find the optimal physical region in the SN, we also designed a metaheuristic optimization method based on SMO to find the highest potential $R(G_v)$ in a parallel manner.

In the second group, Fig. 15 shows that VNE_MR achieves a lower average computational time compared to EE_CTA because there is only a small population that is used to search in the SN at the beginning, and they converge to a potential physical area quickly. Only when the regeneration stagnates do we expand the population size by combining the group into a larger population to explore and repeat the previous search phases until we find the near-optimal solution. The time complexity of our algorithm is $o(|N_v| + |L_v| + |L_{R(G_v)}| + |N_{R(G_v)}|^2) \cdot |P^1|$, where $|P^1| = \sum_{i=1,2,\dots,MG} P_i$, while the time complexity of EE_CTA is $o(|N_v| + |L_v| + |L_{R(G_v)}| + |N_{R(G_v)}|^2) \cdot |P^2|$, where $|P^2|$ is a constant. It can be

Figure 15 Average processing time

Figure 16 Average number of open nodes

Figure 17 Average number of open links

Figure 18 Average node utilization

Figure 19 Average bandwidth utilization

Figure 20 Average energy consumption

Figure 21 Average revenue-cost ratio

seen that the size of our population gradually increases to a scale similar to that of the comparative experiment, which has a direct impact on the computational time. VNE_MR also achieves this result by pruning the redundant SN information to reduce the size of the search space. Therefore, we only concentrate on the physical resources in $R(G_v)$, while EE-CTA explores the appropriate physical resources in the SN; obviously, the former's order of magnitude is much smaller than that of the latter. In particular, although our computational time is sharply reduced, it does not cause the $R(G_v)$ to skip the global minima in the long run, as the VNE_MR achieved good results in terms of energy savings and revenue-cost ratio compared with EE-CTA. As shown in Fig. 20, the average energy consumption of VNE_MR is reduced by 7,000 W compared with that of EE-CTA. Fig. 21 shows that the revenue-cost ratio of the VNE_MR algorithm is approximately 12% higher than that of EE-CTA. Figs. 16–19 also confirm this result; VNE_MR opens fewer nodes and links than EE-CTA and obtains greater utilization than EE-CTA.

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Authors' contributions

ZX is in charge of the major theoretical analysis, algorithm design, numerical simulations, draft of the article, and critical revisions. LZ, YS, LJ and WL provided critical revision and final approval of the version to be published. All authors read and approved the final manuscript.