Information Exposure Maximization Problem with Heuristic and Evolutionary Algorithms

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Abstract—This paper explores the Information Exposure Maximization Problem (IEMP) as it applies to reducing echo chambers and filter bubbles in social networks. By representing the network as a graph, we examine the effectiveness of heuristic and evolutionary algorithms in balancing exposure to diverse information. The heuristic approach employs a Monte Carlobased greedy selection strategy, while the evolutionary algorithm uses a genetic algorithm to optimize influence spread. Results indicate that the heuristic method consistently outperforms the evolutionary algorithm in efficiency and evaluation value, especially for larger datasets. However, the evolutionary algorithm with a halt condition shows promise in reducing computational time. This study underscores the utility of heuristic methods in influence maximization and suggests potential improvements in evolutionary approaches for similar applications.

Keywords—evolutionary algorithm, genetic algorithm, heuristic algorithm, optimization

I. INTRODUCTION

The rise of social networks has led to increased occurrences of echo chambers and filter bubbles, where users predominantly interact with like-minded individuals and viewpoints. This phenomenon restricts exposure to diverse information and contributes to polarization. Addressing this, the Information Exposure Maximization Problem (IEMP) aims to enhance exposure to multiple viewpoints in a network by selecting sets of users that maximize the reach of competing campaigns, thus diversifying the information spread within the network.

This project examines IEMP within a graph-based model of social networks, where users and their connections are represented as nodes and edges. By identifying optimal user sets for two distinct campaigns, we aim to achieve a balanced exposure across the network, potentially countering echo chambers. This report investigates two primary algorithmic approaches: a Monte Carlo-based heuristic strategy and a genetic algorithm-based evolutionary strategy. The purpose of this study is to evaluate the effectiveness and efficiency of these approaches and to assess their suitability for real-world applications in influence maximization and information diversification.

II. PRELIMINARY

The IEMP is framed as an optimization algorithmic problem aimed at reducing the effects of echo chambers and filter bubbles, where users are frequently limited to similar viewpoints, isolating them within their own information spheres. This study examines a social network represented as a graph G = (V, E) where V represents users (nodes) and E represents social connections (edges) between them. The objective in the IEMP is to identify two sets of users that achieve the highest balanced exposure to diverse information.

The following sections provide essential definitions of terminology and notations used, and a formal definition of the problem.

A. Terminology and Notations

- Social Network G = (V, E) represents the network graph, where V = {v₁, v₂, ..., v_n} represents the set of nodes, and E ⊆ V×V represents the edges between nodes
- Campaigns $C = \{c_1, c_2\}$ represents two campaigns; each campaign holds its own viewpoint.
- Initial Seed Set $-I_i \subseteq V$, $i \in \{1, 2\}$ represents the initial seed set for campaign c_i .
- Balanced Seed Set $-S_i \subseteq V$, $i \in \{1, 2\}$ represents the target seed set needed to find for each campaign c_i .
- **Budget** k represents the size's upper bound of the two balanced seed sets; that is, $|S_1| + |S_2| \le k$.
- **Diffusion Probability** $-P_i = \{p^i_{(u,v)} \mid (u,v) \in E\}, i \in \{1, 2\}$ represents the edge weight associated with campaign c_i where $p^i_{(u,v)}$ represents the probability of node u activating node v under each campaign c_i .
- **Diffusion Model** M represents the stochastic process for the seed $U_i = I_i \cup S_i$, which initiates the spread of information on graph G. It is assumed that information from the two campaigns spreads through the network following the independent cascade (IC) model, with each campaign's message spreading independentlythis is known as heterogeneous propagation. The diffusion process for the first campaign (with the second campaign following the same process) proceeds as follows:
 - i. At step t = 0, nodes in the seed set U_I are activated, while all the other nodes remain inactive.
 - ii. In each step t, any active user u for campaign c_l attempts to activate each of its inactive outgoing neighbors v with a probability $p^l_{(u, v)}$. This activation is like flipping a coin with a probability of heads equal to $p^l_{(u, v)}$: if heads, v is activated; if tails, v remains inactive. Each active user u has only one chance to activate each neighbor for campaign c_l , after which u stays active but ceases further activation efforts.
 - iii. The diffusion process ends when no more nodes can be activated.
- **Exposed Node Set** $-r_i(U)$ represents the set of vertices that can be reached from U through the cascade

process mentioned above for campaign c_i , for a given seed set U. Note that in a single propagation step, $r_i(U)$ includes not only nodes that were successfully activated by U but also nodes that were targeted for activation by U but remained inactive. Since the diffusion process is stochastic, $r_i(U)$ is a random variable.

B. Problem Formulation

Given a social network G = (V, E), two sets of I_1 and I_2 of the initial seed sets for the two campaigns, and a budget k, the goal of the IEMP is to identify two sets S_1 and S_2 such that $|S_1| + |S_2| \le k$ and to maximize the expected number of vertices that are either influenced by both campaigns or remain unaffected by both; that is,

```
\max \Phi(S_I, S_I) = \max \mathbb{E}[|V \setminus (r_I (I_I \cup S_I) \triangle r_2 (I_2 \cup S_2))|]
such that |S_I| + |S_2| \le k and S_I, S_2 \in V
```

III. METHODOLOGY

As introduced earlier, IEMP's diffusion model is based on a stochastic process. The two algorithms outlined below will leverage this diffusion model within their respective procedures. To facilitate this, it is essential to define an algorithm that accurately represents the diffusion process, which can be accomplished through the application of a breadth-first search (BFS) approach.

ALGORITHM 1: INFLUENCE DIFFUSION VIA BFS

```
Input: graph(G), seed(U_i = I_i \cup S_i), campaign
    (campaign index)
   Output: active (set of active nodes), exposed (set of
   exposed nodes)
   Initialize queue as an empty queue
   Initialize active as an empty set
   Initialize exposed as an empty set
   for node in seed do
5
       queue.Add(node)
6
      active.Add(node)
7
      exposed.Add(node)
8
   while queue is not emoty do
      node \leftarrow queue.Remove()
10
      for neighbor, probab in
       graph.GetAdjList(node) do
11
          if not exposed. Contains (neighbor) then
12
             exposed.Add(neighbor)
13
          if not active. Contains (neighbor) then
14
              if probab \ge RandomValue() then
15
                 active.Add(neighbor)
16
                 queue.Add(neighbor)
17
   return active, exposed
```

A. Heuristic Search

This section outlines a heuristic approach for solving the IEMP using a Monte Carlo-based greedy selection strategy.

The main idea of this heuristic algorithm is to expand the node with the highest incremental influence value, h(v), to maximize exposure for two campaigns.

To achieve a better computational accuracy, the algorithm applies Monte Carlo simulation to approximate each node's influence impact by averaging results from multiple simulated influence spreads. Furthermore, to improve efficiency, the algorithm shall avoid recalculating the entire influence spread from scratch. It should only compute the changes in active and exposed nodes for each candidate node. This is achieved by diffusing influence to only a depth deeper and incrementally updates active and exposed sets.

In this proposed influence diffusion process, each incremental computation is conducted to a depth of one, focusing solely on evaluating the immediate neighbors of the current source node. This approach does not lead to issues with activation or exposure, as the algorithm will subsequently iterate through all nodes, prompting each node to activate its own neighbors. Consequently, there is no need to explore deeper from the initial node, as doing so would introduce unnecessary redundancy in the checks performed.

ALGORITHM 2: INCREMENTAL INFLUENCE DIFFUSION

Input: graph, src (source node), active, exposed, campaign

Output: active_increment (set of newly activated nodes), exposed_increment (set of newly exposed nodes)

- Initialize active_increment as an empty set
- ² Initialize *exposed_increment* as an empty set
- ³ **if** not *active*.contains(*src*) **then**
 - active increment.Add(src)
- ⁵ **if** not *exposed*.contains(*src*) **then**
- exposed_increment.Add(src)
- **for** *neighbor*, *probab* in *graph*.GetAdjList(*src*) **do**
- if not exposed.Contains(neighbor) then

 | exposed_increment.Add(neighbor) |
 | if not active.Contains(neighbor) then
 | if probab ≥ RandomValue() then
- 12 | active_increment.Add(neighbor)
 13 return active_increment, exposed_increment

By limiting the depth to one at this stage, we streamline the process and enhance computational efficiency while ensuring comprehensive coverage of the network in subsequent iterations. Nonetheless, the aforementioned influence diffusion with exhaustive BFS is still utilized in our final algorithm, but only once for each campaign.

ALGORITHM 3: IEMP MONTE CARLO GREEDY HEURISTIC

Input: graph, initial1 (I_1), initial2 (I_2), budget (k), rep (number of simulations)

Output: balanced1 (S_1), balanced2 (S_2)

- $num_nodes \leftarrow graph.GetNumNodes()$
- ² Initialize *balanced1* as an empty set
- ³ Initialize *balanced2* as an empty set
- while balanced1.Length() + balanced2.Length() < budget do</p>
- Initialize $h1_rec$ as an empty array
- 6 Initialize *h2_rec* as an empty array

```
7
       for j \leftarrow 0 to rep-1 do
 8
           union1 \leftarrow initial1 \cup balanced1
           union2 \leftarrow initial2 \cup balanced2
10
           active1, exposed1 \leftarrow
           InfluenceDiffusionViaBFS(graph, union1, 1)
11
           active2, exposed2 \leftarrow
           InfluenceDiffusionViaBFS(graph, union2, 2)
12
           phi_s1_s2 \leftarrow ComputePhi(num\_nodes,
           exposed1, exposed2)
13
           for i \leftarrow 0 to num nodes-1 do
14
               if not balanced1.contains(i) and not
               balanced2.contains(i) then
15
                   active1_increment, exposed1_increment
                   ← IncrementalInfluenceDiffusion(
                   graph, i, active1, exposed1, 1)
16
                   active2 increment, exposed2 increment
                   ← IncrementalInfluenceDiffusion(
                   graph, i, active2, exposed2, 2)
17
                   phi_s1vi_s2 \leftarrow
                   ComputePhi(num_nodes, exposed1 ∪
                   exposed1 increment, exposed2)
18
                   phi_s1_s2vi \leftarrow
                   ComputePhi(num_nodes, exposed1,
                   exposed2 \cup exposed2\_increment)
19
                   h1 \ rec[i] \leftarrow h1 \ rec[i] + phi \ s1vi \ s2
                   phi_s1_s2
20
                   h2\_rec[i] \leftarrow h2\_rec[i] + phi\_s1\_s2vi -
                  phi s1 s2
        for j \leftarrow 0 to num\_nodes-1 do
21
22
           h1\_rec[j] \leftarrow h1\_rec[j] / rep
23
           h2\_rec[j] \leftarrow h2\_rec[j] / rep
24
        new_v1 \leftarrow IndexOfMaxElement(h1\_rec)
25
        new_v2 \leftarrow IndexOfMaxElement(h2\_rec)
26
        if new_v1 \ge new_v2 then
27
           balanced1.Add(new v1)
28
        else
29
           balanced2.Add(new_v2)
30
    return balanced1, balanced2
```

B. Evolutionary Approach

In this section, we address the IEMP by employing a genetic algorithm (GA) to identify optimal intervention sets S_I and S_2 that maximize the expected influence spread within a network. Given the combinatorial complexity and large search space of this problem, optimization techniques like genetic algorithms provide an effective solution method. GAs are well-suited for this task due to their capability to explore complex, multimodal search spaces and to find near-optimal solutions within reasonable computational time. The GA approach is structured as follows:

- **Solution Representation**: Each candidate solution is represented by a binary vector, $x = \{x_1, x_2, ..., x_{|V|}, x_{|V+I|}, x_{|V+2|}, ..., x_{|V+V|}\}$ where $x_i \in \{0, 1\}$.
- **Fitness Function**: To guide the search towards optimal solutions, a fitness function is defined that

evaluates each candidate solution based on the influence spread it achieves. The fitness function distinguishes between feasible and infeasible solutions.

$$fitness(S_1, S_2) = \begin{cases} \Phi(S_1, S_2) \text{ if } |S1| + |S2| \le k \\ -\Phi(S_1, S_2) \text{ otherwise} \end{cases}$$

- **Genetic Algorithm Operations**: In our approach, diversity is cherished; hence, the GA operations below is full of randomness, which makes this study quite experimental.
 - i. **Initialization**: The initial population is generated in three distinct groups to balance diversity and quality: The random population is created with random binary vectors where nodes are randomly added to intervention sets *S1* or *S2*. The controlled random population has nodes selected within a controlled range, increasing the probability of feasible solutions within budget. The ideal population is generated with exactly the budgeted number of nodes, creating solutions that strictly satisfy constraints.
 - Selection ii. Survivor and **Offspring Generation**: Survivor selection is relatively simple. It involves choosing the topperforming individuals and a few mid-range and lower-performing solutions to maintain genetic diversity. Choosing pairs for breeding is done through random pair selection, except some controlled pairs for breeding among the best parents. Offspring generation is achieved through two-point crossover and mutation, where selected individuals exchange segments of their binary vectors to produce new solutions. Flip-bit mutation is applied probabilistically to encourage diversity, with a defined mutation rate applied to randomly chosen bits in the binary vectors of offspring solutions. This technique introduces further variation into the population and prevents premature convergence.

ALGORITHM 4: IEMP GENETIC ALGORITHM

Input: graph, initial1, initial2, budget

Output: balanced1, balanced2

HALT ← 0.98

num_nodes ← graph.GetNumNodes()

gen0 ← InitGen0(graph, budget)

Evaluate fitness of gen0

Sort gen0 based on fitness value

current_best_solution, current_best_solution_val ←

RetrieveBestSolution(gen0)

current_gen ← gen0

for generation ← 0 to gen-2 do

if current_best_solution_val / num_nodes ≥

HALT then

| break

 $12 \qquad next_gen \leftarrow GenerateOffSpring(parent_gen)$

 $parent gen \leftarrow ChooseSurvivor(current gen)$

10

11

13 Evaluate fitness of next_gen 14 Sort next gen based on fitness value 15 candidate_best_solution, $candidate_best_solution_val \leftarrow$ RetrieveBestSolution(next_gen) 16 **if** $candidate_best_solution_val \ge$ current best solution val then 17 $best_solution \leftarrow candidate_best_solution$ 18 $best_solution_val \leftarrow$ candidate_best_solution_val balanced1, $balanced2 \leftarrow$ ConvertBinaryRepresentationToSetRepresentation(best solution)

Note that we may terminate the algorithm as soon as the halt condition is satisfied, which is defined as checking if $\Phi(S_I, S_I)$ is 98% as big as the total number of nodes. This allows us to improve the computational time by accepting results that are sufficiently close to optimal.

return balanced1, balanced2

IV. EXPERIMENTS

A. Setup

20

In our experiment, we conducted Monte Carlo simulations three times within our heuristic algorithm for each test case to account for stochastic variability and improve result reliability. For the evolutionary algorithm, we configured a population size of 90 and limited the evolutionary generations to 50, balancing computation time with solution quality. Additionally, we ran the evolutionary algorithm both with and without the halt condition to explore how closely it could approach optimal results with and without this constraint.

We tested our approach on three datasets of varying scale and structure, as shown in Table I. Dataset 1 provided a baseline with a smaller network size, while Dataset 2 introduced a moderate level of complexity. Dataset 3, the largest, presented additional challenges due to its scale and influence probabilities skewed towards lower values, which could limit spread and increase unpredictability. For each dataset, we ran the tests three times and computed the average of the final evaluation values based on the outputted balanced seed sets for a better accuracy in performance comparison.

TABLE I. DATASETS' DETAILS

Dataset	Number of Nodes	Number of Edges
1	475	13289
2	13984	17319
3	36742	49248

The testing environment for this experiment was configured on macOS Sequoia version 15.0.1, running on an Apple M3 Pro processor with 18GB of RAM and a 1TB storage capacity. The programming language used for the algorithm implementation was Python 3.12, with essential libraries such as Numpy 2.1 for handling numerical computations.

B. Results

To obtain evaluation values, we ran independent cascade to simulate the influence spread on each outputted balanced seed set for 300 times. We also recorded the running time in the process.

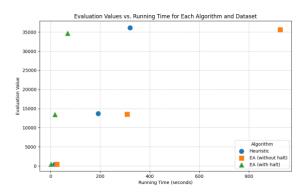


Fig. 1. Evaluation values and running time for each algorithm across different datasets

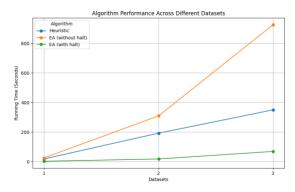


Fig. 2. Running time across different datasets

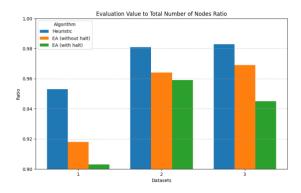


Fig. 3. Evaluation value to total number of nodes ratio across different datasets

In the halt condition of the evolutionary algorithm, we only ran independent cascade 10 times for each instance, which resulted in a high but less accurate evaluation value. This allowed the halt condition to satisfy and terminate the algorithm early, even though the actual ratio upon running 300 simulations is not that large as shown in Figure 3.

C. Analysis

In analyzing the results presented in the grouped bar chart (Figure 3), we observe that the Heuristic method consistently

outperforms both Evolutionary Algorithms (EA) across all datasets in terms of the evaluation value to total number of nodes ratio. Specifically, the Heuristic achieves ratios of 0.953, 0.981, and 0.983 for Datasets 1, 2, and 3, respectively. In comparison, the EA without halt shows a slight drop in performance, with ratios of 0.918, 0.964, and 0.969, while the EA with halt presents the lowest ratios, indicating its relatively diminished efficiency with values of 0.903, 0.959, and 0.945. These results suggest that while all methods demonstrate high effectiveness, the Heuristic method is notably superior, especially in scenarios demanding optimal results.

The running time performance depicted in Figure 2 reveals a stark contrast between the Heuristic method and the EAs. While the Heuristic maintains a significantly lower running time across all datasets—recording 18.11 seconds for Dataset 1, 192.35 seconds for Dataset 2, and 350.03 seconds for Dataset 3—the EAs exhibit a marked increase in time consumption, particularly the EA without halt, which takes 25.27 seconds for Dataset 1 and skyrockets to 925.81 seconds for Dataset 3. In contrast, the EA with halt presents the most favorable time performance for larger datasets, showing a drastic reduction in running time, down to 68.82 seconds for Dataset 3. This indicates that while the Heuristic method excels in efficiency, the halting mechanism of the EA may optimize performance and be able to achieve a close-to-optimal results in significantly less time.

Finally, the scatter plot presented in Figure 1 provides a comprehensive visualization of the relationship between evaluation values and running times across algorithms. It illustrates that the Heuristic consistently achieves high evaluation values with relatively low running times. Notably, the data points for the Heuristic cluster towards the higher evaluation values, contrasting sharply with the EAs, which, despite having longer runtimes, exhibit a less consistent

performance in evaluation values. The EA with halt, although showing improvements in runtime, still fails to reach the evaluation levels of the Heuristic. This analysis emphasizes that while running time is an essential factor in algorithm selection, the balance between efficiency and effectiveness is crucial for achieving optimal results in practical applications, which makes Heuristic our overall best approach.

V. CONCLUSION

In this study, the heuristic and evolutionary algorithms were applied to the IEMP to maximize information exposure within a social network. The heuristic method consistently showed superior performance in terms of efficiency and effectiveness across all datasets. It achieved higher evaluation values and required significantly less computational time compared to the evolutionary algorithms. However, the evolutionary algorithm with a halt condition provided a reasonable compromise by reducing runtime without severely compromising results, making it potentially suitable for larger networks where time constraints are critical.

While the heuristic approach proves robust, further exploration into other evolutionary algorithms, such as particle swarm optimization or differential evolution, may provide improved solutions. Additionally, hybrid models that combine heuristic elements with genetic algorithms could offer a balance between accuracy and computation time. This study demonstrates the potential of heuristic algorithms in tackling complex optimization problems within influence maximization, with the evolutionary approach presenting an adaptable framework for broader applications in network analysis.

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