

Political Districting via Discrete Particle Swarm Optimization

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Abstract—The abstract goes here.

I. INTRODUCTION

A. Problem Importance

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III. PROBLEM FORMULATION AND MODELING

The Political Districting problem can be modeled as an assignment problem with specific constraints. The goal is to assign a fixed set of tracts to a fixed number of districts, where each district is contiguous, while minimizing a certain cost function. The inputs provided to our algorithms are:

- T , The set of tracts
- D , the set of Districts
- $P_T(x)$ which gives the population of a given tract
- $A_T(x)$ which gives the area of a given tract
- $Borders(x)$ which, for a given tract, returns the set of tracks adjacent to it paired with the lengths of those borders

We define a solution as a mapping $s : T \rightarrow D$ where every unit in T is mapped to a single district in D , and every district in D has at least one unit in T assigned to it. The encoding method does not make any effort to guarantee district contiguity, therefore it is left to the algorithm to only generate valid solutions.

IV. PROPOSED SOLUTION

Particle Swarm Optimization is an optimization technique which models particles moving through a continuous n-dimensional search space, where each particle represents a potential solution to the problem. In the classical PSO algorithm, a set of particles are initialized to a certain solution, then at each iteration, each particle computes its next position via the following equations of motion:

$$v_{t+1}^{id} = w \cdot v_t^{id} + c_1 r_1^{id} (lbest_t^{id} - x_t^{id}) + c_2 r_2^{id} (gbest_t^{id} - x_t^{id}) \quad (1)$$

$$x_{t+1}^{id} = x_t^{id} + v_{t+1}^{id} \quad (2)$$

Where v is the velocity of particle id , w is the inertia weight, c_1, c_2 are acceleration coefficients, r_1, r_2 are randomly generated numbers in $[0, 1]$, x is the position of the particle,

$lbest$ is the best solution the particle has traversed, $gbest$ is the best solution traversed across all particles, t is the iteration number, and i and d are the particle number and dimension. While this algorithm was meant for problems on a continuous space, it can also be used for discrete optimization problems. Our solution, inspired by the work of Kang-Ping Wang et al. [1] on the Traveling Salesman Problem, proposes a novel method of applying PSO to the discrete problem of Political Districting.

A. Initial Solution

In order to generate a valid initial solution, we proceed as in Vickrey [6] and Bozkaya [3]. Initially, all tracts are not yet assigned to any district. The goal of this algorithm is to generate an initial solution such that all tracts are assigned to exactly $|D|$ contiguous districts.

For the purpose of this algorithm, we define a target population for the districts in the initial solution:

$$\bar{P} = \frac{\sum_{x \in T} P_T(x)}{|D|}$$

We create our districts by picking an unassigned seed tract, and then gradually add randomly picked unassigned adjacent tracts to the district until the population of the district exceeds \bar{P} for the first time, or there are no unassigned adjacent tracts available. We do this until all the tracts are assigned to a district.

The result of this initial assignment could result in k districts that may or may not be equal to $|D|$.

If $k > |D|$, we iteratively pick the district with the smallest population, and merge it with its smallest neighbour to preserve contiguity. In the case where $k < |D|$, we iteratively increase the number of districts by splitting the district with the largest population into two districts while still preserving contiguity. At the end of this process, there will be exactly $|D|$ valid districts that can be used as an initial solution.

B. Cost Function

While there are many factors that can be considered for political districting, for the sake of efficiency and simplicity, our fitness function only focuses on the two main goals of political districting which are population equality and district compactness. These measures do not assess the contiguity of a solution, as the algorithms used are designed to only evaluate valid solutions. A constant coefficient is introduced to each term to allow us to tune how much each factor contributes to the overall fitness.

$$f(x) = c_{pop} \cdot f_{pop}(x) + c_{shape} \cdot f_{shape}(x)$$

1) *Population Equality*: In an ideal solution, each district would have the same population, which would equal the average population across the districts. We therefore define the measure of population equality for a set of districts D where $P_D(d)$ is the total population of district $d \in D$ as

$$f_{pop}(x) = \sum_{d \in D} P_D(d) - \bar{P}$$

2) *Compactness*: To evaluate compactness we use the Schwartzberg Index [?], which is the perimeter of a district squared divided by its area, due to it being simple to compute. The perimeter of a district is computed by summing up the outward border lengths of the tracts on its border, and the area of a district is computed as the sum of the tract areas.

$$f_{shape} = \sum_{d \in D} \frac{(P_D(d))^2}{A_D(d)}$$

C. Neighborhood

The neighborhood of a given solution is defined as all possible solutions reachable by moving one tract from district i to district j such that the resulting solution remains contiguous.

D. Algorithm Description

In order to apply Particle Swarm Optimization on our problem we must define sufficient terminology and mathematical operators to be able to perform the update function 1.

1) *Velocity*: For the discrete solution space of district assignment, we define "velocity" to be an ordered sequence of Swap Operators. A given Swap Operator $SO(t, d)$ is defined as taking tract t belonging to district d_o and being reassigned to district d , where $d_o \neq d$.

2) *Initial Velocity and Inertia*: Initial particle velocities are generated by performing random swaps that still generate valid solutions. The v term in 1 is also generated as a random sequence of swaps, where the length of this sequence is a constant parameter. As a result, the concept of "inertia" is not present in our solution.

3) *Addition of Velocity to a Solution*: Given a solution X and a velocity $V = (SO_1, SO_2, \dots, SO_m)$, $X + V$ is the solution resulting from the application of swaps $SO_1 \dots SO_m$ on X in ascending order. In order to assure that the algorithm only produces valid solutions, the resulting state must be contiguous.

4) *Subtraction of Solutions from Solutions*: Given two valid solutions A and B , the subtraction operator $A - B = V$ is a velocity described as a sequence of swaps required to transform solution B into A . The resulting velocity V is guaranteed to be valid when applied to solution A , as B is assumed to be valid. In our implementation, this sequence is determined by iterating through districts in B and applying $SO(t, d)$ if tract t is assigned to district d in solution A but not in solution B , and the result remains contiguous.

5) *Multiplication of Real Numbers and Velocities*: Given $V = (SO_1, \dots, SO_m)$ we define $c \cdot V, c \in [0, 1]$ as the sequence constructed from the first $\lceil c \cdot m \rceil$ swap operators in V . Any values of c outside of $[0, 1]$ are considered invalid.

6) *Modified Update*: To perform equation 2, changes must be made to 1 to guarantee that $x_t^{id} + v_{t+1}^{id}$ results in a contiguous solution. Naively performing 1 would result in the $w \cdot v_t^{id}$, $c_1 r_1^{id}(lbest_t^{id} - x_t^{id})$, and $c_2 r_2^{id}(gbest_t^{id} - x_t^{id})$ being valid on x_t^{id} , however in order to add these resulting velocities to calculate v_{t+1}^{id} , we cannot simply append the swap sequences. Even if swap sequences v_1 and v_2 are both valid on solution x , v_2 is not guaranteed to be valid on $x + v_1$. We therefore use a modified position update method

$$\begin{aligned} X_1 &= x_t^{id} + w \cdot v_t^{id} \\ X_2 &= X_1 + r_1 \cdot c_1 \cdot (lbest_t^{id} - X_1) \\ x_{t+1}^{id} &= X_2 + r_2 \cdot c_2 \cdot (gbest_t^{id} - X_2) \end{aligned} \quad (3)$$

Since the $r_1 \cdot c_1 \cdot (lbest_t^{id} - X_1)$ term of X_2 is generated from X_1 , adding the first i elements of that result to X_1 is guaranteed to be a contiguous solution. The same reasoning can be applied to conclude that x_{t+1}^{id} is a valid solution

This method of addition differs from the classical vector addition, however it achieves the same goal of having a particle move in the direction of a velocity, in the direction of the local best, and in the direction of the global best, each to some variable degree. One important difference to note, is that if each coefficient was 1, even though each velocity is applied "equally", X_2 would be equivalent to $lbest$, and x_{t+1} would be equivalent to $gbest$. This detail can be handled by tuning the constant terms.

7) *Implementation Details*: The computation of $A - B$ where A and B are solutions is expensive, as the process of getting from A to B may require a very large number of swaps. We therefore introduce a constant parameter which describes the maximum number of swaps to generate when calculating $A - B$. Another issue with a naive implementation is that when generating a random velocity, certain swaps may undo previous swaps. To avoid this, a tabu list is maintained, which disallows assigning a cell to a district it had previously been assigned to during that swap sequence.

8) *Possible Improvements*: There are certain ideas which we did not have time to implement, however may theoretically result in improvements in performance and solution quality.

While our current implementation has no concept of "inertia", we could define "inertia" to be related to which districts grew or shrunk during the previous iteration. To create the swap sequence $v_t = SO_{t,1} \dots SO_{t,m}$ that is valid on x_t , and inherits the inertia of $v_{t-1} = SO_{t-1,1} \dots SO_{t-1,m}$, we would ensure that for $SO_{t-1,i} = (u_{t-1}, d_{t-1})$ and $SO_{t,i} = (u_t, d_t)$, $d_{t-1} = d_t$ and $s(u_{t-1}) = s(u_t)$. This means that every swap in each swap sequence will take a tract from district i and move it to district j , the only difference between the sequences is which tract was swapped. Another possible improvement would be to establish certain "centers" for each district, similar to what was done in the work by Federica Ricca et al. [2]. Each district would have one "center" tract which is always assigned to that district. This means that between two solutions, each district d will be in the same general area. This trait would theoretically help the efficiency of the algorithm immensely, as in the current implementation, two equivalent districting solutions could be seen as radically different simply because the specific identifiers assigned to the districts are different.

E. Parameter Values

We tested arbitrary combinations of values for w , c_1 , and c_2 where $w, c_1, c_2 \leq 1$ because of the constraints set by how we defined a particle's velocity. A parameter value greater than 1 would not have made any sense given our velocity model. The swarm size was limited by our computing power to 5 particles.

F. Example Run-through

We will demonstrate two iterations of the three algorithms we have chosen to implement using a simple example input. The input's tracts are conceptually a grid of tracts, where each tract shares a border with its vertical and horizontal neighboring tracts.

The population of each tract is 1 for simplicity, and the solution will be a grouping of tracts such that there are exactly two groups that represent two districts.

1) Simulated Annealing:

Figure 1 shows the first two iterations of the SA implementation. We start with a randomly generated initial solution of two districts with a fitness cost of 19.3 (calculated using the previously defined cost function).

The neighborhood of the initial solution is defined by the set of possible tract swaps (switching a tract from one district to another) without breaking continuity.

The algorithm will randomly pick one of these neighbors, and will choose it as the new solution if its fitness is better than the current one, or if the acceptance probability given the current temperature allows for worse solutions.

The algorithm does not stop in the first two iterations since the stopping conditions have not been met. Since the temperature is quite high in the beginning, the first two iterations also end up accepting solutions that have a worse fitness than the previous solution.

-----	Cost: 19.3	-----	
1	1	1	1
1	1	1	1
1	1	1	1
1	1	1	1
-----	Cost: 21.1	-----	
1	1	1	1
1	1	1	1
1	1	1	1
1	1	1	1
-----	Cost: 26.3	-----	
1	1	1	1
1	1	1	1
1	1	1	1
1	1	1	1

Fig. 1. Example run using Simulated Annealing

2) Tabu Search:

Figure 2 shows the first three iterations of a sample run using our tabu search implementation. Similarly to our SA algorithm, we start with a randomly generated initial solution of two districts. In this particular run, the initial solution has a fitness cost of 22.5.

Throughout the algorithm, we maintain a tabu table, which has the districts as its "x-axis" and the tracts as its "y-axis". Adding a value to cell, say $[i][j]$ of the tabu table means that we have chosen to swap tract j from its old district to district i . In every iteration, we first subtract all the non-zero values by 1 before updating with the new swap.

In every iteration, we first find the neighboring solutions to the current solution. These neighboring solutions are found in the same way as done for our SA algorithm. We then iterate through these neighboring solutions to find the one with the best (lowest) fitness cost. We choose these solutions with the constraints that they are not tabu (the solution's value in the tabu table is equal to 0), and that the resulting fitness cost is better than the current solution's fitness cost. The algorithm will stop when the current solution has no more neighbors, or no more viable solutions within its neighbors.

In the specific run shown in Figure 2, the chosen tract to be swapped is the 2nd one of the 1st row. Our tabu tenure value has been set to 5, resulting in the value 5 in our tabu table that can be seen in the next iteration.

The next iteration now has a lower cost of 14.7. The chosen tract to be swapped is the 2nd one in the 3rd row. Again this results in the value 5 in the following iteration's tabu table print-out, but before updating the new solution's tabu value to 5, we decrement the existing non-zero values

----- Cost: 22.5 -----															
1	1	1	1												
1	1	1	1												
1	1	1	1												
1	1	1	1												

0	0	0	0		0	0	0	0							
0	0	0	0		0	0	0	0							
0	0	0	0		0	0	0	0							
0	0	0	0		0	0	0	0							
----- Cost: 14.7 -----															
1	1	1	1												
1	1	1	1												
1	1	1	1												
1	1	1	1												

0	5	0	0		0	0	0	0							
0	0	0	0		0	0	0	0							
0	0	0	0		0	0	0	0							
0	0	0	0		0	0	0	0							
----- Cost: 10.3 -----															
1	1	1	1												
1	1	1	1												
1	1	1	1												
1	1	1	1												

0	4	0	0		0	0	0	0							
0	0	0	0		0	0	0	0							
0	5	0	0		0	0	0	0							
0	0	0	0		0	0	0	0							

Fig. 2. Example run using Tabu Search

by 1, resulting in the 4 that is seen in the next iteration's tabu table.

3) PSO:

Figure 3 is an example of a particle's current position X , it's personal best, and the global best of the swarm. We assume that the values of w , c_1 , c_2 , r_1 , and r_2 are all 1 for simplicity in this example. Swaps for the random velocity are capped at 10 swaps, and 4 for the other two velocities for the reasons mentioned in implementation details. This example will only demonstrate one iteration of a single particle for brevity. We show the swaps represented by a velocity as a table of R 's and B 's as shown in Figure 4, where the letter represents the district color that the tract in the letter's position should be swapped to.

In order to move the particle to a new position during a

X															
1	1	1	1												
1	1	1	1												
1	1	1	1												
1	1	1	1												
Personal Best															
1	1	1	1												
1	1	1	1												
1	1	1	1												
1	1	1	1												
Global Best															
1	1	1	1												
1	1	1	1												
1	1	1	1												
1	1	1	1												

Fig. 3. Example PSO particle, personal best, and global best

Velocity															
B	B	B	B												
-	R	R	-												
R	R	R	-												
R	-	-	-												
X + Velocity															
1	1	1	1												
1	1	1	1												
1	1	1	1												
1	1	1	1												

Fig. 4. Example random velocity

delta PBest															
R	-	-	-												
-	B	B	-												
-	-	B	-												
-	-	-	-												

X + Velocity + delta PBest															
1	1	1	1												
1	1	1	1												
1	1	1	1												
1	1	1	1												

Fig. 5. Swaps to get from (X + Velocity) to Personal Best

single iteration, we first generate a random valid velocity for the particle's current position (limited to a maximum of 10

delta GBest				
-	R	-	-	-
-	-	-	-	-
B	B	-	-	-
B	-	-	-	-

X + Velocity + delta PBest + delta GBest				
1	1	1	1	1
1	1	1	1	1
1	1	1	1	1
1	1	1	1	1

Fig. 6. Swaps to get from (X + Velocity + Pbest) to Global Best

swaps) and add this velocity to X . This new position is shown in Figure 4.

We then calculate the swaps required from the previous position to the particle's personal best (limited to a maximum of 4 swaps) and add this to the previous position. The resulting position is shown in Figure 5.

Finally, we calculate the swaps required from the previous position to the swarm's global best (also limited to a maximum of 4 swaps) and add this to the previous position. This will result in the new position of the particle (X') for this iteration as shown in Figure 6.

V. DATA

Initially, the proposed solutions are run on experimental data in a grid shape, known as the Grid Problem. We wanted to apply the solutions on real world data to observe whether the results are meaningful. Inspired by literature on Tabu Search implementation, we decided to use data from the state of Iowa to perform our solution and called it the Iowa Problem. Another motivation to use Iowa was because of its relatively normal county subdivision without too many abnormally-shaped counties.

The relevant data for the state of Iowa was fetched in several stages due to the difficulty of finding consolidated data in one place. From the US census bureau, we fetched the data for the county shapefiles, which includes information on the county land/water area, name, shape points and relevant identifiers such as GeoID. However, in order for the algorithm to work we also needed information about individual county populations. Fortunately, there is up-to-date county population data available through Iowa Demographics [8], which is scraped into a text file. Lastly, the algorithm also requires data on county adjacencies and the border length between them. The US census surprisingly provides data on county adjacencies on a nationwide level [7], and identifies neighbouring counties by GeoIDs. The border length can be calculated by combining this information with the relevant shape points of the neighbouring counties.

Since each county contains a vast collection of shape points to describe its shape, it becomes a computationally difficult task to find intersecting points and lines between all counties

within the state. Therefore, we applied approximation on the shape points by taking the absolute top right and bottom left points, and approximating the shapes into rectangles. Even after the approximation process we still have an accurate topological mapping of Iowa. By doing so, we can easily compute the border lengths and create visualizations to describe the congressional district division of a solution.

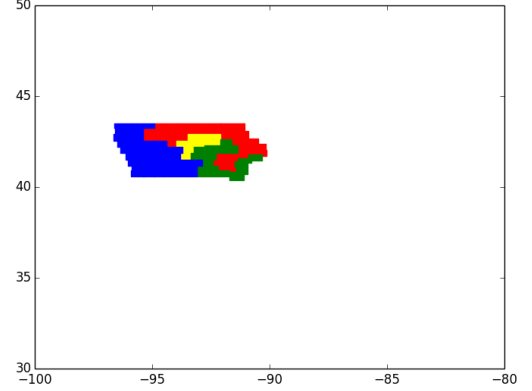


Fig. 7. Visualization of the final congressional district division in the state of Iowa using particle swarm solution from one iteration

VI. PERFORMANCE EVALUATION

Each algorithm was implemented in Python and tests were run on real data for the state of Iowa from data provided by the U.S. government census data in 2016. For each algorithm we ran a series of tests with 2, 10 and 100 iterations with fixed parameters. The initial parameters for each algorithm were chosen at random within reason. For example, the cooling rate selected for simulated annealing was 0.003. Selecting a faster cooling rate would cause the algorithm to converge to a solution more quickly, which is undesirable. For each algorithm, the results for each run was recorded. With this data, the minimum, maximum, median and standard deviation were computed for each of the evaluation criteria: compactness, population equality and overall fitness. These criteria were calculated as discussed in Section IV (Proposed Solution). We assumed that all our solutions met the contiguity criteria since this was a hard constraint implemented in each algorithm (i.e. moves or operations were not selected if they resulted in discontinuities). The statistics for each algorithm were then exported to a csv file and graphed, which can be viewed in the tables and figures below.

A. Compactness Results

Since our results show the values of cost functions, the lower the value, the better the solution is. Based on the numbers in Table I as well as Figure 7, it is evident that simulated annealing had the most unfavourable compactness results by a significant amount. The other 3 algorithms had similar performance in regards to compactness, but based on our results, PSO generally performed the best, followed by hill climbing and then tabu search.

TABLE I
COMPACTNESS RESULTS FOR 100 ITERATIONS ON IOWA

Algorithm	Compactness			
	min	max	median	stdev
SA	19.255	587.016	495.146	244.594
PSO	57.213	229.580	109.956	33.041
Hill Climbing	52.983	264.8	122.061	41.088
Tabu Search	66.946	257.239	127.855	39.242

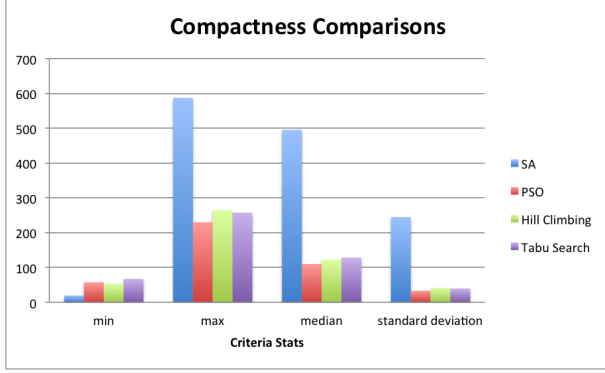


Fig. 8. Compactness Results for 100 Iterations on Iowa

B. Population Equality Results

TABLE II
POPULATION EQUALITY RESULTS FOR 100 ITERATIONS ON IOWA

Algorithm	Population Equality			
	min	max	median	stdev
SA	1841655	4648349.5	3098654	787073.096
PSO	2463	94699	32829.25	15648.72
Hill Climbing	1749	218283	9859.75	28231.930
Tabu Search	36195.5	9319.25	5647.456	5315.779

Since the simulated annealing solution had significantly poorer values for population equality as seen in Table II, they have been excluded in the corresponding graph representation below to better visualize the other algorithm results.

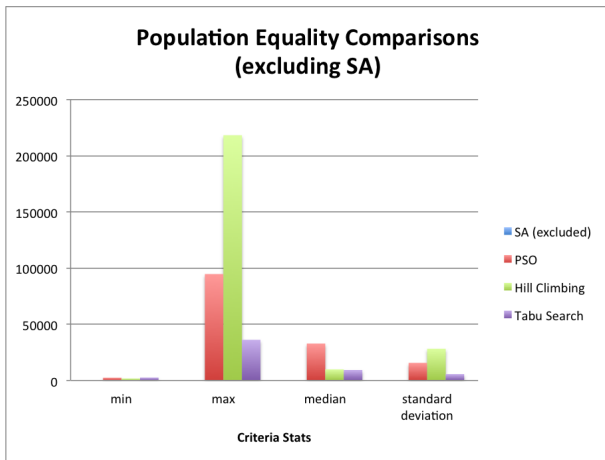


Fig. 9. Population Equality Results for 100 Iterations on Iowa (excluding SA)

Looking at Figure 8, we could deduce that hill climbing, although did relatively well in average cases based on the

median value, did significantly poorer in certain edge cases as seen from the maximum value. This is most likely as a result of being stuck at a local minima. Looking at the median values, PSO did the most poorly on average, while hill climbing did quite a bit better, and tabu search performed the best. Overall, tabu search had the best (lowest) average population equality cost value, the least deviation, and the smallest maximum cost value. Thus, we can conclude from our results that tabu search performed the best in terms of population equality.

C. Overall Fitness Results

TABLE III
OVERALL FITNESS RESULTS FOR 100 ITERATIONS ON IOWA

Algorithm	Fitness			
	min	max	median	stdev
SA	3683810.325	9296718.255	6197845.196	1573915.146
PSO	5082.820	189513.077	65748.985	31296.346
Hill Climbing	3644.275	436675.159	19839.867	56463.519
Tabu Search	5315.779	72478.096	18799.833	11289.642

Looking at the values in Table III, it is evident that simulated annealing performed significantly worse in comparison to the other 3 algorithms, so it has again been excluded in this corresponding graph representation below for better visualization of the other algorithm results.

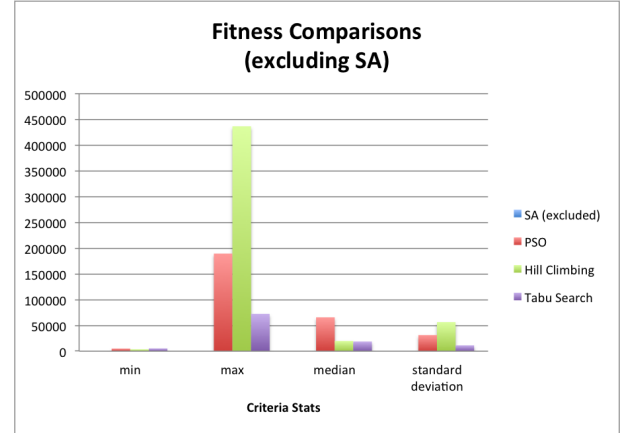


Fig. 10. Overall Fitness Results for 100 Iterations on Iowa (excluding SA)

Figure 9 shows the results for the overall fitness costs for the algorithms we tested, excluding simulated annealing. The conclusions that can be made are similar to that of the results for population equality. Hill Climbing appeared to work well on average cases, but did significantly poorly in certain edge cases. PSO performed the worst on average, looking at the median values. Tabu Search again performed the best on average cases (it has the lowest median value), had the lowest maximum cost, and the lowest deviation. Based on these results, it can be concluded that for the Iowa state dataset of 2016, tabu search would give the best performance for the Political Districting problem.

VII. CONCLUSIONS & RECOMMENDATIONS

Our results show that tabu search provides solutions that have the best overall fitness (the least cost) in comparison to

the other algorithms we tested. This means that when the tabu search algorithm is run on the data set, it produces solutions that are generally more compact and equal in population than what hill climbing, simulated annealing and particle swarm optimization would produce. Looking at the compactness results from each algorithm, we observe that on average, particle swarm optimization provides better solutions compared to hill climbing, tabu search. Meanwhile, simulated annealing produces large fitness values which are not ideal compared to the other algorithms. Comparing the population equality results we see that simulated annealing again does significantly worse compared to the other algorithms and tabu search did relatively well.

As a result of our experiments and observations of results, it is recommended that further research and parameter tuning effort should be put towards the tabu search algorithm. Tabu search performed the best on average and relatively well when considering each evaluation criteria for political districting. A second recommendation would also be to further research particle swarm optimization methods applied to the political districting problem. Particle swarm optimization performed the second best after observing the results of the overall fitness function.

Given more time, we would have further tuned the parameters for each algorithm. Doing so would allow us to compare the optimal solutions that each algorithm produces. This would give us more accurate insights into which algorithm produces the best results based on the criteria. Running a large number of iterations for certain algorithms was very costly in time and limited us to run at most 100 iterations for each algorithm. In the previous papers we found while researching, we found that the number of iterations done in the different experiments varied from 30 iterations [5] to 20,000 iterations [4]. To justify a reasonable accuracy with 100 iterations, we compared the results between running 10 and 100 iterations, and confirmed that the results were relatively similar.

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