

# THE TRAVELLING SALESMAN PROBLEM

By Group M

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Repo URL: [mawande-johnpaul/TSP: Comparison of SOM and Exhaustive Algorithms](https://github.com/mawande-johnpaul/TSP: Comparison of SOM and Exhaustive Algorithms)

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## DEFINITION AND INTRODUCTION

The Travelling Salesman Problem (TSP) is a real-world combinatorial optimization problem that involves finding the shortest possible route that visits each city exactly once and returns to the starting point. The objective is to minimize the total travel distance or cost while ensuring computational efficiency.

**Input:** Data structure

- Points are stored as a dictionary where the;
  - ❖ Keys: All the individual points.
  - ❖ Values: Dictionaries containing the keys as points adjacent to the current point and their respective distances from the current point (node).

```
adjacency_list = {  
    1: {2: 12, 3: 10, 7: 12},  
    2: {1: 12, 3: 8, 4: 12},  
    3: {1: 10, 2: 8, 4: 11, 5: 3, 7: 9},  
    4: {2: 12, 3: 11, 5: 11, 6: 10},  
    5: {3: 3, 4: 11, 6: 6, 7: 7},  
    6: {4: 10, 5: 6, 7: 9},  
    7: {1: 12, 3: 9, 5: 7, 6: 9}  
}
```

## Effectiveness

- It is easy to check validity of a path by checking if the next node is in the adjacent points of that point.
- Every adjacent point has a distance from the current node tied to it so a direct reference can be made for rapid computation.

**Output:** A sequence of numbers representing the most optimal route.

## Solution;

- Two types of solutions (algorithms) are available;
  - Exact (non-heuristic) algorithms

These guarantee an optimal solution.

- ❖ Brute-force (Exhaustive Search)
- ❖ Dynamic Programming (Held-Karp)
- ❖ Branch and Bound

- Heuristic (Approximate)

These do not guarantee the best path or a valid one.

- ❖ Nearest-Neighbor
- ❖ Genetic Algorithm
- ❖ Simulated Annealing
- ❖ Christofides Algorithm (Best approximation technique)
- ❖ Self-Organizing-Maps

## Comparison of solutions

- Algorithms are classified mainly by complexity (number of steps), performance and how much of a guarantee they offer for an optimal solution.

**Table showing comparison in complexity, pros and cons of each algorithm.**

Algorithm	Description	Complexity	Suitability	Pros	Cons
Dynamic Programming	Solves optimally by breaking the problem into subproblems.	$O(n^2 \cdot 2^n)$	20-30 cities	Optimal solution, avoids recompilation.	Memory-heavy, computationally expensive for large n
Branch-and-Bound	Systematically explores the solution space with pruning.	$O(e^n)$	Problems requiring optimal solution.	Guarantees optimal solution if fully explored Prunes parts of the search space to avoid unnecessary checks.	Still computationally expensive, impractical for large datasets.
Nearest Neighbor	Greedy heuristic. Picks nearest city next.	$O(n^2)$	Large problems needing quick solutions.	Fast, simple to implement.	No optimality guarantee, can get stuck in local minima.
Self-Organizing Map (SOM)	Neural network clustering approach, useful for approximation.	Varies (depends on grid size and iterations)	Very large problems.	Scalable, good for large datasets higher dimension, captures patterns.	No guarantee of optimality, solution quality depends on map configuration.
Brute-force	Tests all possible solutions and picks the best.	$O(n!)$	< 10 points	Guarantees optimal path, Easy to implement.	factorial time complexity increases extremely rapidly with n.
Christofides Algorithm	The direct distance between two cities is always less than or equal to the sum of the distances through a third city.	$O(n^3)$	Where the triangle inequality holds.	Guarantees a 1.5 times approximation to the optimal. Easy to implement.	No guarantee of optimal solution.
Genetic Algorithm	Depends on population size and generations.	Depends on population size and generations	When an exact solution is not needed.	Provides a near optimal approximation.	No guarantee of optimality, solution Implementation is hectic.
Simulated Annealing	Probabilistic and depends on temperature decay and iterations.	Depends on temperature decay and iterations.	When an exact solution is not needed.	Less computational since they path with highest probability is considered.	No guarantee of optimality, solution Implementation is hectic.

**Example showing complexity versus number of points.**

<b>N (Number of Cities)</b>	<b>Brute Force <math>O(n!)</math></b>	<b>Dynamic Programming <math>O(n^2 \cdot 2^n)</math></b>	<b>Nearest Neighbor <math>O(n^2)</math></b>	<b>Christofides <math>O(n^3)</math></b>	<b>Branch and Bound (<math>e^n</math>)</b>
<b>1</b>	1	1	1	1	1
<b>2</b>	2	4	4	8	2
<b>3</b>	6	36	9	27	6
<b>4</b>	24	384	16	64	24
<b>5</b>	120	4,800	25	125	120
<b>6</b>	720	61,440	36	216	720
<b>7</b>	5,040	823,680	49	343	5,040
<b>8</b>	40,320	10,485,760	64	512	40,320
<b>9</b>	362,880	134,217,728	81	729	362,880
<b>10</b>	3,628,800	1,717,986,918,400	100	1,000	3,628,800

**According to the statistical facts Exhaustive Search (Brute-Force) will be used because;**

- It guarantees an exact, optimal path.
- Easy to implement.
- Practical for this dataset  $n = 7$  points.
- Can be optimized to use up less memory easily.

## FINDINGS

### How it works

- Rule 1: Visit all points.
  - Find all numbers with n-1 digits without 1 in them.
- Rule 2: Visit each point exactly once.
  - Eliminate all numbers with any repeating digits e.g. 223457
- Rule 3: Start and end at the same point.
  - Add 1 at the beginning and end of each number obtained to obtain all the possible paths.
- Rule 4: Shortest distance possible.
  - Use the adjacency list to calculate the distance of the current digit (current node) from the next digit in the sequence (next node) while checking that the next node is adjacent to the current node. If not, that path is discarded.

```
from itertools import permutations # Library to carry out permutations

# Data structure to represent the points.
adjacency_list = {
    1: {2: 12, 3: 10, 7: 12},
    2: {1: 12, 3: 8, 4: 12},
    3: {1: 10, 2: 8, 4: 11, 5: 3, 7: 9},
    4: {2: 12, 3: 11, 5: 11, 6: 10},
    5: {3: 3, 4: 11, 6: 6, 7: 7},
    6: {4: 10, 5: 6, 7: 9},
    7: {1: 12, 3: 9, 5: 7, 6: 9}
}

# Function to generate permutations.
def generate_valid_sequences():
    """
    Generate all valid permutations of nodes 2-7 and prepend + append node 1
    to ensure cycles start and end at 1. Return a list of all the permutations.
    """
    nodes = "234567"
    perm_list = ['1' + ''.join(p) + '1' for p in permutations(nodes)]
    return perm_list
```

```

sequences = generate_valid_sequences() # Assign the permutations list to sequences variable
results = {} # Dictionary to store all the paths and their respective distances.

# Compute distances for each valid route
for sequence in sequences: # Repeat this block for all sequences
    total_distance = 0 # Initialize the distance covered in the path
    valid = True # Track validity of the path based on adjacency

    for i in range(len(sequence) - 1): # Repeat this block for all digits in a sequence
        current_node = int(sequence[i]) # The point we are currently on
        next_node = int(sequence[i + 1]) # The next to be visited in the sequence

        # Check if the next point is adjacent to the current point.
        if next_node in adjacency_list[current_node]:
            # Update the distance with distance to next point.
            total_distance += adjacency_list[current_node][next_node]
        else: #
            valid = False # If the next point in the sequence is not adjacent to the current point.
            break # Stop checking if an invalid connection is found

    if valid: # Adds sequence to valid results if it has passed all validity tests
        results[sequence] = total_distance # The value of the sequence key is its distance travelled

# Find best routes
best_route = min(results, key=results.get) # Min function to return key of minimum value in dictionary

# Print the optimal sequence (route) and its distance
print(f"Best Route: {best_route} with distance {results[best_route]}")

```

## SELF-ORGANISING-MAPS (SOM'S) / KOHONEN MAPS

- Unsupervised neural network that clusters high dimensional data through creating a discretized version of its input space(map) in lower dimensions.
- Created by Teuvo Kohonen.
- They use competitive learning instead of error-correcting learning by using a function related to neighborhoods to preserve the topological properties of the map.
- They consist of the input layer, the weights and the output (Kohonen/feature/competitive layer)
- They are great for dimensional reduction, clustering, natural language processing, geology, astronomy and finance.
- Weights – Coordinates of each output neuron.
- Output neurons must compete to be activated, hence become the winner.

Before that algorithm must be used, the adjacency list is converted into an adjacency matrix, using `np.zeros()` then filling the distances. It is then saved in a .tsp file (standard file format for SOM input.)

```
NAME : City data
COMMENT : 7 locations from adjacency list
COMMENT : Generated using NetworkX
TYPE : TSP
DIMENSION : 7
EDGE_WEIGHT_TYPE : EUC_2D
NODE_COORD_SECTION
1 46.6385 71.1307
2 43.1025 60.1363
3 37.7358 66.6288
4 31.8756 57.3015
5 40.4456 65.3949
6 35.2848 66.1870
7 42.3950 59.4070
EOF
```



### **The learning scheme is comprised of;**

- Initialization
  - Data structures, samples and training data is loaded
- Competition
  - Winner is determined
- Cooperation
  - Topological neighbor of winner is determined
- Adaptation
  - The weights are updated to increase importance of neurons is updated
- Continuation
  - Algorithm is repeated for a given number of iterations

### **Phases of the learning process**

- Initialization
  - Generate weights from the input map from equation through sampling.

$$W_{ij} = X_{ij}$$

Where W is a weight and X is a point selected randomly.

- Sampling
  - Go through dataset and select samples randomly in each iteration from map.

- Matching

- Finding winner (with weight vector closest (shortest Euclidian distance,  $d$ ) to the input sample) and its topological neighbors.

$$d = \|X - W\| = \sqrt{\sum ((X_i - W_i)^2)}$$

- The winner's index is calculated from;

$$I(x) = \operatorname{argmin}(d), \text{ a 'NumPy' function.}$$

- Its topological neighborhood is calculated from;

$$H(i^*, i) = \exp(-\|R_{i^*} - R_i\|^2 / (2 * \sigma^2))$$

Where  $\sigma$  is the standard deviation of the weights calculated from;

$\sigma = \text{np.std}(\text{array of weights})$  where  $\text{np.std}()$  is a 'NumPy' function

- Topological neighborhood: Radius around winner in which all neurons' weights are updated.

- Updating

- Changing neuron weights of the winner and its topological neighbors based on a specific equation;

$$\Delta W_{ij} = \alpha * H(i^*, i) * (X_j - X_{ij}) \quad \text{Where;}$$

$$\text{Learning rate, } \alpha = \alpha(0) * \exp(-t / T) \quad \text{Where;}$$

$t$  is the current iteration and increases by one each iteration

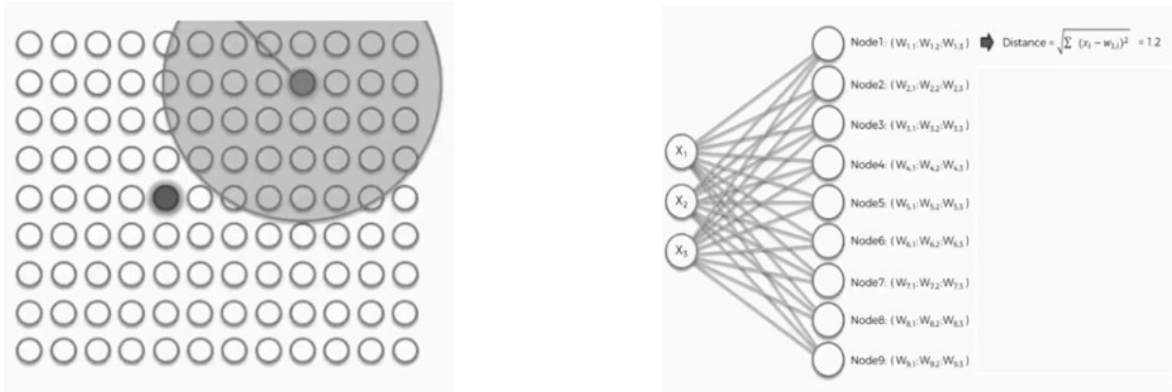
$T$  is the annealing constant that controls how fast the learning rate decays.

- $T$  can be any value of choice e.g. with  $T=100$ , decay is slower than  $T=1000$ .
- Learning rate: Refers to how fast the weights are updated.
- Topological neighborhood and learning rate decay as iterations increase.

- Continuation / Convergence

- Repeat the same algorithm until the feature map remains the same or until the number of iterations (epoch  $t$ ) is reached.

**Illustration of feature(output) map showing winning neuron and its topological neighborhood and mapping.**



- The weights are similar to the inputs of the input space.
- Step by step, the winning neuron is mapped onto the feature map until the whole map is completely covered or all neurons are mapped.

#### PSEUDO CODE

Convert the data into x,y coordinates using automatic mapping with Networkx library.

Initialize a 2D grid of neurons representing possible city positions

Set learning parameters (learning rate, neighborhood radius, number of iterations)

Randomly distribute neurons in a continuous space that covers the cities

FOR each iteration:

FOR each city in the adjacency list:

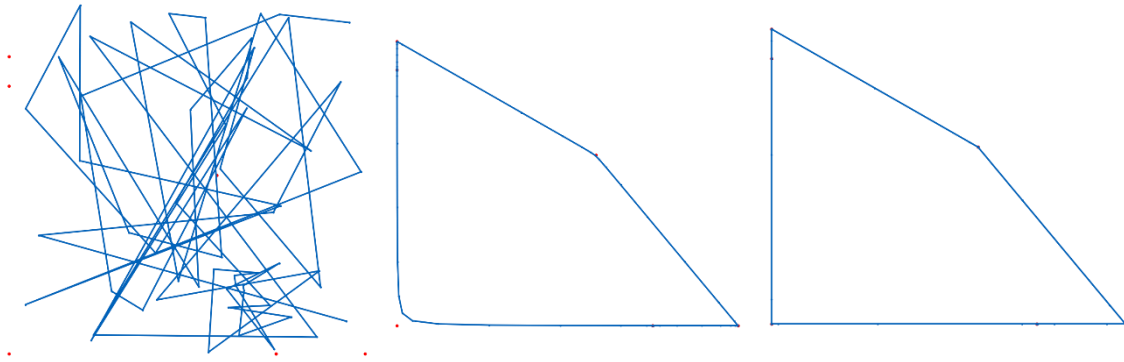
- Find the winning neuron/Best Matching Unit (BMU): The neuron closest to the current city.
- Update the BMU and its neighbors:
- Move the BMU towards the city.
- Adjust neighboring neurons based on the neighborhood radius.
- Reduce the learning rate and neighborhood radius over time.

Sort the neurons based on their positions to form a closed-loop path

Map the cities to the nearest neurons to create a tour sequence

Calculate the total distance of the generated tour using the adjacency list

## Training results



## Limitations and difficulties

- Data must be converted to standard (.tsp) format for input into the algorithm.
- Implementation is not straight forward.
- Slow and computationally expensive (not suitable for small datasets).

## COMPARISON

### Rate quality

- Exhaustive Search is shorter to implement, run and adopt as opposed to SOM that needs multiple files, equations, functions and processes. Speed can be reduced by reducing number of iterations

### With imports

```
[Running] python -u "c:\Users\LENOVO\Documents\TSP\Exhaustive Search\src\with imports.py"
Best Route: 12467531 with distance 63

[Done] exited with code=0 in 0.28 seconds
```

### Without imports

```
[Running] python -u "c:\Users\LENOVO\Documents\TSP\Exhaustive Search\src\without imports.py"
Best Route: 12467531 with distance 63
Worst Route: 12345671 with distance 69

[Done] exited with code=0 in 0.254 seconds
```

### With optimizations

```
[Running] python -u "c:\Users\LENOVO\Documents\TSP\Exhaustive Search\with imports and optimization.py"
Best Route: [1, 2, 4, 6, 7, 5, 3] with distance 63

[Done] exited with code=0 in 0.192 seconds
```

## Complexity

- Exhaustive search is more complex with steps equal to  $7! = 5040$  while SOM complexity depends on the number of iterations used.

## Time complexity

- high-level discussion of the computational cost of the SOM approach (number of iterations, updates per iteration, compare with exhaustive).

**Consider a scenario of a mailman charged with delivering mail to 9 locations in Kampala.**

- The problem size ( $n=9$ ) is **small enough** for an exact algorithm.
- Since a mailman **must** find the best route, approximate methods may lead to inefficiencies.
- **SOMs require parameter tuning**, whereas brute-force guarantees correctness.
- **Exhaustive Search is easier to verify**, while SOMs rely on probabilistic learning.
- **Brute-Force** can act as a Benchmark when trying another method.

**Extensions and optimizations for SOMs.**

- While SOMs can approximate TSP solutions, they can be improved using:
  - **Adaptive Learning Rate:** Decay learning rate dynamically to refine weight updates.
  - **Neighborhood Function Tuning:** Modify the function to prevent early convergence to suboptimal paths.
  - **Hybrid Approaches:** Combine SOMs with local search techniques (e.g., 2-opt) to refine solutions.
  - **Parallel Processing:** Use GPU acceleration to speed up large-scale computations.
  - **Edge Weight Adjustments:** Dynamically adjust weights based on distance constraints.

**Optimizations for Brute-Force Method.**

- Although brute-force is infeasible for large datasets, it can be optimized using:
  - **Branch and Bound:** Prune unnecessary paths early based on cost constraints.
  - **Memoization:** Store intermediate results to prevent redundant calculations.
  - **Parallel Processing:** Utilize multi-threading or distributed computing to explore paths simultaneously.
  - **Bit masking and Dynamic Programming:** Reduce redundant re-computations in Held-Karp algorithm implementations.
  - **Preprocessing:** Sort adjacency lists to prioritize the shortest edges first.