



# New geometric approaches to the map-matching problem

T. Akamatsu, G. Gress, K. Huneycutt, S. Omura

Academic Mentor: Dr. Kano Industry Mentor: Dr. Yamazaki

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# What is g-RIPS?

Not sure yet if I should do this yet; ask Suito-sensei

# \_

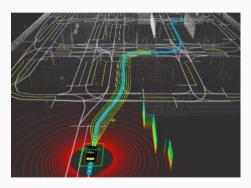
**Introduction to map-matching** 

# Map-matching

Given GPS trajectory data and a road map, **map-matching** is the process of determining the route on the map that corresponds to the trajectory data.

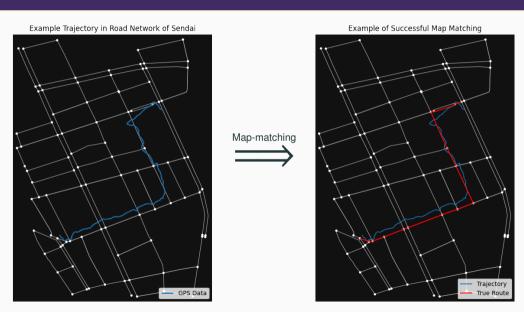


Web mapping services



Autonomous Vehicles [H]

# **Map-matching**



# How do we model this mathematically?

#### **Mathematical Formulation of the Problem**

Let us fix  $N \in \mathbb{N}$ ,  $N \ge 2$ , but almost everywhere we consider the case N = 2.

# **Definition (Trajectory)**

A **trajectory** Tr is a sequence  $\mathbf{p} = (p_1, p_2, \dots, p_n)$  of points in  $\mathbb{R}^N$  equipped with

- a sequence  $t(\mathbf{p}) = (t_1, \dots, t_n)$  of positive numbers satisfying  $t_1 < t_2 < \dots < t_n$ , called the **timestamp** of  $\mathbf{p}$ ,
- a sequence spd(p) = (spd<sub>1</sub>,..., spd<sub>n</sub>) of positive numbers called the speed of p (optional),
- a sequence  $u(\mathbf{p}) = (u_1, \dots, u_n)$  of unit vectors in  $\mathbb{R}^N$ , called the **direction** of  $\mathbf{p}$  (optional).

#### **Mathematical Formulation of the Problem**

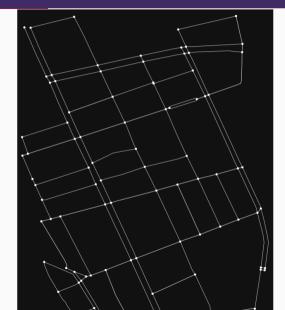
## **Definition (Road Network)**

A **road network** (also known as a map) is a directed graph G=(V,E) consists of the set V (resp. E) of vertices (resp. edges) with an embedding  $\phi:|G|\to\mathbb{R}^N$  of the geometric realization |G| of G. We will identify G and the image  $\phi(|G|)$  by  $\phi$  as long as there is no confusion.

## **Definition (Local Road Network)**

A **local road network** is a directed connected subgraph of G = (V, E).

# (Local) Road Network



#### **Problem Statement**

## **Definition (Route)**

A **route** r on a road network G = (V, E) is a sequence of connected edges  $(e_1, e_2, \dots, e_n) \subset E$ , i.e. the head of  $e_i$  coincides with the tail of  $e_{i+1}$  for each  $i = 1, 2, \dots, n-1$ . Let R denote the set of all routes.

**Definition (Candidate Routes)** For the local road network graph as H of the road network G, we define

 $C\mathcal{R}_H = C\mathcal{R} := \{ \text{routes on a local road network graph } H \},$ 

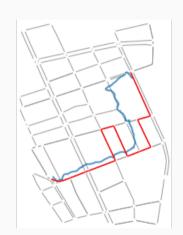
# Route



# **Candidate Routes**



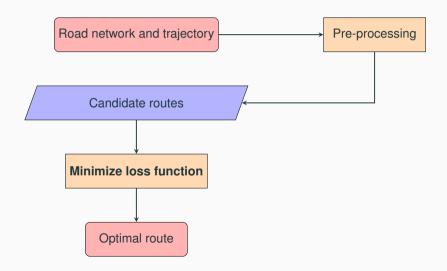




#### **Problem Statement**

**Definition (Map-Matching)** Given a road network G=(V,E) and a trajectory Tr, the map-matching,  $\mathcal{MR}_G(Tr)$ , is the route that is the argument of the minimum of some function  $L: \mathcal{CR} \to \mathbb{R}^+$ , called the loss function.

# **Map-matching Pipeline**



#### **Mathematical Formulation**

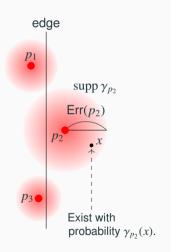
# **Assumption**

• Give the **GPS error** as  $\operatorname{Err}: \mathbf{p} \to \mathbb{R}_{\geq 0}$  and assume that the *spherically-symmetric probability measure*  $\gamma_p$  (e.g. *Gaussian measure*) is given such that

$$\operatorname{supp} \gamma_p = B(p; \mathsf{Err}(p)) := \big\{ x \in \mathbb{R}^N \ \big| \ d_{\mathbb{R}^N}(x,p) \le \mathsf{Err}(p) \big\}.$$

We assume that  $p \in \mathbf{p}$  is truly located at x with probability  $\gamma_p(x)$ .

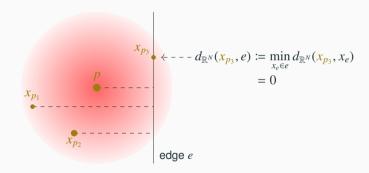
 Suppose that there is NO error with respect to the speed and direction information.



# **Definition (The "distance" with error between** $p \in \mathbf{p}$ and $e \in E$ )

• Define the "distance" with error  $d_{Err}$  between  $p \in p$  (with errors) and  $e \in E$  (without errors) as

$$\mathsf{d}_{\mathsf{Err}}(p,e) \coloneqq \int_{x_p \in B\left(p; \mathsf{Err}(p)\right)} d_{\mathbb{R}^N}(x_p,e) \, \mathrm{d}\gamma_p(x_p).$$



# Wasserstein method

#### Wasserstein method

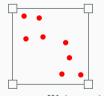
Definition (( $L^1$ -)Wasserstein distance (*review of mid-term presentation*)) Let (X,d) be a complete and separable metric space. For probability measures  $\mu,\nu$  with finite supports, we define  $W_1$  distance between  $\mu$  and  $\nu$  as

$$W_1(\mu,\nu) := \min_{\pi \in \Pi(\mu,\nu)} \sum_{x \in X} \sum_{y \in X} d(x,y)\pi(x,y),$$

where  $\pi \in \Pi(\mu, \nu)$  : $\Leftrightarrow$  for any  $x, y \in X$ ,  $\sum_{v \in X} \pi(x, y) = \mu(x)$ ,  $\sum_{x \in X} \pi(x, y) = \nu(y)$ .

Definition (Prob. meas. associated w/  $\mathbf{p}$  and  $A \in C\mathcal{R}$ )

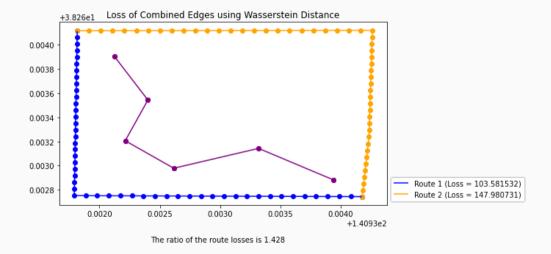
- For the trajectory **p**, define  $\mu_{\mathbf{p}} := (1/n) \sum_{p \in \mathbf{p}} \delta_p$ .
- $\triangleright$  Devide each  $A \in C\mathcal{R}$  into m+1 equal parts and V(A,m) denotes the set of m threshold points.
  - ▶ Define  $\nu_A := (1/m) \sum_{a \in V(A,m)} \delta_a$ .

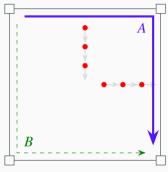




Compare  $W_1(\mu_{\mathbf{p}}, \nu_A)$  with  $W_1(\mu_{\mathbf{p}}, \nu_B)$ .

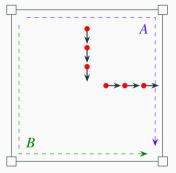
# **Wasserstein Distance: Proof of Concept**





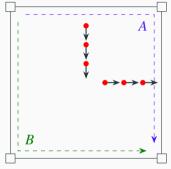
- Each  $p \in \mathbf{p}$  is located on the vertical or parallel bisectors.
- spd(p), Err(p) are the same at each p, respectively.

- · Location only.
- $\triangleright W_1(\mu_{\mathbf{p}}, \nu_A) < W_1(\mu_{\mathbf{p}}, \nu_B).$
- We should select the route B.
- Introduce probability measures  $\mu_{\mathbf{p},A}^{\varepsilon}$ ,  $\nu_{A}^{\varepsilon}$  and  $\nu_{B}^{\varepsilon}$  that include speed and direction information.
- > Compare  $W_1(\mu_{\mathbf{p},A}^{\varepsilon}, \nu_A^{\varepsilon})$  with  $W_1(\mu_{\mathbf{p},B}^{\varepsilon}, \nu_B^{\varepsilon})$
- The effect of location is still strong.
- Normalization



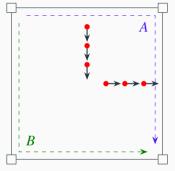
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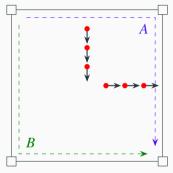
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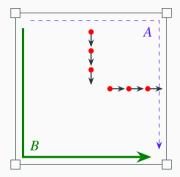
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- Normalization.
- $\quad \triangleright \ \, \text{Compare} \,\, \frac{W_1(\mu_{\mathbf{p},A}^\varepsilon, \nu_A^\varepsilon)}{W_1(\mu_{\mathbf{p}}, \nu_A)} \,\, \text{with} \,\, \frac{W_1(\mu_{\mathbf{p},B}^\varepsilon, \nu_B^\varepsilon)}{W_1(\mu_{\mathbf{p}}, \nu_B)}$



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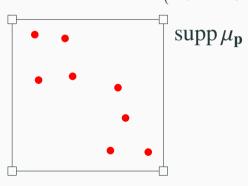
$$\Rightarrow \frac{W_1(\mu_{\mathbf{p},A}^{\varepsilon}, \nu_A^{\varepsilon})}{W_1(\mu_{\mathbf{p}}, \nu_A)} > \frac{W_1(\mu_{\mathbf{p},B}^{\varepsilon}, \nu_B^{\varepsilon})}{W_1(\mu_{\mathbf{p}}, \nu_B)}.$$

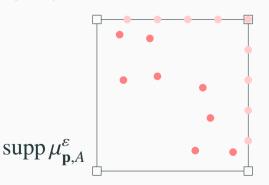
We can select the route B.

# Our strategy: Perturb $\mu_{\mathbf{p}}$ and each $\nu$ only by $\varepsilon$ according to speed and direction.

• Let  $0 < \varepsilon \ll 1$ . For each  $p \in \mathbf{p}$ ,  $A \in C\mathcal{R}$ ,  $a \in V(A, m)$  and  $x \in V(A, m)$ , define

$$\mu^{\varepsilon}_{\mathbf{p},A}(x) \coloneqq \begin{cases} (1-\varepsilon)/n & (x=p), \\ \varepsilon \cdot (\text{our weight including } S \& D) & (x \in V(A,m)), \end{cases}$$

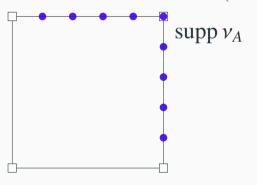


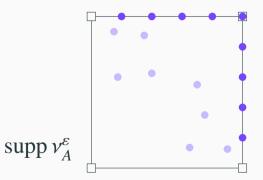


# Our strategy: Perturb $\mu_{\mathbf{p}}$ and each $\nu$ only by $\varepsilon$ according to speed and direction.

• Let  $0 < \varepsilon \ll 1$ . For each  $p \in \mathbf{p}$ ,  $A \in C\mathcal{R}$ ,  $a \in V(A, m)$  and  $x \in V(A, m)$ , define

$$\nu_A^{\varepsilon}(x) := \begin{cases} (1-\varepsilon)/m & (x \in V(A,m)), \\ \varepsilon/n & (x \in \mathbf{p}). \end{cases}$$





# **Summary & Future problem**

# **Summary**

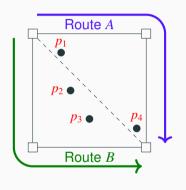
- Quantify the distance between  $p \in \mathbf{p}$  and each  $A \in C\mathcal{R}$  by making  $\mu_{\mathbf{p}}$  and  $\nu_A$   $\varepsilon$ -pertubation according to speed and direction information.
- Conclude that the route A with the smallest  $W_1(\mu_{\mathbf{p},A}^{\varepsilon}, \nu_A^{\varepsilon})/W_1(\mu_{\mathbf{p}}, \nu_A)$  is the true route.

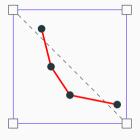
## Future problem (from a theoretical point of view)

- Is this method also effective when CR is dense?
- Formulation of  $\mu_{\mathbf{p}}$  with  $(\gamma_p)_{p \in \mathbf{p}}$ .

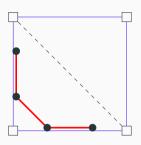
# "Physical" method

# "Electric" Method : Review of Mid-presentation





**Figure 1:** Connecting traj. pts. and giving charges.



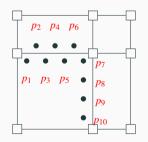
**Figure 2:** Moving to "closer" route.

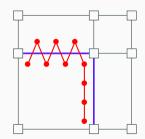
• Considering not only trajectory points, but also the entire polyline.

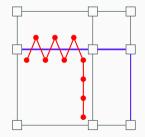
# "Electric" Method: Problems

#### **Problems:**

- × Not taking into account speed and direction information.
- $\times$  Divergence problem :  $\int_{\text{polyline}} \int_{\text{route}} r^{-2}$
- Even if  $\int_{\text{polyline}} \int_{\text{route}} (r^2 + \varepsilon)^{-1}$ , affects of intersection point is too large.







# "Electric" Method : Strategy to Solve the Problems

# Strategy:

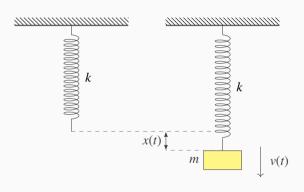
Replace "maximizing inverse square" to "minimizing square";

$$\max_{A \in \mathcal{CR}} \frac{1}{r^2} : \text{``electric''} \text{ method}$$



 $\min_{A \in CR} r^2$ : "harmonic oscillator" method

# "Harmonic Oscillator" Method : General Setting

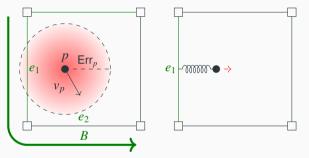


- m:mass,
- *k* :spring constant,
- v(t) :verocity of mass point,
- x(t):displacement from natural length of spring
- Lagrangian

$$L(t) = \frac{1}{2}mv(t)^2 + \frac{1}{2}kx(t)^2$$

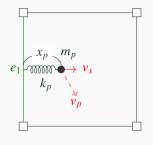
• Action  $Act = \int L(t) dt = \int \left\{ \frac{1}{2} m v(t)^2 + \frac{1}{2} k x(t)^2 \right\} dt$ 

# "Harmonic Oscillator" Method : Settings in Map-matching Problem



- $p \in \mathbb{R}^N$ : trajectory point,
- $v_p = \operatorname{spd}_p u_p \in \mathbb{R}^N$ : speed at p,
- $\operatorname{Err}(p) \in \mathbb{R}$ : error of p,
- $B \in C\mathcal{R}$ ,  $e \in B$
- S(p,e): score of edge e $:= \langle v, \vec{e} \rangle_{\mathbb{P}^N} \exp(-\mathsf{d}_{\mathsf{Frr}}(p,e))$
- Connect trajectory point to the highest score edge by "spring".
- Define "Lagrangian" of this system.

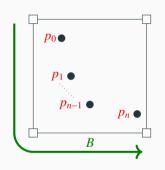
# "Harmonic Oscillator" Method : Settings in Map-matching Problem

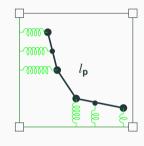


- $v_s$ : spring direction component of  $v_p$
- $x_p = d_{\text{Err}_p}(p, e)$ : "displacement" of p
- $m_p = \frac{1}{1 + \mathsf{Err}(p)}$ : "mass" of p,
- $k_p = \exp(-\text{Err}(p))$ : "spring constant" w.r.t. p,
- $M(p) = m_p \left\| \frac{v_s}{\log(1+|v_p|)} \right\|^2$ : "momentum" of p,
- $P(p) = k_p x_p^2$ : "potential" of p,

$$L := M(p) + P(p)$$
: "Lagrangian".

# "Harmonic Oscillator" Method : Settings in Map-matching Problem





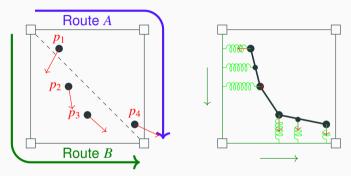
- $\mathbf{p} = (p_i)_{0 \le i \le n}$ : traj. points,
- $v_{\mathbf{p}} = (v_i = \operatorname{spd}_i u_i)_{0 \le i \le n}$ : velocity,
- Err :  $\mathbf{p} \to \mathbb{R}$  : error,
- $B \in CR$ : route.

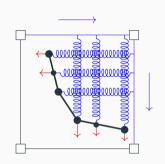


- $l_{\mathbf{p}}:[0,1]\to\mathbb{R}^N$ : polyline,
- $v_{l_{\mathbf{p}}}:[0,1] \to \mathbb{R}^N$ : velocity,
- $\mathsf{Err}_{l_{\mathbf{p}}}:[0,1] \to \mathbb{R}:\mathsf{error},$

$$L(t) := M(l_{\mathbf{p}}(t)) + P(l_{\mathbf{p}}(t)), \quad Act_{\mathbf{p}}(B) := \int_{[0,1]} L(t) dt : \text{``action'' of } \mathbf{p} \text{ on } B.$$

# "Harmonic Oscillator" Method : Settings in Map-matching Problem





- Calculate  $Act_{\mathbf{p}}(A)$  and  $Act_{\mathbf{p}}(B)$ .
- Choose the route that minimize action.

### "Harmonic Oscillator" Method: Strength/Weakness and Future Problem

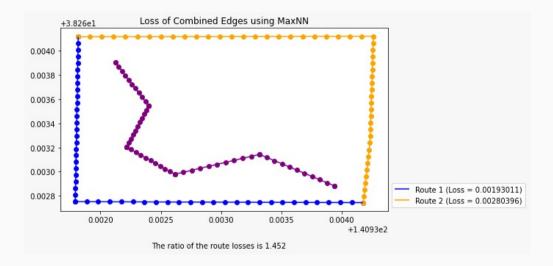
### Strength/weakness:

- √ No divergence problem,
- √ Taking into account speed and direction naturally,
- × Insufficient consideration of the entire route.

### **Future problems:**

- Consider more appropriate way to define mass and spring constant.
- Connect to all edge of a route with appropriate weight to consider the whole route.

### "Electric Method" and "Harmonic Oscillator": Proof of Concept



# Numerical Results

### Implementing metric-based methods

```
from algorithms import metric and, fam him
                                                                                                                                                                       而水业土甲ョ
Wfrom fmm import FastMapMatchConfig
### Define was matching configurations
radius = 0.803
ons error - 8 8885
#fmm config = FastMapMatchConfig(k,radius.gps error)
cfn file . None
## Least squares functions
Is ri = lambda distarray: np.square(distarray) # The function applied directly to the distances from the candidate route to the k-NW GPS coords
ls ro = lambda distarray: 1*(1/np.size(distarray) * np.sum(distarray)) # This is where we 'integrate' over the distances, and if we need to do anything else, we do it
ls qi = lambda distarray: np.square(distarray) # The function applied directly to the distances from the GPS coords to the k-NW candidate route nodes
ls go = lambda distarray: 1*(1/np.size(distarray) * np.sum(distarray))
## Inverse squares function ('Electrical method')
is ri = lambda distarray: np.power(np.square(distarray) + eps. -1) # We need eps to prevent singularities, i.e. r = 0
is ro = Lambda distarray: -1*(1/np.size(distarray) * np.sum(distarray)) # We sup. and then multiply by -1 to turn the minimizing process into a maximizing process
is ni = lambda distarray: nn.nover(nn.square(distarray) + ens. -1)
is go = lambda distarray: -1*(1/np.size(distarray) * np.sum(distarray))
def wrapper_f(ri, ro, gi, go): # This should return a function composed from the basic functions, that can then be applied onto route and gps data.
    return lambda route, gps : 1*ro(ri(route)) + 1*go(gi(gps))
ls loss function - wrapper f(ls ri, ls ro, ls gi, ls go)
is loss function = wrapper f(is_ri, is_ro, is_gi, is_go)
def wasserstein(routeloss, gpsloss):#gpsloss,n,m
    #the (i, i)th entry of the apsloss matrix is the distance from the ith point of the trajectory to the ith point on the candiste route
    # n is the number of points along the trajectory
    # m is the number of points on the candidate route
    n = qpsloss.shape[0]
    n = qpsloss.shape[1]
    #Create equality constraints
    b = [1/n for i in range(0,n)]+ [1/n for i in range(0.m)]
    rowl = [i for i in range(0.n) for i in range(0.m)]
    row2 = [n+i for i in range (0,m) for i in range(0,n)]
    row = np.appending.matrix.flatten(np.array(row1)) np.matrix.flatten(np.array(row2)))
    col1 = [list(range(0.n+m))]
    col2 - [ism*k for i in range[0.m) for k in range[0.n)]
    col = np. appending.matrix.flatten(np.array(col1)).np.matrix.flatten(np.array(col2)))
    data = np.ones(n*n*2)
    A = csr matrix((data, (row, col)), shape = (n+m, n+m)), toarray()
    A = A[:-1]
    h = hf:-11
    #solve the linear program
    res = linprog(np.matrix,flatten(gpsloss),Nome, Nome,A.b)
    loss = res.fun
    return loss
```

### Implementing metric-based methods

Why are we defining each of the metric-based functions here, instead of in a separate Class? Because metric\_mm is designed to accommodate any distance-based loss function, allowing for simple and customizable simulator creations.

```
# The metric_mm algorithm allows you to either directly pass in a loss function, or to pass it all the individual pieces and wrap it itself
# This may be useful in case you wish to utilize the individual functions of a sim elsewhere.

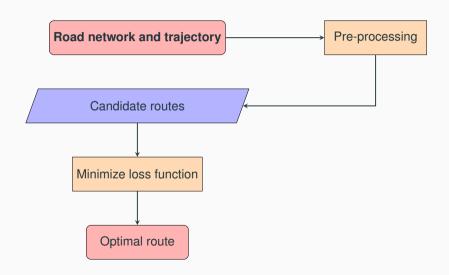
sim1 = fmm_bin.FMM(cfg = fmm_config)
sim2 = metric_mm.Sim(ls_ri, ls_ro, ls_gi, ls_go, wrapper_f) # Least squares metric-based
sim3 = metric_mm.Sim(loss_function = is_loss_function)
sim4 = metric_mm.Sim(loss_function = wasserstein)

## If you have the ground truth, load it here
ground_truth = db.read_text('map-matching-dataset/*route.geojson').map(json.loads).map(gpd.GeoDataFrame.from_features)
```

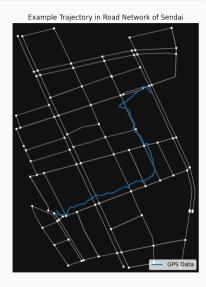
### **Sendai Map Revisited**

Now we revisit the Sendai map case.

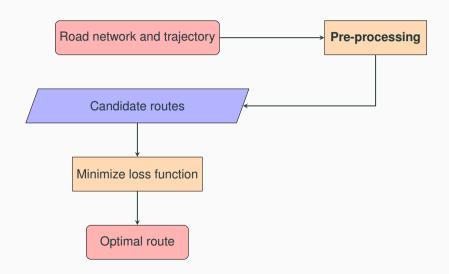
### **Map-matching Pipeline**



# **Sendai Map Revisited**



### **Map-matching Pipeline**



### Sendai Map Revisited

```
Obtaining Candidate Routes (Dijsktra's Algorithm)
```

CPU times: user 45.8 s, sys: 49.7 ms, total: 45.9 s

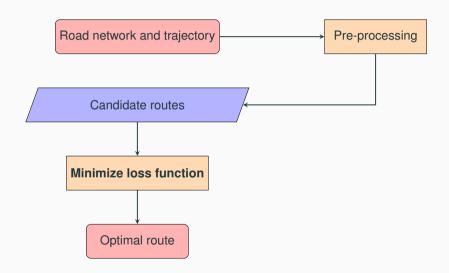
Wall time: 45.9 s

Preprocessing Candidate Routes

CPU times: user 12.7 s, sys: 103 ms, total: 12.8 s

Wall time: 12.8 s

### **Map-matching Pipeline**



### Sendai Map Revisited

```
Least Squares Runtime
```

CPU times: user 1.53 s, sys: 29.9 ms, total: 1.56 s

Wall time: 1.55 s

### Inverse Squares Runtime

CPU times: user 14.2 s, sys: 392 ms, total: 14.6 s

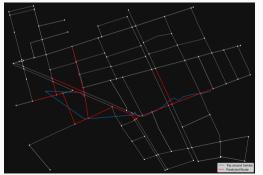
Wall time: 10.2 s

#### Wasserstein Runtime

(Computation times vary wildly based on parameters- anywhere from 20 seconds, to 20 minutes, to 2.5 hours)

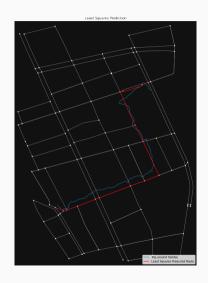
### Sendai Map Revisited

Unfortunately, due to FMM relying on outdated libraries, we could not produce an image demonstrating FMM's performance on the new trip data. Our preliminary findings demonstrated that FMM still performed poorly on the dataset. For the sake of comparison, recall FMM's results on the older GPS dataset:



We also note that FMM runs in roughly 5 seconds on our current dataset.

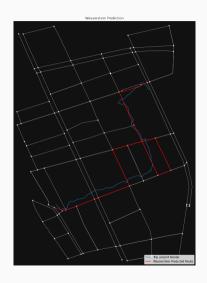
# Sendai Map: Least Squares (Harmonic Oscillator) Result



# Sendai Map: Inverse Squares (Electrical Method) Result

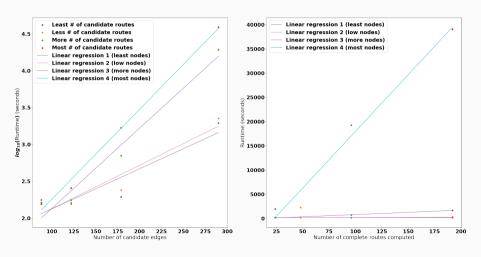


### **Sendai Map: Wasserstein Method Result**



### **Computational Complexity (Preprocessing)**

#### Preprocessing (Dijkstra) Runtime

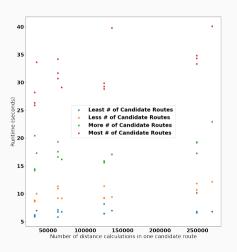


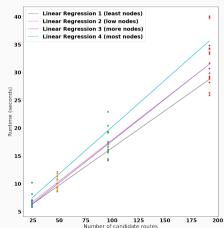
# **Computational Complexity (Metric Calculations)**

The following scatter plots demonstrates how computation grows as a function of input nodes.

### **Computational Complexity (Metric Calculations)**

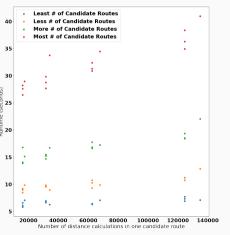


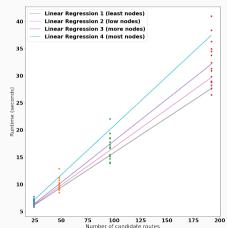




### **Computational Complexity (Metric Calculations)**







### **Computational Complexity**

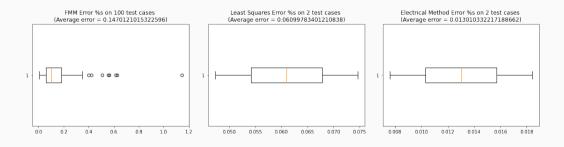
This suggests that simple distance methods (like inverse squares and least squares) have a computational complexity of O(n). Due to processing power, we could not obtain enough data to infer the computational complexity of the Wasserstein method. Because it relies on linear programming, we hypothesize that the growth rate is a polynomial of degree > 1.

Some aspects of these algorithms can be improved upon—simple parallelization techniques offer a noticeable increase. However, other aspects are inherent to the method and are difficult to improve.

### **Preliminary Numerical Results**

Due to limits in processing power, we were unable to test our algorithms against all of the data in the dataset. However, we demonstrate a few cases here.

### **Preliminary Numerical Results**



### **Future Work (Implementation)**

- While a naive Dijkstra method generates a good set of candidate routes, it has
  many shortcomings. It is still relatively slow, it is conditional on the given
  parameters, and it cannot handle stranger routes (traversing an edge more than
  once). Each of these individually can be addressed; or it may be more suitable to
  find an alternative candidate route generation method.
- Investigate Wasserstein method inconsistencies more thoroughly
- Incomplete road networks
- Implement IMU-based map matching approaches

### **Thank You! And References**

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- C. Yang and G. Gidófalvi, Fast map matching, an algorithm for integrating a hidden 50/50