

Density Approximation for Kinetic Groups

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Groups



Photos by T.R. Shankar Raman and M.M. Karim

Groups

Finding groups in...

Static point sets:

K-means clustering

DBSCAN

...

Moving point sets:

[Grouping structure](#) [Buchin et al., 2013]



Group density

Group density contains information about underlying behavior

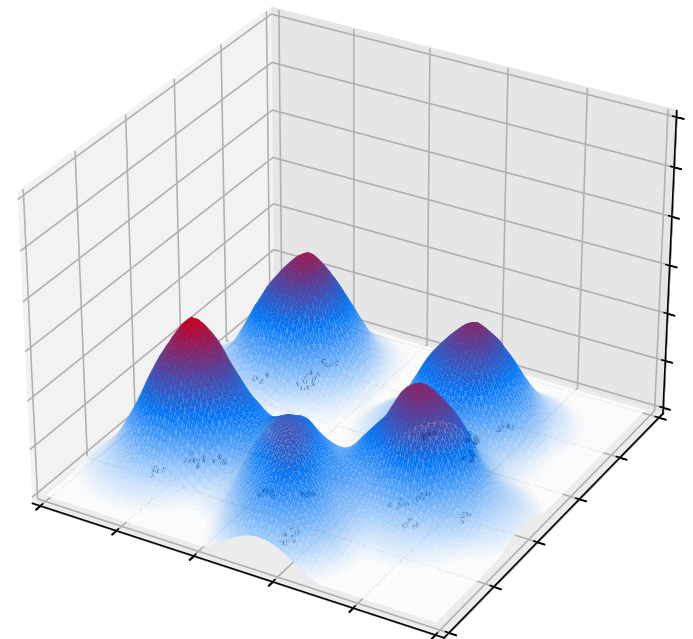
Predator proximity, feeding sites, environmental factors, ...

Function $f(x, y)$ that estimates the density at (x, y)

Goal: Maintain $f(x, y)$ as the points move

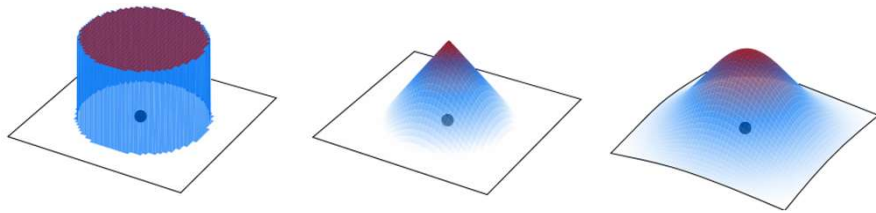


Maintain approximation of $f(x, y)$



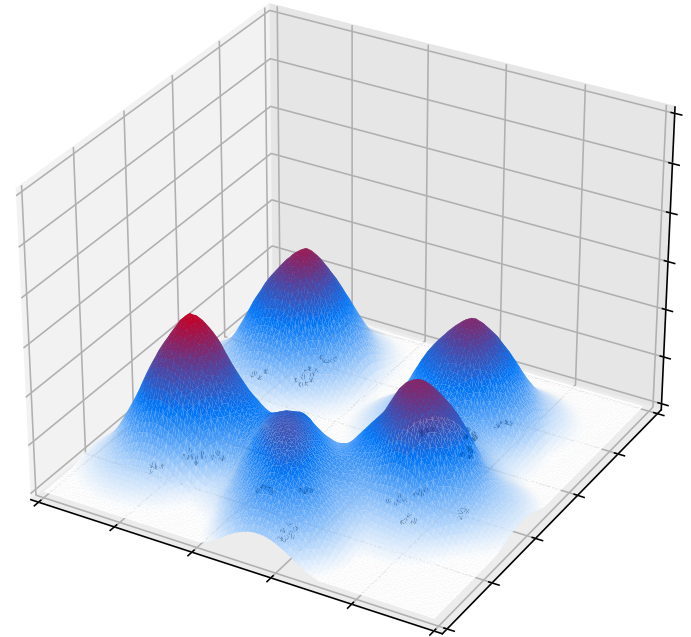
Kernel Density Estimation

Estimate the density function using **kernel function K**



The **density function** is then estimated by:

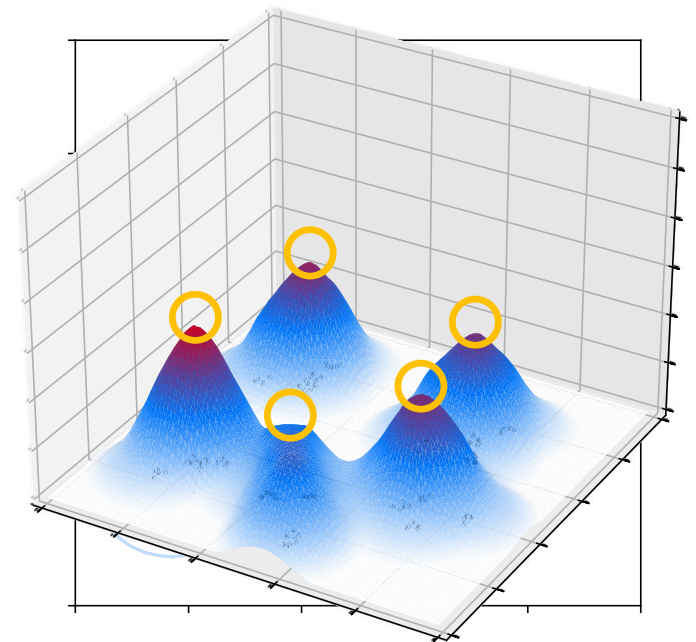
$$KDE_p(x, y) = \frac{1}{n} \sum_{p \in P} K(x - x(p), y - y(p))$$



Kernel Density Estimation

Use $KDE_p(x, y)$ to describe **group characteristics**:

- Local maxima → **Dense clusters**
- Contour lines → **Group shape**
- Kernel radius → **Spatial scales**



Kernel Density Estimation

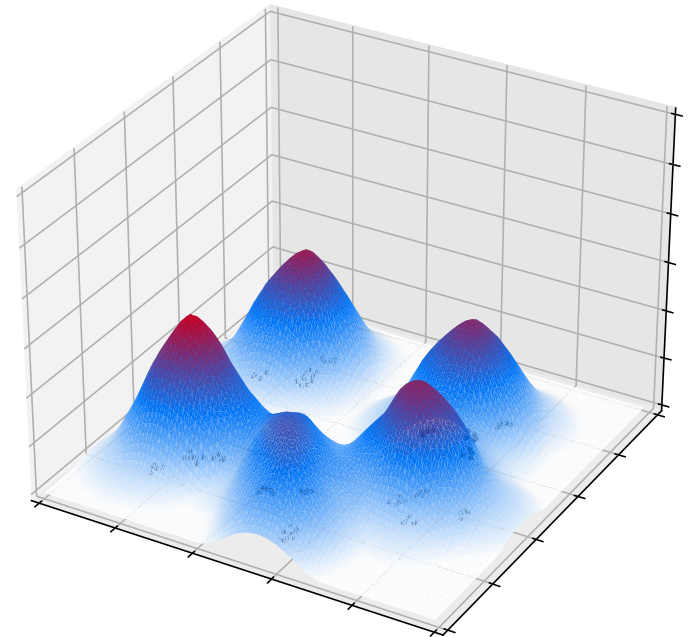
Use $KDE_p(x, y)$ to describe **group characteristics**:

Useful in practice...

but difficult to give general **theoretical guarantees**

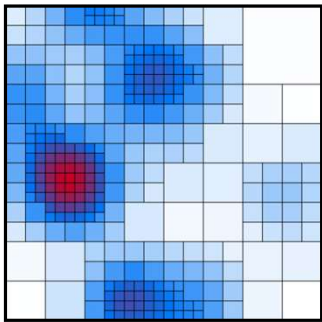
Assumptions:

- Trajectories of points are **known**
- **Piece-wise linear** movement
- Within a **bounding box**

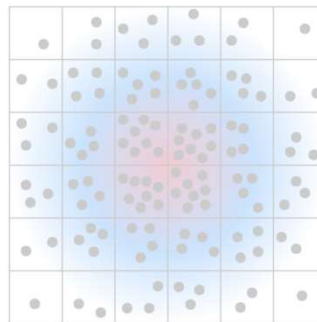


Overview

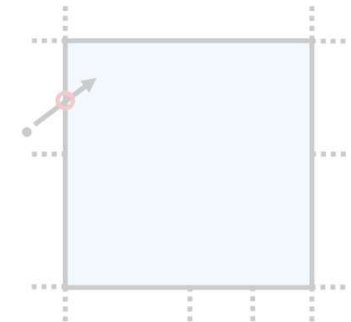
Volume-based quadtree



Discretization

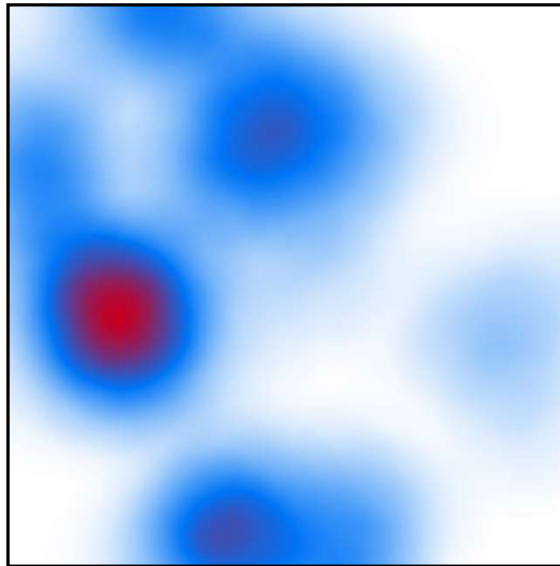


Kinetic Data Structure



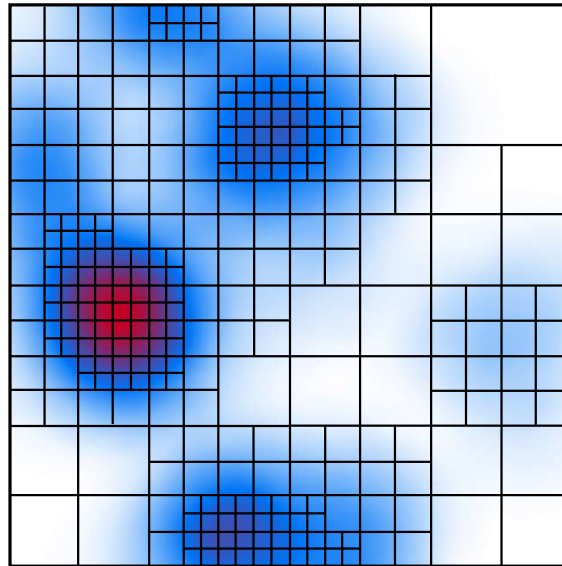
Volume-based Quadtree

Construct quadtree on **volume** under KDE_p



Volume-based Quadtree

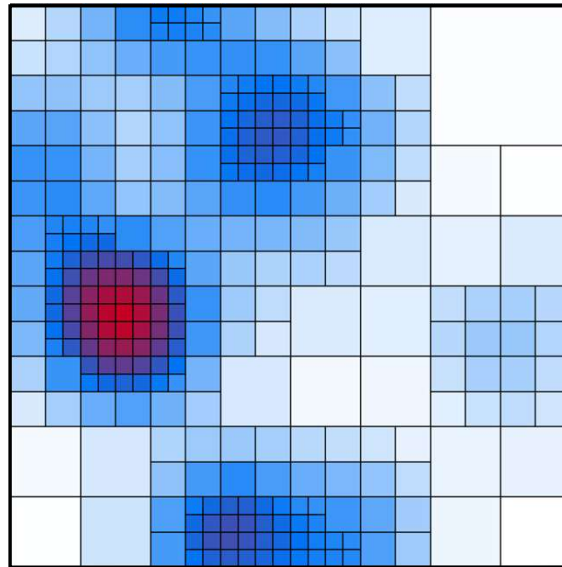
Subdivide cell v if the volume under KDE_p in v exceeds threshold $\rho > 0$



Volume-based Quadtree

Subdivide cell v if the volume under KDE_p in v exceeds threshold $\rho > 0$

Value of density approximation in v equals average value of KDE_p in v



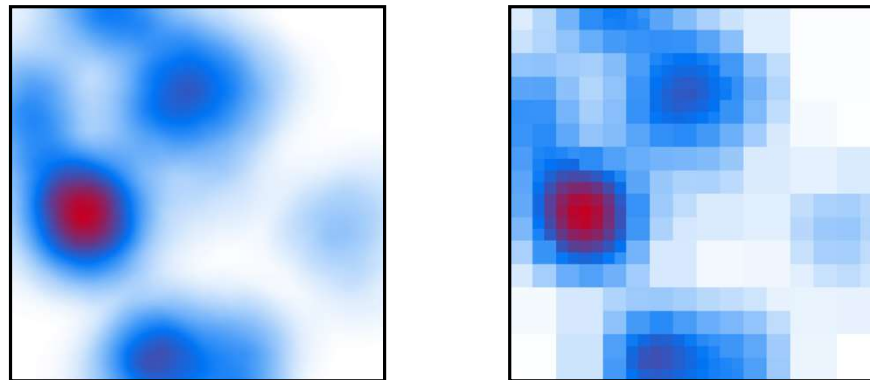
Volume-based Quadtree

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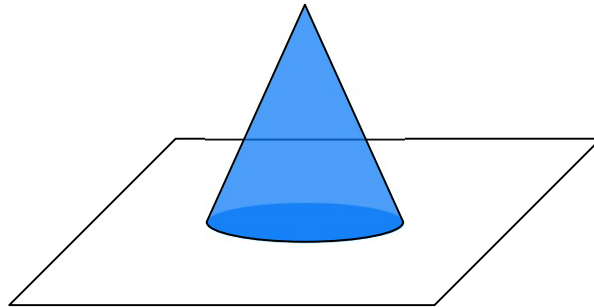
Value of density approximation in v equals average value of KDE_p in v

Measure the quality of approximation f_T of KDE_p as

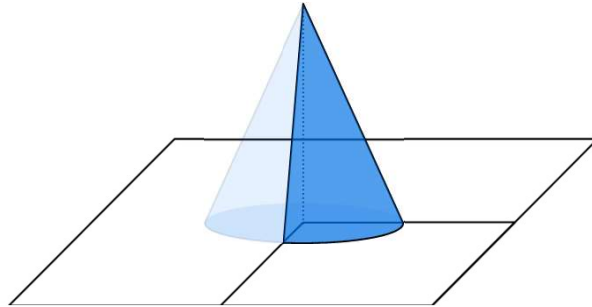
$$|KDE_p(x, y) - f_T(x, y)| \text{ for all } (x, y) \in \mathcal{D}$$



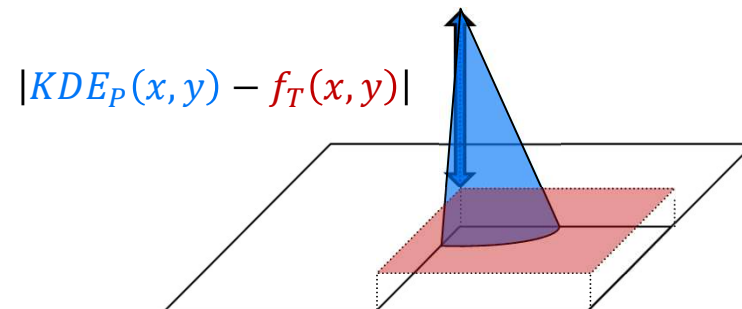
Volume-based Quadtree



Volume-based Quadtree



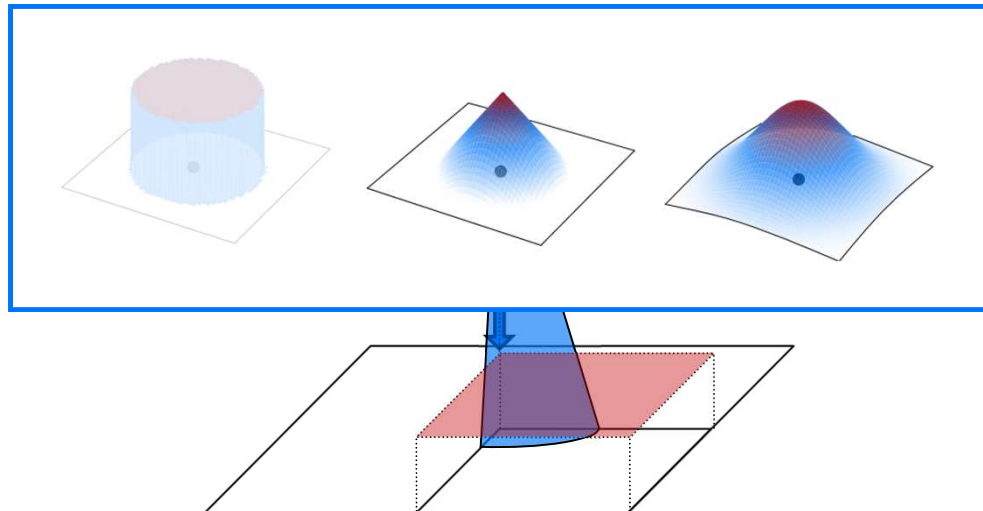
Volume-based Quadtree



Volume-based Quadtree

Bound the maximum **slope** and **height** of KDE_p

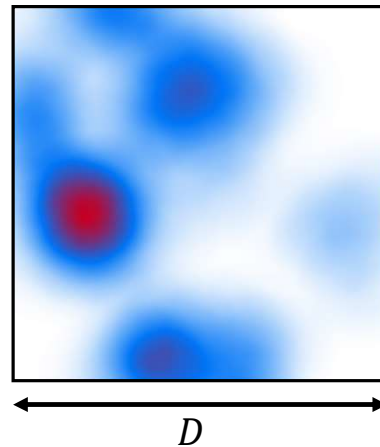
- **Lipschitz constant** λ , maximum absolute slope in any direction
- **Maximum height** z^*



Volume-based Quadtree

Bound the maximum **slope** and **height** of KDE_P

- **Lipschitz constant** λ , maximum absolute slope in any direction
- **Maximum height** z^*
- **Domain size** D

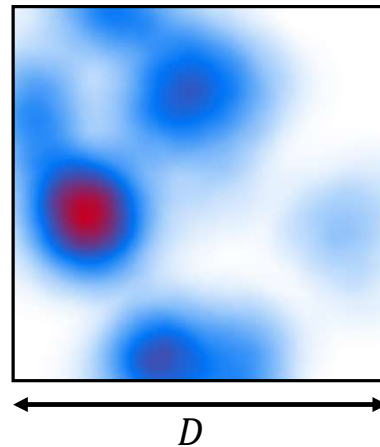


Volume-based Quadtree

For Lipschitz constant λ , maximum height z^* and domain size D :

$$|KDE_p(x, y) - f_T(x, y)| < \varepsilon \text{ for } \rho = O(\varepsilon^3)$$

with polynomial bounds on the size of the quadtree



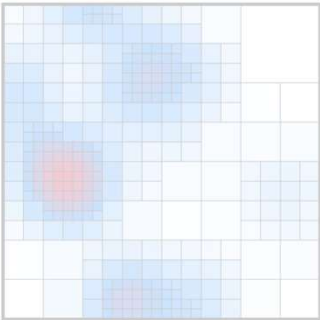
Towards moving points

Difficult to maintain as underlying points [move...](#)

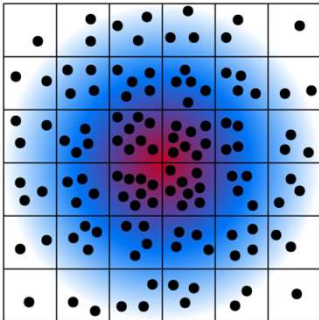
[Solution](#): Approximate volume under KDE_p using a large set of [moving points](#)

Overview

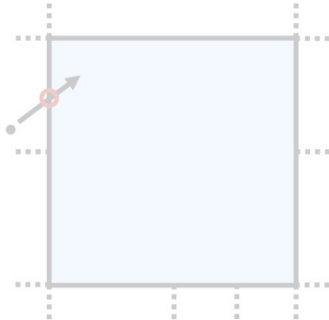
Volume-based quadtree



Discretization



Kinetic Data Structure



Discretization

Construct **moving point set** \mathcal{S} as a stand-in for the volume under KDE_p

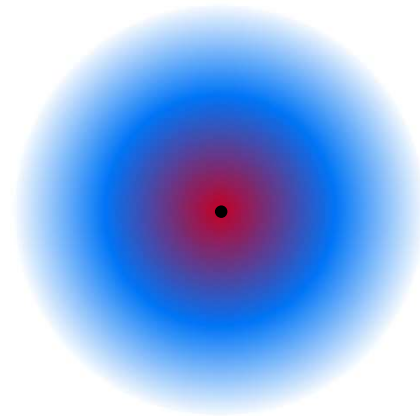
Such that at any **time** and for any **quadtree cell** v :

$$\text{volume under } KDE_p \text{ in } v \approx \text{number of points from } \mathcal{S} \text{ in } v$$

Discretization

Construct **moving point set** \mathcal{S} as a stand-in for the volume under KDE_p :

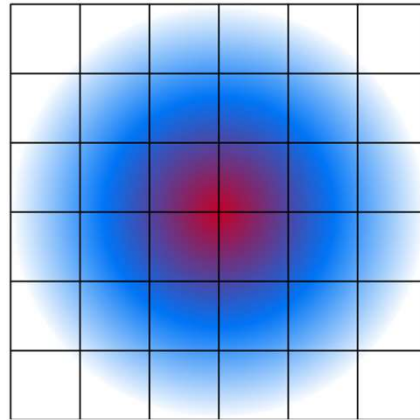
1. Approximate a **single kernel** K with a set of points \mathcal{S}



Discretization

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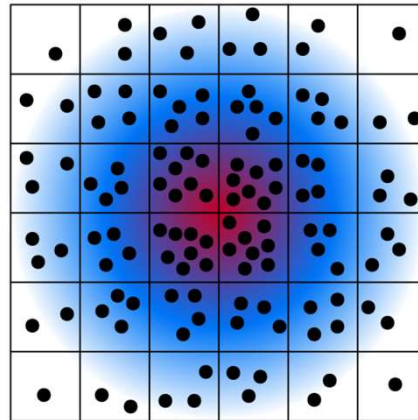
1. Approximate a **single kernel** K with a set of points \mathcal{S}
 - a. Overlay $r \times r$ **grid**



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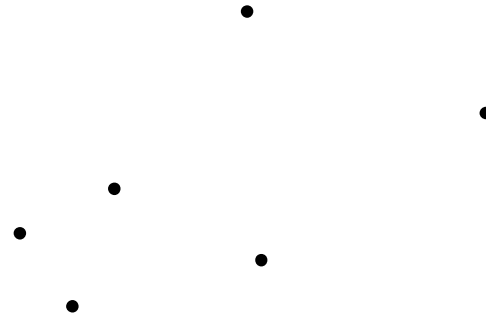
1. Approximate a **single kernel** K with a set of points \mathcal{S}
 - a. Overlay $r \times r$ **grid**
 - b. Sample **random points** in each grid cell proportional to **kernel value**



Discretization

Construct **moving point set** \mathcal{S} as a stand-in for the volume under KDE_p :

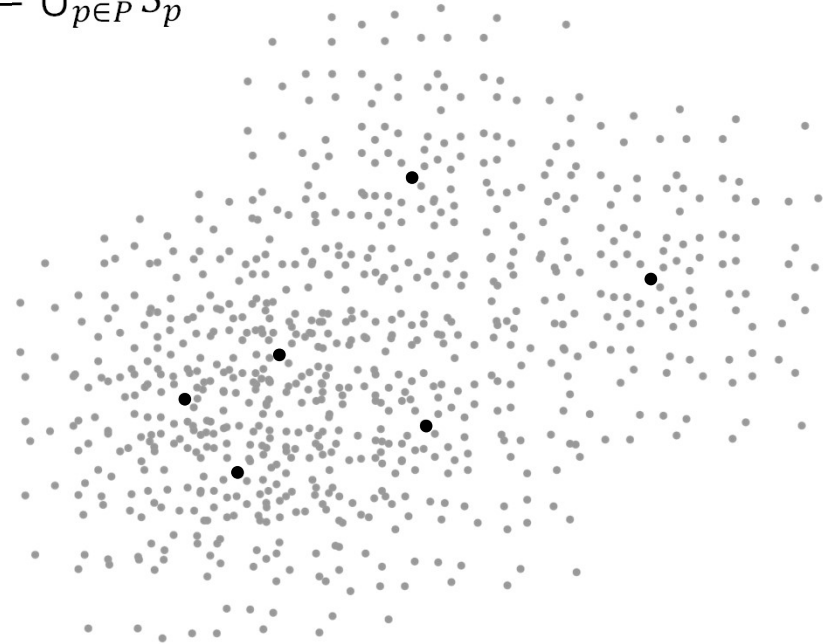
1. Approximate a **single kernel** K with a set of points \mathcal{S}
2. Approximate KDE_p with the union of kernels $\mathcal{S} = \bigcup_{p \in P} \mathcal{S}_p$



Discretization

Construct **moving point set** \mathcal{S} as a stand-in for the volume under KDE_p :

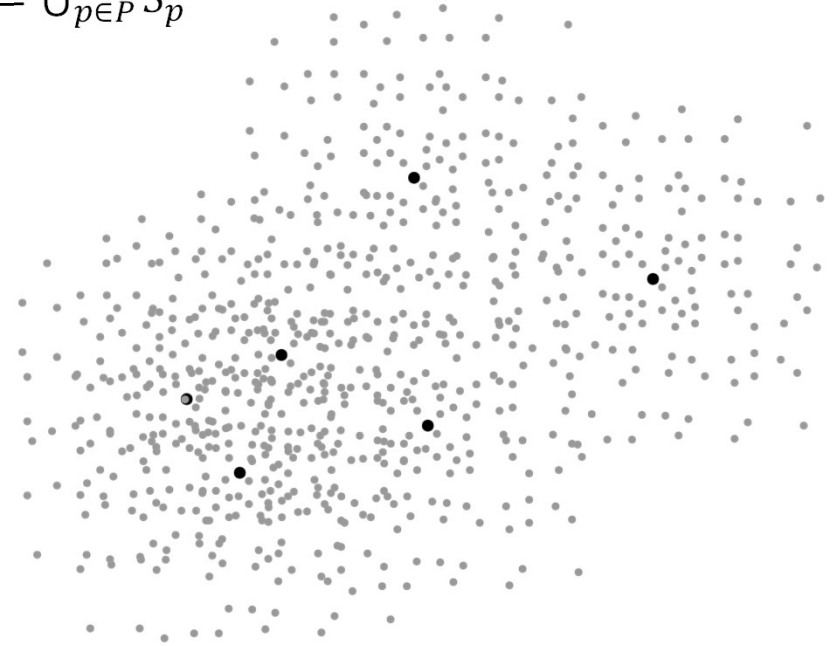
1. Approximate a **single kernel** K with a set of points S
2. Approximate KDE_p with the union of kernels $\mathcal{S} = \bigcup_{p \in P} S_p$
 - a. Assign each **input point** $p \in P$ a set S_p



Discretization

Construct **moving point set** \mathcal{S} as a stand-in for the volume under KDE_p :

1. Approximate a **single kernel** K with a set of points S
2. Approximate KDE_p with the union of kernels $\mathcal{S} = \bigcup_{p \in P} S_p$
 - a. Assign each **input point** $p \in P$ a set S_p
 - b. S_p copies the **movement** of p



Discretization

Construct **moving point set** \mathcal{S} as a stand-in for the volume under KDE_p :

1. Approximate a **single kernel** K with a set of points \mathcal{S}
2. Approximate KDE_p with the union of kernels $\mathcal{S} = \bigcup_{p \in P} \mathcal{S}_p$
3. **Reduce** the size of \mathcal{S}
 - a. Take **coreset** of \mathcal{S} [Agarwal *et al.*, 2005]

Volume-based Quadtree

Construct **volume-based** quadtree on KDE_p :

Subdivide cell v if the **volume** under KDE_p in v exceeds **threshold** $\rho > 0$

Value of **density approximation** in v equals average value of KDE_p in v

Weight-based Quadtree

Construct **volume-based** quadtree on KDE_p :

Subdivide cell v if the **volume** under KDE_p in v exceeds **threshold** $\rho > 0$

Value of **density approximation** in v equals average value of KDE_p in v

Construct **weight-based** quadtree on \mathcal{S} :

Subdivide cell v if the **number of points** from \mathcal{S} in v exceeds **threshold** $\rho > 0$

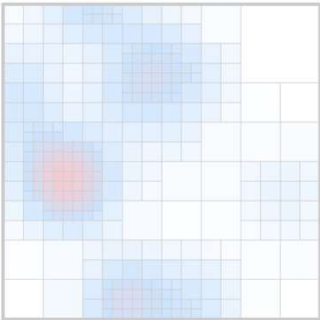
Value of **density approximation** in v proportional to the **number of points** in v

$$|KDE_p(x, y) - f_{\bar{\tau}}(x, y)| < \varepsilon \text{ for } \rho = O(\varepsilon^3)$$

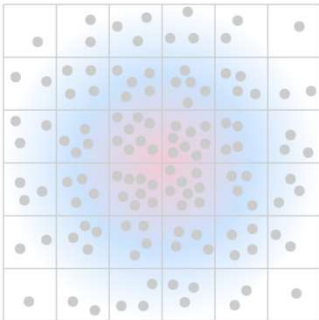
$$\text{with } |\mathcal{S}| = O\left(\frac{1}{\varepsilon^8} \log\left(\frac{1}{\varepsilon}\right)\right)$$

Overview

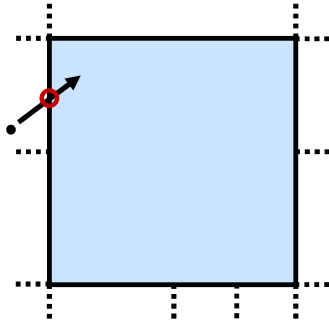
Volume-based quadtree



Discretization



Kinetic Data Structure

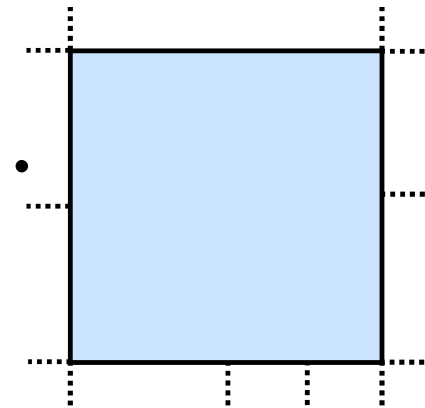


Kinetic Data Structure

Compute **weight-based quadtree** \tilde{T}

Goal: Maintain correctness of \tilde{T} as the points **move**

Observe: \tilde{T} depends only on **distribution** of P into **cells**

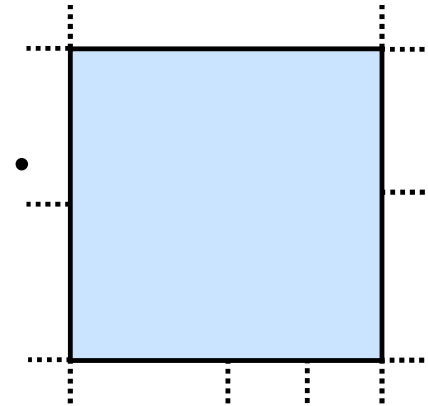


Kinetic Data Structure

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Observe: \tilde{T} changes only when a point from \mathcal{S} **crosses a cell boundary**



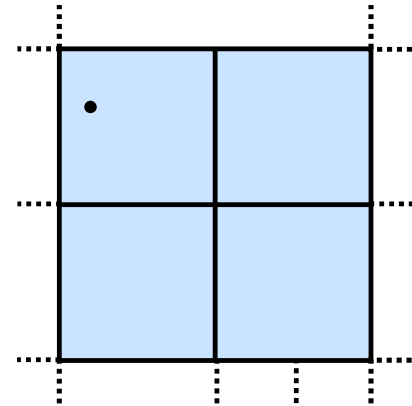
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- Update cell **boundaries**
- Update cell **heights**



Kinetic Data Structure

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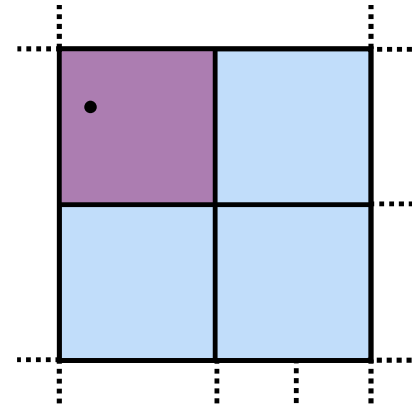
Goal: Maintain correctness of \tilde{T} as the points **move**

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- Update cell **boundaries**
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- Update **local maxima**

Sufficient to maintain \tilde{T} ...

But what about its **local maxima**?



Kinetic Data Structure

Compute **weight-based quadtree** \tilde{T}

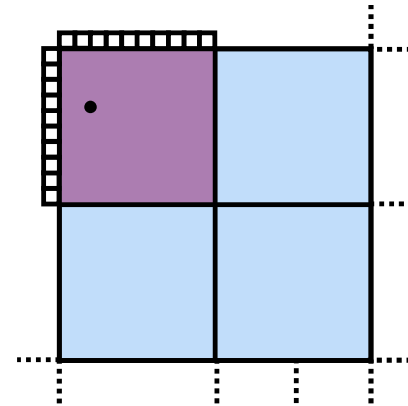
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Kinetic Data Structure

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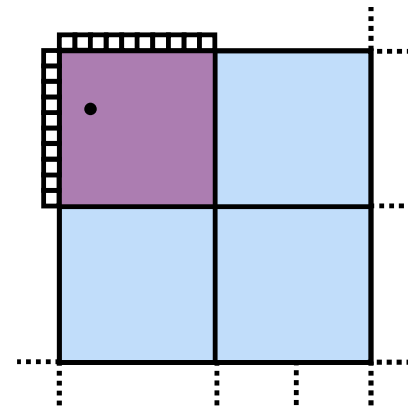
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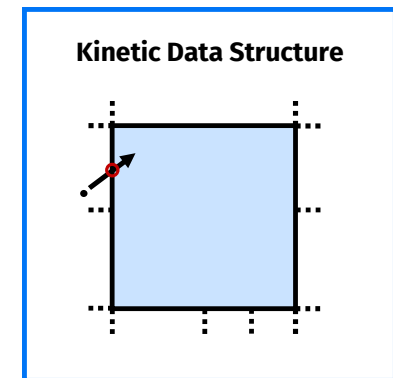
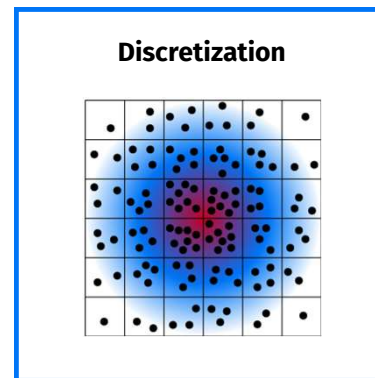
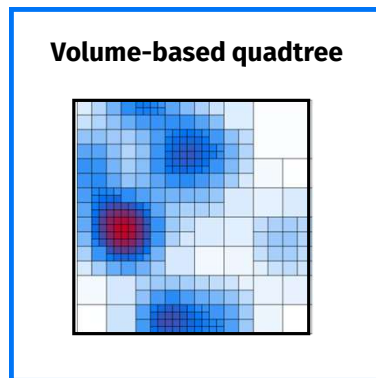
Small cells must have **high values** ...

Local maxima cannot have **small neighbors**



Summary

Theorem. Let $f = KDE_P$ be a KDE function on a set P of n linearly moving points in \mathbb{R}^2 . For any $\varepsilon > 0$, there exists a kinetic data structure that maintains an ε -approximation of f .



Summary

Theorem. *Let $f = KDE_P$ be a KDE function on a set P of n linearly moving points in \mathbb{R}^2 . For any $\varepsilon > 0$, there exists a kinetic data structure that maintains an ε -approximation of f .*

Future work:

- Engineer **practical** approach
 - Deal with assumptions
 - (known trajectories, piece-wise linear, bounding box)
- Use density surface to find **other characteristics**
 - Group shape, clustering, ...