Computational optimal transport: recent speed-ups and applications

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Who am I?

Background in **mathematics** and **data sciences**:

- **2012–2016** ENS Paris, mathematics.
- **2014–2015** M2 mathematics, vision, learning at ENS Cachan.
- **2016–2019** PhD thesis in **medical imaging** with Alain Trouvé at ENS Cachan.
- **2019–2021 Geometric deep learning** with Michael Bronstein at Imperial College.
 - **2021+ Medical data analysis** in the HeKA INRIA team (Paris).

HeKA: a translational research team for public health

Hôpitaux

Inria Inserm

Universités



My main motivation

Develop **robust and efficient** software that **stimulates other researchers**:

- 1. Speed up **geometric machine learning** on GPUs:
 - ⇒ **pyKeOps** library for distance and kernel matrices, 600k+ downloads.
- 2. Scale up **pharmacovigilance** to the full French population:
 - ⇒ **survivalGPU**, a fast re-implementation of the R survival package.
- 3. Ease access to modern statistical **shape analysis**:
 - ⇒ **GeomLoss**, truly scalable optimal transport in Python.
 - ⇒ **scikit-shapes**, alpha release now available.

Today's talk - assuming that you would enjoy some nice simulations

- 1. A quick heads up on **fast geometric methods**.
- 2. Efficient discrete optimal transport solvers.
- 3. New applications for systems of **incompressible particles**.

How to code a N-body simulation?

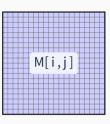
Scientific computing libraries represent most objects as tensors

Context. Constrained **memory accesses** on the GPU:

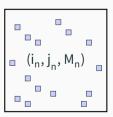
- Long access times to the registers penalize the use of large dense arrays.
- Hard-wired contiguous memory accesses penalize the use of sparse matrices.

Challenge. In order to reach optimal run times:

- **Restrict** ourselves to operations that are supported by the constructor: convolutions, FFT, etc.
- Develop new routines from scratch in C++/CUDA (FAISS, KPConv...): several months of work.



Dense array



Sparse matrix

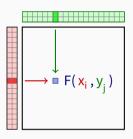
The KeOps library: efficient support for symbolic matrices

Solution. KeOps-www.kernel-operations.io:

- For PyTorch, NumPy, Matlab and R, on **CPU and GPU**.
- Automatic differentiation.
- Just-in-time compilation of optimized C++ schemes, triggered for every new reduction: sum, min, etc.

If the formula "F" is simple (\leqslant 100 arithmetic operations): "100k \times 100k" computation \rightarrow 10ms – 100ms, "1M \times 1M" computation \rightarrow 1s – 10s.

Hardware ceiling of 10^{12} operations/s. \times **10 to** \times **100 speed-up** vs standard GPU implementations for a wide range of problems.



Symbolic matrix Formula + data

- Distances d(x_i,y_j).
 Kernel k(x_i,y_i).
- Numerous transforms.

A first example: efficient nearest neighbor search in dimension 50

Create large point clouds using **standard PyTorch syntax**:

import torch N, M, D = 10**6, 10**6, 50 x = torch.rand(N, 1, D).cuda() # (1M, 1, 50) array y = torch.rand(1, M, D).cuda() # (1, 1M, 50) array

Turn **dense** arrays into **symbolic** matrices:

```
from pykeops.torch import LazyTensor
x_i, y_j = LazyTensor(x), LazyTensor(y)
```

Create a large **symbolic matrix** of squared distances:

```
D_{ij} = ((x_i - y_j) ** 2).sum(dim=2) # (1M, 1M) symbolic
```

Use an .argmin() **reduction** to perform a nearest neighbor query:

```
indices_i = D_ij.argmin(dim=1) # -> standard torch tensor
```

The KeOps library combines performance with flexibility

Script of the previous slide = efficient nearest neighbor query, on par with the bruteforce CUDA scheme of the FAISS library... And can be used with any metric!

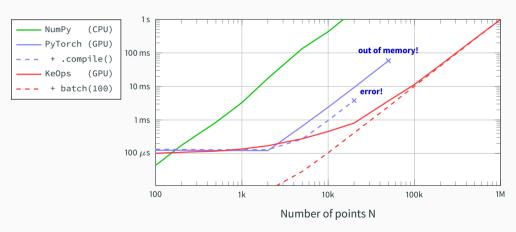
```
D_ij = ((x_i - x_j) ** 2).sum(dim=2)  # Euclidean
M_ij = (x_i - x_j).abs().sum(dim=2)  # Manhattan
C_ij = 1 - (x_i | x_j)  # Cosine
H_ij = D_ij / (x_i[...,0] * x_j[...,0])  # Hyperbolic
```

KeOps supports arbitrary **formulas** and **variables** with:

- Reductions: sum, log-sum-exp, K-min, matrix-vector product, etc.
- **Operations:** +, \times , sqrt, exp, neural networks, etc.
- Advanced schemes: batch processing, block sparsity, etc.
- Automatic differentiation: seamless integration with PyTorch.

KeOps lets users work with millions of points at a time

Benchmark of a Gaussian **convolution** $a_i \leftarrow \sum_{j=1}^N \exp(-\|x_i - y_j\|_{\mathbb{R}^3}^2) \, b_j$ between **clouds of N 3D points** on a A100 GPU.



Yet another ML compiler?

Many impressive tools out there (Numba, Triton, Halide...):

- Focus on **generality** (software + hardware).
- Increasingly easy to use via e.g. PyTorch 2.0.

KeOps fills a different niche (a bit like cuFFT, FFTW...):

- Focus on a **single major bottleneck**: geometric interactions.
- Agnostic with respect to Euclidean / non-Euclidean formulas.
- Fully compatible with PyTorch, NumPy, R.
- Can actually be used by mathematicians.

KeOps is a **bridge** between geometers (with a maths background) and compiler experts (with a CS background).

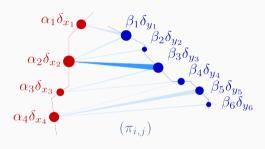
How to solve the OT problem?

Duality: central planning with NM variables \simeq outsourcing with N + M variables

$$\begin{array}{lll} \mathsf{OT}(\mathsf{A},\mathsf{B}) \ = \ \min_{\pi} \ \langle \, \pi \,,\, \mathsf{C} \, \rangle, \ \ \mathsf{with} \ \mathsf{C}(\pmb{x_i},y_j) \ = \ \frac{1}{p} \|\pmb{x_i} - y_j\|^p & \longrightarrow \ \ \mathsf{Assignment} \\ & \mathsf{s.t.} \ \pi \, \geqslant \, 0, \quad \pi \, \mathbf{1} \ = \ \mathsf{A}, \quad \pi^\mathsf{T} \mathbf{1} \ = \ \mathsf{B} \end{array}$$



$$\sum_{i,j} \pi_{i,j} \operatorname{C}({\color{black} \boldsymbol{x_i}}, y_j)$$

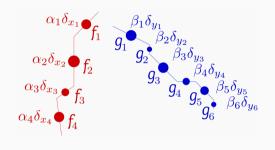


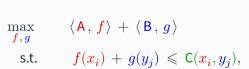
Duality: central planning with NM variables \simeq outsourcing with N + M variables

$$\begin{split} \mathsf{OT}(\mathsf{A},\mathsf{B}) \ = \ \min_{\boldsymbol{\pi}} \ \langle \, \boldsymbol{\pi} \,,\, \mathsf{C} \, \rangle, \ \ \mathsf{with} \, \mathsf{C}(\pmb{x_i},y_j) \ = \ \tfrac{1}{p} \|\pmb{x_i} - y_j\|^p \qquad \longrightarrow \ \ \mathsf{Assignment} \\ \mathsf{s.t.} \ \ \boldsymbol{\pi} \ \geqslant \ 0, \quad \boldsymbol{\pi} \, \mathbf{1} \ = \ \mathsf{A}, \quad \boldsymbol{\pi}^\mathsf{T} \mathbf{1} \ = \ \mathsf{B} \end{split}$$



$$\sum_{i,j} \pi_{i,j} \, \mathsf{C}(\pmb{x_i}, \pmb{y_j})$$







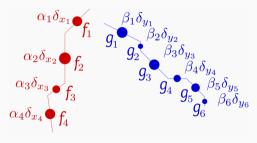
$$rac{1}{i} lpha_i f_i + \sum_j eta_j g_j \
ightarrow ext{FedEx}$$

Duality: central planning with NM variables \simeq outsourcing with N + M variables

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$$\sum_{i,j} \pi_{i,j} \, \mathsf{C}(\boldsymbol{x_i}, y_j)$$



$$= \max_{f,\,g} \qquad \langle \mathsf{A},\,f
angle + \langle \mathsf{B},\,g
angle \ ext{s.t.} \qquad f(x_i) + g(y_j) \leqslant \mathsf{C}(x_i,y_j),$$



Being too greedy... doesn't work!

$$\begin{aligned} \text{OT}(\alpha,\beta) &= & \max_{\substack{(f_i) \in \mathbb{R}^{\text{N}} \\ (g_j) \in \mathbb{R}^{\text{M}}}} \sum_{i=1}^{\text{N}} \alpha_i f_i + \sum_{j=1}^{\text{M}} \beta_j g_j \\ &\text{s.t. } \forall i,j, \ f_i + g_j \leqslant \mathbf{C}(x_i,y_j) \end{aligned}$$

Algorithm 3.1: Naive greedy algorithm

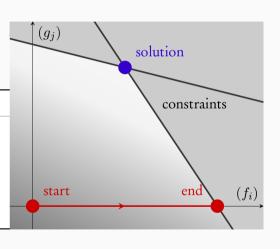
1:
$$f_i, g_j \leftarrow \mathbf{0}_{\mathbb{R}^{\mathrm{N}}}, \mathbf{0}_{\mathbb{R}^{\mathrm{M}}}$$

2: repeat

3:
$$f_i \leftarrow \min_{j=1}^{\mathrm{M}} \left[\mathbf{C}(x_i, y_j) - g_j \right]$$

4:
$$g_j \leftarrow \min_{i=1}^{N} \left[\mathbf{C}(x_i, y_j) - f_i \right]$$

- 5: until convergence.
- 6: return f_i , g_j



The auction algorithm: take it easy with a slackness $\, arepsilon > 0 \,$

$$\begin{aligned} \text{OT}(\alpha,\beta) &= & \max_{\substack{(f_i) \in \mathbb{R}^{\mathrm{N}} \\ (g_j) \in \mathbb{R}^{\mathrm{M}}}} \sum_{i=1}^{\mathrm{N}} \alpha_i f_i + \sum_{j=1}^{\mathrm{M}} \beta_j g_j \\ &\text{s.t. } \forall i,j, \ f_i + g_j \ \leqslant \ \mathbf{C}(x_i,y_j) \end{aligned}$$

Algorithm 3.2: Pseudo-auction algorithm

1:
$$f_i, g_j \leftarrow \mathbf{0}_{\mathbb{R}^N}, \mathbf{0}_{\mathbb{R}^M}$$

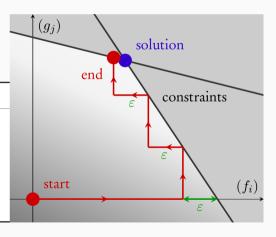
2: repeat

3:
$$f_i \leftarrow \min_{j=1}^{M} \left[\mathbf{C}(x_i, y_j) - g_j \right] - \varepsilon$$

4:
$$g_j \leftarrow \min_{i=1}^{N} [\mathbf{C}(x_i, y_j) - f_i]$$

5: **until** $\forall i, \exists j, f_i + g_j \geq \mathbf{C}(x_i, y_j) - \varepsilon$.

6: return f_i , g_j



The Sinkhorn algorithm: use a softmin, get a well-defined optimum

$$\begin{aligned} \text{OT}(\alpha,\beta) &= & \max_{\substack{(f_i) \in \mathbb{R}^N \\ (g_j) \in \mathbb{R}^M}} \sum_{i=1}^N \alpha_i f_i + \sum_{j=1}^M \beta_j g_j \\ &- \varepsilon \log \left\langle \alpha_i \otimes \beta_j, \exp \frac{1}{\varepsilon} [f_i \oplus g_j - \mathbf{C}_{ij}] \right\rangle \end{aligned}$$

Algorithm 3.3: Sinkhorn or "soft-auction" algorithm

1:
$$f_i, g_j \leftarrow \mathbf{0}_{\mathbb{R}^{\mathrm{N}}}, \mathbf{0}_{\mathbb{R}^{\mathrm{M}}}$$

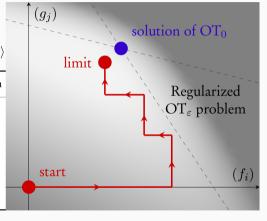
2: repeat

3:
$$f_i \leftarrow -\varepsilon \log \sum_{j=1}^{M} \beta_j \exp \frac{1}{\varepsilon} [g_j - \mathbf{C}(x_i, y_j)]$$

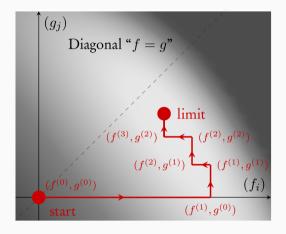
4:
$$g_j \leftarrow -\varepsilon \log \sum_{i=1}^{N} \alpha_i \exp \frac{1}{\varepsilon} [f_i - \mathbf{C}(x_i, y_j)]$$

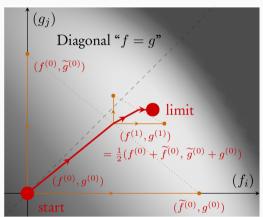
5: **until** convergence up to a set tolerance.

6: return f_i , g_j



The symmetric Sinkhorn algorithm: stay close to the diagonal if A \simeq B





Remark 1: a streamlined algorithm

One key operation – the soft, **weighted distance transform**:

$$\forall i \in [1, \mathsf{N}], \ f(x_i) \leftarrow \min_{y \sim \beta} \left[\mathsf{C}(x_i, y) - g(y) \right] = -\varepsilon \log \sum_{j=1}^{\mathsf{M}} \beta_j \exp \tfrac{1}{\varepsilon} \left[g_j - \mathsf{C}(x_i, y_j) \right].$$

Similar to the chamfer distance transform, convolution with a Gaussian kernel... Fast implementations with **pyKeOps**:

- If $\mathsf{C}(x_i,y_j)$ is a closed formula: **bruteforce** scales to N, M \simeq 100k in 10ms on a GPU.
- If A and B have a low-dimensional support:
 use a clustering and truncation strategy to get a x10 speed-up.
- If A and B are supported on a 2D or 3D grid and $C(x_i, y_j) = \frac{1}{2} ||x_i y_j||^2$: use a **separable** distance transform to get a second x10 speed-up. (N.B.: FFTs run into numerical accuracy issues.)

Remark 2: annealing works!

The **Auction/Sinkhorn** algorithms:

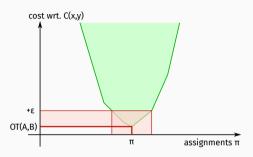
- Improve the dual cost by at least ε at each (early) step.
- Reach an ε -optimal solution with $(\max C) / \varepsilon$ steps.

Simple heuristic: run the optimization with **decreasing values** of ε .

 ε -scaling

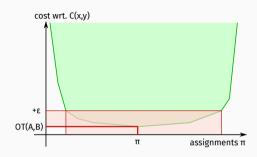
- simulated annealing
- = multiscale strategy
- = divide and conquer

Remark 3: the curse of dimensionality



In low dimension:

- ||x y|| takes large and small values.
- The OT objective is **peaky** wrt. π .
- ε -optimal solutions are **useful**.
- OT(discrete samples) ≃
 OT(underlying distributions)



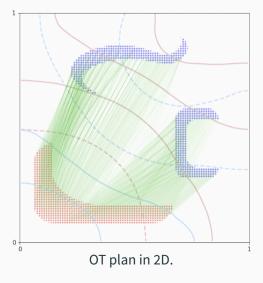
In **high dimension**:

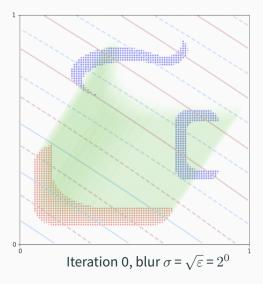
- $\|x-y\|$ gets closer to a constant.
- The OT objective is **flat** wrt. π .
- ε -optimal solutions are **random**.
- OT(discrete samples) \neq OT(underlying distributions)

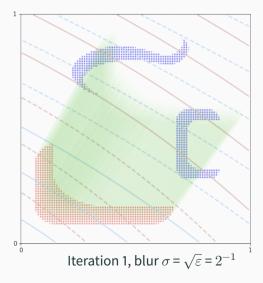
To recap 80+ years of work...

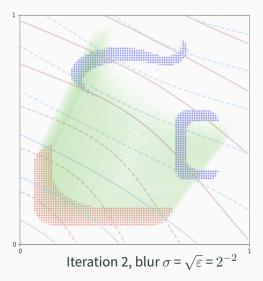
Key dates for discrete optimal transport with N points:

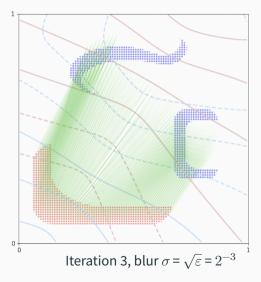
- [Kan42]: **Dual** problem of Kantorovitch.
- [Kuh55]: **Hungarian** methods in $O(N^3)$.
- [Ber79]: **Auction** algorithm in $O(N^2)$.
- [KY94]: **SoftAssign** = Sinkhorn + simulated annealing, in $O(N^2)$.
- [GRL+98, CR00]: **Robust Point Matching** = Sinkhorn as a loss.
- [Cut13]: Start of the GPU era.
- [Mér11, Lév15, Sch19]: **multi-scale** solvers in $O(N \log N)$.
- Solution, today: Multiscale Sinkhorn algorithm, on the GPU.
 - \Longrightarrow Generalized **QuickSort** algorithm.

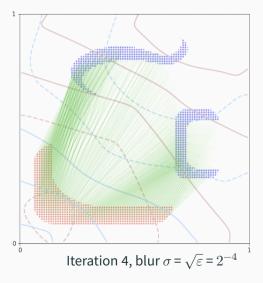


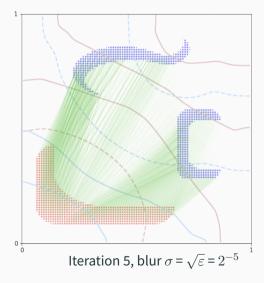


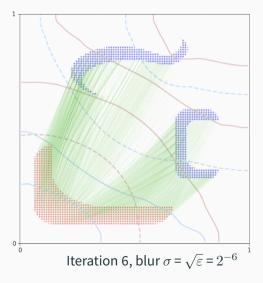


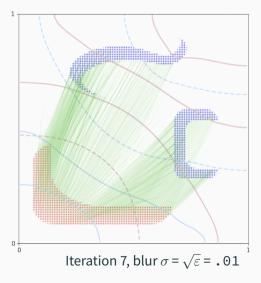


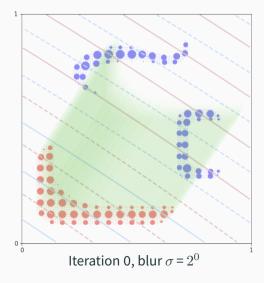


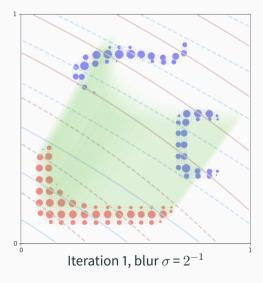


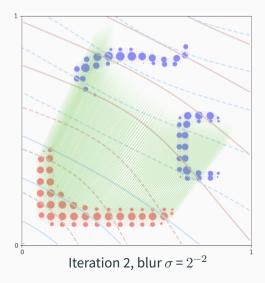


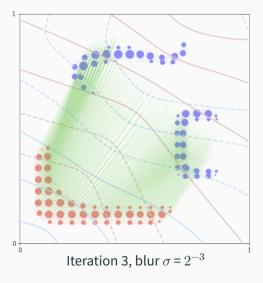


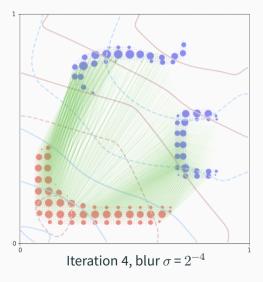


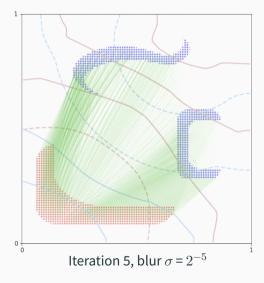


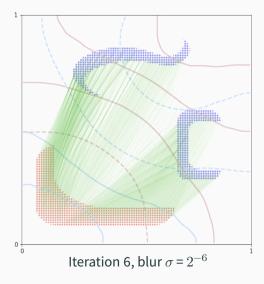


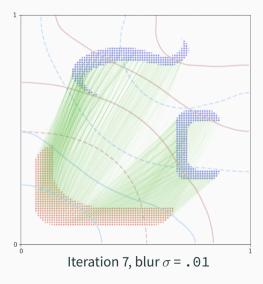












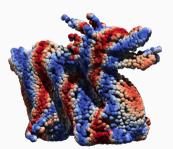
Scaling up optimal transport to anatomical data

Progresses of the last decade add up to a $\times 100 - \times 1000$ acceleration:

$$Sinkhorn~GPU \xrightarrow{\times 10} + KeOps \xrightarrow{\times 10} + Annealing \xrightarrow{\times 10} + Multi-scale$$

With a precision of 1%, on a modern gaming GPU:



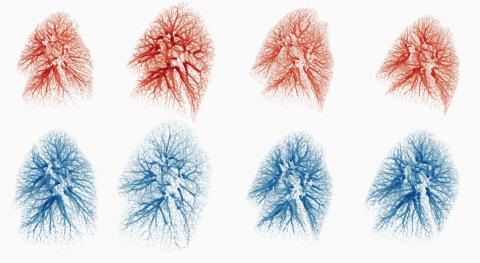


10k points in 30-50ms



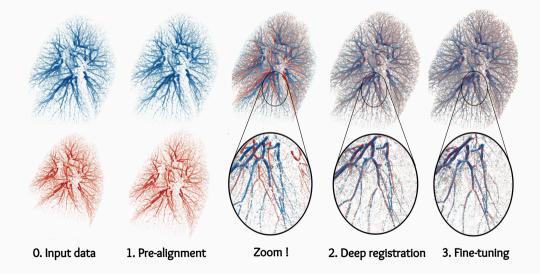
100k points in 100-200ms

A typical example in anatomy: lung registration "Exhale - Inhale"



Complex deformations, high **resolution** (50k–300k points), high **accuracy** (< 1mm).

Three-steps registration



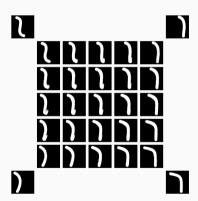
Wasserstein barycenters [AC11]

Barycenter
$$\mathbf{A}^* = \arg\min_{\mathbf{A}} \sum_{i=1}^{4} \lambda_i \operatorname{Loss}(\mathbf{A}, \mathbf{B}_i)$$
.



Euclidean barycenters.

$$\mathsf{Loss}(\mathsf{A},\mathsf{B}) = \|\mathsf{A} - \mathsf{B}\|_{L^2}^2$$



Wasserstein barycenters.

$$Loss(A, B) = OT(A, B)$$

Wasserstein barycenters

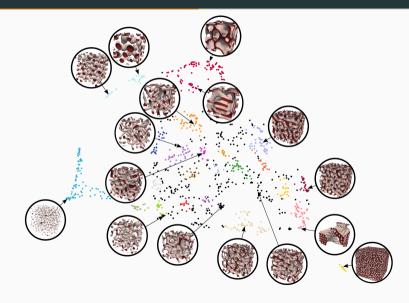
From a computational perspective:

- The problem is **convex** (easy) wrt. the weights.
- The support of the barycenter lies in the **convex hull** of the input distributions.

The curse of dimensionality hits hard:

- In high dimension, identifying the support can become **NP-hard**.
- In dimensions 2 and 3, we can just use a grid and recover super fast algorithms.
 Computing OT distances and barycenters between density maps is a solved problem.
 - ⇒ We can now **easily** do manifold learning with e.g. UMAP in Wasserstein spaces of **2D and 3D** distributions.

An example: Anna Song's exploration of 3D shape textures [Son22]





Particle systems

Two very talented postdocs

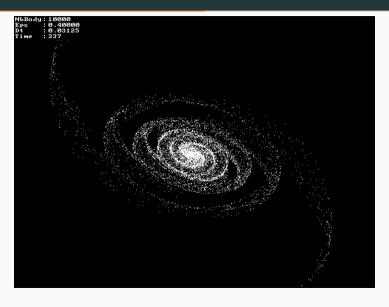


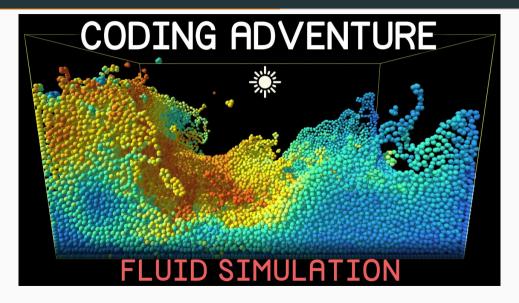
Maciej Buze Heriot-Watt University



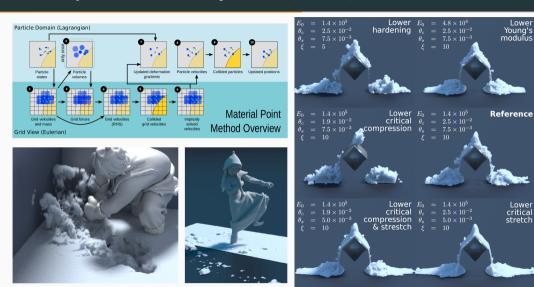
Antoine Diez Kyoto University

Original motivation: the N-body problem [Pri11]

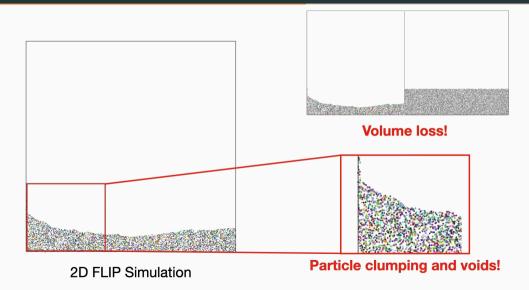




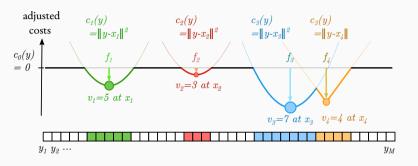
The material point method: Disney's Frozen [SSC+13]

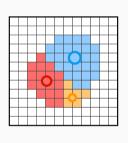


How can we enforce a volume preservation constraint? [QLDGJ22]



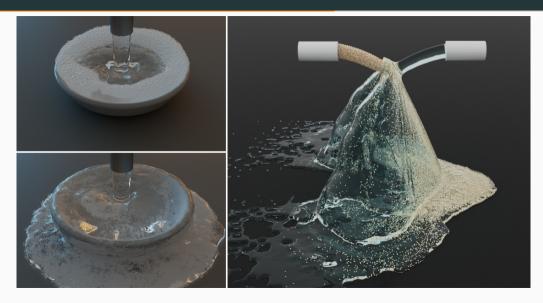
Use power diagrams i.e. semi-discrete optimal transport



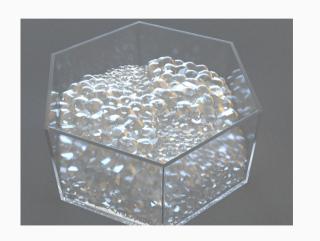


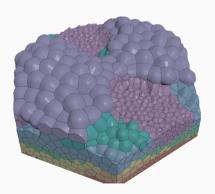
- The f_i 's maximize the dual objective $\sum_{i=1}^N v_i f_i + \int_{v \in \Omega} \min_{i=0}^N [\, c_i(y) f_i\,] \, \mathrm{d}y.$
- Optimality conditions \iff Vol(Cell_i) = v_i .
- To **compute the cells**, the objective and its gradient:
 - If $c_i(y) = \|y x_i\|^2$ for all cells, use a clever **grid-free** algorithm.
 - Otherwise, just use **KeOps**.

Power plastics [QLY⁺23]



Power plastics [QLY+23] – without the eye candy





Main numerical ingredients

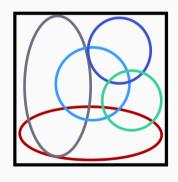
These simulations alternate between:

- 1. **Moving the particles** according to your favorite N-body model.
- 2. Computing Laguerre cells with the correct volumes:
 - (Multiscale) Sinkhorn for tolerance > 5%.
 - (Quasi-)Newton for tolerance < 1%.
- 3. **Correcting** the particle positions to enforce the volume-preservation constraint:
 - Jump to the centroid of the cell.
 - Or add a spring for smoother trajectories.

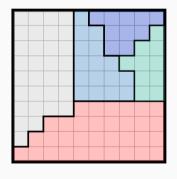
See e.g. Thomas Gallouët for a rigorous analysis with Mérigot, Lévy, etc.

But today: new applications with **custom cost functions** (thanks KeOps).

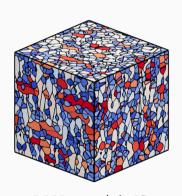
Anisotropic power diagrams let us model polycrystalline metals [BFR+24]



Ellipsoids.

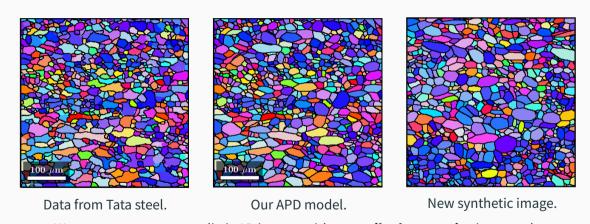


Pixel cells.



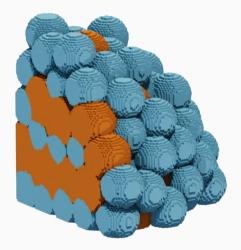
5,000 crystals in 3D.

Fit to real EBSD scan of low-carbon steel [BFR⁺24]

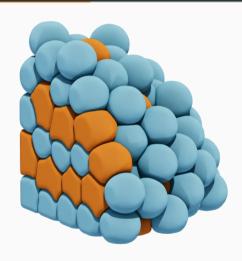


We can generate new, realistic 3D images with **prescribed properties** in seconds.

Change the cost function to simulate hard (blue) and soft (orange) cells [DF24]

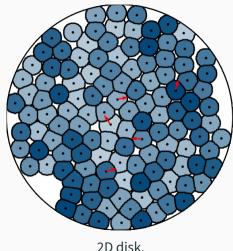


The **raw** 100x100x100 pixel grid...



with some Hollywood **makeup**.

Run-and-tumble motion [DF24]

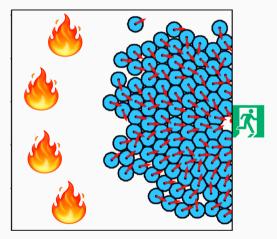


2D disk.

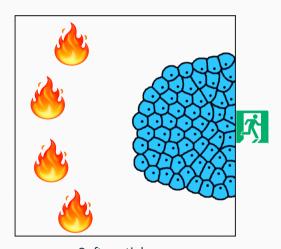


3D cube.

Fire alarm! [DF24]

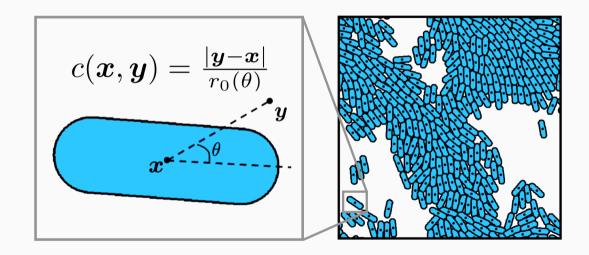


Hard particles **burn**.

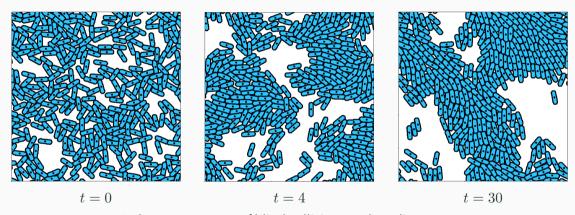


Soft particles **escape**.

Self-organizing swarms of blind, incompressible swimmers [DF24]

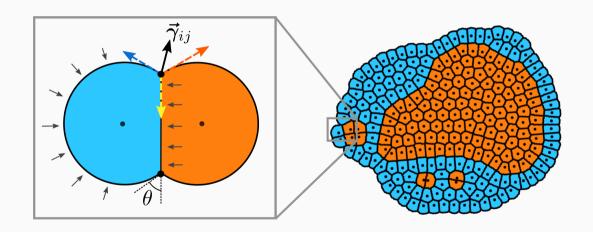


Self-organizing swarms of blind, incompressible swimmers [DF24]

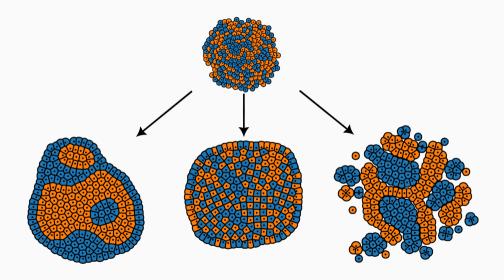


Order emerges out of blind collisions and re-alignments.

Surface tension [DF24]



Surface tension [DF24] – playing with the energy parameters





Conclusion

Genuine team work



Benjamin Charlier



Joan Glaunès



Thibault Séjourné



F.-X. Vialard



Gabriel Peyré



Alain Trouvé



Marc Niethammer



Shen Zhengyang



Olga Mula



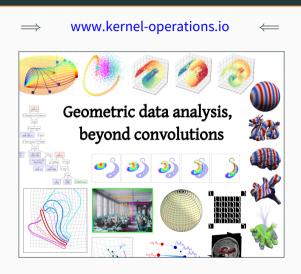
Hieu Do

Key points

- Optimal Transport = volume preservation = generalized sorting :
 - → Super-fast solvers on **simple domains**, especially 2D/3D spaces.
 - → **Fundamental tool** at the intersection of geometry and statistics.
- "Video-game physics" is great for modelling:
 - → **Expressive**, real-time simulations that you can implement without being a Finite Elements guru.
- GPUs are more versatile than you think.
 - Ongoing work to provide **fast GPU backends** to researchers, going beyond what Google and Facebook are ready to pay for.

2026 target for scientific Python: **interactive**, **web-based** simulations à la ShaderToy.

Documentation and tutorials are available online



www.jeanfeydy.com/geometric_data_analysis.pdf



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