Fast geometric libraries for vision and data sciences

.

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Who am I?

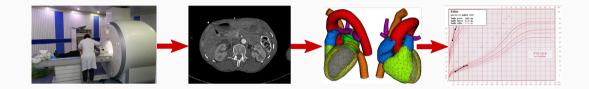
Background in **mathematics** and **data sciences**:

- 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).

Close ties with **healthcare**:

- 2015 Image denoising with Siemens Healthcare in Princeton.
- 2019+ MasterClass Al–Imaging, for radiology interns in the University of Paris.
- 2020+ Colloquium on Medical imaging in the AI era at the Paris Brain Institute.

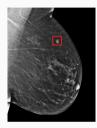
My motivation: medical data analysis



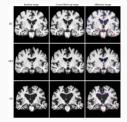
Three main characteristics:

- Heterogeneous data: patient history, images, etc.
- Small stratified samples: 10 1000 patients per group.
- Dealing with **outliers** and the **heavy tails** of our distributions is a priority.

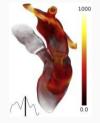
Computational anatomy [CSG19, LSG⁺18, CMN14]



Detect a pattern.



Analyze a variation.



Register a model.

Some characteristics, in the wider context of computer vision research:

- Standard acquisitions, without occlusions.
- Precision work (at millimeter scale).
- Need for **guarantees** of robustness and regularity.

Target. Design models that combine medical **expertise** with modern **datasets**.

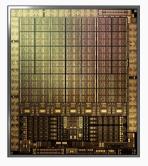
Challenge. The advent of Graphics Processing Units (GPU):

• Incredible value for money:

1 000€ \simeq 1 000 cores \simeq 10¹² operations/s.

• Bottleneck: constraints on register usage.

"User-friendly" Python ecosystem, consolidated around a **small number of key operations**.



7,000 cores in a single GPU.

Solution. Expand the standard toolbox in data sciences to deal with the challenges of the healthcare industry.

Ease the development of **advanced models**, without compromising on numerical performance.

In-depth work, numerical foundations \longrightarrow high-level applications:

- 1. Efficient manipulation of "symbolic" matrices (distances, kernel, etc.).
- 2. Optimal transport: generalized sorting methods.
- 3. Geometric deep learning and biomedical applications.

Future of these tools and clinical perspectives.

1. Symbolic matrices

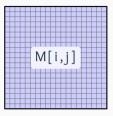
Computing libraries represent most objects as tensors

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

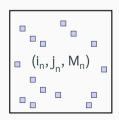
- Long access times to the registers penalize the use of large dense arrays.
- Hard-wired **contiguous** memory accesses penalize the use of **sparse** matrices.

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



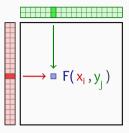
The KeOps library: efficient support for symbolic matrices

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 (\leq 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 matric 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!

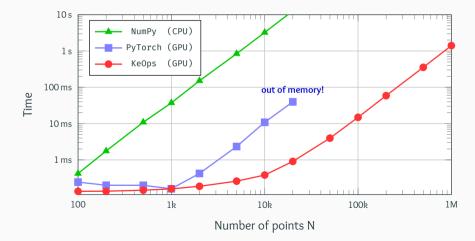
D_ij	= ((x_i - x_j) ** 2).sum(dim=2)	#	Euclidean
M_ij	= (x_i - x_j).abs().sum(dim=2)	#	Manhattan
C_ij	= 1 - (x_i x_j)	#	Cosine
H_ij	= $D_{ij} / (x_{i[,0]} * x_{j[,0]})$	#	Hyperbolic

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.

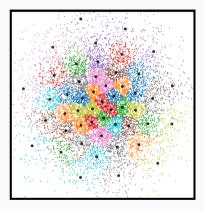
KeOps lets users work with millions of points at a time

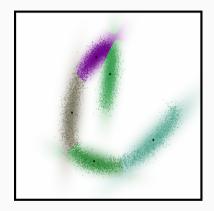
Benchmark of a Gaussian **convolution** between **clouds of N 3D points** on a RTX 2080 Ti GPU.



Applications

KeOps is a good fit for machine learning research



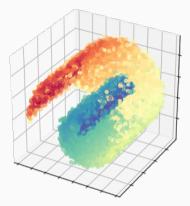


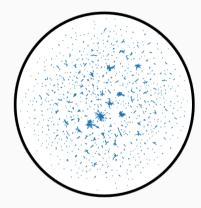
K-Means.

Gaussian Mixture Model.

Use any kernel, metric or formula you like!

KeOps is a good fit for machine learning research





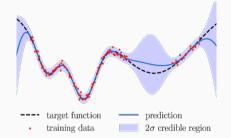
Spectral analysis.

UMAP in hyperbolic space.

Use any kernel, metric or formula you like!

Applications to Kriging, spline, Gaussian process, kernel regression

A standard tool for regression [Lec18]:



Under the hood, solve a kernel linear system:

$$(\lambda \operatorname{Id} + K_{xx}) a = b \qquad \text{i.e.} \qquad a \leftarrow (\lambda \operatorname{Id} + K_{xx})^{-1} b$$

where $\ \lambda \geqslant 0 \ \ {\rm et} \ \ (K_{xx})_{i,j} = k(x_i,x_j)$ is a positive definite matrix.

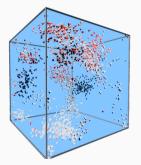
KeOps symbolic tensors $(K_{xx})_{i,j} = k(x_i, x_j)$:

- Can be fed to **standard solvers**: SciPy, GPyTorch, etc.
- GPytorch on the 3DRoad dataset (N = 278k, D = 3): $\label{eq:generalized_states} \begin{array}{c} \textbf{7h avec 8 GPUs} & \rightarrow & 15mn avec 1 GPU. \end{array}$

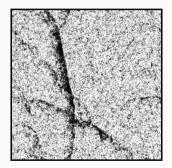
 Provide a fast backend for research codes: see e.g. Kernel methods through the roof: handling billions of points efficiently, by G. Meanti, L. Carratino, L. Rosasco, A. Rudi (2020).

KeOps lets you focus on your models, results and theorems

Some applications to **dynamical systems** [DM08, DFMAT17] and **statistics** [CDF19] with A. Diez, G. Clarté et P. Degond:



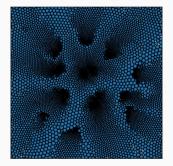
3D Vicsek model with orientation, interactive demo with 2k **flyers**.



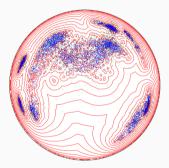
2D Vicsek model on the torus, in real-time with 100k **swimmers**.

KeOps lets you focus on your models, results and theorems

 \implies Scale up to **millions/billions** of agents with Python scripts.



Packing problem in 2D with 10k repulsive balls.



Collective Monte Carlo **sampling** on the hyperbolic Poincaré disk.

2. Optimal transport

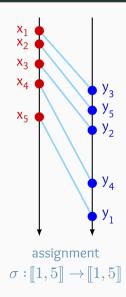
Optimal transport (OT) generalizes sorting to spaces of dimension $\mathsf{D}>1$

Context. If $A = (x_1, ..., x_N)$ and $B = (y_1, ..., y_N)$ are two clouds of N points in \mathbb{R}^D , we define:

$$\mathsf{OT}(\mathbf{A},\mathbf{B}) \;=\; \min_{\sigma\in\mathcal{S}_{\mathsf{N}}}\; \frac{1}{2\mathsf{N}}\sum_{\mathsf{i}=1}^{\mathsf{N}}\|\,\mathbf{x}_{\boldsymbol{i}}-\mathbf{y}_{\sigma(\boldsymbol{i})}\|^2$$

Generalizes **sorting** to metric spaces. **Linear problem** on the permutation matrix P:

$$\begin{aligned} \mathsf{OT}(\mathsf{A},\mathsf{B}) \ &= \ \min_{\mathsf{P}\in\mathbb{R}^{\mathsf{N}\times\mathsf{N}}} \ \frac{1}{2\mathsf{N}} \sum_{i,j=1}^{\mathsf{N}} \mathsf{P}_{i,j} \cdot \| \, \mathsf{x}_{i} - \mathsf{y}_{j} \|^{2} \,, \\ \text{s.t.} \quad \mathsf{P}_{i,j} \ &\geqslant \ \mathsf{O} \quad \underbrace{\sum_{j} \mathsf{P}_{i,j} \ = \ \mathsf{1}}_{\text{Each source point...}} \ \underbrace{\sum_{i} \mathsf{P}_{i,j} \ = \ \mathsf{1}}_{\text{is transported onto the target.}} \end{aligned}$$



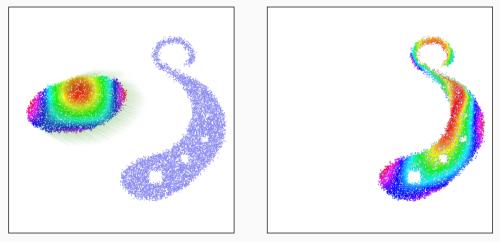
Key properties of this distance "up to permutations"

The Wasserstein distance $\sqrt{OT}(A, B)$ is:

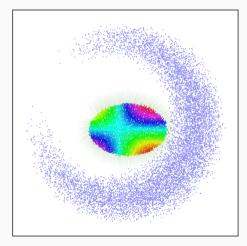
- Symmetric: OT(A, B) = OT(B, A).
- Positive: $OT(A, B) \ge 0$.
- Definite: $OT(A, B) = 0 \iff A = B$.
- Translation-aware: $OT(A, Translate_{\vec{v}}(A)) = \frac{1}{2} \| \vec{v} \|^2$.
- More generally, OT retrieves the unique gradient of a convex function ${\bf T}=\nabla\phi$ that maps A onto B:

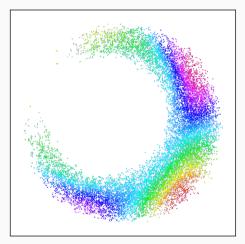
 $\begin{array}{ll} \text{In dimension 1,} & (\mathbf{x}_{i} - \mathbf{x}_{j}) \, \cdot \, (\mathbf{y}_{\sigma(i)} - \mathbf{y}_{\sigma(j)}) & \geqslant \, \mathbf{0} \\ \\ \text{In dimension D,} & \langle \, \mathbf{x}_{i} - \mathbf{x}_{j} \, \ , \, \, \mathbf{T}(\mathbf{x}_{i}) - \mathbf{T}(\mathbf{x}_{j}) \, \rangle_{\mathbb{R}^{D}} \, \geqslant \, \mathbf{0} \, . \end{array}$

 \implies Appealing generalization of an **increasing mapping**.

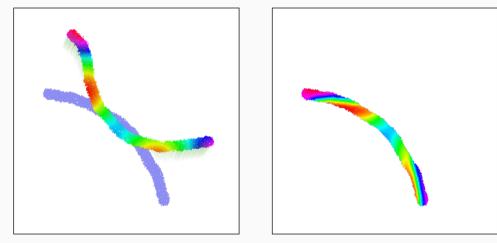




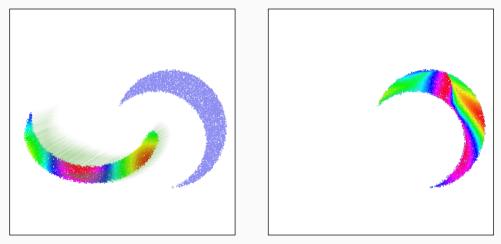




Before









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 $O({\rm N}\log{\rm N}).$
- Solution, today: Multiscale Sinkhorn algorithm, on the GPU.

 \implies Generalized **QuickSort** algorithm.

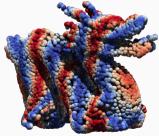
Scaling up optimal transport to anatomical data

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:

pip install geomloss + modern GPU (1000€)



10k points in 30-50ms



100k points in 100-200ms

3. Geometric deep learning

Design task-specific trainable models

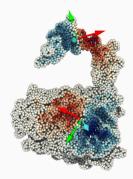
Context. Trainable models on **non-Euclidean domains** (point clouds, surfaces, graphs, etc.), beyond 2D/3D images.

Challenge. In spite of growing interest in the industry, these models still **lack support** on the numerical side. C++/CUDA is (often) required to reach top performance.

Solution. Using KeOps, with a few lines of Python:

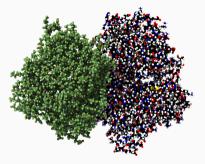
- Local interactions: K-nearest neighbors.
- Global interactions: generalized convolutions.

Modelling **freedom** \implies **Domain-specific** priors.

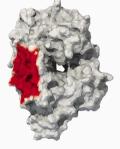


Quasi-geodesic convolution on a protein surface.

Applications to protein sciences [SFCB20]



(a) Raw protein data.

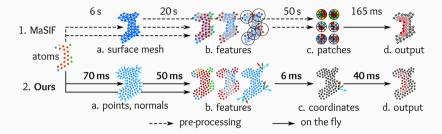


(b) Interface.



(c) Prediction.

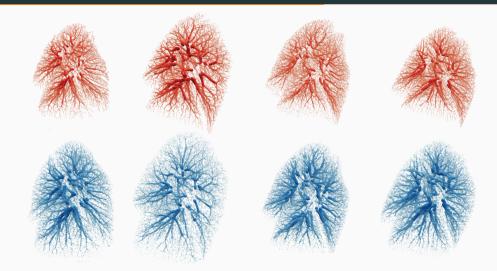
Fast end-to-end learning on protein surfaces





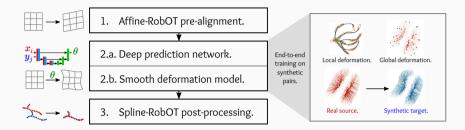
 $\times 100$ - $\times 1{,}000$ faster, lighter and fully differentiable.

Lung registration "Exhale – Inhale"



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

Three-steps registration

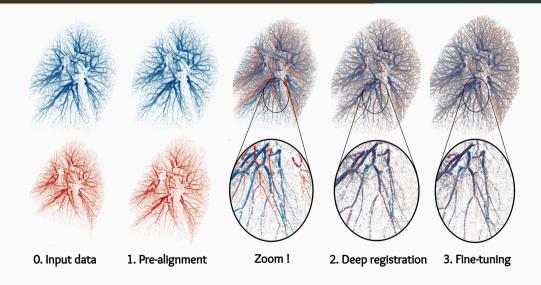


This **pragmatic** method:

- Is easy to train on synthetic data.
- Scales up to high-resolution: 100k points in 1s.
- Excellent results: KITTI (outdoors scans) and DirLab (lungs).

Accurate point cloud registration with robust optimal transport, Shen, Feydy et al., NeurIPS 2021, already on ArXiv.

Three-steps registration



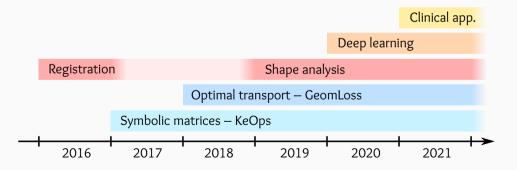
Conclusion

Key points

- Symbolic matrices are to geometric ML what
 - sparse matrices are to graph processing:
 - \longrightarrow KeOps: **x30 speed-up** vs. PyTorch, TF et JAX.
 - $\longrightarrow~$ Useful in a wide range of settings.
- Optimal Transport = generalized sorting :
 - \longrightarrow Geometric gradients.
 - \longrightarrow Super-fast $O(N \log N)$ solvers.
- These tools open **new paths** for geometers and statisticians:
 - \longrightarrow GPUs are more **versatile** than you think.

Two major evolutions:

- "Big" geometric problem: N > 10k \longrightarrow N > 1M.
- Optimal transport: linear problem + generalized quicksort.



Genuine team work



Alain Trouvé



Thibault Séjourné



F.-X. Vialard



Gabriel Peyré



Benjamin Charlier



Joan Glaunès





Freyr Sverrisson

Shen Zhengyang

+ Marc Niethammer, Bruno Correia, Michael Bronstein...

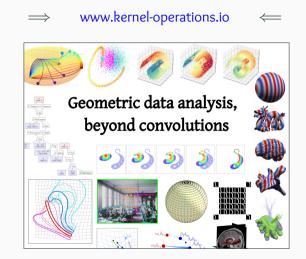
KeOps and GeomLoss are:

- Fast : 10 to 1,000 speedup vs. standard GPU implementations.
- Memory-efficient: O(N), not $O(N^2)$.
- Versatile, with a transparent interface: freedom!
- Powerful and well-documented: research-friendly.
- Slow with large variables of dimension D > 100.

Coming soon:

- ightarrow Approximation strategies (Nyström, etc.) in KeOps.
- $\rightarrow\,$ Wasserstein <code>barycenters</code> and <code>grid</code> <code>images</code> in GeomLoss.

Documentation and tutorials are available online



www.jeanfeydy.com/geometric_data_analysis.pdf

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