

Partial-Hessian Strategies for Fast Learning of Nonlinear Embeddings

Max Vladymyrov and Miguel Á. Carreira-Perpiñán



Electrical Engineering and Computer Science
University of California, Merced

<https://eecs.ucmerced.edu>

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Introduction

We focus on graph-based dimensionality reduction techniques:

- ▶ Input is a (sparse) affinity matrix.
- ▶ Objective function is a minimization over the location of the latent points.
- ▶ Examples:
 - Spectral methods: Laplacian Eigenmaps (LE), LLE;
 - ✓ have a closed-form solution;
 - ✗ results are often not satisfactory.
 - Nonlinear methods: SNE, s-SNE, t -SNE, elastic embedding (EE);
 - ✓ produce good quality embedding;
 - ✗ notoriously slow to train, limited to small data sets.

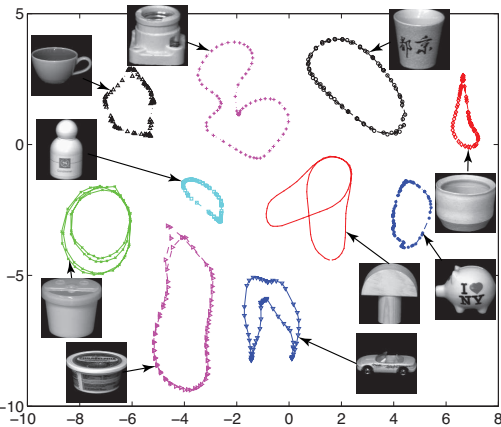
One reason for slow training is inefficient optimization algorithms that take many iterations and move very slowly towards a solution.

COIL-20 Dataset

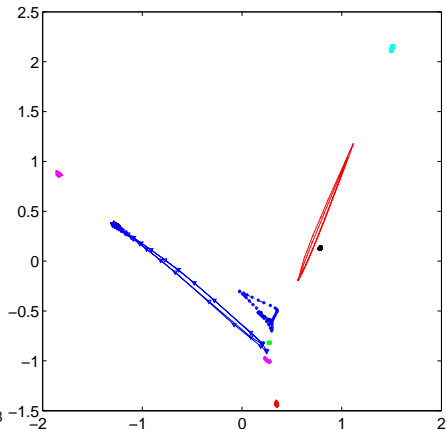
Rotations of 10 objects every 5° ; input is grayscale images of 128×128 .



Elastic Embedding



Laplacian Eigenmaps



Teaser

We are proposing a new training algorithm that:

- ▶ generalizes over multiple algorithms (s-SNE, t -SNE, EE);
- ▶ fast (1-2 orders of magnitude compared to current techniques);
- ▶ allows deep, inexpensive steps;
- ▶ scalable to larger datasets;
- ▶ intuitive and easy to implement.

General Embedding Formulation (Carreira-Perpiñán 2010)

For $\mathbf{Y} = (\mathbf{y}_1, \dots, \mathbf{y}_N) \in \mathcal{R}^{D \times N}$ matrix of high-dimensional points and $\mathbf{X} = (\mathbf{x}_1, \dots, \mathbf{x}_N) \in \mathcal{R}^{d \times N}$ matrix of low-dimensional points, define an objective function:

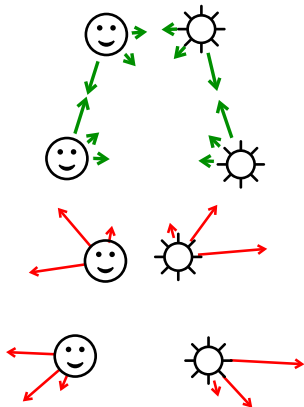
$$E(\mathbf{X}, \lambda) = E^+(\mathbf{X}) + \lambda E^-(\mathbf{X}) \quad \lambda \geq 0$$

E^+ is the *attractive term*:

- ▶ often quadratic,
- ▶ minimal with coincident points;

E^- is the *repulsive term*:

- ▶ often very nonlinear,
- ▶ minimal with points separated infinitely.



Optimal embeddings balance both forces.

Example: SNE (Hinton & Roweis 2003)

Define P_n and Q_n as distributions for each data point over the neighbors in high- and low-dimensional spaces respectively:

$$p_{nm} = \frac{\exp(-\frac{\|\mathbf{y}_n - \mathbf{y}_m\|^2}{\sigma^2})}{\sum_{k=1, k \neq n}^N \exp(-\frac{\|\mathbf{y}_n - \mathbf{y}_m\|^2}{\sigma^2})}; \quad q_{nm} = \frac{\exp(-\|\mathbf{x}_n - \mathbf{x}_m\|^2)}{\sum_{k=1, k \neq n}^N \exp(-\|\mathbf{x}_n - \mathbf{x}_m\|^2)}$$

The goal is to position points \mathbf{X} such that P_n matches the Q_n for every n :

$$\begin{aligned} E_{\text{SNE}}(\mathbf{X}) &= \sum_{n=1}^N D(P_n \| Q_n) \\ &= \sum_{n,m=1}^N p_{nm} \log \frac{p_{nm}}{q_{nm}} = - \sum_{n,m=1}^N p_{nm} \log q_{nm} + C \\ &= \sum_{n,m=1}^N p_{nm} \|\mathbf{x}_n - \mathbf{x}_m\|^2 + \sum_{n=1}^N \log \sum_{m \neq n} \exp(-\|\mathbf{x}_n - \mathbf{x}_m\|^2) + C \\ &= E^+(\mathbf{X}) + \lambda E^-(\mathbf{X}) \quad (\text{In this formulation } \lambda = 1) \end{aligned}$$

General Embedding Formulation: Other Special Cases

	$E^+(\mathbf{X})$	$E^-(\mathbf{X})$
SNE: (Hinton&Roweis,'03)	$\sum_{n,m=1}^N p_{nm} \ \mathbf{x}_n - \mathbf{x}_m\ ^2$	$\sum_{n=1}^N \log \sum_{m=1}^N e^{-\ \mathbf{x}_n - \mathbf{x}_m\ ^2}$
s-SNE: (Cook et al,'07)	$\sum_{n,m=1}^N p_{nm} \ \mathbf{x}_n - \mathbf{x}_m\ ^2$	$\log \sum_{n,m=1}^N e^{-\ \mathbf{x}_n - \mathbf{x}_m\ ^2}$
<i>t</i>-SNE: (van der Maaten & Hinton,'08)	$\sum_{n,m=1}^N p_{nm} \log (1 + \ \mathbf{x}_n - \mathbf{x}_m\ ^2)$	$\log \sum_{n,m=1}^N (1 + \ \mathbf{x}_n - \mathbf{x}_m\ ^2)^{-1}$
EE: (Carreira-Perpiñán,'10)	$\sum_{n,m=1}^N w_{nm}^+ \ \mathbf{x}_n - \mathbf{x}_m\ ^2$	$\sum_{n,m=1}^N w_{nm}^- e^{-\ \mathbf{x}_n - \mathbf{x}_m\ ^2}$
LE & LLE: (Belkin & Niyogi,'03) (Roweis & Saul,'00)	$\sum_{n,m=1}^N w_{nm}^+ \ \mathbf{x}_n - \mathbf{x}_m\ ^2$ s.t. constraints	0

w_{nm}^+ and w_{nm}^- are affinity matrices elements

Optimization Strategy

Look for a search direction \mathbf{p}_k at iteration k as a solution of a linear system $\mathbf{B}_k \mathbf{p}_k = -\mathbf{g}_k$, where \mathbf{g}_k is the current gradient and \mathbf{B}_k is a partial Hessian matrix.

$\mathbf{B}_k = \mathbf{I}$ (grad. descent) $\xrightarrow[\text{faster convergence rate}]{\text{more Hessian information}}$ $\mathbf{B}_k = \nabla^2 E$ (Newton's method)

We want \mathbf{B}_k :

- ▶ contain as much information about the Hessian as possible;
- ▶ positive definite (pd);
- ▶ fast to solve the linear system and scale up to larger N .

After \mathbf{p}_k is obtained, a line search algorithm finds the step size α for the next iteration $\mathbf{X}_{k+1} = \mathbf{X}_k + \alpha \mathbf{p}_k$. We used backtracking line search.

Structure of the Hessian of the Generalized Embedding

Given a symmetric matrix of weights \mathbf{W} , we can always define its degree matrix $\mathbf{D} = \text{diag} \left(\sum_{n=1}^N w_{nm} \right)$ and its graph Laplacian $\mathbf{L} = \mathbf{D} - \mathbf{W}$.

\mathbf{L} is positive semi-definite (psd) when entries of \mathbf{W} are non-negative.

The $Nd \times Nd$ Hessian can be written in terms of certain graph Laplacians:

$$\nabla^2 E = 4\mathbf{L} \otimes \mathbf{I}_d$$

$$+ 8\mathbf{L}^{\text{xx}}$$

$$- 16\lambda \text{vec}(\mathbf{X}\mathbf{L}^q) \text{vec}(\mathbf{X}\mathbf{L}^q)^T$$

$\mathbf{L} = \mathbf{L}^+ - \lambda\mathbf{L}^-$; $\nabla^2 E^+(\mathbf{X}) = \mathbf{L}^+ \otimes \mathbf{I}_d$
 \mathbf{L}^+ is psd and data-independent
for Gaussian kernel.

data-dependent, overall not definite,
but has psd diagonal blocks.[†]

always negative definite.[†]

[†]exact expressions for \mathbf{L}^{xx} and \mathbf{L}^q are in the paper.

Thus, there are several choices for psd parts of the Hessian:

- ▶ The best choice depends on the problem.
- ▶ We focus in particular on the one that does generally well.

The Spectral Direction (definition)

$$\nabla^2 E = 4\mathbf{L} \otimes \mathbf{I}_d + 8\mathbf{L}^{xx} - 16\lambda \text{vec}(\mathbf{X}\mathbf{L}^q) \text{vec}(\mathbf{X}\mathbf{L}^q)^T$$

\downarrow

$$\mathbf{L}^+ - \lambda \mathbf{L}^-$$

$\mathbf{B}_k = 4\mathbf{L}^+ \otimes \mathbf{I}_d$ is a convenient Hessian approximation:

- ▶ equal to the Hessian of the spectral methods: $\nabla^2 E^+(\mathbf{X})$;
- ▶ always psd \Rightarrow global convergence under mild assumptions;
- ▶ block-diagonal and has d blocks of $N \times N$ graph Laplacian $4\mathbf{L}^+$;
- ▶ **constant** for Gaussian kernel. For other kernels we can fix it at some \mathbf{X} ;
- ▶ “bends” the gradient of the nonlinear E using the curvature of the spectral E^+ ;

The Spectral Direction (computation)

We need to solve a linear system $\mathbf{B}_k \mathbf{p}_k = \mathbf{g}_k$ efficiently for every iteration (naively $\mathcal{O}(N^3 d)$).

- ▶ Cache the (also sparse) Cholesky factor of \mathbf{L}^+ in the first iteration. Now, there are just two triangular systems for each iteration.
- ▶ For scalability, we can make \mathbf{W}^+ even more sparse than it was with a κ -NN graph ($\kappa \in [1, M]$ is a user parameter). This affects only the runtime, convergence is still guaranteed.
- ▶ \mathbf{B}_k is psd \Rightarrow add small constant μ to the diagonal elements.

	Cost per iteration
Objective function	$\mathcal{O}(N^2 d)$
Gradient	$\mathcal{O}(N^2 d)$
Spectral direction	$\mathcal{O}(N \kappa d)$

This strategy adds almost no overhead when compared to the objective function and the gradient computation.

The Spectral Direction (pseudocode)

SpectralDirection(\mathbf{X}_0 , \mathbf{W}^+ , κ)

(optional) Further sparsify \mathbf{W}^+ with κ -NN graph

$\mathbf{L}^+ \leftarrow \mathbf{D}^+ - \mathbf{W}^+$

Compute graph Laplacian $\mathcal{O}(N)$

$\mathbf{R} \leftarrow \text{chol}(\mathbf{L}^+ + \mu \mathbf{I})$

compute Cholesky decomposition $\mathcal{O}(N^2\kappa)$

$k \leftarrow 1$

repeat

 Compute E_k and \mathbf{g}_k

 Objective function and the gradient $\mathcal{O}(N^2d)$

$\mathbf{p}_k \leftarrow -\mathbf{R}^{-T}(\mathbf{R}^{-1}\mathbf{g}_k)$

 Solve two triangular systems $\mathcal{O}(N\kappa d)$

$\alpha \leftarrow$ backtracking line search

$\mathbf{X}_k \leftarrow \mathbf{X}_{k-1} + \alpha \mathbf{p}_k$

$k \leftarrow k + 1$

until stop

return \mathbf{X}

Experimental Evaluation: Methods Compared

- Gradient descent (**GD**),
(Hinton&Roweis,'03)

$$\mathbf{B}_k = \mathbf{I}$$

- Diagonal methods:

- ▶ fixed-point iterations (**FP**),
(Carreira-Perpiñán,'10)

$$\mathbf{B}_k = 4\mathbf{D}^+ \otimes \mathbf{I}_d$$

- ▶ the diagonal of the Hessian (**DiagH**);

$$\mathbf{B}_k = 4\mathbf{D}^+ \otimes \mathbf{I}_d + 8\lambda\mathbf{D}^{xx}$$

- Our methods:

- ▶ spectral direction (**SD**);
- ▶ partial Hessian **SD-**,
solve linear system with conjugate gradient;

$$\mathbf{B}_k = 4\mathbf{L}^+ \otimes \mathbf{I}_d$$

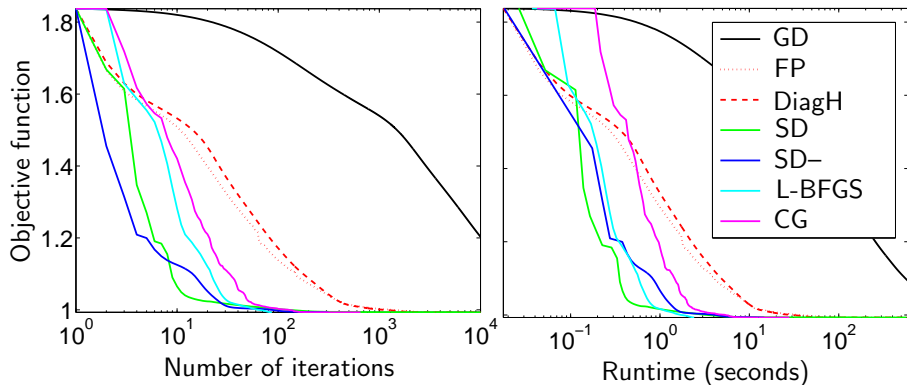
$$\mathbf{B}_k = 4\mathbf{L}^+ \otimes \mathbf{I}_d + 8\lambda\mathbf{L}_{i^*,i^*}^{xx}$$

- Standard large-scale methods:

- ▶ nonlinear Conjugate Gradient (**CG**);
- ▶ **L-BFGS**.

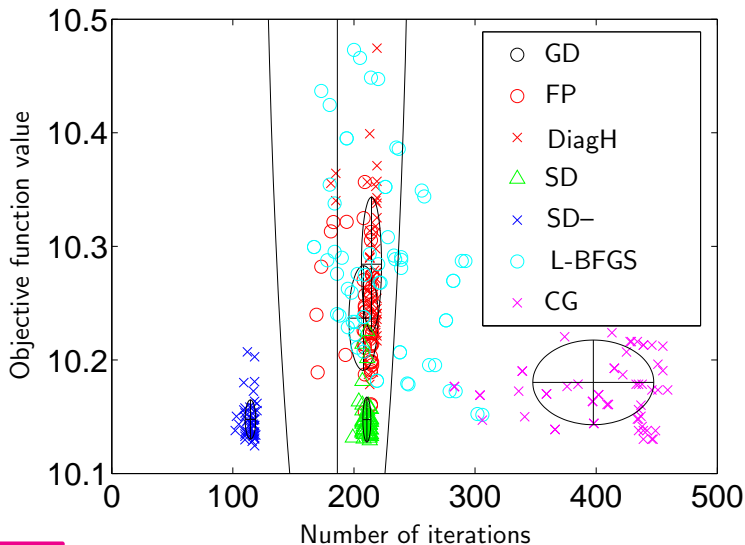
COIL-20. Convergence to the same minimum, EE

Initialize \mathbf{X}_0 close enough to \mathbf{X}_∞ so that all methods have the same initial and final points.



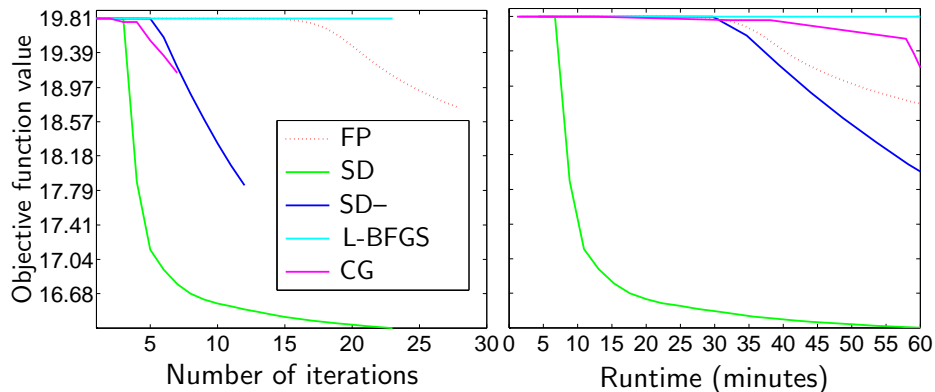
COIL-20. Convergence from random initial \mathbf{X} , s-SNE

Run the algorithms 50 times for 20 seconds each with different initialization.



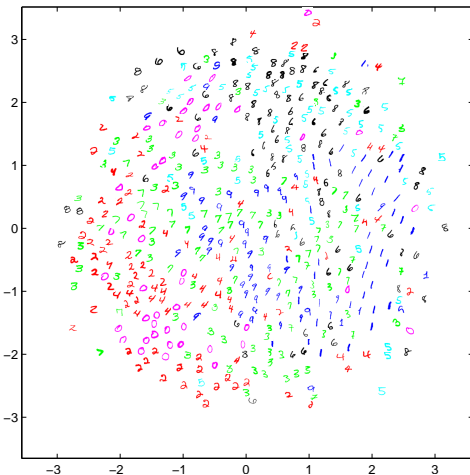
MNIST. t -SNE

$N = 20\,000$ images of handwritten digits (each a 28×28 pixel grayscale image, $D = 784$). 1 hour of optimization.

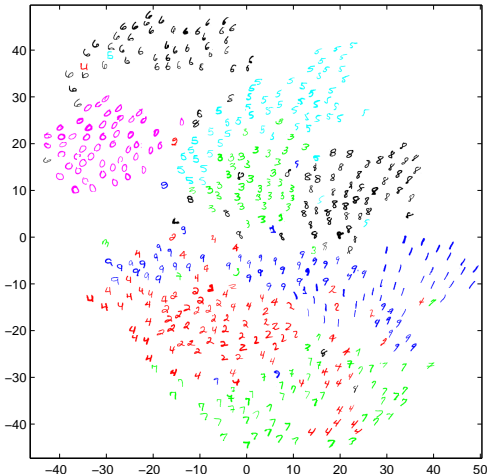


MNIST. Embedding after 1 hour of t -SNE optimization

Fixed-point iteration



Spectral direction



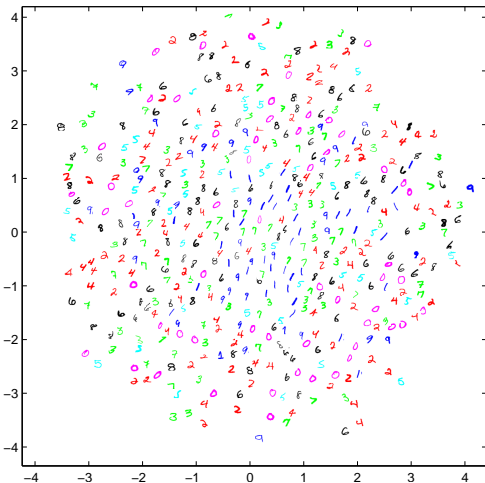
Conclusions

- ▶ We presented a common framework for many well-known dimensionality reduction techniques.
- ▶ We showed the role of graph Laplacians in the Hessian and derived several partial Hessian optimization strategies.
- ▶ We presented the **spectral direction**: a new simple, generic and scalable optimization strategy that runs one to two orders of magnitude faster compared to traditional methods.
- ▶ The evaluation of E and ∇E remains the bottleneck ($\mathcal{O}(N^2d)$) that can be addressed in the future works (e.g. with Fast Multipole Methods).
- ▶ Matlab code (very soon): <http://eecs.ucmerced.edu/>.

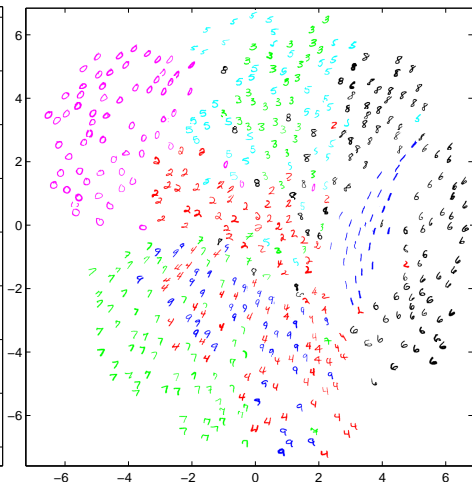
Partially supported by NSF CAREER award IIS-0754089.

MNIST. Embedding after 20 min of EE optimization

Fixed-point iteration



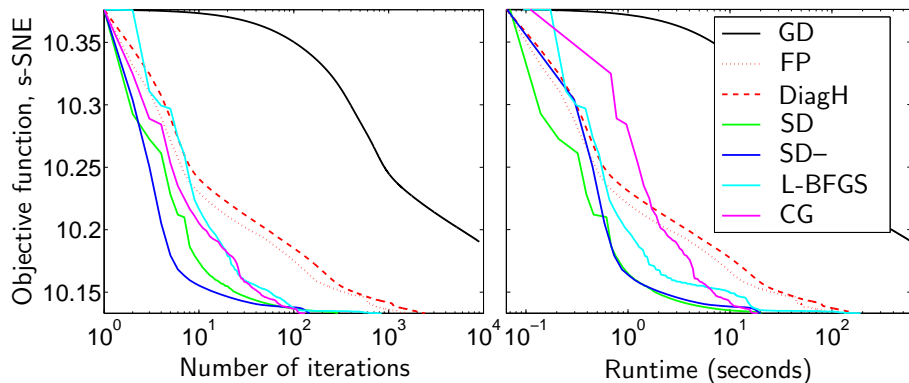
Spectral direction



Animation

COIL-20. Convergence to the same minimum, s-SNE

We initialized \mathbf{X}_0 close enough to \mathbf{X}_∞ so that all methods have the same initial and final points.



COIL-20: Homotopy optimization for EE

Start with small λ where E is convex and follow the path of minima to desired λ by minimizing over \mathbf{X} as λ increases. We used 50 log-spaced values from 10^{-4} to 10^2 .

