

Probabilistic Low-Rank Matrix Completion with Adaptive Spectral Regularization Algorithms

Adrien Todeschini[†], François Caron^{*}, Marie Chavent[†]

[†]Inria Bordeaux, ^{*}Univ. Oxford

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Summary

- 1 Introduction
- 2 Complete case
 - Rank penalty
 - Nuclear norm penalty
 - Hierarchical adaptive spectral penalty
 - EM algorithm for MAP estimation
- 3 Matrix completion
 - EM algorithm for MAP estimation
- 4 Experiments
 - Simulated data
 - Collaborative filtering examples
- 5 Conclusion and perspectives

Summary

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Matrix completion

Objective

- Complete a matrix of potentially large dimension based on a small (and potentially noisy) subset of its entries [Srebro et al., 2005, Candès and Plan, 2010].

Popular application: collaborative filtering

- To build automatic recommender systems, where the rows correspond to users, the columns to items and entries may be ratings or binaries (like/dislike).
- The objective is then to predict user preferences from a subset of the entries.
- e.g. Netflix, Amazon, Google...

Model

- Z an $m \times n$ unknown matrix of preferences
- Low rank assumption:

$$\underbrace{Z}_{m \times n} \simeq \underbrace{A}_{m \times k} \underbrace{B^T}_{k \times n}$$

with $k \ll \min(m, n)$.

- Likelihood: we typically observe a noisy version X_{ij} of some entries $(i, j) \in \Omega$ where $\Omega \subset \{1, \dots, m\} \times \{1, \dots, n\}$.

$$X_{ij} = Z_{ij} + \varepsilon_{ij}, \quad \varepsilon_{ij} \stackrel{\text{iid}}{\sim} \mathcal{N}(0, \sigma^2) \quad \forall (i, j) \in \Omega, \quad (1)$$

Further notations

- Frobenius norm:

$$\|X\|_F^2 = \sum_{(i,j)} X_{ij}^2$$

- Subset operators:

$$P_{\Omega}(X)(i,j) = \begin{cases} X_{ij} & \text{if } (i,j) \in \Omega \\ 0 & \text{otherwise} \end{cases}$$

$$P_{\Omega}^{\perp}(X)(i,j) = \begin{cases} 0 & \text{if } (i,j) \in \Omega \\ X_{ij} & \text{otherwise} \end{cases}$$

- $r = \min(m, n)$
- $X = \tilde{U}\tilde{D}\tilde{V}^T$ is the singular value decomposition (SVD) of X with $\tilde{D} = \text{diag}(\tilde{d}_1, \dots, \tilde{d}_r)$ and $\tilde{d}_1 \geq \tilde{d}_2 \geq \dots \geq \tilde{d}_r \geq 0$
- Nuclear norm: $\|X\|_* = \sum_{i=1}^r \tilde{d}_i$

Optimization problem

$$\begin{aligned} & \underset{Z}{\text{minimize}} && \text{rank}(Z) && (2) \\ & \text{subject to} && \frac{1}{2\sigma^2} \sum_{(i,j) \in \Omega} (X_{ij} - Z_{ij})^2 \leq \delta \end{aligned}$$

$$\Leftrightarrow \underset{Z}{\text{minimize}} \underbrace{\frac{1}{2\sigma^2} \|P_{\Omega}(X) - P_{\Omega}(Z)\|_F^2}_{\text{- log likelihood}} + \underbrace{\lambda \text{rank}(Z)}_{\text{penalty}} \quad (3)$$

- Rank penalty: non convex problem
- Computationally hard for general subset Ω
- Nuclear norm penalty: convex relaxation
[Fazel, 2002, Candès et al., 2008, Mazumder et al., 2010]

$$\underset{Z}{\text{minimize}} \frac{1}{2\sigma^2} \|P_{\Omega}(X) - P_{\Omega}(Z)\|_F^2 + \lambda \|Z\|_* \quad (4)$$

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Rank penalty

- Non convex problem

$$\underset{Z}{\text{minimize}} \quad \frac{1}{2\sigma^2} \|X - Z\|_F^2 + \lambda \text{rank}(Z) \quad (5)$$

- Global solution given by a hard-thresholded (truncated) SVD

$$\hat{Z} = \mathbf{H}_{\lambda\sigma^2}(X) \quad (6)$$

where $\mathbf{H}_\lambda(X) = \tilde{U}\tilde{D}^\lambda\tilde{V}^T$ with $\tilde{D}^\lambda = \text{diag}((\tilde{d}_1)_{\lambda+}, \dots, (\tilde{d}_r)_{\lambda+})$

and $t_{\lambda+} = \begin{cases} t & \text{if } t \geq \lambda \\ 0 & \text{otherwise} \end{cases}$.

Nuclear norm penalty

- Convex relaxation

$$\underset{Z}{\text{minimize}} \quad \frac{1}{2\sigma^2} \|X - Z\|_F^2 + \lambda \|Z\|_* \quad (7)$$

- Global solution given by a soft-thresholded SVD [Cai et al., 2010, Mazumder et al., 2010]

$$\hat{Z} = \mathbf{S}_{\lambda\sigma^2}(X)$$

where $\mathbf{S}_\lambda(X) = \tilde{U}\tilde{D}_\lambda\tilde{V}^T$ with $\tilde{D}_\lambda = \text{diag}((\tilde{d}_1 - \lambda)_+, \dots, (\tilde{d}_r - \lambda)_+)$ and $t_+ = \max(t, 0)$.

- The solution to (7) can be interpreted as the Maximum A Posteriori (MAP) estimate

$$\hat{Z} = \arg \max_Z [\log p(X|Z) + \log p(Z)]$$

under the likelihood (1) and prior

$$p(Z) \propto \exp(-\lambda \|Z\|_*)$$

MAP interpretaton

Assuming $Z = UDV^T$, with $D = \text{diag}(d_1, d_2, \dots, d_r)$ this can be further decomposed as

$$p(Z) = p(U)p(V)p(D)$$

where

- U and V follow a uniform Haar prior distribution on the unitary matrices
- the singular values d_i follow an exponential distribution

$$p(D) = p(d_1, \dots, d_r) = \prod_{i=1}^r \text{Exp}(d_i; \lambda) \quad (8)$$

The exponential distribution has a mode at 0, hence favoring sparse solution.

Hierarchical adaptive spectral penalty

[Todeschini et al., 2013]

- Idea: to bridge the gap between the nuclear norm and the rank penalty
- We consider the following hierarchical prior for the low rank matrix Z .

$$p(d_1, \dots, d_r | \gamma_1, \dots, \gamma_r) = \prod_{i=1}^r p(d_i | \gamma_i) = \prod_{i=1}^r \text{Exp}(d_i; \gamma_i)$$

$$p(\gamma_1, \dots, \gamma_r) = \prod_{i=1}^r p(\gamma_i) = \prod_{i=1}^r \text{Gamma}(\gamma_i; a, b)$$

- Marginal distribution over d_i :

$$p(d_i) = \int_0^{\infty} \text{Exp}(d_i; \gamma_i) \text{Gamma}(\gamma_i; a, b) d\gamma_i = \frac{ab^a}{(d_i + b)^{a+1}} \quad (9)$$

It is a Pareto distribution with heavier tails than exponential distribution

Hierarchical adaptive spectral penalty

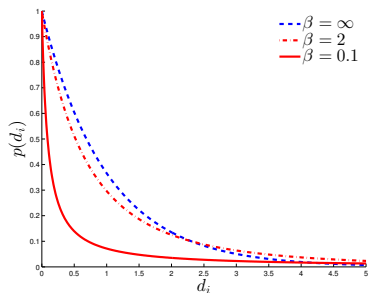
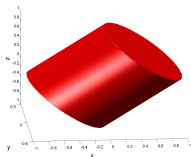


Figure : Marginal distribution $p(d_i)$ with $a = b = \beta$

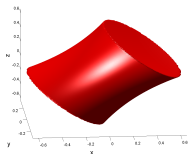
- HASP penalty: admits as special cases the nuclear norm penalty $\lambda \|Z\|_*$ when $a = \lambda b$ and $b \rightarrow \infty$.

$$\text{pen}(Z) = -\log p(Z) = -\sum_{i=1}^r \log(p(d_i)) = \sum_{i=1}^r (a+1) \log(b+d_i) \quad (10)$$

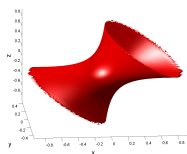
Hierarchical adaptive spectral penalty



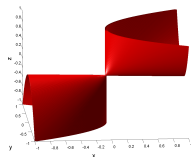
(a) Nuclear norm



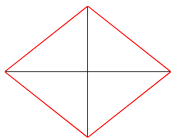
(b) HASP ($\beta = 1$)



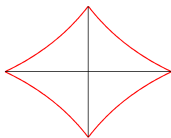
(c) HASP ($\beta = 0.1$)



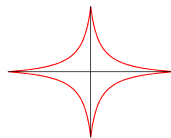
(d) Rank penalty



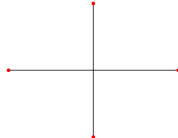
(e) ℓ_1 norm



(f) HAL ($\beta = 1$)



(g) HAL ($\beta = 0.1$)



(h) ℓ_0 norm

Figure : Top: Manifold of constant penalty, for a symmetric 2×2 matrix $Z = [x, y; y, z]$ for (a) the nuclear norm, hierarchical adaptive spectral penalty with $a = b = \beta$ (b) $\beta = 1$ and (c) $\beta = 0.1$, and (d) the rank penalty. Bottom: contour of constant penalty for a diagonal matrix $[x, 0; 0, z]$, where one recovers the classical (e) lasso, (f-g) hierarchical lasso and (h) ℓ_0 penalties.

EM algorithm for MAP estimation

We derive an Expectation Maximization (EM) algorithm to obtain a MAP estimate

$$\hat{Z} = \arg \max_Z [\log p(X|Z) + \log p(Z)]$$

i.e. to minimize

$$L(Z) = \frac{1}{2\sigma^2} \|X - Z\|_F^2 + \sum_{i=1}^r (a + 1) \log(b + d_i) \quad (11)$$

EM algorithm for MAP estimation

- Latent variables: $\gamma = (\gamma_1, \dots, \gamma_r)$
- E step:

$$\begin{aligned}Q(Z, Z^*) &= \mathbb{E} [\log(p(X, Z, \gamma)) | Z^*, X] \\ &= C - \frac{1}{2\sigma^2} \|X - Z\|_F^2 - \sum_{i=1}^r \mathbb{E}[\gamma_i | d_i^*] d_i \\ &= C - \frac{1}{2\sigma^2} \|X - Z\|_F^2 - \sum_{i=1}^r \omega_i d_i\end{aligned}$$

where $\omega_i = \mathbb{E}[\gamma_i | d_i^*] = \frac{a+1}{b+d_i^*}$.

EM algorithm for MAP estimation

- M step:

$$\underset{Z}{\text{minimize}} \quad \frac{1}{2\sigma^2} \|X - Z\|_F^2 + \sum_{i=1}^r \omega_i d_i \quad (12)$$

(12) is an adaptive spectral penalty regularized optimization problem, with weights $\omega_i = \frac{a+1}{b+d_i^*}$.

$$\begin{aligned} d_1^* &\geq d_2^* \geq \dots \geq d_r^* \\ \Rightarrow 0 &\leq \omega_1 \leq \omega_2 \leq \dots \leq \omega_r \end{aligned} \quad (13)$$

Given condition (13), the solution is given by a weighted soft-thresholded SVD [Gaïffas and Lecué, 2011]

$$\hat{Z} = \mathbf{S}_{\sigma^2\omega}(X) \quad (14)$$

where $\mathbf{S}_{\omega}(X) = \tilde{U}\tilde{D}_{\omega}\tilde{V}^T$ with $\tilde{D}_{\omega} = \text{diag}((\tilde{d}_1 - \omega_1)_+, \dots, (\tilde{d}_r - \omega_r)_+)$.

EM algorithm for MAP estimation

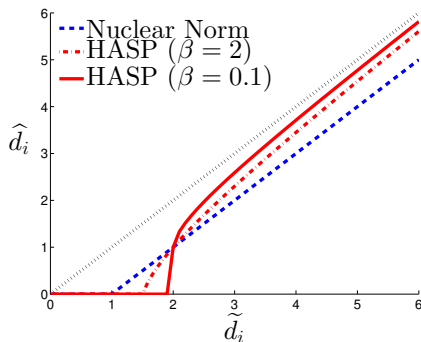


Figure : Thresholding rules on the singular values \tilde{d}_i of X

The weights will penalize less heavily higher singular values, hence reducing bias.

HAST algorithm

Hierarchical Adaptive Soft Thresholded (HAST) algorithm for low rank estimation of complete matrices

Initialize $Z^{(0)}$. At iteration $t \geq 1$

- For $i = 1, \dots, r$, compute the weights $\omega_i^{(t)} = \frac{a+1}{b+d_i^{(t-1)}}$
- Set $Z^{(t)} = \mathbf{S}_{\sigma^2 \omega^{(t)}}(X)$
- If $\frac{L(Z^{(t-1)}) - L(Z^{(t)})}{L(Z^{(t-1)})} < \varepsilon$ then return $\hat{Z} = Z^{(t)}$

This algorithm admits the soft-thresholded SVD operator as a special case when $a = b\lambda$ and $b = \beta \rightarrow \infty$.

Settings

- Parametrization:
 - ▶ We set $b = \beta$ and $a = \lambda\beta$ where λ and β are tuning parameters that can be chosen by cross-validation.
 - ▶ It is possible to estimate σ within the EM algorithm. In our experiments, we have found the results not very sensitive to the setting of σ , and set it to 1.
- Initialization:
 - ▶ As λ is the mean value of the regularization parameter γ_i , we initialize the algorithm with the soft thresholded SVD with parameter $\sigma^2\lambda$.

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Matrix completion

- Only a subset $\Omega \subset \{1, \dots, m\} \times \{1, \dots, n\}$ of the entries of the matrix X is observed.
- Relies on imputing missing values
- Assuming the same prior (9), the MAP estimate is obtained by minimizing

$$L(Z) = \frac{1}{2\sigma^2} \|P_{\Omega}(X) - P_{\Omega}(Z)\|_F^2 + (a + 1) \sum_{i=1}^r \log(b + d_i) \quad (15)$$

EM algorithm for MAP estimation

- Latent variables: γ and $P_{\Omega}^{\perp}(X)$
- E step:

$$\begin{aligned} Q(Z, Z^*) &= \mathbb{E} [\log(p(P_{\Omega}(X), P_{\Omega}^{\perp}(X), Z, \gamma)) | Z^*, P_{\Omega}(X))] \\ &= C_2 - \frac{1}{2\sigma^2} \left\{ \|P_{\Omega}(X) + P_{\Omega}^{\perp}(Z^*) - Z\|_F^2 \right\} - \sum_{i=1}^r \mathbb{E}[\gamma_i | d_i^*] d_i \end{aligned}$$

- M step:

$$\underset{Z}{\text{minimize}} \frac{1}{2\sigma^2} \|X^* - Z\|_F^2 - \sum_{i=1}^r \omega_i d_i \quad (16)$$

where $\omega_i = \mathbb{E}[\gamma_i | d_i^*]$ and $X^* = P_{\Omega}(X) + P_{\Omega}^{\perp}(Z^*)$ is the observed matrix, completed with entries in Z^* .

We now have a complete matrix problem whose solution is obtained with a weighted soft-thresholded SVD.

EM algorithm for MAP estimation

Hierarchical Adaptive Soft Impute (HASI) algorithm for matrix completion

Initialize $Z^{(0)}$. At iteration $t \geq 1$

- For $i = 1, \dots, r$, compute the weights $\omega_i^{(t)} = \frac{a+1}{b+d_i^{(t-1)}}$
- Set $Z^{(t)} = \mathbf{S}_{\sigma^2 \omega^{(t)}} (P_{\Omega}(X) + P_{\Omega}^{\perp}(Z^{(t-1)}))$
- If $\frac{L(Z^{(t-1)}) - L(Z^{(t)})}{L(Z^{(t-1)})} < \varepsilon$ then return $\hat{Z} = Z^{(t)}$

- HASI algorithm admits the Soft-Impute algorithm of [Mazumder et al., 2010] as a special case when $a = \lambda b$ and $b = \beta \rightarrow \infty$. In this case, one obtains at each iteration $\omega_i^{(t)} = \lambda$ for all i .
- On the contrary, when $\beta < \infty$, our algorithm adaptively updates the weights so that to penalize less heavily higher singular values.

Initialization

- The objective function (15) is in general not convex and different initializations may lead to different modes.
- As in the complete case, we suggest to set $a = \lambda b$ and $b = \beta$ and to initialize the algorithm with the Soft-Impute algorithm with regularization parameter $\sigma^2 \lambda$.

Scaling

- Similarly to the Soft-Impute algorithm, the computationally demanding part of HASI is $\mathbf{S}_{\sigma^2\omega^{(t)}} (P_{\Omega}(X) + P_{\Omega}^{\perp}(Z^{(t-1)}))$ which requires calculating a low rank truncated SVD.
- For large matrices, one can resort to the PROPACK algorithm [Larsen, 2004]. This sophisticated linear algebra algorithm can efficiently compute the truncated SVD of the “sparse + low rank” matrix

$$P_{\Omega}(X) + P_{\Omega}^{\perp}(Z^{(t-1)}) = \underbrace{P_{\Omega}(X) - P_{\Omega}(Z^{(t-1)})}_{\text{sparse}} + \underbrace{Z^{(t-1)}}_{\text{low rank}}$$

and can thus handle large matrices.

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Simulated data

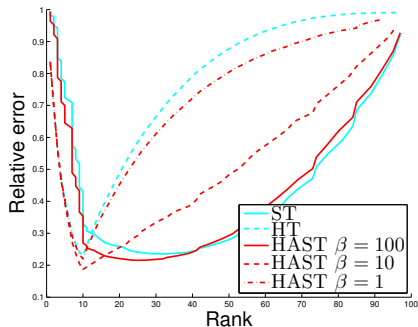
Procedure

- We generate Gaussian matrices A and B respectively of size $m \times q$ and $n \times q$, $q \leq r$ so that the matrix $Z = AB^T$ is of low rank q . A Gaussian noise of variance σ^2 is then added to the entries of Z to obtain the matrix X .
- The signal to noise ratio is defined as $\text{SNR} = \sqrt{\frac{\text{var}(Z)}{\sigma^2}}$.
- We set $m = n = 100$ and $\sigma = 1$.
- We run all the algorithms with a precision $\epsilon = 10^{-9}$ and a maximum number of $t_{\max} = 200$ iterations (initialization included for HASI).
- For the HASP penalty, we set $a = \lambda\beta$ and $b = \beta$.
- We compute the solutions over a grid of 50 values of the regularization parameter λ linearly spaced from λ_0 to 0, where $\lambda_0 = \|P_{\Omega}(X)\|_2$ is the largest singular value of the input matrix X , padded with zeros. This is done for three different values $\beta = 1, 10, 100$.
- We compute err , the relative error between the estimated matrix \hat{Z} and the true matrix Z in the complete case, and $err_{\Omega^{\perp}}$ in the incomplete case, where

$$err = \frac{\|\hat{Z} - Z\|_F^2}{\|Z\|_F^2} \quad \text{and} \quad err_{\Omega^{\perp}} = \frac{\|\hat{P}_{\Omega^{\perp}}(\hat{Z}) - P_{\Omega^{\perp}}(Z)\|_F^2}{\|P_{\Omega^{\perp}}(Z)\|_F^2}$$

Simulated data

Complete case



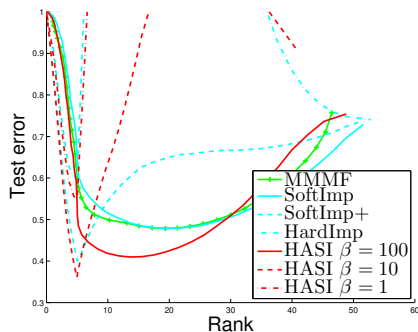
(a) SNR=1; Complete; rank=10

Figure : Test error w.r.t. the rank obtained by varying the value of the regularization parameter λ .

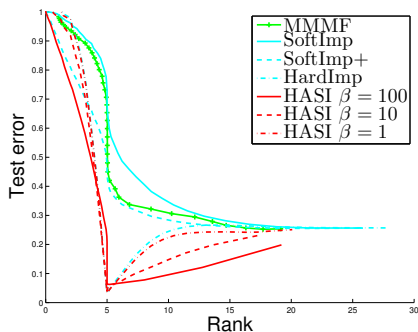
- The HASP penalty provides a bridge/tradeoff between the nuclear norm and the rank penalty.
- For example, value of $\beta = 10$ show a minimum at the true rank $q = 10$ as HT, but with a lower error when the rank is overestimated.

Simulated data

Incomplete case



(a) SNR=1; 50% missing; rank=5



(b) SNR=10; 80% missing; rank=5

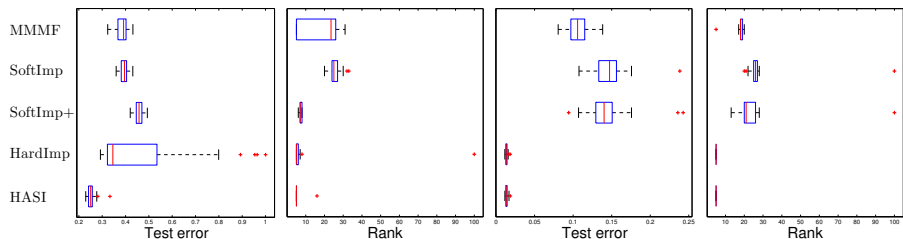
Figure : Test error w.r.t. the rank obtained by varying the value of the regularization parameter λ , averaged over 50 replications.

- Similar behavior is observed, with the HASI algorithm attaining a minimum at the true rank $q = 5$.

Simulated data

Incomplete case

We then remove 20% of the observed entries as a validation set to estimate the regularization parameters. We use the unobserved entries as a test set.



(a) SNR=1; 50% miss. (b) SNR=1; 50% miss. (c) SNR=10; 80% miss. (d) SNR=10; 80% miss.

Figure : Boxplots of the test error and ranks obtained over 50 replications.

- For 50% missing data, HASI is shown to outperform the other methods.
- For 80% missing data, HASI and Hard Impute provide the best performances.
- In both cases, it is able to recover very accurately the true rank of the matrix.

Collaborative filtering examples (Jester)

Procedure

- We randomly select two ratings per user as a test set, and two other ratings per user as a validation set to select the parameters λ and β .
- The results are computed over four values $\beta = 1000, 100, 10, 1$.
- We compare the results of the different methods with the Normalized Mean Absolute Error (NMAE)

$$\text{NMAE} = \frac{\frac{1}{\text{card}(\Omega_{\text{test}})} \sum_{(i,j) \in \Omega_{\text{test}}} |X_{ij} - \hat{Z}_{ij}|}{\max(X) - \min(X)}$$

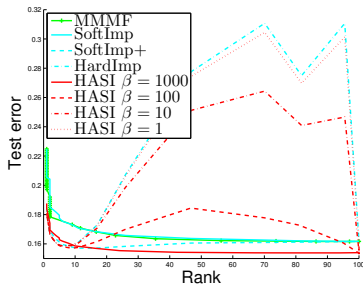
Collaborative filtering examples (Jester)

Table : Results on the Jester datasets, averaged over 10 replications

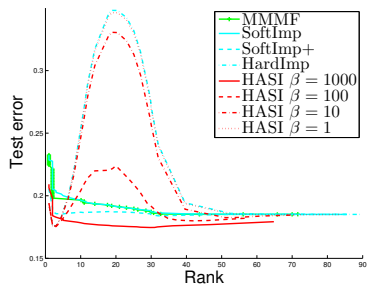
Method	Jester 1 24983 × 100 27.5% miss.		Jester 2 23500 × 100 27.3% miss.		Jester 3 24938 × 100 75.3% miss.	
	NMAE	Rank	NMAE	Rank	NMAE	Rank
MMMF	0.161	95	0.162	96	0.183	58
Soft Imp	0.161	100	0.162	100	0.184	78
Soft Imp+	0.169	14	0.171	11	0.184	33
Hard Imp	0.158	7	0.159	6	0.181	4
HASI	0.153	100	0.153	100	0.174	30

- The HASI algorithm provides very good performance on the different Jester datasets, with lower NMAE than the other methods.

Collaborative filtering examples (Jester)



(a) Jester 1



(b) Jester 3

Figure : NMAE w.r.t. the rank obtained by varying the regularization parameter λ .

- Low values of β exhibit a bimodal behavior with two modes at low rank and full rank.
- High value $\beta = 1000$ is unimodal and outperforms Soft-Impute at any particular rank.

Collaborative filtering examples (MovieLens)

Procedure

- We randomly select 20% of the entries as a test set, and the remaining entries are split between a training set (80%) and a validation set (20%).
- For all the methods, we stop the regularization path as soon as the estimated rank exceeds $r_{max} = 100$.
- For the larger MovieLens 1M dataset, the precision, maximum number of iterations and maximum rank are decreased to $\epsilon = 10^{-6}$, $t_{max} = 100$ and $r_{max} = 30$.

Collaborative filtering examples (MovieLens)

Table : Results on the MovieLens datasets, averaged over 5 replications

Method	MovieLens 100k 943 × 1682 93.7% miss.		MovieLens 1M 6040 × 3952 95.8% miss.	
	NMAE	Rank	NMAE	Rank
MMMF	0.195	50	0.169	30
Soft Imp	0.197	156	0.176	30
Soft Imp+	0.197	108	0.189	30
Hard Imp	0.190	7	0.175	8
HASI	0.187	35	0.172	27

- For the MovieLens 100k dataset, HASI provides better NMAE than the other methods with a low rank solution.
- For the MovieLens 1M dataset, MMMF provides the best NMAE at maximum rank. HASI provides the second best performances with a slightly lower rank.

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Conclusion and perspectives

- Conclusion:

- ▶ The proposed class of methods has shown to provide good results compared to several alternative low rank matrix completion methods.
- ▶ It provides a bridge between nuclear norm and rank regularization algorithms.
- ▶ Although the related optimization problem is not convex, experiments show that initializing the algorithm with the Soft-Impute algorithm of [Mazumder et al., 2010] provides very satisfactory results.

- Perspectives:

- ▶ Investigate a fully Bayesian approach and derive a Gibbs sampler or variational algorithm to approximate the posterior distribution.
- ▶ Application to larger Netflix dataset

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