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## PREPRINT August 11, 2013

#### LSSC

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#### Abstract

- 1 Overview
- 2 Theoritical Description

#### 2.1 Notations

In order to keep coherence with [3], the notations used are the same. A picture of n pixels is seen as a column vector in  $\mathbb{R}^n$ . The noisy picture is noted  $\mathbf{y}$  and the denoised one  $\mathbf{x}$ . The i-th pixel of  $\mathbf{x}$  is noted  $\mathbf{x}[i]$  and the patch centered in  $\mathbf{x}[i]$  and of size m is noted  $\mathbf{x}_i$ .

## 2.2 Learned Sparse Coding

The idea behind this method is to assume that the denoised picture is a signal that can be approximated by a sparse linear combinations of elements from a basis set. The basis set is called a dictionary  $\mathbf{D} \in \mathbb{R}^{m \times k}$  and is composed of k elements. The denoised patches are then computed from  $\mathbf{D}$  with:

$$min_{\boldsymbol{\alpha} \in \mathbb{R}^k} ||\boldsymbol{\alpha}||_p \quad s.t. \quad ||\mathbf{y}_i - \mathbf{D}\boldsymbol{\alpha}||_2^2 \leqslant \epsilon$$
 (1)

 $\mathbf{D}\alpha$  is the estimate of the denoised patch and  $||\alpha||_p$  is a regularization term that impose sparsity for  $\alpha$ .

p is usually 0 or 1. Eq. 1 becomes NP-hard to solve when p=0 and a greedy algorithm such as Orthogonal Matching Pursuit [5] can give an approximation. With p=1, the problem is convex and solved efficiently with the LARS algorithm [6]. Experimental observations [7] have shown that the learning part is better with p=1 and the recomposition part with p=0.

 $\epsilon$  can be chosen according to the value of the estimated standard deviation of the noise.

#### 2.3 Simultaneous Sparse Coding

# 3 Algorithm Description

### 3.1 Algorithm Overview

The first part consists in initializing a dictionary that will denoise roughly the picture.

Once the picture is denoised a first time, a clustering is made in order to regroup similar patches for further treatment.

Then, iteratively for each clusters, the dictionary is updated using simultaneous sparse coding and the cluster is denoised.

## 3.2 Dictionnary Initialization

The initial dictionnary is first learned offline on the 10 000 images of the PASCAL VOC'07 database using the online dictionary learning procedure of [2]. This procedure is then used on the noisy picture in order to improve the dictionary efficiency.

In fact, only a fixed number T of patches in the picture are used to update the dictionary. However they are chosen so that they are independently and identically distributed in the picture.

The algorithm corresponds then to the minimization of eq.1 on the T patches with the  $l_1$  norm using the LARS [6] algorithm.

**Input**: number of iterations T, i.i.d. sampling of T patches of the noisy picture  $Y_T$ , initial dictionary  $\mathbf{D}^0 \in \mathbb{R}^{m \times k}$ , regularization parameter  $\lambda$ 

Output: learned dictionary D

Initialization:  $\mathbf{A}^0 \in \mathbb{R}^{k \times k} \leftarrow 0, \, \mathbf{B}^0 \in \mathbb{R}^{m \times k} \leftarrow 0$ 

for t = 1..T do

$$\mathbf{y}_t = \mathbf{Y}_T[t]$$

Sparse coding: compute with LARS algorithm:

$$\alpha^t = \operatorname{argmin}_{\alpha \in \mathbb{R}^k} ||\alpha||_1 \text{ s.t. } ||\mathbf{y}_t - \mathbf{D}^{t-1}\alpha||_2^2 \le \lambda$$

$$\mathbf{A}^t \leftarrow \mathbf{A}^{t-1} + \boldsymbol{\alpha}^t \boldsymbol{\alpha}^{tT} \\ \mathbf{B}^t \leftarrow \mathbf{B}^{t-1} + \mathbf{y}_t^t \boldsymbol{\alpha}^{tT}$$

$$\mathbf{B}^t \leftarrow \mathbf{B}^{t-1} + \mathbf{v}_{\perp}^t \boldsymbol{\alpha}^{tT}$$

Update dictionary from  $\mathbf{D}^{t-1}$  to  $\mathbf{D}^t$  so that:

$$\mathbf{D}^{t} = \operatorname{argmin}_{\mathbf{D} \in \mathbb{R}^{m \times k}} \frac{1}{t} \sum_{i=1}^{t} (\frac{1}{2} |\mathbf{y}_{t}^{i} - \mathbf{D} \boldsymbol{\alpha}^{i}||_{2}^{2} + \lambda ||\boldsymbol{\alpha}^{i}||_{1})$$

end return  $\mathbf{D}^T$ 

**Algorithm 1**: Online Dictionary Learning

Input :input dictionary  $\mathbf{D} = [\mathbf{d}^1, \dots, \mathbf{d}^k] \in \mathbb{R}^{m \times k}$ 

 $\mathbf{A} = [\mathbf{a}^1, \dots, \mathbf{a}^k] \in \mathbb{R}^{k \times k}, \, \mathbf{B} = [\mathbf{b}^1, \dots, \mathbf{b}^k] \in \mathbb{R}^{m \times k}$ 

Output: updated dictionary D

repeat

for 
$$j = 1..k$$
 do update the  $j^{th}$  column: if  $\mathbf{A}(j,j) = 0$  then 
$$\begin{vmatrix} \mathbf{d}^j \leftarrow 0 \\ \mathbf{e} \mathbf{n} \mathbf{d} \\ \mathbf{e} \mathbf{l} \mathbf{s} \mathbf{e} \mathbf{d} \end{vmatrix} = \mathbf{d}^j \leftarrow \mathbf{d}^j + \mathbf{d}^j \mathbf{d}^j \leftarrow \mathbf{d}^j + \mathbf{d}^j$$

until convergence Marc: apparemment, dans le code on a une boucle sur params.updateIteration = 1. Est-ce qu'il ne vaudrait mieux pas calculer l'argmin à chaque boucle et s'arrêter lorsque c'est plus petit qu'une certaine valeur?;

return D Algorithm 2: Dictionary Update updateDictionary Marc: Algo validé

```
Input: Gram matrix of the dictionary G = \mathbf{D}^T \mathbf{D} \in \mathbb{R}^{k \times k} where \mathbf{D} = [\mathbf{d}^1, \dots, \mathbf{d}^k] \in \mathbb{R}^{m \times k},
noisy patch \mathbf{y}, constraint \lambda
 Output : sparse vector \boldsymbol{\alpha} \in \mathbb{R}^k
 —INITIALIZATION—
Marc : \mathcal{A} n'est pas définie, et \mathbf{D}_{\mathcal{A}} non plus.
normX \in \mathbb{R} \leftarrow ||\mathbf{y}||_2^2
coeffs \in \mathbb{R}^k \leftarrow \mathbf{0}
 Most correlated element
\hat{\mathbf{c}} \in \mathbb{R}^k \leftarrow \mathbf{D}^T \mathbf{y}
C \in \mathbb{R}^+ \leftarrow \max_{i=1..m}(|\hat{\mathbf{c}}_i|)
currentInd \leftarrow j s.t. \hat{\mathbf{c}}_i = C
if normX > \lambda \ Marc : pourquoi > et \ pas < ? then
   return 0
end
newAtom \leftarrow True
for i = 1..k do
                  --NEW ATOM---
             if newAtom then
                           \mathcal{A}[i] \leftarrow \text{currentInd}
                           G_A[i^{th} \text{ line}] \leftarrow G[\mathcal{A}[i]^{th} \text{ line}]
                           G_S \leftarrow \boldsymbol{D}_{\mathcal{A}}^T \boldsymbol{D}_{\mathcal{A}}
                         UPDATE G_S^-
              end
              —VARIABLES UPDATES—
             \mathbf{u} \leftarrow G_S^{-1}(\operatorname{sgn}(\hat{\mathbf{c}}_j))_{j \in \mathcal{A}} \operatorname{Marc} : \operatorname{qu'est-ce} \operatorname{que} \operatorname{que} \operatorname{que} \operatorname{que} : \operatorname{qu'est-ce} \operatorname{que} \operatorname{que} : \operatorname{qu'est-ce} \operatorname{que} : \operatorname{qu'est-ce} \operatorname{que} : \operatorname{qu'est-ce} : \operatorname{qu'est-ce}
            ratio \leftarrow \left(-\frac{\text{coeffs}[j]}{u_j}\right)
              stepMAX \leftarrow min^+(ratio) (min^+ means its the minimum between positive values)
              criticalInd ← j s.t. ratio[j] = stepMAX Marc : criticalInd n'est jamais plus utilisé par la
              suite??!!
              C \leftarrow \hat{\mathbf{c}}[0] \text{ Marc} : \text{ ca ne devrait pas être } C = \max \hat{\mathbf{c}}[j] ?
             \gamma \leftarrow \min^+ \left( \frac{C \pm \hat{\mathbf{c}}_j}{1 \pm (G_A u)[j]} \right)_{j \notin \mathcal{A}} \text{Marc} : \text{ça fait 2 possibilités ou bien 4 avec les } \pm ?
              —POLYNOMIAL REŚOLUTION—
             \begin{aligned} a &\leftarrow \sum_{j \in \mathcal{A}} \operatorname{sgn}(\hat{\mathbf{c}}[\mathcal{A}[j]]) u[j] \\ b &\leftarrow \sum_{j \in \mathcal{A}} \hat{\mathbf{c}}[\mathcal{A}[j]] u[j] \\ c &\leftarrow \operatorname{normX} - \lambda \end{aligned}
              \Delta \leftarrow b^2 - ac
             \mathrm{stepMAX2} \leftarrow \min(\tfrac{b-\sqrt{\Delta}}{c}, C)
              —FINAL STEP & BŘEAK-
              \gamma \leftarrow \min(\gamma, \text{stepMAX}, \text{stepMAX2})
              coeffs \leftarrow coeffs + \gamma u Marc : a priori coeffs et u ne sont pas de la même taille
              \hat{\mathbf{c}} \leftarrow \hat{\mathbf{c}} - \gamma G_A u
              normX \leftarrow normX + a\gamma^2 - 2b\gamma
              if |\gamma| < 1e^{-15} || \gamma = stepMAX2 || normX < 1e^{-15} || normX - \lambda < 1e^{-15} then
                        break Marc: ici on travaille en float, donc je pense qu'on peut mettre 1e^{-6}
              end
             if \gamma = stepMAX then
                           DOWNDATE G_s^{-1}
                           newAtom \leftarrow False
                         i \leftarrow i-2 Marc c'est possible que i devienne négatif? Ça risque de poser problème non?
              else
                                                                                                                                                                                5
                \perp newAtom \leftarrow True
             end
end
```

ant/oraffa A) (gented coefficient with respect to the increasing order of A

```
Input: Gram matrix G_S \in \mathbb{R}^{i \times i}, and its former inverse to update G_S^{-1} \in \mathbb{R}^{i-1 \times i-1}
    Output: updated G_S^{-1} \in \mathbb{R}^{i \times i}
    if i = 1 then
     return \frac{1}{G_S}
    end
    u \leftarrow G_S^{-1}G_S[i^{th} \text{ line}]

\sigma \leftarrow \frac{1}{G_s(i,i) - u.G_S[i^{th} \text{ line}]} 

G_s^{-1}(i,i) \leftarrow \sigma 

G_s^{-1}[i^{th} \text{ line}] \leftarrow -\sigma u

return G_s^{-1} \leftarrow G_s^{-1} + \sigma u u^T

Algorithm 4: Update invert algorithm updateGram Marc: Algo correspondent, mais à valider.
    Input: pseudo-Gram matrix G_A \in \mathbb{R}^{i \times i} Gram matrix G_S \in \mathbb{R}^{i \times i}, and its inverse G_S^{-1} \in \mathbb{R}^{i \times i},
    active indexes list A \in \mathbb{R}^i, sparse coefficient list coeffs \in \mathbb{R}^k, criticalInd \in [1; k], current
    iteration i
    Output: downdated matrices G_S, G_S^{-1}, G_A \in \mathbb{R}^{i-1 \times i-1}, downdated lists A \in \mathbb{R}^{i-1}, coeffs \in \mathbb{R}^k
    \sigma \leftarrow \frac{1}{G_S^{-1}(\text{criticalInd},\text{criticalInd})}
    u \leftarrow G_S^{-1}[criticalInd<sup>th</sup> line] without its criticalInd<sup>th</sup> coefficient
    for j = criticalInd:i do
         G_A[j^{th} \text{ line}] \leftarrow G_A[(j+1)^{th} \text{ line}]
          for k=1:criticalInd-1 do
               G_S(j,k) \leftarrow G_S(j+1,k)
               G_s^{-1}(j,k) \leftarrow G_s^{-1}(j+1,k)
          end
          for k = criticalInd:i do
               G_S(j,k) \leftarrow G_S(j+1,k+1)
               G_S^{-1}(j,k) \leftarrow G_S^{-1}(j+1,k+1)
         \mathcal{A}[j] \ \leftarrow \ \mathcal{A}[j+1]
         coeffs[j] \leftarrow coeffs[j+1]
    end
    coeffs[i] \leftarrow 0
 G_s^{-1} \leftarrow G_s^{-1} - \sigma u u^T

Algorithm 5: Downdate invert algorithm downdateGram Marc: Algo ne correspondant pas, j'ai
```

Algorithm 5: Downdate invert algorithm downdateGram Marc: Algo ne correspondant pas, j'ai créé une nouvelle fonction downdateGramBis dont je pense est une meilleure version. Il faudra la tester et la valider.

## Glossary

## **Image Credits**

POL (there's no need to credit this image, here is used as an example.)

## 4 References

#### References

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