

This is just a quick document for describing the differences between vanilla gradient descent, iterative thresholding and AMP.

1 Gradient Descent

Suppose we observe a vector $y \in \mathbb{R}^m$, through an operator $A \in \mathbb{R}^{m \times n}$, i.e. we have the model

$$y = Ax \quad (1.0.1)$$

$x \in \mathbb{R}^n$. One way to guess x , is to minimise the mean squared error:

$$J = \frac{1}{2} \|y - Ax\|_2^2 = (y - Ax)^T (y - Ax) \quad (1.0.2)$$

We can do this by gradient descent, as

$$\frac{\partial J}{\partial x} = A^T (y - Ax) \quad (1.0.3)$$

So we have an iterative algorithm:

$$x^{t+1} = x^t + A^T (y - Ax^t)$$

Putting it in a more suggestive form:

$$\begin{aligned} x^{t+1} &= x^t + A^T z^t \\ z^t &= y - Ax^t \end{aligned}$$

2 Iterative Thresholding

The algorithm from the previous section will perform poorly, as we've not used any prior information about x . We assume that x is sparse, so we now seek solutions to:

$$J = \frac{1}{2} \|y - Ax\|_2^2 + \lambda \|x\|_1 \quad (2.0.4)$$

The presence of the ℓ_1 norm, is inconvenient: J in this case has no formal gradient. In this case we can pretend that J is differentiable by doing a step in the direction of the gradient, and correcting for any error using the soft thresholding operator: $S_\gamma(x)_i = \text{sign}(x_i) (|x_i| - \gamma)^+$.

So our algorithm becomes:

$$\begin{aligned} x^{t+1} &= S_\lambda (x^t + A^T z^t) \\ z^t &= y - Ax^t \end{aligned}$$

The soft-thresholding operator can be derived by considering the MAP estimate of the following model:

$$y = x + w \quad (2.0.5)$$

where x is some (sparse) signal, and w is additive white Gaussian noise. We seek

$$\hat{x} = \arg \max_x \mathbb{P}_{x|y}(x|y) \quad (2.0.6)$$

This can be recast in the following form by using Bayes rule, noting that the denominator is independent of x and taking logarithms:

$$\hat{x} = \arg \max_x [\log \mathbb{P}_w(y - x) + \log \mathbb{P}(x)] \quad (2.0.7)$$

The term $\mathbb{P}_n(y - x)$ arises because we are considering $x + w$ with w zero mean Gaussian, with variance σ_n^2 . So, the conditional distribution of y (given x) will be a Gaussian centred at x .

We will take $\mathbb{P}(x)$ to be a Laplacian distribution:

$$\mathbb{P}(x) = \frac{1}{\sqrt{2}\sigma} \exp -\frac{\sqrt{2}}{\sigma}|x| \quad (2.0.8)$$

Note that $f(x) = \log \mathbb{P}_x(x) - \frac{\sqrt{2}}{\sigma}|x|$, and so by differentiating $f'(x) = -\frac{\sqrt{2}}{\sigma}\text{sign}(x)$
Taking the maximum of 2.0.7 we obtain:

$$\frac{y - \hat{x}}{\sigma_n^2} - \frac{\sqrt{2}}{\sigma}\text{sign}(x) = 0 \quad (2.0.9)$$

Which leads the soft thresholding operation defined earlier, with $\gamma = \frac{\sqrt{2}\sigma_n^2}{\sigma}$ as (via rearrangement):

$$y = \hat{x} + \frac{\sqrt{2}\sigma_n^2}{\sigma}\text{sign}(x)$$

or

$$\hat{x}(y) = \text{sign}(y) \left(y - \frac{\sqrt{2}\sigma_n^2}{\sigma} \right)_+$$

i.e $S_\gamma(y)$.

3 AMP

Approximate message passing further corrects this idea of pretending our optimisation objective is differentiable and correcting, by making a quadratic approximation to the likelihood of the model.

The algorithm becomes

$$\begin{aligned} x^{t+1} &= S_\lambda(x^t + A^T z^t) \\ z^t &= y - Ax^t + b_t z^t \end{aligned}$$

which is similar to the iterative thresholding algorithm, but with an additional 'momentum' term added. A good choice of b is:

$$\frac{1}{m} \|x^t\|_0 \quad (3.0.10)$$

where $\|t\|_0$ is the number of non-zero elements of t .

4 DISTA

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1: procedure DADMM( $y_j, M_j, \varepsilon$ )
2:    $x^0 = 0, z^0 = 0, \theta^0 = 0, \eta^0 = 0,$ 
3:    $Q = (M_j^T M_j + (\rho D_J + 1)I)^{-1}, w_j = M_j^T y_j$ 
4:   while  $\|z^{k+1} - z^k\| \leq \varepsilon$  do
5:     for  $c = 1, \dots, C$  do
6:        $x^{k+1} \leftarrow Q(w_j + z^k - \theta^{kT} - \nu^{kT})$ 
7:        $z^{k+1} \leftarrow S_{\beta/\rho}(x_j^{k+1})$ 
8:        $\theta^{k+1} \leftarrow \theta_j^k + \rho(x^{k+1} - z^{k+1})$ 
9:     end for
10:  Each node transmits  $x^{k+1}$  in  $\mathcal{N}_j$  and calculates
11:     $\nu_j^{k+1} \leftarrow \nu_j^k + \rho(\sum_{m \in \mathcal{N}_j} z_m^k - z_j^k)$ 
12:  end while
13:  return  $z^{k+1}$ 
14: end procedure

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Figure 4.1: DADMM at Node j

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1: procedure DAMP( $y_j, M_j, \varepsilon$ )
2:    $x^0 = 0, \alpha,$ 
3:
4:   while  $\|x_j^{k+1} - x_j^k\| \leq \varepsilon$  do
5:  Each node transmits  $x^k$  in  $\mathcal{N}_j$  and calculates
6:     $x_j^{k+1/2} \leftarrow \frac{1}{D_p+1} \sum_{j \in \mathcal{N}_j} x_j^k$ 
7:    for  $c = 1, \dots, C$  do
8:       $x^{k+1} \leftarrow x_j^{k+1/2} - \alpha S_\beta(M_j^T x_j^{k+1/2})$ 
9:    end for
10:  end while
11:  return  $x^{k+1}$ 
12: end procedure

```

Figure 4.2: DAMP at Node j

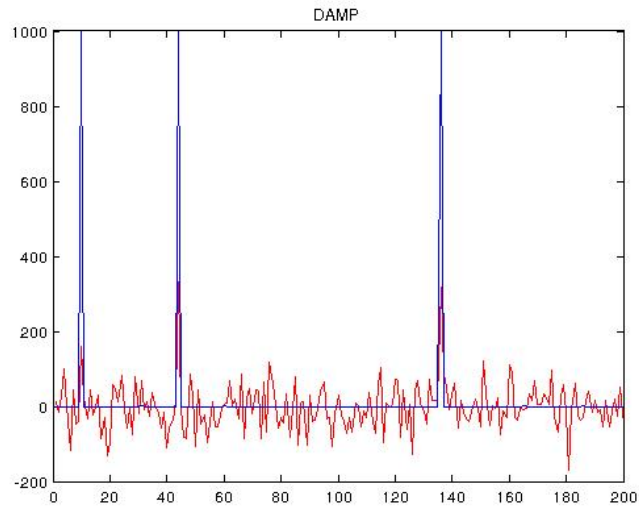


Figure 4.3: DAMP estimate after 1000 iterations

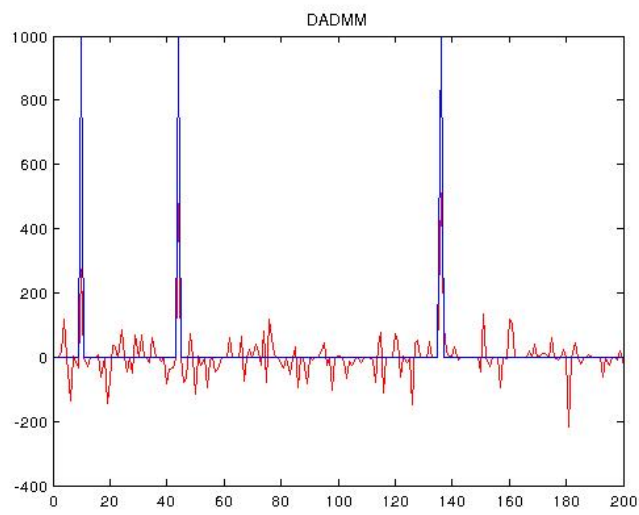


Figure 4.4: DADMM estimate after 1000 iterations

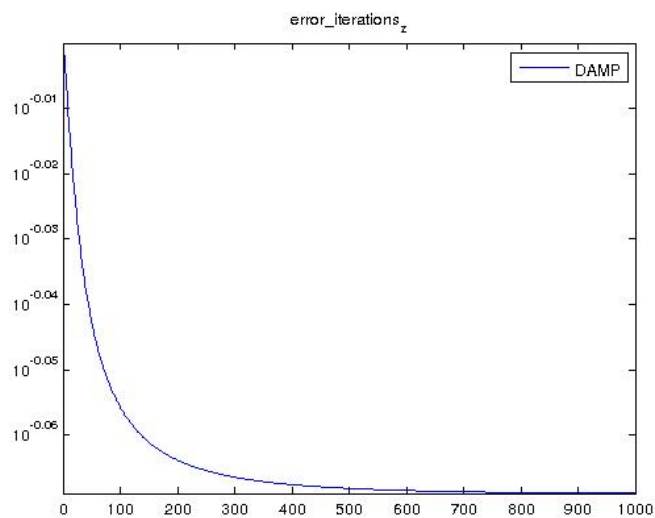


Figure 4.5: DAMP progress

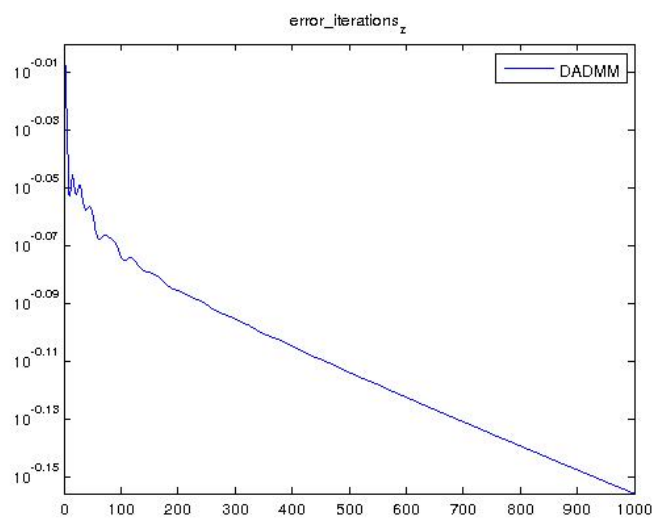


Figure 4.6: DADMM progress

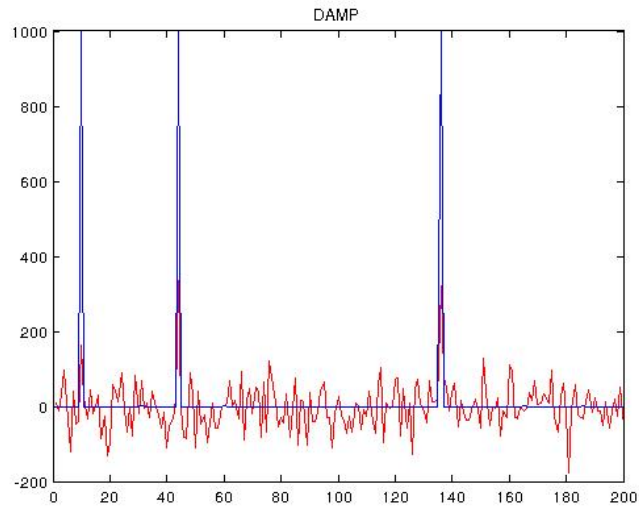


Figure 4.7: DAMP estimate after 5000 iterations

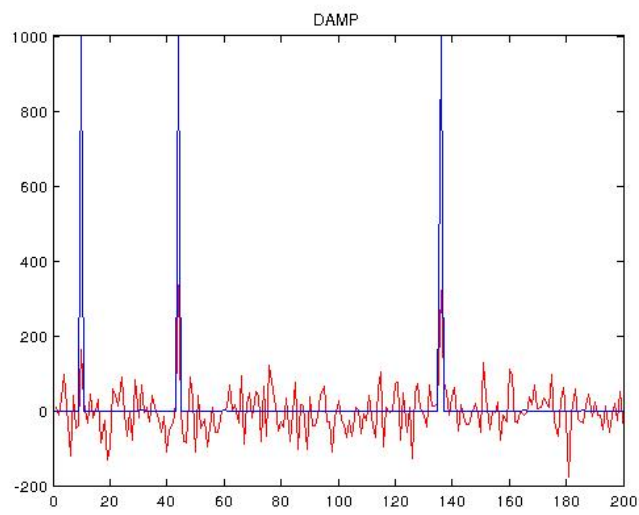


Figure 4.8: DADMM estimate after 5000 iterations

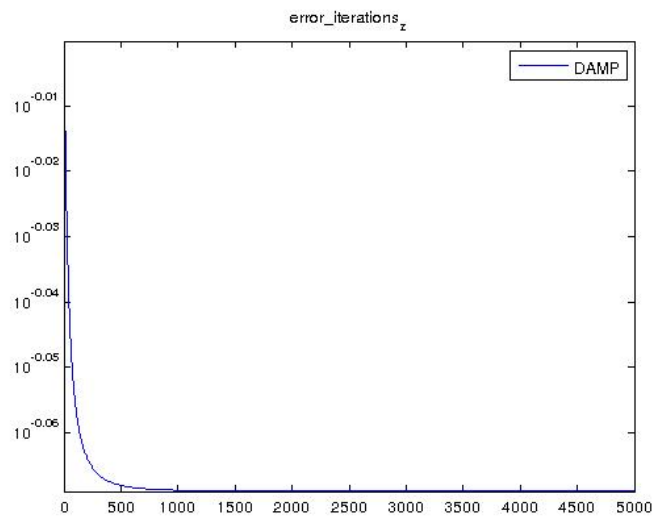


Figure 4.9: DAMP progress

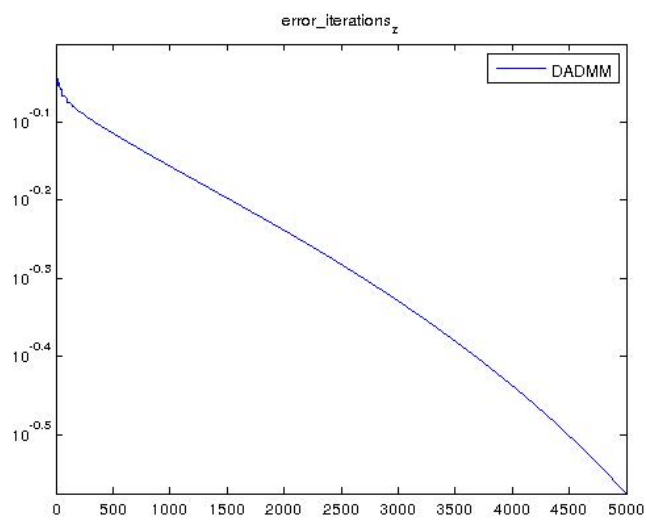


Figure 4.10: DADMM progress