# Three $l_1$ based nonconvex methods in constructing sparse mean reverting portfolios

Dedicated to the memory of our good friend and colleague Ernie Esser

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Abstract We study the problem of constructing sparse and fast mean reverting portfolios. The problem can be motivated by convergence trading and formulated as a generalized eigenvalue problem with a cardinality constraint [6]. We use a proxy of mean reversion coefficient, the direct Ornstein-Uhlenbeck (OU) estimator, which can be applied to both stationary and nonstationary data. In addition, we introduce three different methods to enforce the sparsity of the solutions. One method uses the ratio of  $l_1$  and  $l_2$  norms and the other two use  $l_1$  norm. We analyze various formulations of the resulting nonconvex optimization problems and develop efficient algorithms to solve them for portfolio sizes as large as hundreds. By adopting a simple convergence trading strategy, we test the performance of our sparse mean reverting portfolios on both synthetic and historical real market data. In particular, the  $l_1$  regularization method, in combination with quadratic program formulation as well as

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difference of convex function and least angle regression treatment, gives fast and robust on large out-of-sample data set.

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#### 1 Introduction

Convergence trade is a trade designed to benefit from the phenomenon that the difference of the prices of two assets may fluctuate around a certain level. The deviations from this level are temporary and thus investors can build an appropriate trading strategy when they observe the deviations and expect to profit by the amount of convergence. Mean reversion rate is a well-known indicator of predictability in financial markets. Constructing a fast mean reverting portfolio from a set of assets is a challenging problem. Classical methods include cointegration [9] and canonical correlation analysis [13].

Cointegration is a statistical property of a set of time series. It tests whether we could build a stationary time series as a linear combination of a set of nonstationary time series. In convergence trade, investors want the values of the portfolio to be stationary since it is easy to determine trading signals for the trading strategy. Canonical correlation analysis aims at maximizing the correlation of the linear combinations on two sets of random vectors. This can help investors find portfolios that share a common stochastic drift. Their difference will be a desired mean reverting portfolio.

The risk of convergence trade is that the expected convergence does not happen, or that it takes too long such that it possibly diverges before converging. Therefore, it is important to quantify how fast the portfolio will converge and how an optimal portfolio could be constructed based on this criterion. In addition, a sparse portfolio may be preferred in convergence trade since sparsity typically means less transaction costs. However, there will be a trade-off between the sparsity and the convergence rate.

An optimization framework for constructing sparse mean reverting portfolios is proposed in [6]. The author uses the idea of predictability [2] as a proxy for the rate of mean reversion. By presetting the desired cardinality, an optimization problem is formulated. In [5], the authors replace the cardinality constraint by the variance constraint in order to improve the profits during arbitrage opportunities, while in [11], the authors discuss more details about parameter estimation and trading strategies. In [27], the authors propose a method called DC-PCA which solves for sparse generalized eigenvectors using the difference of convex algorithms. However, the authors only test their algorithm for principal component analysis. Based on [6], the sparse optimization problem can be formulated as

$$\min_{x} x^{T} A x / x^{T} B x$$

$$s.t. \quad ||x||_{0} \le k$$

$$||x||_{2} = 1,$$
(1)

where  $||x||_0$  is the number of the non-zero entries of x and A and B are both positive definite matrices.

Without the cardinality constraint, the problem is a generalized eigenvalue problem. The goal is to find the unit vector x, such that

$$Bx = \lambda Ax$$

where  $\lambda$  has the largest possible value. When the matrix A is an identity matrix, then the problem reduces to finding the first principal component. Sparse principal component analysis problems are well-studied, see for instance [7,17,18,26,33]. In [26], the authors discuss 8 different formulations of the sparse PCA problem and introduce alternating maximization (AM) method to solve them efficiently.

In our work, we take the approach set forth in [6] as a starting point and develop three optimization methods that use  $l_1$  based norms to enforce the sparsity of the portfolio.

The first method extends the AM method to the generalized eigenvalue problem. We uses the  $l_1$  constraint to enforce the sparsity. The problem formulation is

$$\max_{x} x^{T} B x$$

$$s.t. \quad ||x||_{1} \le \sqrt{s}$$

$$x^{T} A x \le 1,$$

The second method uses the ratio of  $l_1$  and  $l_2$  norms as a penalty function assuming limited prior information of the assets. Such a penalty term arises in nonnegative matrix factorization (NMF), blind deconvolution, and sparse representation in coherent dictionaries [10,14,16,19,31]. For example, the nonnegative least squares (NNLS) problem under such penalty takes the following form [10]:

$$\min_{x \ge 0} \|\mathbf{X}x - \mathbf{Y}\|_2^2 + \gamma P(x)$$

where  $P(x) = \frac{\|x\|_1}{\|x\|_2}$ , **X** is an  $m \times n$  matrix and **Y** is an  $m \times 1$  vector. For our problem, we formulate the following minimization:

$$\min_{x \neq 0} \frac{x^T A x}{x^T B x} + \gamma \frac{\|x\|_1}{\|x\|_2} \tag{2}$$

where A and B are positive definite matrices and  $\gamma$  is a nonnegative tuning parameter. The first term of the objective function of (2) models predictability and the second term promotes sparsity. The details are in section 5. The  $l_1$ 

norm regularization technique for variational problems under nonconvex (orthogonality/unit ball) constraints has been actively developed lately to construct compactly supported eigen-functions [1,20,24,25]. In fact (1.2) is simply imposing  $l_1$  regularization on the unit ball in  $l_2$ .

Due to the nonconvexity of (2), finding a global minimum is challenging. To deal with this aspect of global optimization, we shall incorporate a recent variant of simulated annealing, the so called intermittent diffusion method with discontinuous diffusion coefficient [4]. The combined local minimization of (2) and random search for global minimum is however expensive for large size portfolio computation.

The third method uses  $l_1$  norm and our partial knowledge on the collection of assets. We reformulate the problem as a quadratic program:

$$f(r) = \max_{x_i = 1, \|x\|_1 \le m} x^T B x - r x^T A x$$
 (3)

where A and B are both positive definite matrices, r is a nonnegative number,  $x_i$  is the weight of the ith asset, and the choice of (i, m) will be addressed later. The ratio minimization problem:

$$\min_{x_i=1, \|x\|_1 \le m} \frac{x^T A x}{x^T B x}$$

will be shown to be equivalent to finding the positive root of f(r).

Nonconvexity still exists and a special algorithm designed to minimize a difference of convex functions [12] can be applied to overcome this difficulty in combination with convex algorithms such as the least angle regression [8]. The resulting algorithm is faster than the above methods and can handle portfolios with hundreds of assets with only limited prior knowledge. The combined difference of convex function and least angle regression method for root finding of f(r) is our main contribution for computing large size (over 100's assets) portfolios with a significant speedup over other nonconvex methods.

The paper is organized as follows. In section 2, we restate the problem of sparse and fast mean reverting portfolio proposed in [6]. Section 3 discusses two proxies of the mean reversion coefficient and one of them is the direct OU estimator which we found to work better. In section 4, we present the optimization problem in finding the desired portfolios and give a brief introduction to existing methods for solving it. In sections 5, 6 and 7, we introduce three new approaches in constructing portfolios. We formulate new optimization problems and develop the algorithms to solve them. Numerical experiments are presented in section 8. We perform in-sample tests and out-of-sample tests on both synthetic data and historical market data. The in-sample and out-of-sample performances curves for the mean reversion coefficients and the trading opportunities have a similar performance trend as functions of cardinality or  $l_1$  penalty level.

Before we introduce the portfolio problem, let us summary some notations.

$$- \langle x, y \rangle = x^T y;$$

- For 
$$x \in \mathbb{R}$$
,  $(x)_{+} = max(x, 0)$ ;  
-  $\chi_{A}(\mathbf{x}) = \begin{cases} 0 : \mathbf{x} \in A \\ \infty : else \end{cases}$ ;  
- If  $v \in \mathbb{R}^{n}$ ,  $v_{i}$  is the *i*th entry of  $v$ ;

## 2 Sparse and fast mean reverting portfolio problem

We briefly review the formulation in [6]. Suppose that  $S_{ti}$  is the value at time t of the ith asset with i = 1, 2..., n and t = 1, 2, ..., l, we want to form a portfolio  $P_t$  of these assets with coefficients  $x_i$ , and assume it follows an Ornstein-Uhlenbeck process given by:

$$dP_t = \lambda(\mu - P_t)dt + \sigma dW_t$$
 with  $P_t = \sum_{i=1}^n x_i S_{ti}$ 

where  $\lambda > 0$ ,  $\sigma > 0$  and  $\mu$  are parameters and  $W_t$  is a standard Brownian motion.

The objective here is to maximize the mean reversion coefficient  $\lambda$  of  $P_t$  by adjusting the portfolio weights  $x_i$  under the normalization  $\sum_{i=1}^n x_i^2 = 1$ . In addition, we want to limit the number of assets in the portfolio, i.e. we want to balance the number of nonzeros of  $x_i$ 's and the mean reversion coefficient.

# 3 Proxy of the mean reversion coefficients

In [5], the authors discuss three different criteria to measure how fast a portfolio is mean-reverting. They are predictability, the portmanteau statistics and the crossing statistics. In our paper, we mainly consider two proxies. One is predictability and the other one is called the direct OU estimator.

## 3.1 Predictability

The idea of predictability of a time series is first derived in [2]. They consider a stationary vector autoregressive (VAR) model:

$$S_t^T = S_{t-1}^T \beta + Z_t^T \tag{4}$$

where the column vector  $S_t = (S_{t1}, ..., S_{tn})^T$ ,  $\beta \in \mathbb{R}^{n \times n}$ ,  $Z_t$  is a vector of i.i.d Gaussian noise with zero mean and a covariance matrix  $\Sigma$ , independent of  $S_{t-1}$ .

In the univariate case,

$$\mathbf{E}[S_t^2] = \mathbf{E}[(S_{t-1}\beta)^2] + \mathbf{E}[Z_t^2]$$

which can be rewritten as  $\sigma_t^2 = \sigma_{t-1}^2 + \Sigma$ . Box & Tiao (1977) measure the predictability of stationary series by:

$$\nu = \frac{\sigma_{t-1}^2}{\sigma_t^2}$$

d'Aspremont [6] propose to use this measure of predictability as a proxy for the mean reversion parameter  $\lambda$ .

In multivariate case, consider a portfolio  $P_t = S_t^T x$  with weights  $x \in \mathbb{R}^n$ , then by multiplying both sides of (4) by x, we get

$$S_t^T x = S_{t-1}^T \beta x + Z_t^T x$$

and we can measure its predictability as:

$$\nu_1(x) = \frac{x^T \beta^T \Gamma \beta x}{x^T \Gamma x}$$

where  $\Gamma$  is the covariance matrix of  $S_t$ . Minimizing the predictability  $\nu_1(x)$  corresponds to maximizing  $\lambda$ .

#### 3.2 Direct OU estimator

The Ornstein-Uhlenbeck process plays an important role in constructing sparse and fast mean reverting portfolios. We will use the mean reversion coefficient,  $\lambda$ , to test the performance of a portfolio. Therefore, we will need an estimator for the mean reversion coefficient and will use this estimator as its proxy value.

Consider an Ornstein-Uhlenbeck process:

$$dP_t = \lambda(\mu - P_t)dt + \sigma dW_t \tag{5}$$

We can estimate the parameters of an OU process by linear regression. We refer the readers to [15,32] for more details. By writing (5) in a discrete form, we have

$$P_t - P_{t-1} = \lambda \mu \Delta t - \lambda P_{t-1} \Delta t + \sigma \Delta W_t$$

Note that it can be written as:

$$P_t = \lambda \mu \Delta t + P_{t-1}(1 - \lambda \Delta t) + \sigma \Delta W_t$$

When  $\mu = 0$ , this has the form (4) in the case that n = 1. It is equivalent to the linear regression:

$$y = a + bx + \epsilon_t$$

Therefore, we could estimate all the parameters by regressing  $P_t - P_{t-1}$  on  $P_{t-1}$ . Then we can recover  $\hat{\lambda}$  as  $-\frac{\hat{b}}{\Delta t}$ ,  $\hat{\mu}$  as  $\frac{\hat{a}}{\hat{\lambda} \Delta t}$  and  $\hat{\sigma}$  as  $\frac{\hat{\sigma}(\epsilon_t)}{\sqrt{\Delta t}}$ .

Maximizing the estimated mean reversion coefficient  $\lambda$  corresponds to minimizing the estimated slope of  $\hat{b}$ . Note that

$$b = \frac{\operatorname{cov}(P_t - P_{t-1}, P_{t-1})}{\operatorname{var}(P_{t-1})} = \frac{\operatorname{cov}(P_t, P_{t-1})}{\operatorname{var}(P_{t-1})} - 1$$

Replacing  $P_t$  by  $S_t^T x$ , we have:

$$b = \frac{\text{cov}(S_t^T x, S_{t-1}^T x)}{\text{var}(S_{t-1}^T x)} - 1$$

We define the direct OU estimator as:

$$\nu_2(x) = \frac{\text{cov}(S_t^T x, S_{t-1}^T x)}{\text{var}(S_{t-1}^T x)}$$
 (6)

If  $S_t$  is stationary, then we can rewrite (6) in the following form:

$$\nu_2(x) = \frac{x^T(\text{cov}(S_t, S_{t-1}) + \text{cov}(S_t, S_{t-1})^T)x}{2x^T \text{var}(S_{t-1})x}$$

Minimizing the predictability  $\nu_2(x)$  corresponds to maximizing  $\lambda$ .

This proxy is a speical case of the crossing statistics in [5] when p = 1. The only assumption needed is that the linear combinations of the asset  $S_t^T x$  is stationary. This phenomenon is called cointegration [9]. According to our numerical tests, the portfolios constructed under this proxy gives better trading performance.

Though the estimation of  $cov(S_t, S_{t-1}) + cov(S_t, S_{t-1})^T$  is not guaranteed to be positive definite, our numerical tests show that in most cases it is still true for the market data. Even if this condition fails, we could use the estimation of  $cov(S_t, S_{t-1}) + cov(S_t, S_{t-1})^T + \rho var(S_{t-1})$  as the matrix in the quadratic form of the numerator where  $\rho$  is a tuning parameter. By choosing an appropriate  $\rho$ , the matrix will be positive definite and it is equivalent to the optimization problem when  $\rho = 0$ .

# 4 Sparse optimization problem and existing methods

In the previous section, we discussed two proxies of the mean reversion coefficient. Note that both of them can be written as

$$\frac{x^T A x}{x^T B x}$$

where A and B are  $n \times n$  positive definite matrices and we will make this assumption for the rest of the paper. The estimation procedures for A and B for two proxies are discussed in the Appendix B.

If we do not penalize the cardinality of x, then minimizing these proxies is the same as a generalized eigenvalue problem. In order to obtain a sparse

solution, d'Aspremont in [6] proposed the following sparse optimization problem:

$$\min_{x} x^{T} A x / x^{T} B x$$

$$s.t. \quad ||x||_{0} \le k$$

$$||x||_{2} = 1,$$
(7)

where the  $l_0$  norm of a vector x is the number of nonzero entries of x. This problem has been proved to be NP-hard [22]. When the dimension of the problem is large, we cannot expect to find the optimal solution. Several methods of solving (7) have been proposed in [6,11,27]. We give a brief summary here:

- Exhaustive search method: it tests all  $\frac{n!}{k!(n-k)!}$  possible combinations of assets and find the smallest generalized eigenvalue and eigenvectors. This method gives us an optimal solutions and works very well when n is small. However, it will be extremely slow when n is large.
- Greedy search: denote  $I_k$  as the support of the solution and set  $I_k = \emptyset$  initially. Each time, we pick one asset from  $\{1, 2, ..., n\} \setminus I_k$  such that it has smallest objective among all the other choices. Then we add it to  $I_k$  and repeat this procedure k times. This method gives a sub-optimal solution.
- Truncation method: first we solve the unconstrained problem and find  $x_{opt}$ . Then we find the index set  $J_k$  of the largest k components of  $|x_{opt}| = (|x_1|, ..., |x_n|)^T$ . Next we solve the generalized eigenvalue problem on the set  $J_k$  by taking the corresponding part of matrices A and B. It gives a sub-optimal solution and it is the fastest among all the listed methods since it only requires solving the generalized eigenvalue problem twice.
- Semidefinite relaxation method: this method reformulates the problem (7) as a semidefinite program. This can be considered as an extension of DSPCA in [33]. We found that this relaxiation method can be considered as replacing the  $||x||_0 \le k$  by  $||x||_1/||x||_2 \le \sqrt{k}$ . For more details, we refer the readers to [6] and the Appendix A. The major drawback of this algorithm is that it is not scalable to high-dimensional datasets [27].
- DC-PCA: the problem (7) is formulated as difference of convex functions problem and the authors apply the difference of convex functions algorithm(DCA) to efficiently solve the optimization problem. In [27], this method was proven to be efficient for solving various PCA problems, but the authors do not apply it for the generalized eigenvalue problems.

In the next three sections, we will present three approaches for approximating the optimal solution to the problem (7).

# 5 Optimization problems based with $l_1$ constraint

In this section, we consider using the  $l_1$  constraint in the sparse generalized eigenvalue problem. This can be considered as an extension of the  $l_1$  constraint

sparse principal component analysis problem. The problem is formulated as

$$\max_{x} x^{T} B x$$

$$s.t. \quad ||x||_{1} \le \sqrt{s}$$

$$x^{T} A x \le 1,$$
(8)

One way to tackle this problem is to treat the problem as a difference of convex functions and using the DC (difference of convex functions) algorithm. DC programming is extensively studied in [12,21,28,29]. The method aims to minimize a function g-h where both g and h are convex functions on the whole space. In most of the literature on DC programming, an algorithm to find a local optimum is used. In practice, the local algorithm often yields a global minimum.

For 8, we have  $g(x) = \chi_F(x)$  and  $h(x) = -x^T B x$ , where  $F = \{x : x^T A x \le 1, ||x||_1 < \sqrt{s}\}.$ 

Therefore, we will update x by solving the following problem by the DC algorithm:

$$\max_{x} \langle 2Bx^{(k-1)}, x \rangle$$

$$s.t. \quad ||x||_1 \le \sqrt{s}$$

$$x^T Ax < 1,$$
(9)

where  $x^{(k)}$  is the x in the kth interaction.

In fact, we could also obtain the same formulation by the generalized power iteration [18] and alternating maximization [26]. When the A is an identity matrix, then we reduce to a sparse PCA problem. The explicit solution is discussed in [18,26]. For a general A, the maximizer of (9) is then

$$x^* = \frac{A^{-1}(v-z)}{\sqrt{(v-z)^T A^{-1}(v-z)}}$$

where  $v = 2Bx^{(k-1)}$  and z is the optimizer of the dual problem of (9):

$$\min_{\lambda > 0, \|z\|_{\infty} < \lambda} \lambda \sqrt{s} + \sqrt{(v-z)^T A^{-1} (v-z)}$$

When  $A = (A_{ij})$  is a diagonal matrix, then the above reduces to a closed-form expression:

$$x_i^* = \frac{A_{ii}^{-1} \operatorname{sgn}(v_i)(|v_i| - \lambda)_+}{\sqrt{\sum_i A_{ii}^{-1}(|v_i| - \lambda)_+^2}}$$

## 5.1 Comparison to DC-PCA

DC-PCA in [27] starts with an  $l_0$  penalized optimization problem. The authors approximate the  $||x||_0$  by  $\sum_i \log(x_i + \epsilon)$  and formulate

$$\max_{x} x^{T} B x - \rho \sum_{i} \log(x_{i} + \epsilon)$$
s.t. 
$$x^{T} A x \leq 1,$$
(10)

After applying the DC algorithm, we will need to solve the following problem at each iteration:

$$\max_{x} \langle D(x^{(k-1)})Bx^{(k-1)}, x \rangle - \frac{\rho}{2} ||x||_{1} s.t. \quad x^{T} D(x^{(k-1)})AD(x^{(k-1)})x \le 1,$$
(11)

where D(x) is the diagonal matrix whose diagonal entries  $D_{ii} = x_i$ . The next  $x^{(k)}$  follows  $x_i^{(k)} = x_i^{(k-1)} x_i$ . One thing to note that is if  $x_i^{(k-1)} = 0$  then  $x_i^{(k)} = 0$ . Therefore, the initial  $x^{(0)}$  needs to be chosen carefully.

When  $A = (A_{ij})$  is a diagonal matrix, then we also have an explicit solution to the problem (11):

$$x_i^* = \begin{cases} \frac{(c_i)^{-1} \operatorname{sgn}(v_i)(|v_i| - \rho/2)_+}{\sqrt{\sum_i c_i^{-1}(|v_i| - \rho/2)_+^2}} & x_i^{(k-1)} \neq 0\\ 0 & x_i^{(k-1)} = 0 \end{cases}$$

where 
$$c_i = A_{ii}(x_i^{(k-1)})^2$$
 and  $v = D(x^{(k-1)})Bx^{(k-1)}$ .

## 5.2 Comments

These two algorithms are very efficient if the matrix A is a diagonal matrix due to the existence of the explicit solution. When A is not a diagonal matrix, we have not found them efficient to handle a large size problem.

# 6 Optimization problems based on the ratio of $l_1$ and $l_2$ norms

#### 6.1 Motivation

One popular method in handling a cardinality constraint is to use a norm penalization. Standard choices include  $l_1$  and  $l_p$  (0 < p < 1). In our case, we would prefer using  $\frac{\|x\|_1}{\|x\|_2}$  since our problem (2) is scale invariant. The scale constraint is required for  $\|x\|_1$  and  $\|x\|_p$  penalties otherwise they cannot enforce sparsity. The reason is that we could simply decrease the penalty by decreasing the scale of all the elements of x. For analysis of sparsity promoting properties of ratio of  $l_1$  and  $l_2$  norms, we refer to [31].

In the following sections, we will present two formulations of sparse mean reverting problems (7), discuss the algorithms to solve them and show two theorems of the ratio of  $l_1$  and  $l_2$  norms in enforcing the sparsity.

## 6.2 Formulation 1

We could reformulate the problem (7) in the following way:

$$\min_{x} x^{T} A x / x^{T} B x$$

$$s.t. \quad \frac{\|x\|_{1}}{\|x\|_{2}} \leq m$$

$$x \neq 0,$$
(12)

Using a similar method in [6], we find that we can recover the semidefinite programming problem in [6]. The details can be found in Appendix A.

## 6.3 Formulation 2

We modify the problem (7) in the following way:

$$\min_{x \neq 0} \frac{x^T A x}{x^T B x} + \gamma \frac{\|x\|_1}{\|x\|_2} \tag{13}$$

where  $\gamma$  is a tuning parameter.

Comparing with the problem (7), the problem (13) does not specify the cardinality beforehand. In this way, we could consider the  $l_1$  norm in the numerator as a way to quantify the uniform transaction costs.

## 6.4 Algorithm to solve (13)

The major difficulty comes from the nonconvexity of the problem. Most optimization algorithms will only provide a local minimizer. In our problem, the performance of local minimizers may vary a lot. Therefore, it is important to develop an algorithm to approximate the global minimizer. The intermittent diffusion(ID) algorithm proposed in [4] can be helpful. The main idea of this algorithm is "to add intermittent, instead of continuously diminishing, random perturbations to the gradient flow generated by the objective, so that the trajectories can quickly escape from the trap of one minimizer and then approach others."

# 6.5 Analysis of the ratio of $l_1$ and $l_2$ penalty

In this section, we want to prove some useful properties of the solutions to the problem (13). Theorem 1 indicates that an almost 1-sparse solution can always be recovered if the value of  $\gamma$  is large enough. Theorem 2 shows that the ratio of  $l_1$  and  $l_2$  norms of the solution to the problem (13) decreases as we increase the parameter  $\gamma$ .

We will define  $f(x, \gamma)$ , L(x) and P(x) as:

$$\begin{split} f(x,\gamma) &= \frac{x^T A x}{x^T B x} + \gamma \frac{\|x\|_1}{\|x\|_2} \\ L(x) &= \frac{x^T A x}{x^T B x} \\ P(x) &= \frac{\|x\|_1}{\|x\|_2} \end{split}$$

# Algorithm 1 Solve (13)

Input A, B and  $\gamma$ ;

Let  $\alpha$  be the scale of diffusion strength,  $\kappa$  be the scale for diffusion time and N be the total number of realizations.

Set the initial state  $x_0$  as the minimizer of  $\frac{x^TAx}{x^TBx}$  with the constraint that  $||x_0||_2 = 1$ , i.e. the generalized eigenvector associated with the smallest eigenvalue

Find a local minimizer  $\hat{x}_0$  of problem (13) given  $x_0$  and set  $X_{opt} = \hat{x}_0$ .

for i = 1 to N do

Generate two positive random numbers d, s within [0,1] by uniform distribution and let  $\sigma := \alpha d$  and  $T := \kappa s$ .

Solve the stochastic equation for  $t \in [0, T]$ 

$$dx(t,\omega) = -\nabla^s F(x(t,\omega))dt + \sigma dW(t), \quad x(0,\omega) = X_{opt}$$

where  $\nabla^s F$  is the gradient of the objective (13) and record the final state  $x_T := x(T, \omega)$ .

Find a local minimizer  $\hat{x}_i$  of problem (13) by line search algorithm with starting point  $x_T$ .

$$X_{opt} = \hat{x}_i \text{ if } f(\hat{x}_i) < f(X_{opt}).$$

end for

We denote the set of minimizers of the problem (13) by  $x(\gamma)$  and we will set their  $l_2$  norm to 1. When  $\gamma$  is fixed, the existence of the minima of  $f(x,\gamma)$  is due to its continuity on the unit sphere and the compactness of the unit sphere. Since  $x^T A x / x^T B x$  is scale invariant, any vector on the same direction yields the same value. If there are different directions that yield the same  $f(x,\gamma)$ , then we always prefer those with the smaller ratio of  $l_1$  and  $l_2$  norms. Therefore, the set of the optimizers,  $x(\gamma)$ , has a unique value of the function  $P(\cdot)$  on this set (and then a unique value of  $L(\cdot)$ ). Mathematically, it can be defined in the following way:

$$X(\gamma) = \{ x \in \mathbb{R}^n : ||x||_2 = 1, f(x, \gamma) = \min_{z \neq 0, ||z||_2 = 1} \{ \frac{z^T A z}{z^T B z} + \gamma \frac{||z||_1}{||z||_2} \} \}$$

$$x(\gamma) = \{x \in X(\gamma) : P(x) = \min\{P(y) : y \in X(\gamma)\}\}$$

Under this definition,  $P(x(\gamma))$ ,  $L(x(\gamma))$  and  $f(x(\gamma), \gamma)$  are well defined, since  $P(\cdot)$ ,  $L(\cdot)$  and  $f(\cdot, \gamma)$  are constant on  $x(\gamma)$ .

In the following proofs, we will use the notations:

$$B(x, \delta) \equiv \{x' : ||x' - x||_2 < \delta\}$$

$$S_k \equiv \{x \in \mathbb{R}^n : ||x||_2 = 1, ||x||_0 \le k\}$$

$$S^n \equiv \{x \in \mathbb{R}^n : ||x||_2 = 1\}$$

 $d(U,V) = \min\{\|x - y\|_2 : x \in U, y \in V, U \text{ and } V \text{ are compact sets in } \mathbb{R}^n\}$ 

**Theorem 1** Given  $A = \{a_{ij}\}$  and  $B = \{b_{ij}\}$  both  $n \times n$  positive-definite matrices, then for any  $\epsilon > 0$ , there exists a number  $\gamma(\epsilon)$ , such that for any  $\gamma \geq \gamma(\epsilon), \|x-e\|_2 < \epsilon, \text{ where } x(\gamma) \text{ is the set of optimal solutions of problem}$ (13) given  $\gamma$ ,  $x \in x(\gamma)$ , e is a 1-sparse vector and  $||e||_2 = 1$ .

*Proof* First, note that the problem (13) is scale invariant, therefore, we could constrain the problem on the sphere  $||x||_2 = 1$ . All the vectors in our proof have  $l_2$  norm 1.

Note that the 1-sparse minimizers of L(x) on the sphere ||x|| = 1 are the vectors  $\pm e_i$  with all 0s except for a 1 in the ith coordinate. The i is determined

$$i = \arg\min_{i} \{\frac{a_{ii}}{b_{ii}}\}$$

Without loss of generality, we will assume that i = 1 and this corresponds to a unique minimizer, i.e.

$$\frac{a_{11}}{b_{11}} < \frac{a_{ii}}{b_{ii}}, \quad \text{for all } i \neq 1$$

Note that both  $f(x, \gamma)$  and L(x) are even functions of x. We could further

restrict our region to  $S_+ \equiv \{x = (x_1, ..., x_n)^T : x_1 \ge 0 \text{ and } ||x||_2 = 1\}.$ Since L(x) is continuous on  $S_+$ , then  $\exists \delta_i$ , for any point x in the region

 $\{x: \|x \pm e_i\| < \delta_i\} \cap S_+, L(x) \ge L(e_1) = \frac{a_{11}}{b_{11}}, \text{ for } i = 2, 3, ..., n.$ Denote  $D_i = \{x: \|x \pm e_i\| < \delta_i\} \text{ for } i = 2, 3, ..., n, \text{ then for all the points}$  $x \in S_+ \cap (\bigcup_{i=2}^n D_i), L(x) \ge L(e_1)$  and  $||x||_1/||x||_2 \ge 1$ . Therefore, they cannot beat  $e_1$  for any  $\gamma$ .

For any  $\epsilon > 0$ , let  $D_1 = \{x : ||x - e_1|| < \epsilon\}$ , then  $S_+(\epsilon) \equiv S_+ \setminus (\bigcup_{i=1}^n D_i)$  is a closed and bounded set. Therefore,  $\gamma(\epsilon)$  must satisfy the following inequality:

$$f(e_1, \gamma(\epsilon)) \le f(x, \gamma(\epsilon)), \quad \forall x \in S_+(\epsilon).$$

This is equivalent to

$$\frac{a_{11}}{b_{11}} + \gamma(\epsilon) \le L(x) + \gamma(\epsilon) \frac{\|x\|_1}{\|x\|_2}, \quad \text{for any } x \in S_+(\epsilon).$$

And it implies:

$$\gamma(\epsilon) \ge (\frac{a_{11}}{b_{11}} - L(x))/(\frac{\|x\|_1}{\|x\|_2} - 1), \quad \text{for any } x \in S_+(\epsilon).$$

The last line holds, since for any  $x \in S_+(\epsilon)$ ,  $\frac{\|x\|_1}{\|x\|_2}$  is guaranteed to be greater than 1. Note that the function  $(\frac{a_{11}}{b_{11}} - L(x))/(\frac{\|x\|_1}{\|x\|_2} - 1)$  is well-defined and continuous on  $S_{+}(\epsilon)$ . Therefore, we can set  $\gamma(\epsilon)$  as:

$$\gamma(\epsilon) = \max_{x \in S_+(\epsilon)} (\frac{a_{11}}{b_{11}} - L(x)) / (\frac{\|x\|_1}{\|x\|_2} - 1)$$

**Theorem 2** Suppose  $x(\gamma)$  is the set of minimizers of the problem (13), then  $P(x(\gamma))$  is nonincreasing in  $\gamma$ .

*Proof* For any arbitrary  $\gamma_1$  and  $\gamma_2$ , assume  $\gamma_1 < \gamma_2$ . We want to show that  $P(x(\gamma_1)) \ge P(x(\gamma_2))$ . Since  $x(\gamma_1)$  and  $x(\gamma_2)$  are two minimizers, so we have

$$L(x(\gamma_1)) + \gamma_1 P(x(\gamma_1)) \le L(x(\gamma_2)) + \gamma_1 P(x(\gamma_2)) \tag{14a}$$

$$L(x(\gamma_2)) + \gamma_2 P(x(\gamma_2)) \le L(x(\gamma_1)) + \gamma_2 P(x(\gamma_1))$$
(14b)

(14a) and (14b) implies

$$L(x(\gamma_1)) - L(x(\gamma_2)) \le \gamma_1(P(x(\gamma_2)) - P(x(\gamma_1)))$$

$$L(x(\gamma_1)) - L(x(\gamma_2)) \ge \gamma_2(P(x(\gamma_2)) - P(x(\gamma_1)))$$

Then we have:

$$\gamma_2(P(x(\gamma_2)) - P(x(\gamma_1))) \le L(x(\gamma_1)) - L(x(\gamma_2)) \le \gamma_1(P(x(\gamma_2)) - P(x(\gamma_1)))$$

This implies

$$(\gamma_2 - \gamma_1)(P(x(\gamma_2)) - P(x(\gamma_1))) \le 0$$

By the assumption that  $\gamma_1 < \gamma_2$ , therefore  $P(x(\gamma_1)) \ge P(x(\gamma_2))$ .

# 7 Optimization problems based on the $l_1$ norm and prior knowledge

## 7.1 Motivation

In the previous section, we applied a stochastic algorithm in searching for a global optimizer of the problem (13). The major drawback of the algorithms is that we can not guarantee a global optimizer. Based on the theory of the intermittent diffusion(ID) algorithm, we could increase the probability of finding the global optimizer by increasing the number of realizations N. However, choosing N is problem dependent and it will be a difficult task when the dimension of the problem is large. We have to balance the efficiency and the performance of the solution.

Therefore, an efficient global optimization algorithm is desired. In addition, we may prefer a simpler norm to enforce sparsity, because it will lead to simpler algorithms. We also would like to keep the number of tuning parameters as low as possible.

## 7.2 Formulation

Note that the problem (7) is equivalent to:

$$\max x^T B x / x^T A x$$
s.t.  $||x||_0 \le k$ 

$$||x||_2 = 1$$

Now, we use the  $l_1$  norm to enforce the sparsity and consider the following problem:

$$f(r) = \max_{x_i = 1, \|x\|_1 \le m} x^T B x - r x^T A x$$
 (15)

where i is a predetermined fixed number. We could choose i based on our prior knowledge or our investment need. For example, we could choose i by selecting the asset which has the largest entry in absolute value by solving the unconstraint problem. The constraint  $x_i = 1$  enables using the  $l_1$  norm to enforce the sparsity and also simplifies the problem to a quadratic program.

## 7.3 Equivalence of two formulations

We would like to show that

# Theorem 3 Let

$$f(r) = \max_{x_i = 1, ||x||_1 \le m} x^T B x - r x^T A x$$

where A and B are both positive definite matrices, then

- a For any given R > 0, (15) is continuous on  $0 \le r \le R$ ;
- b f(r) is an nonincreasing function of r;
- c There exists  $r_*$  such that  $f(r_*) = 0$ ;
- d Suppose the optimizer at  $r_*$  is  $x_*$ , then  $x_*$  is the optimizer of the following problem:

$$\max_{x \neq 0} x^T B x / x^T A x$$

$$s.t. \quad x_i = 1$$

$$\|x\|_1 \leq m$$
(16)

*Proof* a. For any given R > 0, since  $g(x,r) = x^T B x - r x^T A x$  is continuous on the closed and bounded set  $\{(x,r), x_i = 1, \|x\|_1 \le m, 0 \le r \le R\}$ , all the optimizers are obtainable. Due to the uniform continuity, f(r) is continuous.

b. Suppose  $r_1 > r_2$ , then for any feasible x, we must have

$$(x^T B x - r_1 x^T A x) - (x^T B x - r_2 x^T A x)$$
  
=  $(r_2 - r_1) x^T A x < 0$ 

Since A is positive definite and  $x \neq 0$ . Therefore, it is nonincreasing in r.

- c. Notice that f(0) > 0. Since A is positive definite, there must exist an R > 0 such that B RA is negative definite. Therefore, f(R) < 0. Due to the continuity of f(r), there exists  $r_*$  such that  $f(r_*) = 0$ .
  - d. Suppose there exists a feasible u such that

$$r_1 := \frac{u^T B u}{u^T A u} > r_*$$

Then we must have:

$$0 = u^{T}Bu - r_{1}u^{T}Au$$

$$= u^{T}Bu - r_{*}u^{T}Au - (r_{1} - r_{*})u^{T}Au$$

$$\leq x_{*}^{T}Bx_{*} - r_{*}x_{*}^{T}Ax_{*} - (r_{1} - r_{*})u^{T}Au$$

$$= -(r_{1} - r_{*})u^{T}Au < 0$$

Therefore, the result is shown by contradiction.

This theorem tells us that if we can find a root of f(r) and the corresponding maximizer, then we have found the global maximizer of problem (16).

The conversion of the ratio minimization problem to a sequence of difference minimization problems has been proposed in solving the trace ratio optimization problem arising in machine learning and high dimensional data analysis [23]. Here in Theorem 6.1, we considered the additional  $l_1$  constraint.

## 7.4 Algorithm for finding $r_*$

We used a binary search algorithm.

# **Algorithm 2** Find $r_*$

```
Input the matrices B and A, a threshold \epsilon > 0 for stopping, initial x^{(0)}, r_{min}, r_{max}, i, m and maximum iterations NMAX;

Set N = 1;

while N \leq \text{NMAX do}

Set r_0 = \frac{1}{2}(r_{min} + r_{max});

Solve the quadratic program (15) with r_0 and obtain the maximizer x^{(0)};

if |x^{(0)}|^T Bx^{(0)} - r_0(x^{(0)})^T Ax^{(0)}| < \epsilon OR r_{max} - r_{min} < \epsilon then Stop

end if

if (x^{(0)})^T Bx^{(0)} - r_0(x^{(0)})^T Ax^{(0)} > 0 then r_{min} = r_0

else r_{max} = r_0

end if

end while
```

# 7.5 Algorithm for solving (15)

The difficult part is solving the quadratic program (15). It is a nonconvex optimization problem. Therefore, classical algorithms can not guarantee a global optimizer. In addition, the computational cost is expensive when the dimension of the problem is  $large(\geq 100)$ . We will again apply the DC algorithm to solve the problem.

In order to apply the DC algorithm, we could first reformulate the problem (15) in the following way by plugging in the constraint  $x_i = 1$ :

$$\min_{\|x\|_1 \le m'} rx^T V x + c^T x - x^T U x$$

where U and V are the submatrices of B and A without the ith row and column and  $x \in \mathbb{R}^{n-1}$ .

In addition, by a standard method in DC algorithm, we could change it to an unconstraint problem:

$$\min_{x} rx^{T} V x + c^{T} x - x^{T} U x + \chi_{\|x\|_{1} \le m'}(x)$$
(17)

Let

$$g(x) = rx^{T}Vx + c^{T}x + \chi_{\|x\|_{1} \le m'}(x)$$
  $h(x) = x^{T}Ux$ 

then the objective is a difference of g and h. We now can use the following DC algorithm:

# **Algorithm 3** Solve (17) for a given r

Choose  $x^{(0)}$  in  $\mathbb{R}^{n-1}$ ;

#### repeat

Set  $y^{(k)} = 2Ux^{(k)}$ ;

Solve the optimizer  $x^{k+1}$  of the convex program:

$$\inf\{rx^T V x + c^T x - x^T y^{(k)} + \chi_{\|x\|_1 \le m'}(x), x \in \mathbb{R}^{n-1}\}$$
 (18)

until convergence

## 7.6 Algorithms for solving (18)

The problem (18) can be considered as a quadratic program with  $l_1$  constraint.

$$\min_{x} rx^{T}Vx + c^{T}x + x^{T}y^{(k)} 
s.t. ||x||_{1} \le m'$$
(19)

It can also be rewritten as a least squares optimization with  $l_1$  constraint. This problem has been well studied in the literature. The major difficulty is handling the  $l_1$  constraint. We apply the least angle regression method to solve this problem.

The idea is that we first reformulate (19) as a LASSO problem [30]. Notice that the objective of a LASSO problem is

$$\|\mathbf{X}x - \mathbf{Y}\|_2^2 + \lambda |x|_1 = x^T \mathbf{X}^T \mathbf{X}x - 2\mathbf{Y}^T \mathbf{X}x + \mathbf{Y}^T \mathbf{Y} + \lambda |x|_1$$

where **X** is a matrix of predictor variables, **Y** is an outcome vector and  $\lambda$  is a tuning parameter.

Therefore, by solving the following system for X and Y, we retrieve a LASSO type problem from (19):

$$rV = \mathbf{X}^T \mathbf{X}$$
  $c + y = -2\mathbf{X}^T \mathbf{Y}$ 

The first equation can be solved by Cholesky decomposition and then the second equation is easy to solve. Finally, we could apply the least angel regression algorithm (LARS) [8].

## 8 Numerical tests

Our procedures are implemented in Matlab R2015a. We used the optimization package YALMIP. Computations are performed on a Dell desktop with 8G RAM and 3.4 GHz i7 CPU. We have used two historical data sets. One is the U.S. daily swaps data for maturities 1Y, 2Y, 3Y, 4Y, 5Y, 7Y, 10Y and 30Y from July 3, 2000 until July 15, 2005. The data are obtained from **www.Economagic.com**. The total number of data is  $1257 \times 8$ . The data are in percent with two digits after the decimal point. The other one is the daily closed prices of S&P 500 companies. The data are collected from Yahoo finance. In order to obtain a large sample set, we only select those companies that have been in the list since July 2005. After this preselection procedure, we have 458 companies left. We used the stock prices of the first 100 companies according to the alphabetical order of the ticker symbols. The data size in our numerical test is  $2000 \times 100$ .

In the following sections, we performed several tests:

- Scalability and performance tests of DC-PCA and l<sub>1</sub>-constraint DC-PCA.
   Details are in section 8.1;
- In-sample and out-of-sample tests for the performance of the portfolios constructed under two mean reverting proxies: predictability and direct OU estimator on synthetic data. Details are in section 8.2;
- The impact of the size of training set on out-of-sample performance. Details are in section 8.3;
- In-sample and out-of-sample tests for the ratio of  $l_1$  and  $l_2$  norms approach on the U.S. swaps data. Details are in section 8.4;
- In-sample tests for the  $l_1$  norm with prior knowledge approach on the historical daily closing prices of 100 stocks. Details are in section 8.5.1;
- In-sample and out-of-sample tests for the  $l_1$  norm with prior knowledge approach on the U.S. swaps data. Details are in section 8.5.2;
- In-sample and out-of-sample tests for the  $l_1$  norm with prior knowledge approach on synthetic data with 100 assets. Details are in section 8.5.3;

## 8.1 Scalability and performance tests of DC-PCA and $l_1$ -constraint DC-PCA

In [27], the numerical tests show that the DC-PCA is preferred over SPCA[7] and DSPCA[33] on sparse PCA problem. The solution of DC-PCA gives larger

explained variance and smaller computational costs relatively. Now we want to see how well our proposed method and DC-PCA perform on the sparse generalized eigenvalue problem.

We have performed two tests on synthetic data. The first test is a scalability test. We apply these two method on randomly chosen problems of size n ranging from 200 to 2500 for 6 different values of  $\rho$  and s and compute the average time costs. The matrices A and B are both positive definite and A is a diagonal matrix. The second test is to compare the explained variances of two methods. We generate the two positive definite matrices A and B of size  $3000 \times 3000$  where A is a diagonal matrix. We compute a series of solutions with different sparse parameters  $\rho$  and s and compare the explained variances of the solutions with the same numbers of non-zero loadings.

The results are shown in 1. The left shows that the plot of average CPU time vs. n(the number of rows of the matrix) for these two methods with the empirical complexity growing as  $O(n^p)$ , where p=1.61 for our proposed  $l_1$  constraint DC-PCA and p=2.40 for DC-PCA. We could see that the  $l_1$  constraint formulation is more efficient than the DC-PCA. The right plot shows the comparative performance (explained variance vs. sparsity) of two methods for the first generalized eigenvector. The amount of variance explained by the first generalized eigenvector is 23%. In this case, two methods are very close and DC-PCA is slightly better than our proposed method.

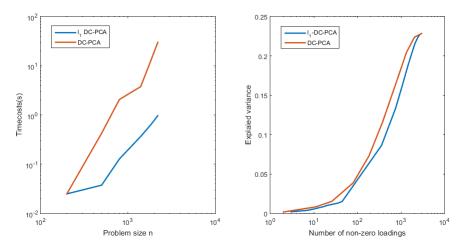


Fig. 1 Comparison between  $l_1$  constraint DC-PCA and DC-PCA in DC-PCA in scalability and performance.

From this test, two methods are very efficient when the matrix A is diagonal. The computation can be done in less than one second even if the problem is of size of 2500.

## 8.2 Comparison of two mean reverting proxies

We wanted to compare the performance of portfolio selection via two proxies that we discussed in section 3.

We set the matrix  $\beta$  and the noise covariance matrix  $\Sigma$  in the VAR(1) model (4) as our estimations based on the U.S. swaps data. We generated a data set of size  $350 \times 8$ , so there are 350 observations for each asset and there are 8 assets in total. We used the first  $100 \times 8$  samples as the training set and the rest as the test set. Next we made estimations of parameters and solved for the optimal solutions by the exhaustive search method based on the training set. We repeated our simulation 1000 times and then we compared the average of the estimated mean reversion coefficients of our sparse portfolios on both the training set and the test set.

Next we tested the performance of the sparse portfolios based on a simple convergence trading strategy. In most of the application of convergence trading, investors will consider two parameters  $\mu$  and  $\tau$ , where  $\mu$  is the estimated average asset value and  $\tau$  is the tolerance of mispricing. In [11], the authors developed a strategy that only takes advantages of underpricing of the portfolio. We generalized their strategy and took advantages of both the underpricing and overpricing of the portfolio. For simplicity, we also assumed that we have the ability to buy and sell assets without any transaction costs and we also have the ability to short sell. We will use the number of trading opportunities to measure the performance of the portfolio. Our objective here is using a simple procedure to evaluate portfolio performance but not building an "optimal" convergence trading strategy. The trading opportunity means the observation that the price converges after out-of-tolerance mispricing. We use K to denote the number of trading opportunities and  $P_t$  to denote the portfolio value at time t.

The trading strategy can be summarized as follows:

- If the observed sample  $P_t > \mu + \tau$ , we will sell our portfolio if we already hold one and K = K + 1. We will short this portfolio if we didn't short it before. Otherwise we perform no action.
- If the observed sample  $P_t < \mu \tau$ , we will go long our portfolio if we already short one and K = K + 1. We will buy this portfolio if we didn't hold it before. Otherwise we perform no action.
- If the observed sample  $\mu \tau \leq P_t \leq \mu + \tau$ , we will go long our portfolio if we already short one and K = K + 1. We will sell our portfolio if we already hold one and K = K + 1. Otherwise we perform no action.

Figure 2 will be helpful in understanding the trading strategy. The X-axis shows the time periods are from day 1 to day 60. The Y-axis shows the values of the portfolio. The green dashed line is  $y = \mu$ , the red solid line (overpriced bound) is  $y = \mu + \tau$  and the teal solid line (underpriced bound) is  $y = \mu - \tau$ .

The trading opportunities will increase by 1 if we have a high price or low price before returning to the normal range.

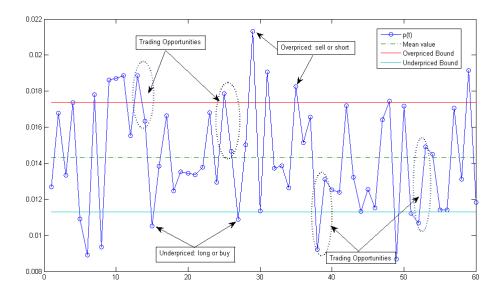


Fig. 2 Trading Opportunities (Not all are marked in the plot.)

The simulation test is similar as before. After we found the optimal solutions to the problem (7) of a given k by the exhaustive search method, we tested the trading strategy only on the test set and counted the trading opportunities. We repeated 1000 times and calculated the averages. The estimation of  $\mu$  and  $\tau$  is based on the training set. We set  $\mu$  as the sample mean and  $\tau$  as the sample standard deviation.

The results are shown in Figure 3. The top left plot is the average in-sample mean reversion coefficients versus cardinality. The top right plot is the average out-of-sample mean reversion coefficients versus cardinality. The bottom plot is the average out-of-sample trading opportunities versus cardinality.

We see that the solutions under direct OU estimator always perform better than that under the predictability. We can also note that the trading opportunities go down somewhat for the portfolios constructed under direct OU estimator when the cardinalities increase from 4 to 8. This could be caused by choice of the parameter  $\tau$ . A higher  $\tau$  may increase the profit at each trading opportunity, but decrease the total number of trades. We remark that in [5] the authors use a variance constraint to improve the profits during arbitrage opportunities.

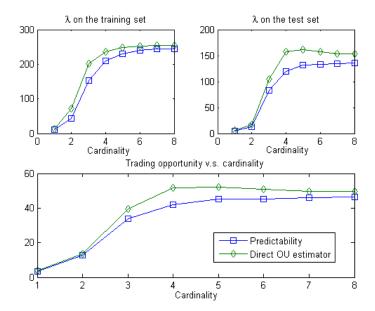


Fig. 3 Comparison of the average performance of sparse mean reverting portfolios solved under different proxies using the exhaustive search method on synthetic data from VAR(1) model (4). Top left: average in-sample mean reversion coefficients versus cardinality; Top right: average out-of-sample mean reversion coefficients versus cardinality; Bottom: average out-of-sample trading opportunities versus cardinality;

## 8.3 The impact of the size of training set on out-of-sample performance

From our numerical results, we notice that the size of the training set will affect the out-of-sample performance. To illustrate this, we perform the following tests.

After presetting proper  $\beta$  and  $\Sigma$ , we generate a data set of 10 assets for 400 observations based on the VAR(1) model (4). We use the last 50, 100, 200 and 400 observations as the training set and estimate A's and B's in (7) respectively. Next, we construct the portfolios under different cardinality constraints by the exhaustive search method by using those A's and B's.

We then generate the test data sets. Each has 400 observations of 10 assets. They follow the same VAR(1) model. The starting values are the final observations of the training data. In total, we have 100 test data sets and we compare the average out-of-sample mean reversion coefficients of the portfolios constructed based on the training set. If we use the last 50 observations of the training set for estimation, then we will only test those portfolios on the first 50 observations of each test set, and so forth.

Figure 4 shows the results. It illustrates impact of the size of training set on out-of-sample performance. We compare the in-sample mean reversion coefficients with the average out-of-sample mean reversion coefficients under different cardinality constraints. The portfolios are constructed by solving the

problem (7) by the exhaustive search method. The top left uses 50 observations of the training set and 50 observations of the test set; the top right uses 100 observations of the training set and 100 observations of the test set; the bottom left uses 200 observations of the training set and 200 observations of the test set; the bottom right uses 400 observations of the training set and 400 observations of the test set. Notice that when the size of the training set increases, the out-of-sample performance becomes closer to the in-sample performance.

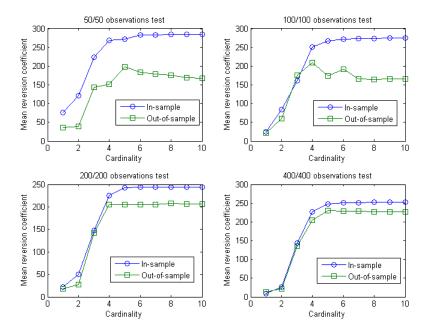


Fig. 4 The tests of the impact of the size of training set on out-of-sample performance based on synthetic data. The portfolios are constructed by the exhaustive search method under different cardinality constraints. Top left: 50 observations of the training set and 50 observations of the test set; Top right: 100 observations of the training set and 100 observations of the test set; Bottom left: 200 observations of the training set and 200 observations of the test set; Bottom right: 400 observations of the training set and 400 observations of the test set;

# 8.4 Tests of the ratio of $l_1$ and $l_2$ norms approach

We performed in-sample and out-of-sample tests on the U.S. swaps data and tried to solve the following problem in section 5:

$$x^T A x / x^T B x + \gamma \frac{\|x\|_1}{\|x\|_2}$$

by using intermittent diffusion algorithm (ID algorithm) [4]. After each iteration of the line search algorithm [3], we fix the  $l_2$  norm to 1.

In our tests of ID algorithm, we let  $\alpha = 20$ ,  $\kappa = 20$  and N = 20. In addition, to satisfy the conditions of ID algorithm so that a global minimizer exists in a bounded set, we also add a penalty function  $p(x, \theta, \xi, \zeta)$  to f(x):

$$p(x, \theta, \xi, \zeta) = \sum_{i} u(x_i, \theta, \xi, \zeta)$$

where

$$u(x_i, \theta, \xi, \zeta) := \begin{cases} \xi(x_i - \theta)^{\zeta}, & x_i > \theta \\ 0, & |x_i| \le \theta \\ \xi(\theta - x_i)^{\zeta}, & x_i < -\theta \end{cases}$$

We set  $\theta = 10$ ,  $\xi = 2$  and  $\zeta = 100$  in our test.

For in-sample tests, we used the entire U.S. swaps data set to estimate the parameters A and B and calculated the solutions for different  $\gamma$ . We found the minimizers to the problem (13) by the algorithm 2 in section 5 for different  $\gamma$ 's between 0 and 1.5 with a step size 0.02. Then we calculated the estimated mean reversion coefficients and counted the trading opportunities of the whole period for all the minimizers. The results are shown in Figure 5. It shows the estimated mean reversion coefficients/trading opportunities versus the values of  $\gamma$  of this in-sample test on the entire swap data set.

For out-of-sample tests, we used every 100 consecutive observations to estimate those parameters and found the minimizers for different  $\gamma$ 's between 0 and 1.5 with a step size 0.02. This time, we calculated the estimated mean reversion coefficients and counted the trading opportunities for both these 100 days and the next 100 trading days. The results are shown in Figure 6. It shows the average estimated mean reversion coefficients/trading opportunities versus the values of  $\gamma$  from in-sample and out-of-sample tests.

From the numerical results, we see that as the  $\gamma$  increases the mean reversion coefficients and the trading opportunities have a decreasing trend for both in-sample and out-of-sample tests. There are big jumps in Figure 5. After checking those portfolios, we found that we only recovered portfolios of cardinality 1, 5, 6, 7 and 8. We failed to recover the portfolios of cardinality 2, 3 and 4. The reason could be that we got trapped in the local minima.

The curves in Figure 6 look smoother, since they are the average performance of portfolios on about 12 data sets. Normally, the in-sample performance is better than the out-of-sample performance. However, if we treat the in-sample performance as a benchmark, we can still maintain about 60% of the performance.

When  $\gamma$  is close to 1.5, the minimizer will be an 1-sparse vector. This shows that the ratio of  $l_1$  and  $l_2$  norms indeed enforces extreme sparsity in our problem

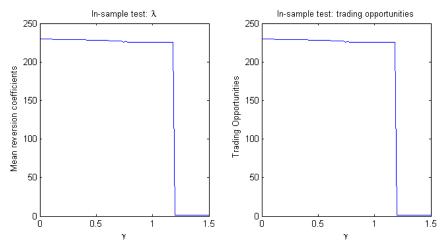


Fig. 5 In-sample tests on the whole data set of the U.S. swaps. Estimated mean reversion coefficients/trading opportunities versus the values of  $\gamma$  using the intermittent diffusion(ID) algorithm with the ratio of  $l_1$  and  $l_2$  norms penalty.

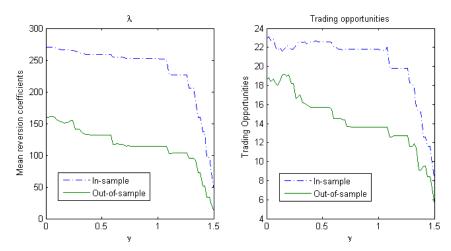


Fig. 6 In-sample and out-of-sample tests on the U.S. swaps. Average estimated mean reversion coefficients/trading opportunities versus the values of  $\gamma$  using the intermittent diffusion(ID) algorithm with the ratio of  $l_1$  and  $l_2$  norms penalty.

# 8.4.1 Comments

The advantages of the ratio of  $l_1$  and  $l_2$  norms approach are: 1. It does not require any prior knowledge of the assets; 2. It does not predetermine the cardinality. This approach also has its disadvantages: 1. It is difficult to recover a portfolio whose cardinalities are intermediate. We could encounter big jumps of the mean reversion coefficients and trading opportunities as we change the tuning parameter  $\gamma$ ; 2. The algorithm is not very efficient. It takes more than 5

minutes in solving a problem of size 100 (100 different assets for us to choose). If we compare this speed with methods of the next section, it is relatively slow.

## 8.5 Tests of the methods in section 6

## 8.5.1 In-Sample tests on 100 stocks

In this test, we applied the method in section 7 to a data set of 100 stocks. These stocks are the first 100 S&P stocks in ticker symbols' alphabetical order from our preselected list of S& P 500 stocks. Therefore, the size of the matrices A and B is  $100 \times 100$ .

When the problem is of this size, we are not able to use the exhaustive search method to get the optimal solution for the middle cardinalities. Therefore, in order to set up a criterion, we will treat the  $l^1$  contraint-free solution as a benchmark. This solution should give us the largest possible mean reversion coefficient based on the data set.

For our data set, the largest possible mean reversion coefficient is 109.97. In addition, we will set the weight of the No. 43 stock (ticker symbol: AMAT) to be 1, since it has the largest weight among all in the densest solution.

Figure 7, 8 and 9 give the numerical results. The  $l_1$  constraints, m, are in the range of 3 to 16 with a step size 1. Figure 7 demonstrates the mean reversion coefficients of portfolios built under different m's. The X-axis shows the level of m and the Y-axis is the mean reversion coefficients. Figure 7 also demonstrates the cardinality of those portfolios and Figure 8 shows the trading opportunities. Figure 9 displays the weights of 100 stocks.

The average computational cost of the method of all these tests is 4.928 seconds. This is a remarkable improvement over the other computational methods to date in terms of efficiency and speed.

## 8.5.2 Out-of-sample tests on the U.S. swaps

We also performed out-of-sample tests. We still worked on the U.S. swap data. Each time, we used 100 days data as the training data and built a sparse mean reverting portfolio. Then we estimated the mean reversion coefficients of the portfolio on these 100 days, next 50 days and next 100 days.

Figure 10 shows the numerical results of the average  $\lambda$  and Figure 11 shows the numerical results of the average trading opportunities. Both the X-axes are the  $l_1$  level. The range is from 1.5 to 2.7 with a step size 0.1. For each level, we performed several tests and the Y-axis is the average of the estimated mean reversion coefficients and the trading opportunities of all these tests.

From Figure 10, we see that as expected the in-sample performance is better. Our portfolios could maintain about 70% of the in-sample mean reversion coefficients during the 50 out-of-sample days and maintain about 65% of the in-sample mean reversion coefficients during the 100 out-of-sample days.

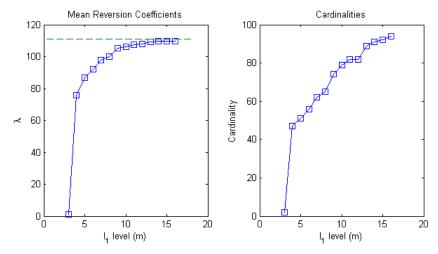


Fig. 7 Mean reversion coefficients/cardinality versus different  $l_1$  constraints using the method of the least angle regression and the difference of convex functions algorithm on 100 stocks. The horizontal line in the left plot shows the largest possible mean reversion coefficient.

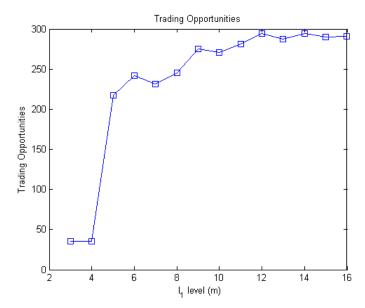


Fig. 8 Trading Opportunities versus different  $l_1$  constraints using the method of the least angle regression and DCA on 100 stocks.

In order to make a fair comparison of the trading opportunities, we performed the following counts. We counted the trading opportunities of the last 50 days of the training period and compare it with the trading opportunities of the next 50 days. We counted the trading opportunities of the last 100 days of

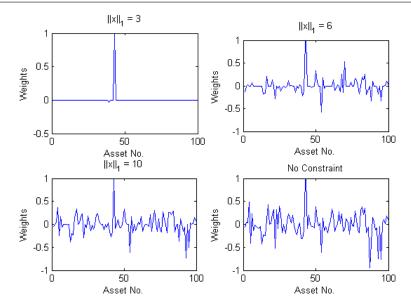


Fig. 9 Solutions under different  $l_1$  constraints using the method of the least angle regression and the difference of convex functions algorithm on 100 stocks.

the training period and compare it with the trading opportunities of the next 100 days. In this experiment, the total length of trading days are identical for in-sample and out-of-sample tests. We find that in both cases our portfolios can maintain about 75% of the in-sample trading opportunities during the out-of-sample period.

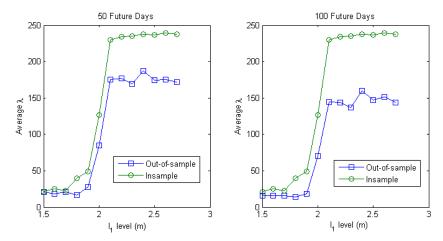


Fig. 10 In-sample and out-of-Sample tests on the U.S. swaps. Average mean reversion coefficients versus different  $l_1$  constraints using the method of the least angle regression and DCA.

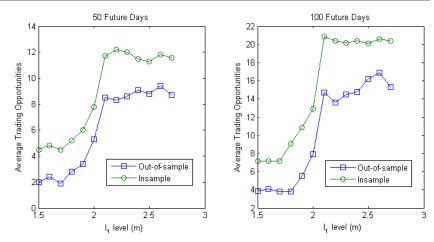


Fig. 11 In-sample and out-of-Sample tests on the U.S. swaps. Average trading opportunities versus different  $l_1$  constraints using the method of the least angle regression and DCA.

## 8.5.3 Out-of-sample tests on high dimensional synthetic data

In this section, we perform in-sample and out-of-sample tests on high dimensional synthetic data.

By presetting the matrix  $\beta$  and the noise covariance matrix  $\Sigma$  in the VAR(1) model (4), we generate a data set of size  $1000 \times 100$ . The eigenvalues of the matrix  $\beta$  lie inside the unit circle. We consider this set as the training set and estimate the matrices A and B based on this set. After picking an appropriate i, we solve for the optimal solutions under different m's (from 1.0683 to 17.0933 with a step size 1.0683) by the least angle regression and the difference of convex functions algorithm.

We also test another two special portfolios. One is a uniform portfolio which has equal weights on 100 assets. The other one is a random portfolio. We set the weight of the same ith asset to 1. For all the other assets, their weights are independently and identically distributed random variables which follows a uniform distribution in the interval [-1,1]. These two portfolios can be considered as two benchmarks. We expect that the portfolios constructed under our algorithms will beat them in out-of-sample performance.

We generate  $1000 \times 100$  observations based on the same VAR(1) model in each trial. These are test sets on which to evaluate our constructed portfolios and those two benchmark portfolios.

We generate 100 different test sets and then we compare the average of the estimated mean reversion coefficients and trading opportunities of these portfolios on both the training set and the test set. When we count the outof-sample trading opportunities, we still use the mean and standard deviation of the training set, since we are not supposed to know the future mean or variance.

The results are shown in Figure 12. The blue curves show the in-sample performance and the green curves show the out-of-sample performance. The red horizontal lines show the level of average mean reversion coefficients/trading opportunities of a random portfolio. The light blue horizontal dashed lines show the level of average mean reversion coefficients/trading opportunities of a uniform portfolio. We notice that it is very hard to maintain a high level of mean reversion coefficients when the dimension of the problem is high. They are about 22% of the in-sample levels. We believe that the estimations of Aand B will impact this performance. When the number of observations increases, we can expect that the out-of-sample performance will be better. The numbers of trading opportunities do not decrease such dramatically. They are about 55% of the in-sample trading opportunities. In fact, this is more important since the profits of convergence trading strategy come directly from those trading opportunities. The out-of-sample performance of the uniform portfolio and the random portfolio are much lower than our sparse and fast mean reverting portfolios.

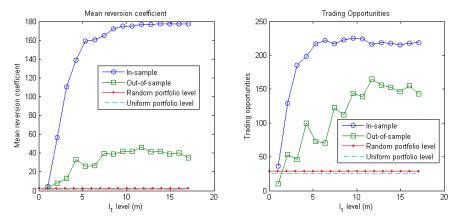


Fig. 12 In-sample and out-of-Sample tests on high dimensional synthetic data from VAR(1) model. Average mean reversion coefficients/trading opportunities versus different  $l_1$  constraints using the method of the least angle regression and the difference of convex functions algorithm. The blue curves show the in-sample performance and the green curves show the out-of-sample performance. The red horizontal line shows the level of average mean reversion coefficients/trading opportunities of a random portfolio. The light blue horizontal dashed line shows the level of average mean reversion coefficients/trading opportunities of a uniform portfolio.

## 9 Conclusion

In this work, we used a proxy of mean reversion coefficient: direct OU estimator. From numerical tests, the portfolios constructed under this proxy perform better in convergence trading than the portfolios constructed under predictability.

We developed several different types of optimization problems for building sparse mean reverting portfolios.

Without any prior knowledge of the assets, we used two  $l_1$ -based constraint/penalty to enforce the sparsity. If we used  $l_1$  norm, the problem could be considered as a simple extension of the sparse PCA. Using the algorithms in [18,26], we could solve the problem efficiently in some special cases. We also studied the properties of the ratio of  $l_1$  and  $l_2$  norms and designed an algorithm in solving the penalized optimization problem.

With prior knowledge of the assets, we adapted the  $l_1$  norm to enforce the sparsity and we simplified the problem to a nonconvex quadratic program. We developed various algorithms for approximating the global minimizer.

In our numerical tests, we applied our methods on both historical market data and synthetic data. We presented the in-sample and out-of-sample performance of the portfolios constructed under different algorithms and different problem settings. We also compared the computation costs of different algorithms for non-trivial sparse generalized eigenvalue problem (neither A not B is a diagonal matrix). The average time costs for a problem of 100-dimensions are shown in the table below.

	$l_1$ DC-PCA	$l_1/l_2$ Regularization	DC-LAR
Time costs(s)	100+	100+	4.9

Our numerical tests suggest that the combination of the least angle regression and the difference of convex functions algorithm is the best choice for non-trivial generalized eigenvalue problems. We carried out efficient computation for portfolios with hundreds of assets. The in-sample and out-of-sample performances of the mean reversion coefficients and the trading opportunities have a similar performance trend as functions of cardinality or  $l_1$  penalty level.

## Acknowledgements

We would like to thank Dr. Wuan Luo for bringing reference [6] to our attention and for helpful communication.

## A Semi-definite relaxiation method derived from $l_1/l_2$

By setting  $X = xx^T$ , then the problem (12) is equivalent to

$$\begin{aligned} \min_{X} & \operatorname{trace}(AX)/\operatorname{trace}(BX) \\ s.t. & & \frac{1^{T}|X|1}{\operatorname{trace}(X)} \leq m^{2} \\ & & rank(X) = 1 \\ & & X \succ 0 \end{aligned}$$

where |X| means we take the absolute value for each entry of X. Then after a change of variables:

$$Y = \frac{X}{\operatorname{trace}(BX)}, \quad z = \frac{1}{\operatorname{trace}(BX)}$$

and dropping the rank constraint, the previous problem can be written as a semidefinite programming problem:

$$\begin{aligned} & \underset{s.t.}{\min_{Y}} & \operatorname{trace}(AY) \\ & s.t. & & 1^{T}|Y|1 \leq m^{2}z \\ & & \operatorname{trace}(Y) - z = 0 \\ & & \operatorname{trace}(BY) = 1 \\ & & Y \succeq 0 \end{aligned} \tag{20}$$

If we set  $Card(x) = k = m^2$ , this is exactly the semidefinite relaxation in [6].

# B Estimation of the matrices A and B

For the estimations below, we assume that each column of the data matrix S represents an asset and its mean is 0. Its size is  $l \times n$ , so we have l observations and n assets.

We define  $S_c$  and  $S_f$  in the following way:

$$S_c = \begin{pmatrix} S_1^T \\ \vdots \\ S_{l-1}^T \end{pmatrix} = \begin{pmatrix} S_{11} & \dots & S_{1n} \\ \vdots & \ddots & \vdots \\ S_{l-1,1} & \dots & S_{l-1,n} \end{pmatrix}$$

$$S_f = \begin{pmatrix} S_2^T \\ \vdots \\ S_l^T \end{pmatrix} = \begin{pmatrix} S_{21} \dots S_{2n} \\ \vdots & \ddots & \vdots \\ S_{l1} \dots S_{ln} \end{pmatrix}$$

where  $S_{ti}$  is the value at time t of the ith asset.

# **B.1** Predictability

In [11], the authors discussed several methods in estimating  $\beta$  and  $\Gamma$  in VAR(1) model (4). In most cases, the number of the observations of assets values l is greater than the number of assets n. Under this case and previous assumptions, we could use the following estimates:

$$\hat{\beta} = (S_c^T S_c)^{-1} (S_c^T S_f))$$
  $\hat{\Gamma} = \frac{1}{l-1} S_c^T S_c$ 

Therefore, the matrices in problem 7 can be estimated as:

$$\hat{A} = \hat{\beta}^T \hat{\Gamma} \hat{\beta} \qquad \hat{B} = \hat{\Gamma}$$

## B.2 Direct OU estimator

Using a similar method as B.1, we estimate the matrices A and B as follows:

$$\hat{A} = \frac{1}{l-1} (S_c^T S_f + S_f^T S_c) \qquad \hat{B} = \frac{2}{l-1} S_c^T S_c$$

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