

# Article Combining Alphas via Bounded Regression

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**Abstract:** We give an explicit algorithm and source code for combining alpha streams via bounded regression. In practical applications, typically, there is insufficient history to compute a sample covariance matrix (SCM) for a large number of alphas. To compute alpha allocation weights, one then resorts to (weighted) regression over SCM principal components. Regression often produces alpha weights with insufficient diversification and/or skewed distribution against, e.g., turnover. This can be rectified by imposing bounds on alpha weights within the regression procedure. Bounded regression can also be applied to stock and other asset portfolio construction. We discuss illustrative examples.

**Keywords:** hedge fund; alpha stream; alpha weights; portfolio turnover; investment allocation; weighted regression; diversification; bounds; optimization; factor models

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# 1. Introduction

With technological advances, there is an ever-increasing number of alpha streams.<sup>1</sup> Many of these alphas are ephemeral, with relatively short lifespans. As a result, in practical applications, typically,

<sup>&</sup>lt;sup>1</sup> Here "alpha", following the common trader lingo, generally means any reasonable "expected return" that one may wish to trade on and is not necessarily the same as the "academic" alpha. In practice, often, the detailed information about how alphas are constructed may not be available, e.g., the only data available could be the position data, so "alpha" then is a set of instructions to achieve certain stock holdings by some times  $t_1, t_2, \ldots$ 

there is insufficient history to compute a sample covariance matrix (SCM) for a large number of alpha streams; SCM is singular. Therefore, directly using SCM in, say, alpha portfolio optimization is not an option.<sup>2</sup>

One approach to circumvent this difficulty is to build a factor model for alpha streams [1]. Because of the utmost secrecy in the alpha business, such factor models must be built in-house; there are no commercial providers of "standardized" factor models for alpha streams. As with factor models for equities, such model building for alphas requires certain nontrivial expertise and time expenditure.

Therefore, in practice, one often takes a simpler path. As was discussed in more detail in [1], one can deform SCM, such that it is nonsingular and, then, use the so-deformed SCM in, say, Sharpe ratio optimization for a portfolio of alphas. For small deformations, this then reduces to a cross-sectional weighted regression of the alpha stream expected returns [1]. The regression weights are the inverse sample variances of the alphas. The columns of the loadings matrix, over which the expected returns are regressed, are nothing but the first K principal components of SCM corresponding to its positive (*i.e.*, non-vanishing) eigenvalues [1].

Regression often produces alpha weights with insufficient diversification and/or skewed distribution against, e.g., turnover. Thus, if some expected returns are skewed, then, despite non-unit regression weights (which suppress more volatile alphas), the corresponding alpha weights can be larger than desired by diversification considerations. Furthermore, the principal components know nothing about quantities, such as turnover.<sup>3</sup> A simple way of obtaining a more "well-rounded" portfolio composition is to set bounds on alpha weights. This is the approach we discuss here.

When individual alpha streams are traded on separate execution platforms, the alpha weights are non-negative. By combining and trading multiple alpha streams on the same execution platform, the framework we adapt here, one saves on transaction costs by internally crossing trades between different alpha streams (as opposed to going to the market).<sup>4</sup> Then, the alpha weights can be negative.

When alpha weights can take both positive and negative values, the bounded regression problem simplifies. It boils down to an iterative algorithm that we discuss in Section 2. This algorithm can actually be derived from an optimization algorithm with bounds (in a factor model context) discussed in [2] by taking the regression limit of optimization. We also give R source code for the bounded regression algorithm in Appendix A. Appendix B contains some legalese. We conclude in Section 3, where we also discuss bounded regression with transaction costs following [3].

## 2. Bounded Regression

#### 2.1. Notations

We have N alphas  $\alpha_i$ , i = 1, ..., N. Each alpha is actually a time series  $\alpha_i(t_s)$ , s = 0, 1, ..., M, where  $t_0$  is the most recent time. Below,  $\alpha_i$  refers to  $\alpha_i(t_0)$ .

<sup>&</sup>lt;sup>2</sup> For a partial list of hedge fund literature, see, e.g., [4–23] and the references therein. For a partial list of portfolio optimization and related literature, see, e.g., [24–58] and the references therein.

<sup>&</sup>lt;sup>3</sup> One approach to rectify this is to add a turnover-based factor to the loadings matrix [1].

<sup>&</sup>lt;sup>4</sup> For a recent discussion, see [59].

Let  $C_{ij}$  be the sample covariance matrix (SCM) of the N time series  $\alpha_i(t_s)$ . If M < N, then only M eigenvalues of  $C_{ij}$  are non-zero, while the remainder have "small" values, which are zeros distorted by computational rounding.<sup>5</sup>

Alphas  $\alpha_i$  are combined with weights  $w_i$ . Any leverage is included in the definition of  $\alpha_i$ , *i.e.*, if a given alpha labeled by  $j \in \{1, ..., N\}$  before leverage is  $\alpha'_j$  (this is a raw, unlevered alpha) and the corresponding leverage is  $L_j$ : 1, then we define  $\alpha_j \equiv L_j \alpha'_j$ . With this definition, the weights satisfy the condition:

$$\sum_{i=1}^{N} |w_i| = 1$$
 (1)

Here, we allow the weights to be negative, as we are interested in the case where the alphas are traded on the same execution platform, and trades between alphas are crossed, so one is actually trading the combined alpha  $\alpha \equiv \sum_{i=1}^{N} \alpha_i w_i$ .

#### 2.2. Weighted Regression

When SCM  $C_{ij}$  is singular and no other matrix (e.g., a factor model) to replace it is available, one can deform SCM, such that it is nonsingular and then use the so-deformed SCM in, say, Sharpe ratio optimization for a portfolio of alphas [1]. For small deformations, this reduces to a cross-sectional weighted regression of the alpha stream expected returns [1]. The regression weights  $z_i$  (not to be confused with the alpha weights  $w_i$ ) are the inverse sample variances of the alphas:  $z_i \equiv 1/C_{ii}$ . The columns of the loadings matrix  $\Lambda_{iA}$ ,  $A = 1, \ldots, K$ , over which the expected returns are regressed, are nothing but the first K principal components of SCM corresponding to its positive (*i.e.*, non-vanishing) eigenvalues. However, for now, we will keep  $\Lambda_{iA}$  general (e.g., one may wish to include other risk factors in  $\Lambda_{iA}$  [1]).

The weights  $w_i$  are given by:

$$w_i = \gamma \ z_i \ \varepsilon_i \tag{2}$$

where  $\varepsilon_i$  are the residuals of the cross-sectional regression of  $\alpha_i$  over  $\Lambda_{iA}$  (without the intercept, unless the intercept is subsumed in  $\Lambda_{iA}$ , see below) with the regression weights  $z_i$ :

$$\varepsilon_i = \alpha_i - \sum_{j=1}^N z_j \,\alpha_j \sum_{A,B=1}^K \Lambda_{iA} \,\Lambda_{jB} \,Q_{AB}^{-1} \tag{3}$$

where  $Q_{AB}^{-1}$  is the inverse of:

$$Q_{AB} \equiv \sum_{i=1}^{N} z_i \Lambda_{iA} \Lambda_{iB} \tag{4}$$

and the overall factor  $\gamma$  in Equation (2) is fixed via Equation (1). Note that we have:

$$\forall A \in \{1, \dots, K\}: \quad \sum_{i=1}^{N} w_i \Lambda_{iA} = 0 \tag{5}$$

<sup>&</sup>lt;sup>5</sup> Actually, this assumes that there are no N/Asin any of the alpha time series. If some or all alpha time series contain N/As in a non-uniform manner and the correlation matrix is computed by omitting such pair-wise N/As, then the resulting correlation matrix may have negative eigenvalues that are not zeros distorted by computational rounding.

Therefore, the weights  $w_i$  are neutral w.r.t. the risk factors defined by the columns of the loadings matrix  $\Lambda_{iA}$ .

## 2.3. Bounds

Since the weights  $w_i$  can have either sign, we will assume that the lower and upper bounds on the weights:

$$w_i^- \le w_i \le w_i^+ \tag{6}$$

satisfy the conditions:

$$w_i^- \le 0 \tag{7}$$

$$w_i^+ \ge 0 \tag{8}$$

$$w_i^- < w_i^+ \tag{9}$$

The last condition is not restrictive: if for some alpha labeled by i, we have  $w_i^- = w_i^+$ , then we can simply set  $w_i = w_i^-$  and altogether exclude this alpha from the bounded regression procedure below. Furthermore, if, for whatever reason, we wish to have no upper/lower bound for a given  $w_i$ , we can simply set  $w_i^{\pm} = \pm 1$ .

The bounds can be imposed for diversification purposes: e.g., one may wish to require that no alpha has a weight greater than some fixed (small) percentile  $\xi$ , *i.e.*,  $|w_i| \leq \xi$ , so  $w_i^{\pm} = \pm \xi$ . One may also wish to suppress the contributions of high turnover alphas, e.g., by requiring that  $|w_i| \leq \tilde{\xi}$  if  $\tau_i \geq \tau_*$ , where  $\tau_i$ is the turnover,<sup>6</sup>  $\tau_*$  is some cut-off turnover, and  $\tilde{\xi}$  is some (small) percentile. Bounds can also be used to limit the weights of low capacity<sup>7</sup> alphas, *etc*.<sup>8</sup>

#### 2.4. Running a Bounded Regression

Therefore, how do we impose the bounds in the context of a regression? There are two subtleties here. First, we wish to preserve the factor neutrality property (5), which is invariant under the simultaneous rescalings  $w_i \rightarrow \zeta w_i$  (where  $\zeta$  is a constant). If we simply set some  $w_i$  to their upper or lower bounds, this generally will ruin the rescaling invariance, so the property (5) will be lost. Second, we must preserve the normalization condition (1). In fact, it is precisely this normalization condition that allows one to meaningfully set the bounds  $w_i^{\pm}$ , as the regression itself does not fix the overall normalization coefficient  $\gamma$  in Equation (2), owing to the rescaling invariance  $w_i \rightarrow \zeta w_i$ .

<sup>&</sup>lt;sup>6</sup> Here, the turnover (over a given period, e.g., daily turnover) is defined as the ratio  $\tau_i \equiv D_i/I_i$  of total dollars  $D_i$  (long plus short) traded by the alpha labeled by *i* over the corresponding total dollar holdings  $I_i$  (long plus short).

<sup>&</sup>lt;sup>7</sup> By capacity  $I_i^*$  for a given alpha, we mean the value of the investment level  $I_i$  for which the P&L  $P_i(I_i)$  is maximized (considering nonlinear effects of impact).

<sup>&</sup>lt;sup>8</sup> Since the regression we consider here is weighted with the regression weights  $z_i = 1/C_{ii}$ , this already controls exposure to alpha volatility, so imposing bounds based on volatility would make a difference only if one wishes to further suppress volatile alphas.

Here, we discuss the bounded regression algorithm. To save space, we skip the detailed derivation as it follows straightforwardly by taking the regression limit of optimization with bounds in the context of a factor model, both of which are discussed in detail in [2].<sup>9</sup>

Let us define the following subsets of the index  $i \in J \equiv \{1, ..., N\}$ :

$$w_i = w_i^+, \quad i \in J^+ \tag{10}$$

$$w_i = w_i^-, \quad i \in J^- \tag{11}$$

$$\overline{J} \equiv J^+ \cup J^- \tag{12}$$

$$\widetilde{J} \equiv J \setminus \overline{J} \tag{13}$$

Further, let:

$$\widetilde{\alpha}_i \equiv \gamma \; \alpha_i \tag{14}$$

$$y_A \equiv \sum_{i \in \widetilde{J}} z_i \, \widetilde{\alpha}_i \, \Lambda_{iA} + \sum_{i \in J^+} w_i^+ \, \Lambda_{iA} + \sum_{i \in J^-} w_i^- \, \Lambda_{iA} \tag{15}$$

where  $\gamma$  is to be determined (see below). Then, we have:

$$w_i = z_i \left( \widetilde{\alpha}_i - \sum_{A,B=1}^K \Lambda_{iA} \, \widetilde{Q}_{AB}^{-1} \, y_B \right), \quad i \in \widetilde{J}$$
(16)

$$\forall i \in J^+: \quad z_i \left( \widetilde{\alpha}_i - \sum_{A,B=1}^K \Lambda_{iA} \, \widetilde{Q}_{AB}^{-1} \, y_B \right) \ge w_i^+ \tag{17}$$

$$\forall i \in J^{-}: \quad z_{i} \left( \widetilde{\alpha}_{i} - \sum_{A,B=1}^{K} \Lambda_{iA} \, \widetilde{Q}_{AB}^{-1} \, y_{B} \right) \leq w_{i}^{-} \tag{18}$$

where  $\widetilde{Q}^{-1}$  is the inverse of the  $K \times K$  matrix  $\widetilde{Q}$ :

$$\widetilde{Q}_{AB} \equiv \sum_{i \in \widetilde{J}} z_i \Lambda_{iA} \Lambda_{iB}$$
<sup>(19)</sup>

Here, the loadings matrix  $\Lambda_{iA}$  must be such that  $\widetilde{Q}$  is invertible.<sup>10</sup> Furthermore, note that  $w_i$ ,  $i \in \widetilde{J}$  given by Equation (16) together with  $w_i = w_i^+$ ,  $i \in J^+$  and  $w_i = w_i^-$ ,  $i \in J^-$  satisfy Equation (5), as they should.

Note that, for a given value of  $\gamma$ , Equation (15) solves for  $y_A$  given  $J^+$  and  $J^-$ . On the other hand, Equations (17) and (18) determine  $J^+$  and  $J^-$  in terms of  $y_A$ . The entire system is then solved iteratively, where at the initial iteration, one takes  $\tilde{J}^{(0)} = J$ , so that  $J^{+(0)}$  and  $J^{-(0)}$  are empty. However, we still need to fix  $\gamma$ . This is done via a separate iterative procedure, which we describe below.

<sup>&</sup>lt;sup>9</sup> The regression limit of optimization essentially amounts to the limit  $\xi_i^2 \equiv \eta \ \tilde{\xi}_i^2, \eta \to 0, \ \tilde{\xi}_i^2 = \text{fixed}$ , where  $\xi_i$  is the specific (idiosyncratic) risk in the factor model with the factor loadings matrix identified with the regression loadings matrix  $\Lambda_{iA}$  (and the  $K \times K$  factor covariance matrix becomes immaterial in the regression limit); see [2] for details.

<sup>&</sup>lt;sup>10</sup> This is the case if the columns of  $\Lambda_{iA}$  are comprised of the first K principal components of SCM  $C_{ij}$  corresponding to its positive eigenvalues. However, as mentioned above, here, we keep the loadings matrix general.

Because we have two iterations, to guarantee (rapid) convergence, the  $J^{\pm}$  iteration (that is, for a given value of  $\gamma$ ) can be done as follows. Let  $\widehat{w}_i^{(s)}$  be such that:

$$\forall i \in J: \quad w_i^- \le \widehat{w}_i^{(s)} \le w_i^+ \tag{20}$$

$$\forall A \in \{1, \dots, K\}: \quad \sum_{i=1}^{N} \widehat{w}_i^{(s)} \Lambda_{iA} = 0 \tag{21}$$

At the (s + 1)-th iteration, let  $w_i^{(s+1)}$  be given by Equation (16) for  $i \in \tilde{J}^{(s)}$ , with  $w_i^{(s+1)} = w_i^{\pm}$  for  $i \in J^{\pm(s)}$ . This solution satisfies Equation (5), but may not satisfy the bounds. Let:

$$q_i \equiv w_i^{(s+1)} - \widehat{w}_i^{(s)} \tag{22}$$

$$h_i(t) \equiv \widehat{w}_i^{(s)} + t q_i, \quad t \in [0, 1]$$
 (23)

Then:

$$\widehat{w}_{i}^{(s+1)} \equiv h_{i}(t_{*}) = \widehat{w}_{i}^{(s)} + t_{*} q_{i}$$
(24)

where  $t_*$  is the maximal value of t, such that  $h_i(t)$  satisfies the bounds. We have:

$$q_i > 0: \quad p_i \equiv \min\left(w_i^{(s+1)}, w_i^+\right) \tag{25}$$

$$q_i < 0: \quad p_i \equiv \max\left(w_i^{(s+1)}, \ w_i^-\right) \tag{26}$$

$$t_* = \min\left(\frac{p_i - \widehat{w}_i^{(s)}}{q_i} \mid q_i \neq 0, \ i \in J\right)$$
(27)

Now, at each step, instead of Equations (17) and (18), we can define  $J^{\pm(s+1)}$  via:

$$\forall i \in J^{+(s+1)}: \quad \hat{w}_i^{(s+1)} = w_i^+ \tag{28}$$

$$\forall i \in J^{-(s+1)}: \quad \widehat{w}_i^{(s+1)} = w_i^-$$
(29)

where  $\widehat{w}_i^{(s+1)}$  is computed iteratively as above, and we can take  $\widehat{w}_i^{(0)} \equiv 0$  at the initial iteration. Unlike Equations (17) and (18), Equations (28) and (29) add new elements to  $J^{\pm}$  one (or a few) element(s) at each iteration.

The convergence criteria are given by:

$$J^{+(s+1)} = J^{+(s)} \tag{30}$$

$$J^{-(s+1)} = J^{-(s)} \tag{31}$$

These criteria are based on discrete quantities and are unaffected by computational (machine) precision effects. However, in practice, the equalities in Equations (28) and (29) are understood within some tolerance (or machine precision); see the R code in Appendix A. We will denote the value of  $\widehat{w}_i^{(s+1)}$  at the final iteration (for a given value of  $\gamma$ ) via  $\widetilde{w}_i$ .

Finally,  $\gamma$  is determined via another iterative procedure as follows (we use superscript *a* for the  $\gamma$  iterations to distinguish them from the superscript *s* for the  $J^{\pm}$  iterations):

$$\gamma^{(a+1)} = \frac{\gamma^{(a)}}{\sum_{i=1}^{N} \left| \widetilde{w}_{i}^{(a)} \right|}$$
(32)

where  $\widetilde{w}_i^{(a)}$  is computed as above for  $\gamma = \gamma^{(a)}$ . To achieve rapid convergence, the initial value  $\gamma^{(0)}$  can be set as follows:

$$\gamma^{(0)} = \frac{1}{\sum_{i=1}^{N} z_i \left|\varepsilon_i\right|} \tag{33}$$

where  $\varepsilon_i$  are the residuals of the weighted regression (without bounds) given by Equation (3). The convergence criterion for the  $\gamma$  iteration is given by:

$$\gamma^{(a+1)} = \gamma^{(a)} \tag{34}$$

understood within some preset computational tolerance (or machine precision).

The R code for the above algorithm with some additional explanatory documentation is given in Appendix A. Note that this code is not written to be "fancy" or optimized for speed or in any other way. Instead, its sole purpose is to illustrate the bounded regression algorithm as it is described above in a simple-to-understand fashion. Some legalese relating to this code is given in Appendix B.

#### 2.5. Application to Stock Portfolios

Above, we discussed the bounded regression algorithm in the context of computing weights for portfolios of alpha streams. However, the algorithm is quite general and, with appropriate notational identifications, can be applied to portfolios of stocks or other suitable instruments. In fact, it can also be applied outside of finance. Here, for the sake of definiteness, we will focus on stock portfolios; in fact, we will assume that they are dollar neutral, so both long and short positions are allowed.<sup>11</sup>

#### 2.5.1. Establishing Trades

Let us first discuss establishing trades, *i.e.*, we start from nil positions and establish a portfolio of N stocks. Instead of alpha streams, our index  $i \in \{1, ..., N\} \equiv J$  now labels the stocks. We will denote the desired dollar (not share) holdings via  $H_i$  and the total dollar investment (long plus short) via I:

$$I \equiv \sum_{i=1}^{N} |H_i| \tag{35}$$

Let  $w_i \equiv H_i/I$ . These are now our stock weights (analogous to the alpha weights). Then, we have the familiar normalization condition:

$$\sum_{i=1}^{N} |w_i| = 1 \tag{36}$$

However, normally, one imposes bounds on  $H_i$ , not on  $w_i$ . For example, in the case of establishing trades, one may wish to cap the positions, such that: (i) not more than a small percentile  $\xi$  of the total dollar investment I is allocated to any given stock; this is a diversification constraint; and (ii) only a

<sup>&</sup>lt;sup>11</sup> Various generalizations are possible, some more straightforward than others.

small percentile  $\tilde{\xi}$  of ADDV (average daily dollar volume)  $V_i$  is traded; this is a liquidity constraint (see below). In this case, we have the following bounds on the dollar holdings  $H_i$ :

$$H_i^- \le H_i \le H_i^+ \tag{37}$$

$$H_i^{\pm} = \pm \min\left(\xi \ I, \ \widetilde{\xi} \ V_i\right) \tag{38}$$

In this case, the upper and lower bounds are symmetrical. In some cases, such as for hard-to-borrow stocks, we may have some  $H_i^- = 0$ . In other cases, one may not wish to have a long position in some stocks, *etc.* We will only assume that  $H_i^- \leq 0$  and  $H_i^+ \geq 0$ , in line with our discussion above for the bounds on the weights, which are then given by:

$$w_i^{\pm} \equiv H_i^{\pm}/I \tag{39}$$

The final touch then is that instead of  $\alpha_i$ , one uses some expected returns  $E_i$  in the case of stocks. The rest goes through exactly as above for a suitably-chosen  $\Lambda_{iA}$ .

#### 2.5.2. Rebalancing Trades

With rebalancing trades, we have the current dollar holdings  $H_i^*$  and the desired dollar holdings  $H_i$ . In this case, one may wish to cap the positions such that: (i) not more than a small percentile  $\xi$  of the total dollar investment I is allocated to any given stock; this is the same diversification constraint as above; (ii) only a small percentile  $\xi$  of ADDV  $V_i$  is traded; this the same liquidity constraint as above; and (iii) not more than a small percentile  $\xi'$  of ADDV  $V_i$  is allocated to any given stock; this is another liquidity constraint stemming from the consideration that, if the portfolio must be liquidated swiftly (e.g., due to an unforeseen event), to mitigate liquidation costs, the positions are capped based on liquidity. Here,  $\xi'$  typically can be several times larger than  $\tilde{\xi}$ ; the portfolio can be built up in stages as long as at each stage, the bounds are satisfied. The bounds on  $H_i$  now read:

$$|H_i| \le \min\left(\xi \ I, \ \xi' \ V_i\right) \tag{40}$$

$$|H_i - H_i^*| \le \xi \, V_i \tag{41}$$

It is more convenient to rewrite these bounds in terms of the traded dollar amounts  $D_i \equiv H_i - H_i^*$ :

$$D_i^- \le D_i \le D_i^+ \tag{42}$$

$$D_i^+ = \min\left(\min\left(\xi \ I, \ \xi' \ V_i\right) - H_i^*, \ \widetilde{\xi} \ V_i\right) \ge 0 \tag{43}$$

$$D_i^- = \max\left(-\min\left(\xi \ I, \ \xi' \ V_i\right) - H_i^*, \ -\widetilde{\xi} \ V_i\right) \le 0$$
(44)

and we are assuming that  $|H_i^*| \leq \min(\xi I, \xi' V_i)$ . Furthermore, we will assume that  $H_i^*$  itself satisfies Equation (5):

$$\forall A \in \{1, \dots, K\}: \quad \sum_{i=1}^{N} H_i^* \Lambda_{iA} = 0$$
(45)

Then, the bounded regression algorithm can be straightforwardly applied to the weights  $w_i$  and  $x_i$  defined as follows:

$$w_i \equiv H_i / I \tag{46}$$

$$x_i \equiv D_i / I \tag{47}$$

In the  $J^{\pm}$  iteration, we now use  $x_i$  instead of  $w_i$ , while in the  $\gamma$  iteration, we still use  $w_i$ . Then, the rest of the algorithm goes through unchanged. Let us note, however, that the source code given in Appendix A is written with alpha weights in mind, so while it can be adapted to the case of stock portfolios in the case of establishing trades, straightforward modifications are required to accommodate rebalancing trades.

#### 2.5.3. Examples: Intraday Mean Reversion Alphas

To illustrate the use of the algorithm, we have employed it to construct portfolios for intraday mean reversion alphas with the loadings matrix  $\Lambda_{iA}$  in the following five incarnations: (i) intercept only (so K = 1); (ii) BICS (Bloomberg Industry Classification System) sectors; (iii) BICS industries; (iv) BICS sub-industries; and (v) the four style factors prc, mom, hlvand vol, of [60] plus BICS sub-industries. The regression weights are the inverse sample variances:  $z_i = 1/C_{ii}$  (see below). In Cases (ii)–(v) above, the intercept is subsumed in the loadings matrix  $\Lambda_{iA}$ . Indeed, we have  $\sum_{A \in G} \Lambda_{iA} \equiv 1$ , where G is the set of columns of  $\Lambda_{iA}$  corresponding to sectors in Case (ii), industries in Case (iii) and sub-industries in Cases (iv) and (v). Consequently, the resultant portfolios are automatically dollar neutral.

The portfolio construction and backtesting are identical to those in [61], where a more detailed discussion can be found; so, to save space, here, we will only give a brief summary. The portfolios are assumed to be established at the open and liquidated at the close on the same day, so they are purely intraday, and the algorithm of Section 2.5.1 for establishing trades applies. The expected returns  $E_i$  for each date are taken to be  $E_i = -R_i$ , where  $R_i \equiv \ln (P_i^{open}/P_i^{close})$ , and for each date,  $P_i^{open}$  is today's open, while  $P_i^{close}$  is yesterday's close adjusted for splits and dividends if the ex-date is today. Therefore, these are intraday mean-reversion alphas.

The universe is top 2000 by ADDV  $V_i$ , where ADDV is computed based on 21-trading day rolling periods. However, the universe is not rebalanced daily, but also every 21 trading days (see [61] for details). The sample variances  $C_{ii}$  are computed based on the same 21-trading day rolling periods and are not applied daily, but also, every 21 trading days, the same as the universe rebalancing (see [61] for details). We run our simulations over a period of five years (more precisely,  $252 \times 5$  trading days going back from 5 September 2014, inclusive). The annualized return-on-capital (ROC) is computed as average daily P&L divided by the total (long plus short) intraday investment level I (with no leverage) and multiplied by 252. The annualized Sharpe ratio (SR) is computed as the daily Sharpe ratio multiplied by  $\sqrt{252}$ . Cents-per-share (CPS) are computed as the total P&L divided by the total shares traded. On each day, the total (establishing plus liquidating) shares traded for each stock are given by  $Q_i = 2|H_i|/P_i^{open}$  (see [61] for details).

For comparison purposes, the results for regressions without bounds are given in Table 1. The results for the bounded regressions, with the bounds on the desired holdings set as:

$$|H_i| \le 0.01 \, V_i \tag{48}$$

so not more than 1% of each stock's ADDV is bought or sold, are given in Table 2, and the corresponding P&Ls are plotted in Figure 1. Thus, as expected, adding the liquidity bounds has the diversification effect on the portfolios, so the Sharpe ratios are substantially improved; as usual, at the expense of (slightly)

lowering paper ROC and CPS. Note that, even with tight liquidity bounds, the four style factors, prc, mom, hlv and vol, of [60] add value, further validating the four-factor model of [60].

**Table 1.** Simulation results for the 5 alphas via regression without bounds discussed in Section 2.5.3. ROC, return-on-capital; SR, Sharpe ratio; CPS, cents-per-share; BICS, Bloomberg Industry Classification System.

Alpha	ROC	SR	CPS
Regression: Intercept only	33.59%	5.59	1.38
Regression: BICS sectors	39.28%	7.05	1.61
Regression: BICS industries	42.66%	8.19	1.75
Regression: BICS sub-industries	45.25%	9.22	1.84
Regression: 4 style factors plus BICS sub-industries	46.60%	9.85	1.90



**Figure 1.** P&L graphs for the intraday alphas discussed in Section 2.5.3, with a summary in Table 2. Bottom-to-top-performing: (i) regression over intercept only; (ii) regression over BICS sectors; (iii) regression over BICS industries; (iv) regression over BICS sub-industries; and (v) regression over the four style factors, prc, mom, hlvand vol, of [60] plus BICS sub-industries. The investment level is \$10 M long plus \$10 M short.

Alpha	ROC	SR	CPS
Regression: Intercept only	29.66%	7.36	1.25
Regression: BICS sectors	35.32%	9.89	1.48
Regression: BICS industries	39.25%	12.00	1.65
Regression: BICS sub-industries	42.23%	14.13	1.75
Regression: 4 style factors plus BICS sub-industries	43.70%	15.54	1.82

**Table 2.** Simulation results for the 5 alphas via bounded regression discussed in Section 2.5.3.

#### 3. Concluding Remarks

One, but not the only, way to think about bounded regression is as an alternative to optimization with bounds when the latter is not attainable. In fact, as mentioned above, bounded regression is a zero specific risk limit of optimization with bounds in the context of a factor model. Therefore, when a factor model is not available, e.g., in the context of alpha streams, bounded regression can be used in lieu of optimization.

In this regard, one can further augment the bounded regression algorithm we discussed above by including linear transaction costs, as in [3]. A systematic approach is to start with optimization with bounds and linear transaction costs in the context of a factor model, as in [2], and to take a zero specific risk limit. Non-linear transaction costs (impact) in the context of alpha weights can be treated using the approximation discussed in [3] using the spectral model of turnover reduction [62].

#### Appendix

#### A. The R Code

Below, we give R (R Package for Statistical Computing, http://www.r-project.org) source code for the bounded regression algorithm we discuss in the main text. The entry function is calc.bounded.lm(), which runs the  $\gamma$  iteration loop and calls the function bounded.lm(), which runs the  $J^{\pm}$  iteration loop. The args () of calc.bounded.lm() are: ret, which is the N-vector of alphas  $\alpha_i$  (or, more generally, some other returns); load, which is the  $N \times K$  loadings matrix  $\Lambda_{iA}$ ; weights, which is the N-vector of the regression weights  $z_i$ ; upper, which is the N-vector of the upper bounds  $w_i^+$ ; lower, which is the N-vector of the lower bounds  $w_i^-$ ; and prec, which is the desired precision with which the output weights  $w_i$ , the N-vector of which calc.bounded.lm() returns, must satisfy the normalization condition (1). Internally, bounded.lm() calls the function calc.bounds(), which computes  $\widehat{w}_i^{(s+1)}$  in Equation (24) at each iteration. The code is straightforwardly self-explanatory. Jp, Jm in bounded.lm() correspond to  $J^{\pm}$ . One subtlety is that, when restricting  $\Lambda_{iA}$  to  $\tilde{J} \subset J$ , in the case of binary industry classification (e.g., when  $\Lambda_{iA}$  corresponds to BICS sub-industries, which can be small), the so-restricted  $\Lambda_{iA}$  may have null columns, which must be omitted, and the code below does just that. For non-binary cases, one may wish to augment the code to ensure that the matrix Q < t(load[Jt, take]) %\*% w.load[Jt, take] is nonsingular (and, if it is, then remove the culprit columns in  $\Lambda_{iA}$  or otherwise modify the latter); however, for non-binary  $\Lambda_{iA}$  and generic regression weights, this should not occur, except for special, non-generic cases.

```
calc.bounded.lm <- function(ret, load, weights, upper, lower, prec = 1e-5)</pre>
{
  reg <- lm(ret \sim -1 + load, weights = weights)
  x <- weights * residuals(reg)</pre>
  ret <- ret / sum(abs(x))</pre>
  repeat{
     x <- bounded.lm(ret, load, weights, upper, lower)</pre>
     if(abs(sum(abs(x)) - 1) < prec)
       break
    ret <- ret / sum(abs(x))</pre>
  }
  return(x)
}
bounded.lm <- function(ret, load, weights, upper, lower, tol = 1e-6)
{
  calc.bounds <- function(z, x)</pre>
  {
     q <- x - z
    p <- rep(NA, length(x))</pre>
     pp <- pmin(x, upper)</pre>
     pm <- pmax(x, lower)</pre>
     p[q > 0] <- pp[q > 0]
     p[q < 0] < - pm[q < 0]
     t <- (p - z)/q
     t <- min(t, na.rm = T)
     z <- z + t * q
     return(z)
  }
  if(!is.matrix(load))
     load <- matrix(load, length(load), 1)</pre>
  n <- nrow(load)</pre>
  k <- ncol(load)</pre>
  ret <- matrix(ret, n, 1)</pre>
```

upper <- matrix(upper, n, 1)</pre>

```
lower <- matrix(lower, n, 1)</pre>
z <- diag(weights)</pre>
w.load <- z %*% load
w.ret <- z %*% ret
J \leq -rep(T, n)
Jp <- rep(F, n)
Jm <- rep(F, n)
z < - rep(0, n)
repeat{
  Jt <- J & !Jp & !Jm
  y <- t(w.load[Jt, ]) %*% ret[Jt, ]</pre>
  if(sum(Jp) > 1)
    y <- y + t(load[Jp, ]) %*% upper[Jp, ]</pre>
  else if(sum(Jp) == 1)
     y <- y + upper[Jp, ] * matrix(load[Jp, ], k, 1)</pre>
  if(sum(Jm) > 1)
    y <- y + t(load[Jm, ]) %*% lower[Jm, ]</pre>
  else if(sum(Jm) == 1)
     y <- y + lower[Jm, ] * matrix(load[Jm, ], k, 1)</pre>
  if(k > 1)
    take <- colSums(abs(load[Jt, ])) > 0
  else
    take <- T
  Q <- t(load[Jt, take]) %*% w.load[Jt, take]</pre>
  Q \ll solve(Q)
  v <- Q %*% y[take]
  xJp <- Jp
  xJm <− Jm
  x <- w.ret - w.load[, take] %*% v</pre>
  x[Jp, ] <- upper[Jp, ]</pre>
  x[Jm, ] <- lower[Jm, ]</pre>
  z <- calc.bounds(z, x)</pre>
  Jp <- abs(z - upper) < tol</pre>
  Jm < -abs(z - lower) < tol
```

```
if(all(Jp == xJp) & all(Jm == xJm))
            break
}
return(z)
}
```

#### **B.** Disclaimers

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## **Conflicts of Interest**

The author declares no conflict of interest.

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