Using the sprintr package

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The sprintr package contains the implementations of a computationally efficient method, called sprinter, to fit large interaction models based on the reluctant interaction selection principle. The details of the method can be found in Yu, Bien, and Tibshirani (2019) *Reluctant interaction modeling*. In particular, sprinter is a multi-stage method that fits the following pairwise interaction model:

$$y = \sum_{j=1}^{p} X_j \beta_j^* + \sum_{\ell \le k} X_\ell X_k \gamma_{\ell k}^* + \varepsilon.$$

This document serves as an introduction of using the package with a simple simulated data example.

Data simulation

We consider the following simple simulation setting, where $X \sim N(\mathbf{0}, \mathbf{I}_p)$. There are two non-trivial main effects $\beta_1 = 1$, $\beta_2 = -2$, and $\beta_j = 0$ for j > 2. The two important interactions are $X_1 * X_3$ with $\gamma_{13} = 3$, and $X_4 * X_5$ with $\gamma_{45} = -4$. With $\varepsilon \sim N(0, 1)$, the following code simulates n = 100 observation from the model above with p = 200.

```
library(sprintr)
set.seed(123)
n <- 100
p <- 200
x <- matrix(data = rnorm(n * p), nrow = n, ncol = p)
y <- x[, 1] - 2 * x[, 2] + 3 * x[, 1] * x[, 3] - 4 * x[, 4] * x[, 5] + rnorm(100)</pre>
```

Using sprinter function

The function sprinter implements the sprinter method (please note that the function name sprinter is different from the package name sprintr), which involves the following three main steps:

- Fit a lasso (with cross-validation) of the response y only on main effects X (if square = FALSE) or with both main effects and squared effects (X, X^2) (if square = TRUE).
- Carry out a screening procedure based on the residual from the previous step. The number of the selected candidate interactions is specified by num_keep.
- With a path of tuning parameter lambda, fit a lasso of the response on main effects, squared effects (if square = TRUE), and selected interactions from the previous step.

There are two tuning parameters: num_keep (used in Step 2) and lambda (used in Step 3). If num_keep is not specified, it will then be set to $n/\lceil \log n \rceil$ (see, e.g., Fan & Lv (2008)). If lambda is not specified, then sprinter would compute its own path of tuning parameter values based on the number of tuning parameters (nlam) and the range of the path (lam_min_ratio).

```
mod <- sprinter(x = x, y = y, square = FALSE, nlam = 100, lam_min_ratio = 0.01)</pre>
```

The output of **sprinter** is a S3 object including several useful components. In particular, it involves a matrix idx that represents the index pairs of all variables considered in Step 3:

```
mod\$idx[(p + 1) : nrow(mod\$idx),]
          index_1 index_2
#>
               96
                       140
    [1,]
#>
    [2,]
                5
                        79
#>
    [3,]
                7
                       113
#>
    [4,]
              113
                       155
#>
    [5,]
               94
                       148
    [6,]
#>
                5
                       173
#>
    [7,]
              108
                       144
#>
    [8,]
                7
                       165
#>
   [9,]
               17
                        77
#> [10,]
              158
                       175
#> [11,]
                       102
                4
#> [12,]
               58
                       108
#> [13,]
                5
                        97
#> [14,]
               30
                       168
#> [15,]
                5
                       182
#> [16,]
                5
                        25
#> [17,]
              135
                       168
#> [18,]
              115
                       175
#> [19,]
                3
                       177
#> [20,]
                        96
                4
#> [21,]
                         3
                1
                         5
#> [22,]
```

Since Step 3 of sprinter always includes the main effects, mod\$idx[(p + 1): nrow(mod\$idx),] contains the indices of all the selected interactions from Step 2. The two columns of this output represents the index pair (ℓ, k) of a selected interaction $X_{\ell} * X_k$, where $\ell \leq k$. Note that here the last two rows are the true interactions $X_1 * X_3$ and $X_4 * X_5$. If the first entry of an index pair is zero, i.e., $(\ell = 0, k)$, then it represents a main effect X_k .

The output mod\$coef is a nrow(mod\$idx)-by-length(mod\$lambda) matrix. Each column of mod\$coef is a vector of estimate of all variable coefficients considered in Step 3 corresponding to one value of the lasso tuning parameter lambda. For example, for the 30-th tuning parameter, we have the corresponding coefficient estiamte:

```
estimate <- mod$coef[, 30]
cb <- cbind(mod$idx, estimate)</pre>
cb[cb[, 3] != 0, ]
#>
        index_1 index_2
                            estimate
#> V1
               0
                       1 0.3062352
               0
                       2 -0.4982855
#> V2
                       3 1.9034944
#> V221
               1
#> V222
                       5 -2.9489912
```

Using cross-validation with cv.sprinter

The function cv.sprinter() performs cross-validation to select the value of lasso tuning parameter lambda used in Step 3, while holding the value of num_keep fixed.

```
mod_cv <- cv.sprinter(x = x, y = y, square = FALSE, nlam = 100, lam_min_ratio = 0.01)</pre>
```

The output of cv.sprinter is a S3 object. The most intersting information is mod_cv\$compact, which is a matrix of three columns. The first two columns show the index pairs of all variables finally selected by the

lasso in Step 3, and the last column is the coefficient estimate corresponding to that selected variable.

```
mod cv$compact
         index_1 index_2 coefficient
#>
#>
    [1,]
               0
                       1 0.988300175
#>
    [2,]
               0
                       2 -1.454164327
#>
    [3,]
               0
                       4 -0.217993186
#>
   [4,]
               0
                      36 0.070174151
#>
   [5,]
               0
                      59 -0.028997635
   [6,]
               0
#>
                      66 -0.076081341
#>
   [7,]
               0
                      87 -0.015292763
#>
  [8,]
               0
                      94 0.071703217
#> [9,]
               0
                     112 0.032026436
#> [10,]
               0
                     123 -0.028004284
#> [11,]
                     157 -0.167206693
               0
#> [12,]
             158
                     175 0.025534472
#> [13,]
             141
                     168 -0.006547523
                       5 -0.155987560
#> [14,]
#> [15,]
              95
                     194 0.049738621
#> [16,]
               5
                      97 0.091145031
#> [17,]
                      29 -0.050334021
#> [18,]
              17
                      77 0.023256339
#> [19,]
              30
                     168 -0.052412220
#> [20,]
             115
                     175 -0.027360353
#> [21,]
               1
                       3 2.471482722
#> [22,]
                       5 -3.583816424
               4
```

We see (from the first two rows and the last two rows) that the fit selected by cross-validation includes all the four important variables in the model, with relatively accurate estimates of their coefficients.

Finally, there is a predict function for the S3 object returned by cv.sprinter that computes the prediction for a new data matrix of main effects:

```
newdata <- matrix(rnorm(20 * p), nrow = 20, ncol = p)
pred <- predict(mod_cv, newdata = newdata)</pre>
```