Solution 1: Multivariate Regression

Consider the multivariate regression setting on $\mathcal{X} \subset \mathbb{R}^p$ without target features, i.e., $\mathcal{Y} = \mathbb{R}$ and $\mathcal{T} = \{1, \dots, m\}$. Furthermore, consider the approach of learning a (simple) linear model $f_j(\mathbf{x}) = \mathbf{a}_j^{\top} \mathbf{x}$ for each target j independently. For this purpose, we would face the following optimization problem:

$$\min_{A} \|Y - \mathbf{X}A\|_F^2,$$

where $\|B\|_F^2 = \sqrt{\sum_{i=1}^n \sum_{j=1}^m B_{i,j}^2}$ is the Frobenius norm for a matrix $B \in \mathbb{R}^{n \times m}$ and

$$\mathbf{X} = \begin{bmatrix} (\mathbf{x}^{(1)})^{\top} \\ \vdots \\ (\mathbf{x}^{(n)})^{\top} \end{bmatrix}, \qquad A = \begin{bmatrix} \mathbf{a}_1 & \cdots & \mathbf{a}_m \end{bmatrix}, \qquad Y = \begin{bmatrix} \mathbf{y}^{(1)} \\ \vdots \\ \mathbf{y}^{(n)} \end{bmatrix}.$$

(a) Show that $\hat{A} = (\mathbf{X}^{\top}\mathbf{X})^{-1}\mathbf{X}^{\top}Y$ is the optimal solution in this case (provided that $\mathbf{X}^{\top}\mathbf{X}$ is invertible).

Solution:

Note that for a matrix $B = (\boldsymbol{b}_1 \dots \boldsymbol{b}_m) \in \mathbb{R}^{n \times m}$, it holds that

$$||B||_F^2 = \sum_{i=1}^n \sum_{j=1}^m B_{i,j}^2 = \sum_{j=1}^m \sum_{i=1}^n B_{i,j}^2 = \sum_{j=1}^m \boldsymbol{b}_j^{\mathsf{T}} \boldsymbol{b}_j.$$

Thus, the function $f(A) = ||Y - \mathbf{X}A||_F^2$ we want to minimize can be written as

$$||Y - \mathbf{X}A||_F^2 = \sum_{j=1}^m (\mathbf{y}_j - \mathbf{X}\mathbf{a}_j)^\top (\mathbf{y}_j - \mathbf{X}\mathbf{a}_j)$$
$$= \sum_{j=1}^m \mathbf{y}_j^\top \mathbf{y}_j - 2\mathbf{y}_j^\top \mathbf{X}\mathbf{a}_j + \mathbf{a}_j^\top \mathbf{X}^\top \mathbf{X}\mathbf{a}_j,$$

where y_j is the j-th column of Y. Therefore, we can write $f(A) = \sum_{j=1}^m f_j(\mathbf{a}_j)$, where

$$f_j(\mathbf{a}) = \mathbf{y}_j^{\mathsf{T}} \mathbf{y}_j - 2\mathbf{y}_j^{\mathsf{T}} \mathbf{X} \mathbf{a} + \mathbf{a}^{\mathsf{T}} \mathbf{X}^{\mathsf{T}} \mathbf{X} \mathbf{a}$$

and we can minimize each f_j separately (w.r.t. to \mathbf{a}_j). We compute the gradient of f_j and set it to $\mathbf{0}$ and solve w.r.t. \mathbf{a} :

$$\nabla f_j(\mathbf{a}) = -2\mathbf{y}_j^{\top} \mathbf{X} + 2\mathbf{X}^{\top} \mathbf{X} \mathbf{a} \stackrel{!}{=} \mathbf{0}$$

$$\Leftrightarrow \mathbf{a} = (\mathbf{X}^{\top} \mathbf{X})^{-1} \mathbf{X}^{\top} \mathbf{y}_j.$$

We check that this is indeed a minimum, by checking that the Hessian matrix is positive semi-definite: The Hessian matrix is

$$\nabla \nabla^{\top} f_i(\mathbf{a}) = 2\mathbf{X}^{\top} \mathbf{X}.$$

It is positive semi-definite, since for any $z \in \mathbb{R}^p$ it holds that

$$z^{\top}(2\mathbf{X}^{\top}\mathbf{X})z = 2z^{\top}\mathbf{X}^{\top}\mathbf{X}z = 2(\mathbf{X}z)^{\top}\mathbf{X}z = 2\|\mathbf{X}z\|_2^2 \ge 0.$$

Consequently, the minimizer of f_j is $\hat{\mathbf{a}}_j = (\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top \mathbf{y}_j$, so that the minimizer of f is

$$\hat{A} = (\hat{\mathbf{a}}_1 \dots \hat{\mathbf{a}}_m) = ((\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top \mathbf{y}_1 \dots (\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top \mathbf{y}_m) = (\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top Y.$$

In particular, the gradient of f w.r.t. matrix $A = (\mathbf{a}_1 \dots \mathbf{a}_m)$ is

$$\nabla f(A) = (\nabla f_1(\mathbf{a}_1) \dots \nabla f_m(\mathbf{a}_m))$$

$$= (-2\mathbf{y}_1^{\mathsf{T}} \mathbf{X} + 2\mathbf{X}^{\mathsf{T}} \mathbf{X} \mathbf{a}_1 \dots -2\mathbf{y}_m^{\mathsf{T}} \mathbf{X} + 2\mathbf{X}^{\mathsf{T}} \mathbf{X} \mathbf{a}_m)$$

$$= -2Y^{\mathsf{T}} \mathbf{X} + 2\mathbf{X}^{\mathsf{T}} \mathbf{X} A.$$

Hence, a gradient descent routine with (fixed) step size α for f would iterate as follows:

$$\hat{A} \leftarrow \hat{A} - 2\alpha \left(-Y^{\top} \mathbf{X} + \mathbf{X}^{\top} \mathbf{X} \hat{A} \right).$$

(b) Assume that the data $(\mathbf{x}^{(i)}, \mathbf{y}^{(i)}) \in \mathcal{X} \times \mathcal{Y}^m$ is generated according to the following statistical model

$$(y_1,\ldots,y_m) = \mathbf{y} = (\mathbf{x}^{(i)})^{\top} A^* + \boldsymbol{\epsilon}^{\top},$$

where $A^* \in \mathbb{R}^{p \times m}$ and $\epsilon \sim \mathcal{N}(\mathbf{0}, \Sigma)$. Show that the maximum likelihood estimate for A^* coincides with the estimate in (a).

Solution:

Under the statistical model it holds that $\mathbf{y}^{(i)} \mid \mathbf{x}^{(i)} \sim \mathcal{N}\left((\mathbf{x}^{(i)})^{\top} A^*, \Sigma\right)$, i.e.²,

$$p(\mathbf{y}^{(i)} \mid \mathbf{x}^{(i)}, A^*) = (2\pi)^{-m/2} |\mathbf{\Sigma}|^{-1/2} \exp\left[-\frac{1}{2}(\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A^*) \mathbf{\Sigma}^{-1} (\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A^*)^{\top}\right]$$

$$\propto \exp\left[-\frac{1}{2} \left(\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A^*\right) \mathbf{\Sigma}^{-1} \left(\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A^*\right)^{\top}\right]. \tag{1}$$

So the log-likelihood for A is

$$l(A \mid \mathcal{D}) = \log \left(\prod_{i=1}^{n} p(\mathbf{y}^{(i)} \mid \mathbf{x}^{(i)}, A) \right)$$

$$\propto \log \left(\exp \left[-\frac{1}{2} \sum_{i=1}^{n} \left(\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A \right) \mathbf{\Sigma}^{-1} \left(\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A \right)^{\top} \right] \right)$$

$$= -\sum_{i=1}^{n} \left(\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A \right) \mathbf{\Sigma}^{-1} \left(\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A \right)^{\top}.$$

So, we want to maximize the function

$$g(A) = -\sum_{i=1}^{n} \left(\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A \right) \mathbf{\Sigma}^{-1} \left(\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A \right)^{\top}$$
$$= -\sum_{i=1}^{n} \mathbf{y}^{(i)} \mathbf{\Sigma}^{-1} (\mathbf{y}^{(i)})^{\top} - 2(\mathbf{x}^{(i)})^{\top} A \mathbf{\Sigma}^{-1} (\mathbf{y}^{(i)})^{\top} + (\mathbf{x}^{(i)})^{\top} A \mathbf{\Sigma}^{-1} A^{\top} \mathbf{x}^{(i)}.$$

We compute the gradient of g and set it to $\mathbf{0}_{p\times m}$ and solve w.r.t. A:

$$\nabla g(A) = \sum_{i=1}^{n} 2\mathbf{x}^{(i)}\mathbf{y}^{(i)}\mathbf{\Sigma}^{-1} - 2\mathbf{x}^{(i)}(\mathbf{x}^{(i)})^{\top}A\mathbf{\Sigma}^{-1} \stackrel{!}{=} \mathbf{0}_{p \times m}$$

$$\Leftrightarrow \mathbf{X}^{\top}Y\mathbf{\Sigma}^{-1} - \mathbf{X}^{\top}\mathbf{X}A\mathbf{\Sigma}^{-1} \stackrel{!}{=} \mathbf{0}_{p \times m}$$

$$\Leftrightarrow A = (\mathbf{X}^{\top}\mathbf{X})^{-1}\mathbf{X}^{\top}Y.$$

where we used for computing the gradient that

$$egin{aligned} rac{\partial oldsymbol{z}^{ op} oldsymbol{B} ilde{z}}{\partial oldsymbol{B}} &= oldsymbol{z} ilde{oldsymbol{z}}^{ op}, & orall oldsymbol{z} \in \mathbb{R}^n, ilde{oldsymbol{z}} \in \mathbb{R}^m, oldsymbol{B} \in \mathbb{R}^{n imes m}, \\ rac{\partial oldsymbol{z}^{ op} oldsymbol{B} oldsymbol{C} oldsymbol{B}^{ op} ilde{z}}{\partial oldsymbol{B}} &= 2oldsymbol{z} ilde{oldsymbol{z}}^{ op} oldsymbol{B} oldsymbol{C}^{ op}, & orall oldsymbol{z} \in \mathbb{R}^m, oldsymbol{ar{z}} \in \mathbb{R}^m, oldsymbol{B} \in \mathbb{R}^{n imes m}, oldsymbol{C} \in \mathbb{R}^{n imes m}. \end{aligned}$$

Moreover, any matrix product $\mathbf{X}^{\top}Y$ can be written as the sum of outer products of the column and row vectors: $\sum_{i=1}^{n} \mathbf{x}^{(i)} \mathbf{y}^{(i)}$.

Thus, the MLE coincides with the OLS in (a).

 $^{^1\}mathrm{Of}$ course, in an iid fashion and the \mathbf{x} 's are independent of the $\boldsymbol{\epsilon}$'s.

²Note that $\mathbf{y}^{(i)} - (\mathbf{x}^{(i)})^{\top} A^*$ is a row vector.

- (c) Write a function implementing a gradient descent routine for the optimization of this linear model.
- (d) Run a small simulation study by creating 20 data sets as indicated below and test different step sizes α (fixed across iterations) against each other and against the state-of-the-art routine for linear models (in R, using the function lm, in Python, e.g., sklearn.linear_model.LinearRegression).
 - Compare the difference in the estimated parameter matrices \hat{A} using the mean squared error, i.e.,

$$\frac{1}{m \cdot p} \sum_{i=1}^{p} \sum_{j=1}^{m} (\mathbf{a}_{i,j}^* - \hat{\mathbf{a}}_{i,j})^2$$

and summarize the difference over all 100 simulation repetitions.

• Compare the estimation also with the James-Stein estimate of A^* , which is given by

$$A^{JS} = \left(\mathbf{a}_1^{JS} \dots \mathbf{a}_m^{JS}\right),\,$$

where

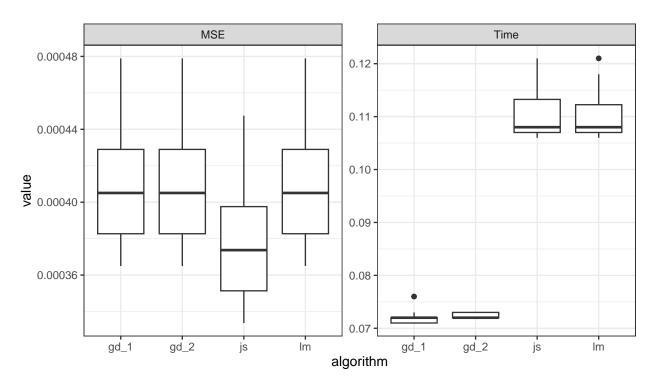
$$\mathbf{a}_{j}^{JS} = \left(1 - \frac{(m-2)\sigma^{2}}{n\|\hat{\mathbf{a}}_{j} - \mathbf{a}_{j}^{*}\|_{2}^{2}}\right)(\hat{\mathbf{a}}_{j} - \mathbf{a}_{j}^{*}) + \mathbf{a}_{j}^{*}, \quad j = 1, \dots, m.$$

and $\hat{\mathbf{a}}_{j}$ is the MLE for the jth target parameter.

```
#' Oparam step_size the step_size in each iteration
#' @param X the feature input matrix X
#' @param Y the score matrix Y
#' @param A the parameter matrix
#' Oparam eps a small constant measuring the changes in each update step.
#' Stop the algorithm if the estimated model parameters do not change
#' more than \code{eps}.
#' @return a set of optimal parameter matrix A
gradient_descent <- function(step_size, X, Y, A = matrix(rep(0,ncol(X)*m),ncol=m),</pre>
                              eps = 1e-8){
  change <- 1 # something larger eps
 XtX <- crossprod(X)</pre>
 XtY <- crossprod(X,Y)</pre>
  while(sum(abs(change)) > eps){
    # Use standard gradient descent:
    change <- + step_size * (XtY - XtX%*%A)
    # update A in the end
    A <- A + change
 return(A)
# make it all reproducible
set.seed(123)
# settings
n <- 10000
p <- 100
```

```
m <- 6
nr_sims <- 20
# define mse
mse <- function(x,y) mean((x-y)^2)</pre>
# create data (only once)
X <- matrix(rnorm(n*p), ncol=p)</pre>
A_truth <- matrix(runif(p*m, -2, 2),ncol=m)
f_truth <- X%*%A_truth
# create result object
result_list <- vector("list", nr_sims)</pre>
js_estimate <- function(A,A_truth){</pre>
A_JS = A
         for(i in 1:ncol(A)){
                  A_JS[,i] = (1-4*(m-2)/(n*sum((A[,i]-A_truth[,i])^2)))*
                    (A[,i]-A_truth[,i])+A_truth[,i]
A_JS
for(sim_nr in seq_len(nr_sims))
  # create response
  Y \leftarrow f_{truth} + rnorm(n*m, sd = 2)
  time_lm <- system.time(</pre>
    coef_{lm} \leftarrow coef(lm(Y^-1+X))
  )["elapsed"]
  time_js <- system.time(</pre>
    coef_js <- js_estimate(coef_lm,A_truth)</pre>
  )["elapsed"]
  time_js = time_js + time_lm
  time_gd_1 <- system.time(</pre>
    coef_gd_1 <- gradient_descent(step_size = 0.0001, X = X, Y = Y)</pre>
  )["elapsed"]
  time_gd_2 <- system.time(</pre>
    coef_gd_2 <- gradient_descent(step_size = 0.00005, X = X, Y = Y)</pre>
  )["elapsed"]
  mse_lm <- mse(coef_lm, A_truth)</pre>
  mse_js <- mse(coef_js, A_truth)</pre>
  mse_gd_1 <- mse(coef_gd_1, A_truth)</pre>
  mse_gd_2 <- mse(coef_gd_2, A_truth)</pre>
  # save results in list (performance, time)
  result_list[[sim_nr]] <- data.frame(mse_lm = mse_lm,</pre>
                                                        mse_js = mse_js,
```

```
mse_gd_1 = mse_gd_1,
                                      mse_gd_2 = mse_gd_2,
                                       time_lm = time_lm,
                                       time_js = time_js,
                                       time_gd_1 = time_gd_1,
                                       time_gd_2 = time_gd_2
library(ggplot2)
library(dplyr)
library(tidyr)
do.call("rbind", result_list) %>%
  gather() %>%
  mutate(what = ifelse(grepl("mse", key), "MSE", "Time"),
         algorithm = gsub("(mse|time))\_(.*)","\2", key)) %>%
  ggplot(aes(x = algorithm, y = value)) +
  geom_boxplot() + theme_bw() +
  facet_wrap(~ what, scales = "free")
```



Solution 2: Conditional Random Fields vs. Structured SVMs

Similar to probabilistic classifier chains, conditional random fields try to model the conditional distribution $\mathbb{P}(\mathbf{y} \mid \mathbf{x})$ by means of

$$\pi(\mathbf{x}, \mathbf{y}) = \frac{\exp(s(\mathbf{x}, \mathbf{y}))}{\sum_{\mathbf{y}' \in \mathcal{Y}^m} \exp(s(\mathbf{x}, \mathbf{y}'))},$$

where $x \in \mathcal{X}$ and $\mathbf{y} \in \mathcal{Y}$ with \mathcal{Y} being a finite set (e.g., multi-label classification), and $s : \mathcal{X} \times \mathcal{Y} \to \mathbb{R}$ being a scoring function. Training of a conditional random field is based on (regularized) empirical risk minimization using the negative log-loss:

$$\ell_{log}(\mathbf{x}, \mathbf{y}, s) = \log \left(\sum_{\mathbf{y}' \in \mathcal{Y}^m} \exp(s(\mathbf{x}, \mathbf{y}')) \right) - s(\mathbf{x}, \mathbf{y}).$$

$$h(\mathbf{x}) = \arg\max_{\mathbf{v} \in \mathcal{V}^m} s(\mathbf{x}, \mathbf{y}). \tag{2}$$

Structured Support Vector Machines (Structured SVMs) are also using scoring functions for the prediction, but use the structured hinge loss for the (regularized) empirical risk minimization approach:

$$\ell_{shinge}(\mathbf{x}, \mathbf{y}, s) = \max_{\mathbf{y}' \in \mathcal{Y}^m} (\ell(\mathbf{y}, \mathbf{y}') + s(\mathbf{x}, \mathbf{y}') - s(\mathbf{x}, \mathbf{y})),$$

where $\ell: \mathcal{Y}^m \times \mathcal{Y}^m \to \mathbb{R}$ is some target loss function (e.g., Hamming loss or subset 0/1 loss).

Show that if we use scoring functions s of the form

$$s(\mathbf{x}, \mathbf{y}) = \sum_{j=1}^{m} s_j(\mathbf{x}, y_j),$$

where $s_j: \mathcal{X} \times \mathcal{Y} \to \mathbb{R}$ are scoring functions for the j-th target, then

(a) conditional random fields are very well suited to model the case, where the distributions of the targets y_1, \ldots, y_m are conditionally independent,

Solution:

The idea of conditional random fields is to model the joint conditional distribution $\mathbb{P}(\mathbf{y} \mid \mathbf{x})$ by means of $\pi(\mathbf{x}, \mathbf{y})$. Thus, it should hold $\mathbb{P}(\mathbf{y} \mid \mathbf{x}) \approx \pi(\mathbf{x}, \mathbf{y})$ and with this,

$$\mathbb{P}(\mathbf{y} \mid \mathbf{x}) \approx \pi(\mathbf{x}, \mathbf{y})$$

$$= \frac{\exp(s(\mathbf{x}, \mathbf{y}))}{\sum_{\mathbf{y}' \in \mathcal{Y}^m} \exp(s(\mathbf{x}, \mathbf{y}'))}$$

$$= \frac{\exp\left(\sum_{j=1}^m s_j(\mathbf{x}, y_j)\right)}{\sum_{\mathbf{y}' \in \mathcal{Y}^m} \exp\left(\sum_{j=1}^m s_j(\mathbf{x}, y_j')\right)}$$

$$= \frac{\prod_{j=1}^m \exp(s_j(\mathbf{x}, y_j))}{\sum_{\mathbf{y}' \in \mathcal{Y}^m} \prod_{j=1}^m \exp(s_j(\mathbf{x}, y_j))}$$

$$= \frac{\prod_{j=1}^m \exp(s_j(\mathbf{x}, y_j))}{\prod_{j=1}^m \sum_{y_j' \in \mathcal{Y}} \exp(s_j(\mathbf{x}, y_j'))}$$

$$= \prod_{j=1}^m \underbrace{\exp(s_j(\mathbf{x}, y_j))}_{\sum_{y_j' \in \mathcal{Y}} \exp(s_j(\mathbf{x}, y_j'))}.$$

$$= \frac{\exp(s_j(\mathbf{x}, y_j))}{\sum_{j=1}^m \exp(s_j(\mathbf{x}, y_j))}.$$

So, if $\pi_j(\mathbf{x}, y_j)$ is interpreted as a model for the marginal conditional distribution $\mathbb{P}(y_j \mid \mathbf{x})$, we see from above

$$\mathbb{P}(\mathbf{y} \mid \mathbf{x}) \approx \prod_{j=1}^{m} \mathbb{P}(y_j \mid \mathbf{x}),$$

i.e., the targets are conditionally independent.

(b) the structured hinge loss corresponds to the multiclass hinge loss for the targets if we use the (non-averaged) Hamming loss for $\ell(\mathbf{y}, \mathbf{y}') = \sum_{j=1}^{m} \mathbb{1}_{[y_j \neq y'_j]}$, i.e.,

$$\ell_{shinge}(\mathbf{x}, \mathbf{y}, s) = \sum_{j=1}^{m} \max_{y'_j \in \mathcal{Y}} \left(\mathbb{1}_{[y_j \neq y'_j]} + s_j(\mathbf{x}, y'_j) - s_j(\mathbf{x}, y_j) \right).$$

Solution:

This can be seen immediately from the definition:

$$\ell_{shinge}(\mathbf{x}, \mathbf{y}, s) = \max_{\mathbf{y}' \in \mathcal{Y}^m} \left(\ell(\mathbf{y}, \mathbf{y}') + s(\mathbf{x}, \mathbf{y}') - s(\mathbf{x}, \mathbf{y}) \right)$$

$$= \max_{\mathbf{y}' \in \mathcal{Y}^m} \left(\sum_{j=1}^m \mathbb{1}_{[y_j \neq y_j']} + s(\mathbf{x}, \mathbf{y}') - s(\mathbf{x}, \mathbf{y}) \right)$$

$$= \max_{\mathbf{y}' \in \mathcal{Y}^m} \left(\sum_{j=1}^m \mathbb{1}_{[y_j \neq y_j']} + \sum_{j=1}^m s_j(\mathbf{x}, y_j') - \sum_{j=1}^m s_j(\mathbf{x}, y_j) \right)$$

$$= \max_{\mathbf{y}' \in \mathcal{Y}^m} \left(\sum_{j=1}^m \mathbb{1}_{[y_j \neq y_j']} + s_j(\mathbf{x}, y_j') - s_j(\mathbf{x}, y_j) \right)$$

$$= \sum_{j=1}^m \max_{y_j' \in \mathcal{Y}} \left(\mathbb{1}_{[y_j \neq y_j']} + s_j(\mathbf{x}, y_j') - s_j(\mathbf{x}, y_j) \right).$$
 (Summands are independent.)