

Linear Classification

Shyue Ping Ong

University of California, San Diego

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Overview

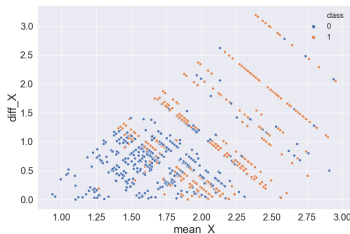
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Preliminaries

- Linear methods can also be used for classification, i.e., decision boundaries are linear.
- These methods are surprisingly effective across a large spectrum of datasets, even compared to more complex ML models.

Metal vs Insulator Dataset

- To demonstrate the use of these methods, we will first discuss the “toy” dataset.
- 2000+ binary (A_xB_y) compounds with experimental band gaps.
- Class 0: metals; Class 1: insulators.
- Using pymatgen, we can generate some simple features. Here, we will create simply features based on the mean and absolute difference in electronegativity between A and B (why?).



Creating the features and classes

```
import pandas as pd
from pymatgen.core import Composition
binaries = pd.read_csv('binary_band_gap.csv')
# We create a column holding the Composition object.
# Note the use of list comprehension in Python.
binaries['composition'] = [Composition(c) for c in binaries['Formula']]
electronegs = [[el.X for el in c] for c in binaries['composition']]
# Create the features mean and difference between electronegativities
binaries['mean_X'] = [np.mean(e) for e in electronegs]
binaries['diff_X'] = [max(e) - min(e) for e in electronegs]
# Label metals (band gap of 0. 1e-5 is used as numerical tolerance) as class 0
# Insulators are labelled as class 1.
binaries['class'] = [0 if eg < 1e-5 else 1 for eg in binaries['Eg (eV)']]
```

Basic concepts

- If there are K classes, we have a $N \times K$ indicator response matrix. Each row is a vector $Y = (Y_1, Y_2, \dots, Y_K)$ where $Y_k = 1$ if the instance belongs to the k th class and all other Y s are 0.

$$\mathbf{Y} = \begin{pmatrix} 0 & 0 & \dots & 1 \\ 1 & 0 & \dots & 0 \\ \dots & & & \\ 0 & 1 & \dots & 0 \end{pmatrix}$$

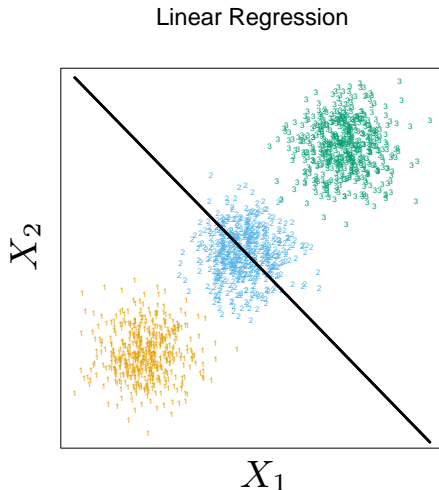
- For the k th response variable, the fitted $\hat{f}_k(x) = \hat{\beta}_{k0} + \hat{\beta}_k^T x$.
- Decision boundary between k and l class is given by $\hat{f}_k(x) = \hat{f}_l(x)$.
- Input is divided into regions.
- Similar to linear regression, we can augment the input space with polynomial (e.g., $X_1^2, X_2^s, X_1 X_2$) and other basis functions, leading to boundaries that are non-linear.

Linear regression of indicator matrix

- Treat each column of \mathbf{Y} as a target. Least squares solution:

$$\hat{\mathbf{Y}} = \mathbf{X}(\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{Y}$$

- For each new observation x , we compute $\hat{f}_k(x) = (1, x^T)(\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{Y}$.
- Find the largest component, and that will result in the classification k ,
 $G(x) = \operatorname{argmax}_{k \in G} \hat{f}_k(x)$.
- Major issue: some categories may be masked for $K \geq 3$.



Discriminant Analysis

- From Bayes rule, we have:

$$P(G = k|X = x) = \frac{f_k(x)\pi_k}{\sum_{l=1}^K f_l(x)\pi_l}$$

- where $f_k(x)$ are the class conditional probability densities ($P(X = x|G = k)$) and π_k are the prior probabilities of being in class k .
- Most common approach - assume Gaussian class densities.

$$f_k(x) = \frac{1}{(2\pi)^{p/2}|\Sigma_k|^{1/2}} \exp -\frac{1}{2}(x - \mu_k)^T \Sigma_k^{-1}(x - \mu_k)$$

Linear Discriminant Analysis

- Assume all classes have a common covariance matrix, i.e., $\Sigma_k = \Sigma$.
- To compare two classes k and l , we can compare the log ratios.

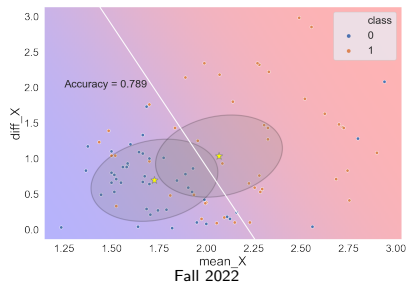
$$\begin{aligned} \log \frac{P(G = k|X = x)}{P(G = l|X = x)} &= \log \frac{f_k(x)}{f_l(x)} + \log \frac{\pi_k}{\pi_l} \\ &= \log \frac{\pi_k}{\pi_l} - \frac{1}{2}(\mu_k + \mu_l)^T \Sigma^{-1}(\mu_k - \mu_l) \\ &\quad + x^T \Sigma^{-1}(\mu_k - \mu_l) \end{aligned}$$

- At the decision boundary, $P(G = k|X = x) = P(G = l|X = x)$, which leads to a linear equation in x .
- Equivalently, we have

$$G(x) = \operatorname{argmax}_k \left\{ \log \pi_k - \frac{1}{2} \mu_k^T \Sigma^{-1} \mu_k + x^T \Sigma^{-1} \mu_k \right\}$$

Linear Discriminant Analysis, contd.

- In general, we do not know the prior distributions and covariance matrix. These are estimated from the data.
 - $\hat{\pi}_k = N_k / N$
 - $\hat{\mu}_k = \sum_{g_i=k} x_i / N$
 - $\hat{\Sigma} = \sum_{k=1}^K \sum_{g_i=k} (x_i - \hat{\mu}_k)^T (x_i - \hat{\mu}_k) / (N - K)$
- Avoids masking problem of linear regression classification.
- For the example data,

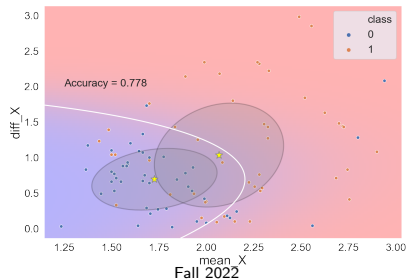


Quadratic Discriminant Analysis

- Covariances are not assumed equal.

$$G(x) = \operatorname{argmax}_k \left\{ \log \pi_k - \frac{1}{2} (x - \mu_k)^T \Sigma_k^{-1} (x - \mu_k) - \frac{1}{2} \log |\Sigma_k| \right\}$$

- No cancellation of terms and decision boundaries are quadratic.
- Covariances must be estimated for each category.
- For the same metal-insulator example,



Discriminant analysis in scikit-learn

```
from sklearn.discriminant_analysis import LinearDiscriminantAnalysis,  
    QuadraticDiscriminantAnalysis  
lda = LinearDiscriminantAnalysis(solver="svd", store_covariance=True)  
X = binaries[["mean_X", "diff_X"]]  
y = binaries["class"]  
model = lda.fit(X, y)  
y_pred = model.predict(X)  
  
qda = QuadraticDiscriminantAnalysis(store_covariance=True)  
y_pred = qda.fit(X, y).predict(X)  
\end{minted}
```

Logistic regression

- Model posterior probabilities with linear function.

$$\log \frac{P(G = 1|X = x)}{P(G = K|X = x)} = \beta_{10} + \beta_1^T x$$

$$\log \frac{P(G = 2|X = x)}{P(G = K|X = x)} = \beta_{20} + \beta_2^T x$$

...

$$\log \frac{P(G = K - 1|X = x)}{P(G = K|X = x)} = \beta_{(k-1)0} + \beta_{k-1}^T x$$

- Results in the following posterior probabilities:

$$P(G = 1|X = x) = \frac{\exp(\beta_{10} + \beta_1^T x)}{1 + \sum_{l=1}^{K-1} \exp(\beta_{l0} + \beta_l^T x)}$$

$$P(G = K|X = x) = \frac{1}{1 + \sum_{l=1}^{K-1} \exp(\beta_{l0} + \beta_l^T x)}$$

Solving for the Logistic Regression Coefficients

- Typically fitted using *maximum likelihood*.

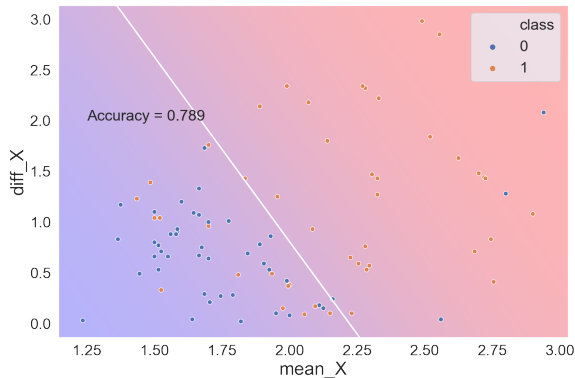
$$l(\beta) = \sum_{i=1}^N \log P(G = k|X = x_i; \beta)$$

- Differentiation and setting $\frac{\partial l}{\partial \beta} = 0$ leads to equations that are non-linear in β .
- These equations are solved using some optimization algorithm (e.g., Newton-Raphson, BFGS, etc.).

Logistic regression on metal/insulator dataset

```
from sklearn.linear_model import LogisticRegression
```

```
clf = LogisticRegression(penalty="none", random_state=0)  
model = clf.fit(X, y)  
y_pred = model.predict(X)
```



Loss functions for binary classification

- Consider a simple binary classification with two levels $(-1, 1)$. The decision boundary is at 0.
- Using the square error does not make sense, since we only care about whether it is > 0 or < 0 .
- *Margin* $yf(x)$ is positive when prediction and actual value is in the same class, and negative if they are in opposite classes.
- Need a loss that penalizes negative values much more than positive values for margins, i.e., monotone decreasing function.
- Exponential loss: $L(y, f(x)) = e^{-yf(x)}$
- Binomial/multinomial loss (can be used for K-classes):

$$L(y, p(x)) = - \sum_{k=1}^K I(y = G_k) f_k(x) + \log \left(\sum_{l=1}^K e^{f_l(x)} \right)$$

Loss functions for binary classification

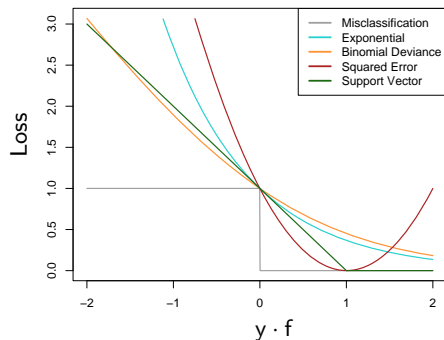


Figure: Loss functions for binary classification. Response: $y = \pm 1$. X-axis is the margin $y \cdot f$.
 Misclassification : $I(\text{sign}(f) \neq y)$; exponential: e^{-yf} ; binomial deviance: $\log(1 + e^{-2yf})$; squared error: $(y - f)^2$; and support vector: $(1 - yf)_+$. Source: [?]

The End