## Lab 2

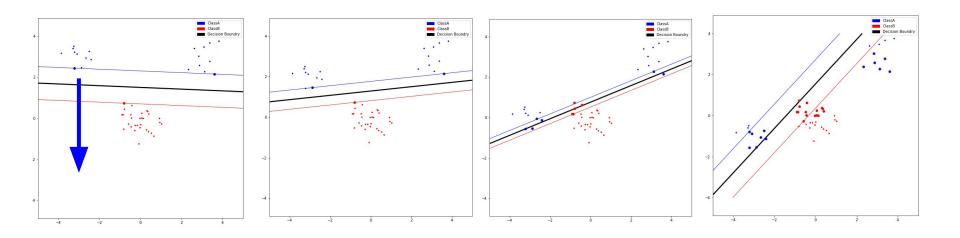
**Quick Summary** 

#### **Outline**

- Move the clusters around and change their sizes to make it easier or harder for the classifier to find a decent boundary. Pay attention to when the optimizer (minimize function) is not able to find a solution at all.
- 2. Implement the two non-linear kernels. You should be able to classify very hard data sets with these.
- 3. The non-linear kernels have parameters; explore how they influence the decision boundary. Reason about this in terms of the biasvariance trade-off.
- 4. Explore the role of the slack parameter C. What happens for very large/small values?
- 5. Imagine that you are given data that is not easily separable. When should you opt for more slack rather than going for a more complex model (kernel) and vice versa?

#### Linearly separable.

#### Not linearly separable.



## Non-linear Kernels

### **RBF Kernel**

$$\mathcal{K}(\vec{x}, \vec{y}) = e^{-\frac{||\vec{x} - \vec{y}||^2}{2\sigma^2}}$$

As sigma is increases, the function smoothness increases

Variance decreases as well, but the bias will be higher

$$\sigma = 1$$

$$\sigma = 2$$

$$\sigma = 3$$

$$\sigma = 5$$

$$\frac{1}{2}$$

$$\frac{$$

## Poly Kernel

$$\mathcal{K}(\vec{x}, \vec{y}) = (\vec{x}^T \cdot \vec{y} + 1)^p$$

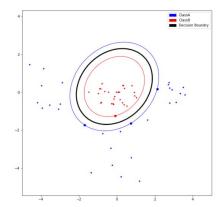
Complexity will increase with a higher order regression polynomial (p)

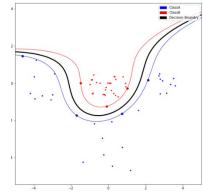
Variance increases and bias decreases when p is higher

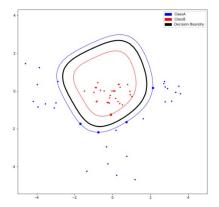
$$p = 2$$

$$p = 3$$

$$p = 4$$





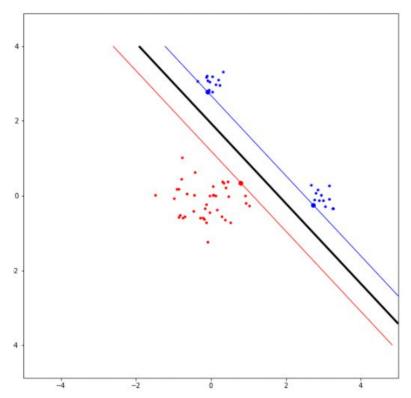


# Compare C

$$C = 100$$

$$\min_{\vec{w},b,\vec{\xi}} ||\vec{w}|| + C \sum_{i} \xi_{i}$$

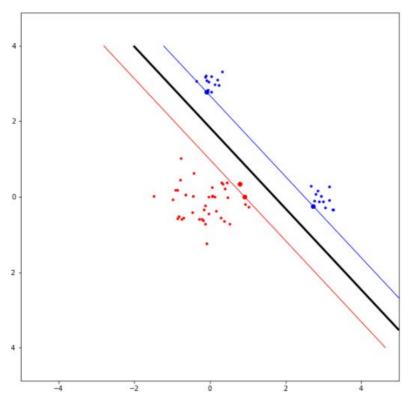
```
#Generate Data
num_points = 40
#Class A1
std_a1 = 0.2
x_a1 = 0.0
y_a1 = 3.0
#Class A2
std_a2 = 0.2
x_a2 = 3.0
y_a2 = 0.0
# #Class A3
# std a3 = 0.5
# x a3 = 0.0
# y a3 = -3.0
#Class B
std_b = 0.8
x_b = 0.0
y_b = 0.0
```



$$C = 1$$

$$\min_{\vec{w}, b, \vec{\xi}} ||\vec{w}|| + C \sum_{i} \xi_{i}$$

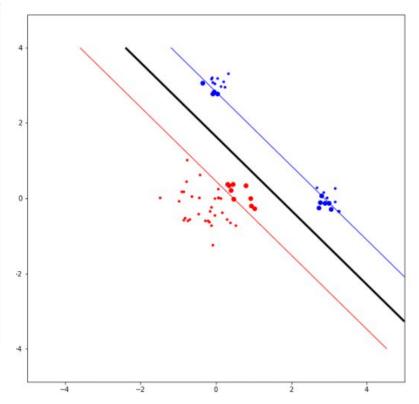
```
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#Class A2
std_a2 = 0.2
x_a2 = 3.0
y_a2 = 0.0
# #Class A3
# std a3 = 0.5
# x a3 = 0.0
# y a3 = -3.0
#Class B
std_b = 0.8
x_b = 0.0
y_b = 0.0
```



$$C = 0.1$$

$$\min_{\vec{w},b,\vec{\xi}} ||\vec{w}|| + C \sum_{i} \xi_{i}$$

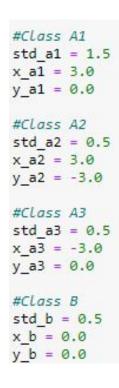
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# std a3 = 0.5
# x a3 = 0.0
# y a3 = -3.0
#Class B
std_b = 0.8
x_b = 0.0
y_b = 0.0
```

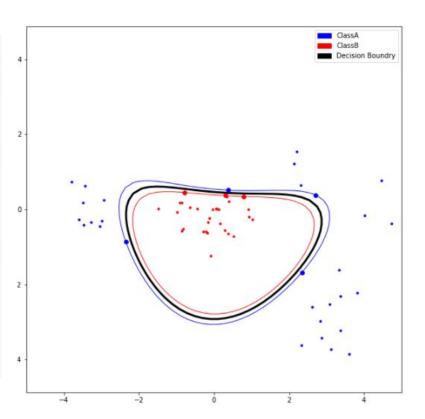


## **RBF Kernel**

$$C = 1000$$

$$\min_{\vec{w},b,\vec{\xi}} ||\vec{w}|| + C \sum_{i} \xi_{i}$$

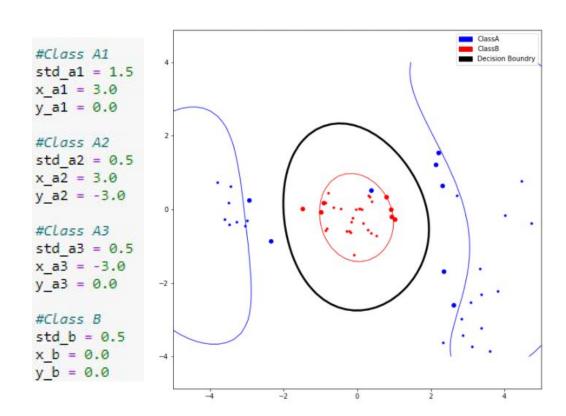




## **RBF Kernel**

$$C = 1$$

$$\min_{\vec{w},b,\vec{\xi}} ||\vec{w}|| + C \sum_{i} \xi_{i}$$



When to address Slack versus Kernels?

It depends on the data set as well.

Opting for more slack is a good idea when there is noise in the data, but the classes are still **linearly separable**. More slack will lead to higher variance and lower bias.

Otherwise, more complex data sets may require more complex model for better return on bias and variance.

