h1

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Generate a dataset of two-dimensional points, and choose a random line in the plane as your target function f, where one side of the line maps to +1 and the other side to -1. Let the inputs $\mathbf{x}_n \in \mathbb{R}^2$ be random points in the plane, and evaluate the target function f on each \mathbf{x}_n to get the corresponding output $y_n = f(\mathbf{x}_n)$. Experiment with the perceptron algorithm in the following settings:

```
[1]: import numpy as np
import matplotlib.pyplot as plt
#import matplotlib as mpl
#mpl.rcParams['figure.dpi'] = 300
%config InlineBackend.figure_format = 'retina'

pColor = 'b'
nColor = 'r'
```

```
[2]: def test_convergence(w, x, y):
    for x_n, y_n in zip(x, y):
        x_n = np.insert(x_n, 0, 1)
        if y_n > 0 and np.dot(w, x_n) < 0:
            return False

    if y_n < 0 and np.dot(w, x_n) >= 0:
        return False

    return True
```

1 a

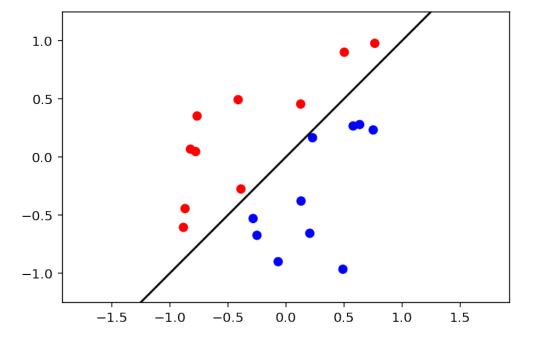
Generate a dataset of size 20. Plot the examples $\{(\mathbf{x}_n, y_n)\}$ as well as the target function f on a plane.

```
[3]: # generate a dataset of points in 2D space from -1 to 1
N = 20
x = (np.random.rand(N,2) * 2) - 1
```

```
[4]: # create a "random" target function
w_t = [0, 0.5, -0.5]
def target(x_n):
    return 1 if np.dot(w_t,np.insert(x_n, 0, 1)) >= 0 else -1
```

```
[5]: # create vector of y values (labels) with target function
y = [target(x_n) for x_n in x]
```

[6]: [-1.25, 1.25, -1.25, 1.25]



2 b

Run the perceptron algorithm on the dataset. Report the number of updates that the algorithm takes before converging. Plot the examples $\{\{(\mathbf{x}_n,y_n)\},$ the target function f, and the final hypothesis g in the same figure.

```
[7]: # randomally initialize weights
w = np.zeros(3)
```

```
while not test_convergence(w, x, y):
    # pick a random sample
    n = np.random.randint(0,x.shape[0]-1)
    x_n = np.insert(x[n,:], 0, 1)
    y_n = y[n]

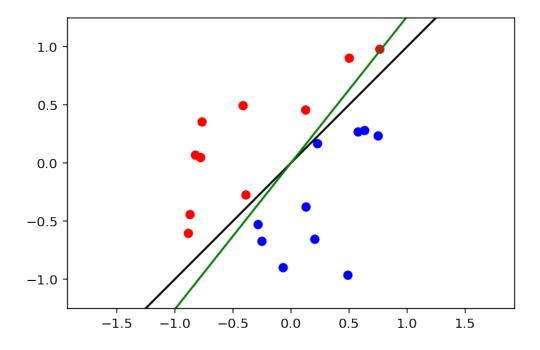
if y_n > 0 and np.dot(w, x_n) < 0:
    w += x_n

if y_n < 0 and np.dot(w, x_n) >= 0:
    w -= x_n

idx += 1

print(f"Converged in {idx} steps with w = {w}")
```

Converged in 17 steps with w = [0. 1.43335106 -1.13837864]



3 c

Repeat everything in b) with another randomly generated dataset of size 20, and compare the result to b).

```
[9]: # generate a dataset of points in 2D space from -1 to 1
N = 20
x = (np.random.rand(N,2) * 2) - 1

# create vector of y values (labels) with target function
y = [target(x_n) for x_n in x]

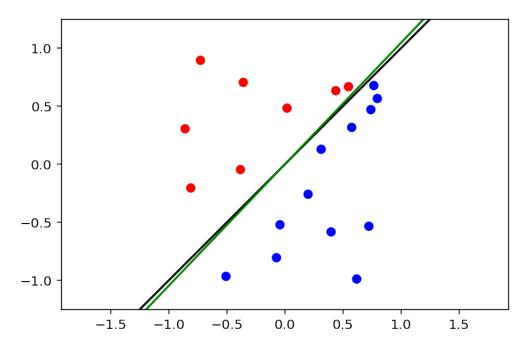
# randomally initialize weights
w = np.zeros(3)
idx = 0

while not test_convergence(w, x, y):
    # pick a random sample
    n = np.random.randint(0,x.shape[0]-1)
    x_n = np.insert(x[n,:], 0, 1)
    y_n = y[n]

if y_n > 0 and np.dot(w, x_n) < 0:
    w += x_n</pre>
```

Converged in 53 steps with w = [0.

1.22477878 -1.17061912]

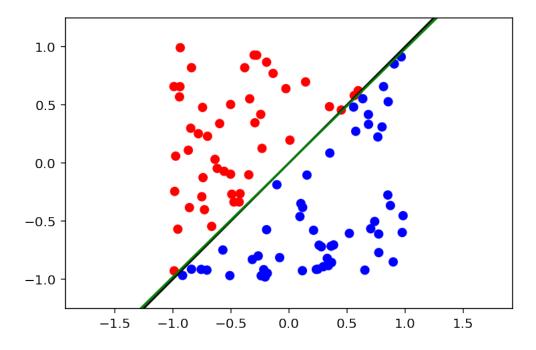


4 d

Repeat everything in b) with another randomly generated dataset of size 100, and compare the result to b).

```
[10]: # generate a dataset of points in 2D space from -1 to 1
      N = 100
      x = (np.random.rand(N,2) * 2) - 1
      # create vector of y values (labels) with target function
      y = [target(x_n) for x_n in x]
      # randomally initialize weights
      w = np.zeros(3)
      idx = 0
      while not test_convergence(w, x, y):
          # pick a random sample
          n = np.random.randint(0,x.shape[0]-1)
          x_n = np.insert(x[n,:], 0, 1)
          y_n = y[n]
          if y_n > 0 and np.dot(w, x_n) < 0:
              w += x_n
          if y_n < 0 and np.dot(w, x_n) >= 0:
              w -= x n
          idx += 1
      print(f"Converged in {idx} steps with w = {w}")
      plt.scatter(x[:,0], x[:,1], c=[pColor if y_n >= 0 else nColor for y_n in y]) #__
      →plot the points in the dataset
      plt.plot(np.linspace(-2,2,50), -(w_t[1]/w_t[2]) * (np.linspace(-2,2,50) -_{\sqcup}
      \hookrightarrow (w_t[0]/w_t[2])), c='k')
      plt.plot(np.linspace(-2,2,50), -(w[1]/w[2]) * (np.linspace(-2,2,50) - (w[0]/w[2])
      \hookrightarrow w[2])), c='g')
      plt.axis('equal')
      plt.axis([-1.25, 1.25, -1.25, 1.25])
      plt.show()
```

Converged in 217 steps with w = [0. 3.47857187 -3.55081271]



5 e

Repeat everything in b) with another randomly generated dataset of size 1000, and compare the result to b).

```
[11]: # generate a dataset of points in 2D space from -1 to 1
N = 1000
x = (np.random.rand(N,2) * 2) - 1

# create vector of y values (labels) with target function
y = [target(x_n) for x_n in x]

# randomally initialize weights
w = np.zeros(3)
idx = 0

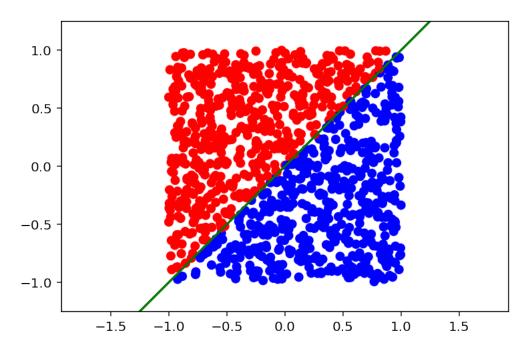
while True:
    # pick a random sample
    n = np.random.randint(0,x.shape[0]-1)
x_n = np.insert(x[n,:], 0, 1)
y_n = y[n]

if y_n > 0 and np.dot(w, x_n) < 0:
    w += x_n</pre>
```

```
if y_n < 0 and np.dot(w, x_n) >= 0:
    # check for convergece every 100 steps
    if idx % 100 == 0:
        if test_convergence(w, x, y):
             break
    idx += 1
print(f"Converged in {idx} steps with w = {w/np.linalg.norm(w, ord=1)}")
plt.scatter(x[:,0], x[:,1], c=[pColor if y_n \ge 0 else nColor for y_n in y]) #_\(\text{u}\)
→plot the points in the dataset
plt.plot(np.linspace(-2,2,50), -(w_t[1]/w_t[2]) * (np.linspace(-2,2,50) -
\hookrightarrow (w_t[0]/w_t[2])), c='k')
plt.plot(np.linspace(-2,2,50), -(w[1]/w[2]) * (np.linspace(-2,2,50) - (w[0]/w[2])
\hookrightarroww[2])), c='g')
plt.axis('equal')
plt.axis([-1.25, 1.25, -1.25, 1.25])
plt.show()
```

Converged in 1800 steps with w = [0.

0.49997488 -0.50002512]



6 f

Modify the experiment such that $\mathbf{x}_n \in \mathbb{R}^{10}$ instead of \mathbb{R}^2 . Run the algorithm on a randomly generated dataset of size 1000. How many updates does the algorithm take to converge?

```
[12]: # generate a dataset of points in 10D space from -1 to 1
      N = 1000
      x = (np.random.rand(N,10) * 2) - 1
      # make a new taget function in R^10
      # create a "random" target function
      w_t = np.random.rand(11)
      def target(x_n):
          return 1 if np.dot(w_t,np.insert(x_n, 0, 1)) >= 0 else -1
      # create vector of y values (labels) with target function
      y = [target(x_n) for x_n in x]
      # initialize weights to 0
      w = np.zeros(11)
      idx = 0
      while True:
          # pick a random sample
          n = np.random.randint(0,x.shape[0]-1)
          x_n = np.insert(x[n,:], 0, 1)
          y_n = y[n]
          if y_n > 0 and np.dot(w, x_n) < 0:
              w += x_n
          if y_n < 0 and np.dot(w, x_n) >= 0:
              w -= x_n
          if idx % 5000 == 0:
              if test_convergence(w, x, y):
                  break
          idx += 1
      print(f"Converged in {idx} steps with w = {w}")
```

```
Converged in 225000 steps with w = [ 3. 24.79573418 1.07691329 29.49325043 17.13808812 23.41487054 25.23573745 25.52347617 30.29276203 14.11818992 33.08671431]
```

7 g

Summarize your conclusions regarding the accuracy and running time of the algorithm as a function of N (the number of data points) and d (the number of dimensions).

As we increase the number of data points, N, we are able to more closely learn the target function. This is shown when we compare the results from the examples with N=20 vs N=100. In the cases where N=20, once we reach convergence, there is often still a significant difference between the true target function and the function we learn. With N=100 we get fairly close to the true target function, and at N=1000, we learn nearly the same function.

We also note, that as we increase N and d, run time, or the steps required for convergence increases. This can result in a significant slow down if we check for convergence by iterating over all data points at every step. For this reason we opt to only check for convergence every k steps (in the larger models), since a single step is much less costly than checking for convergence.