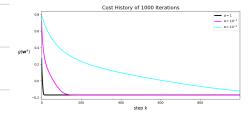
3.5 Try out gradient descent

Run gradient descent to minimize the function

$$g(w) = \frac{1}{50} \left(w^4 + w^2 + 10w \right) \tag{3.44}$$

with an initial point $w^0=2$ and 1000 iterations. Make three separate runs using each of the steplength values $\alpha=1$, $\alpha=10^{-1}$, and $\alpha=10^{-2}$. Compute the derivative of this function by hand, and implement it (as well as the function itself) in Python using NumPy.

Plot the resulting cost function history plot of each run in a single figure to compare their performance. Which steplength value works best for this particular function and initial point?



 α =1 works best for this function, α s:

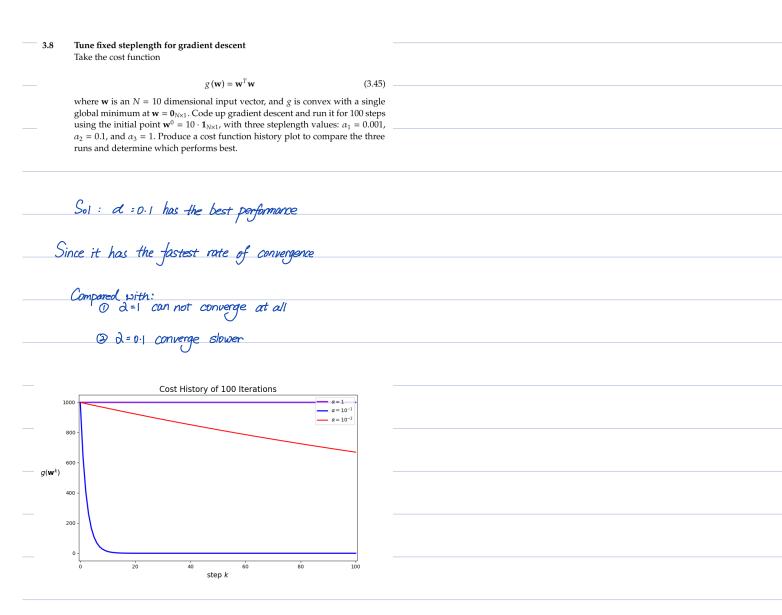
10 converge faster

O Converge without oscillation

• 3-5 Try out gradient descent

from mlrefined libraries.math optimization library import static plotter

```
def gradient decent(G, Gradient, study rate, iteration, w):
   cost = []
   for k in range(1, iteration + 1):
      grad eval = Gradient(w)
      w -= study rate * grad eval
      cost.append(G(w))
   return cost
if __name__ == '__main__':
   iteration = 1000
   w = 2
   G = lambda w: 1 / 50 * (w ** 4 + w ** 2 + 10 * w)
   delta G = lambda w: 1 / 50 * (4 * w ** 3 + w * 2 + 10)
   study rate = [1, 0.1, 0.01]
   cost1 = gradient decent(G, delta G, study rate[0], iteration, w)
   cost2 = gradient decent(G, delta G, study rate[1], iteration, w)
   cost3 = gradient decent(G, delta G, study rate[2], iteration, w)
   plotter = static plotter.Visualizer()
   plotter.plot cost histories(histories=[cost1, cost2, cost3], start=0,
                           labels=[r'$\alpha = 1$', r'$\alpha = 10^{-1}$',
r'\$\alpha = 10^{-2}\$'],
                          title="Cost History of 1000 Iterations"
                           )
```



• 3-8 Tune fixed steplength for gradient descent

```
from mlrefined libraries.math optimization library import static plotter
from autograd import grad
from autograd import numpy as np
plotter = static plotter.Visualizer()
class fixed steplength():
   def __init__(self, function, Dim):
      self.fun = function
      self.ini w = 10 * np.ones((Dim, 1))
   def gradient_descent(self, a):
      w = self.ini w
      max its = 100
      gradient = grad(self.fun)
      weight his = [w]
      cost his = [self.fun(w)]
      for k in range(max its):
          grad eval = gradient(w)
          w = w - a * grad eval
          weight his.append(w)
          cost his.append(self.fun(w))
      return weight his, cost his
if name == '__main__':
   g = lambda w: np.dot(w.T, w)[0][0]
   FSG = fixed steplength(g, 10)
   weight1, cost1 = FSG.gradient descent(a=0.001)
   weight2, cost2 = FSG.gradient_descent(a=0.1)
   weight3, cost3 = FSG.gradient descent(a=1)
   plotter.plot_cost_histories(histories=[cost3, cost2, cost1], start=0,
                           labels=[r'$\alpha = 1$', r'$\alpha = 10^{-1}$',
r'$\alpha = 10^{-2}$'],
                           title="Cost History of 100 Iterations"
                           )
```

Code up momentum-accelerated gradient descent

3.9

Code up the momentum-accelerated gradient descent scheme described in Section A.2.2 and use it to repeat the experiments detailed in Example A.1 using a cost function history plot to come to the same conclusions drawn by studying the contour plots shown in Figure A.3.

Example A.1 Accelerating gradient descent on a simple quadratic

In this example we compare a run of standard gradient descent to the *momentum-accelerated* version using a quadratic function of the form

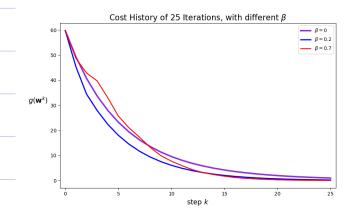
$$g(\mathbf{w}) = a + \mathbf{b}^T \mathbf{w} + \mathbf{w}^T \mathbf{C} \mathbf{w}$$
 (A.9)

where
$$a = 0$$
, $\mathbf{b} = \begin{bmatrix} 0 \\ 0 \end{bmatrix}$, and $\mathbf{C} = \begin{bmatrix} 0.5 & 0 \\ 0 & 9.75 \end{bmatrix}$.

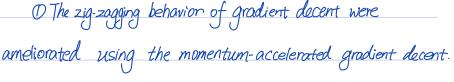
Here we make three runs of 25 steps: a run of gradient descent and two runs of momentum-accelerated gradient descent using two choices of the parameter $\beta \in \{0.2, 0.7\}$. All three runs are initialized at the same point $\mathbf{w}^0 = [10 \ 1]^T$, and use the same steplength $\alpha = 10^{-1}$.

We show the resulting steps taken by the standard gradient descent run in the top panel of Figure A.3 (where significant zig-zagging is present), and the momentum-accelerated versions using $\beta=0.2$ and $\beta=0.7$ in the middle and bottom panels of this figure, respectively. Both momentum-accelerated versions clearly outperform the standard scheme, in that they reach a point closer to the true minimum of the quadratic. Also note that the overall path taken by gradient descent is smoother in the bottom panel, due to the larger value of its corresponding β .

Zig-zagging Behavior for $\beta = 0, 0.2, 0.7$

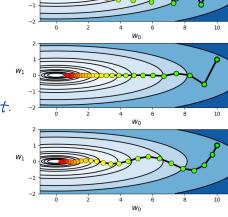


Sol: As illustrated in the history plat below:



 \mathcal{D} $\beta = 0.2.8 \cdot 0.7$, the cost converge closer to 0 when

compared with $\beta = 0$.



MAXXXXXXX

3
$$\beta$$
 = 0-7 converge above than β = 0 & 0.2, that is

because when a large β is chosen, each update will use

less of each subsequent negative gradient direction.

• 3-9 Code up momentum-accelerated gradient descent

from mlrefined libraries.math optimization library import static plotter

```
from autograd import *
from autograd import numpy as np
plotter = static plotter.Visualizer()
class momentum_gradient decent:
   def __init__(self, function, w):
      self.fun = function
      self.ini w = w
   def gradient descent(self, a):
      w = self.ini w
      max its = 100
      gradient = grad(self.fun)
      weight his = [w]
      cost his = [self.fun(w)]
      for k in range(max its):
          grad eval = gradient(w)
          w = w - a * grad eval
          weight his.append(w)
          cost his.append(self.fun(w))
      return weight his, cost his
   def momentum(self, beta):
      w = self.ini w
      max its = 25
      gradient = value and grad(self.fun)
      weight history = []
      cost history = []
      a = 0.1
      h = np.zeros(w.shape)
      for k in range(1, max its + 1):
          cost_eval, grad_eval = gradient(w)
          weight history.append(w)
          cost history.append(cost eval)
          h = beta * h - (1 - beta) * grad eval
          w = w + a * h
      weight history.append(w)
      cost history.append(self.fun(w))
      return weight history, cost history
if __name__ == '__main__':
   g = lambda w: (np.dot(0 * np.ones((2, 1))).T, w) + np.dot(np.dot(w.T,
np.array([[0.5, 0], [0, 9.75]])), w))[0]
```

```
MGD = momentum gradient decent(g, np.array([10.0, 1.0]))
   weight1, cost1 = MGD.momentum(beta=0)
   weight2, cost2 = MGD.momentum(beta=0.2)
   weight3, cost3 = MGD.momentum(beta=0.7)
   his = [weight1, weight2, weight3]
   gs = [g, g, g]
   plotter.two input contour vert plots(title=r"Zig-zagging Behavior for $\beta=0,
0.2, 0.7$", gs=gs, histories=his,
                                  num contours=30, xmin=-1, xmax=11,
                                  ymin=-2.0,
                                  ymax=2)
   plotter.plot_cost_histories(histories=[cost1, cost2, cost3], start=0,
                          labels=[r'$\beta = 0$', r'$\beta = 0.2$', r'$\beta =
0.7$'],
                          title=r"Cost History of 25 Iterations, with different
$\beta$"
                           )
```

3.10

Slow-crawling behavior of gradient descent

In this exercise you will compare the standard and fully normalized gradient descent schemes in minimizing the function

$$g(w_1, w_2) = \tanh(4 w_1 + 4 w_2) + \max(1, 0.4 w_1^2) + 1.$$

Sol: the fully magnitude-normalized gradient decent has steplength $\frac{\alpha}{11 \sqrt{9} (w^{\kappa_1})112}$

(3.46)

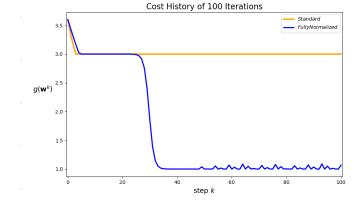
that can adjust itself at each step based on the magnitude of the gradient.

it can ameliorate slow-crawling near minima and saddle point-

As can be illustrated below, after reaching a relative low cost

normalized gradient decent enables the cost converge closer to 0.

W is initialized at [2,2]:



• 3-10 Slow-crawling behavior of gradient descent

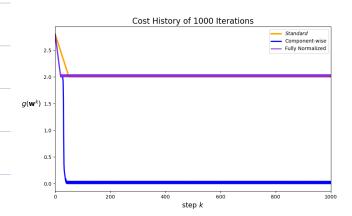
```
from mlrefined libraries.math optimization library import static plotter
from autograd import *
from autograd import numpy as np
plotter = static plotter.Visualizer()
class normalized gradient decent:
   def init (self, function, w):
      self.fun = function
      self.ini w = w
   def gradient descent(self, normalized=False):
      gradient = value and grad(g)
      w = self.ini w
      weight his = []
      cost his = []
      max its = 100
      alpha = 0.1
      for k in range(1, max its + 1):
          cost eval, grad eval = gradient(w)
          weight his.append(w)
          cost his.append(cost eval)
          if normalized:
             grad norm = np.linalg.norm(grad eval)
             if grad norm == 0:
                grad norm += 10 ** -6 * np.sign(2 * np.random.rand(1) - 1)
             grad_eval /= grad_norm
          else:
             grad eval = grad eval
          w = w - alpha * grad_eval
      weight his.append(w)
      cost his.append(self.fun(w))
      return weight_his, cost_his
if name == ' main ':
   g = lambda w: np.tanh(4 * w[0] + 4 * w[1]) + max(0.4 * w[0] ** 2, 1) + 1
   w = np.array([2.0, 2.0])
   print(w)
   NGD = normalized gradient decent(g, w)
   weight1, cost1 = NGD.gradient descent()
   weight2, cost2 = NGD.gradient descent(normalized=True)
   plotter.plot cost histories(histories=[cost1, cost2], start=0,
                           labels=[r'$Standard$', r'$Fully Normalized$'],
                           title="Cost History of 100 Iterations"
```

3.11 Comparing normalized gradient descent schemes

Code up the full and component-wise normalized gradient descent schemes and repeat the experiment described in Example A.4 using a cost function history plot to come to the same conclusions drawn by studying the plots shown in Figure A.6.

Sol: The cost-history for 1000 iterations

Result is illustrated as below:



Our result are completely consistent with

For "fully normalized", its cost function can not coverge

to the minimum, because we is almost 0 everywhere.

Example A.4 Full versus component-normalized gradient descent

In this example we use the function

$$g(w_1, w_2) = \max(0, \tanh(4w_1 + 4w_2)) + |0.4w_1| + 1$$
 (A.25)

to show the difference between full and component-normalized gradient descent steps on a function that has a very narrow flat region along only a single dimension of its input. Here this function - whose surface and contour plots can be seen in the left and right panels of Figure A.6, respectively - is very flat along the w_2 dimension for any fixed value of w_1 , and has a very narrow valley leading towards its minima in the w_2 dimension where $w_1 = 0$. If initialized at a point where $w_2 > 2$ this function cannot be minimized very easily using standard gradient descent or the fully normalized version. In the latter case, the magnitude of the partial derivative in w_2 is nearly zero everywhere, and so fully normalizing makes this contribution smaller, and halts progress. In the top row of the figure we show the result of 1000 steps of fully normalized gradient descent starting at the point $\mathbf{w}^0 = \begin{bmatrix} 2 & 2 \end{bmatrix}^T$, colored green (at the start of the run) to red (at its finale). As can be seen, little progress is made.

In the bottom row of the figure we show the results of using componentnormalized gradient descent starting at the same initialization and employing the same steplength. Here we only need 50 steps in order to make significant progress.

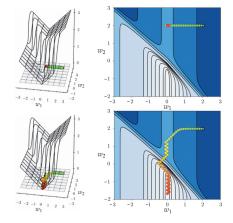


Figure A.6 Figure associated with Example A.4. See text for details.

• 3-11 Comparing normalized gradient descent schemes

```
from mlrefined libraries.math optimization library import static plotter
from autograd import *
from autograd import numpy as np
plotter = static plotter.Visualizer()
class normalized gradient decent:
         def init (self, function, w):
                  self.fun = function
                  self.ini w = w
         def gradient descent(self, normalized=None, max its=100):
                  gradient = value and grad(g)
                  w = self.ini w
                  weight his = []
                 cost his = []
                  alpha = 0.1
                  for k in range(1, max its + 1):
                           cost eval, grad eval = gradient(w)
                           weight his.append(w)
                           cost his.append(cost eval)
                           if normalized == "full":
                                   grad norm = np.linalg.norm(grad eval)
                                    if grad norm == 0:
                                             grad norm += 10 ** -6 * np.sign(2 * np.random.rand(1) - 1)
                                    grad_eval /= grad_norm
                           elif normalized == "Component-wise":
                                    component norm = np.abs(grad eval) + 10 ** (-8)
                                    grad_eval /= component_norm
                           else:
                                   grad eval = grad eval
                           w = w - alpha * grad eval
                  weight his.append(w)
                  cost his.append(self.fun(w))
                  return weight his, cost his
if __name__ == '__main__':
         g = lambda w: np.max(np.tanh(4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[1]), 0) + np.max(np.abs(0.4 * w[0] + 4 * w[0]
w[0]), 0) + 1
         w = np.array([2.0, 2.0])
        max its = 1000
        print(w)
        NGD = normalized gradient decent(g, w)
         weight1, cost1 = NGD.gradient descent(max its=max its)
```