Solutions to hw7 homework on Convex Optimization https://web.stanford.edu/class/ee364a/homework.html

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8.16

9.30

Gradient and Newton methods. Consider the unconstrained problem:

$$\text{minimize} \qquad f(x) = -\sum_{i=1}^{m} log(1 - a_i^T x) - -\sum_{i=1}^{n} log(1 - x_i^2)$$

with variable $x \in R^n$ and $\mathbf{dom} f = \{x \mid a_i^T x < 1,\ i = 1, \dots, m,\ |x_i| < 1,\ i = 1, \dots, n\}$

This is the problem of computing the analytic center of the set of linear inequalities

$$a_i^T x \leq 1, \quad i = 1, \dots, m, \quad |x_i| < 1, \quad i = 1, \dots, n$$

Note that we can choose $x^{(0)} = 0$ as our initial point. You can generate instances of this problem by choosing ai from some distribution on R^n.

(a) Use the gradient method to solve the problem, using reasonable choices for the backtracking parameters, and a stopping criterion of the form $||\nabla f||_2 \leq \nu$. Plot the objective function and step length versus iteration number. (Once you have determined p^* to high accuracy, you can also plot $f - p^*$ versus iteration. Experiment with the backtracking parameters α and β. to see their effect on the total number of iterations required. Carry these experiments out for several instances of the problem, of different sizes.

(b) Repeat using Newton's method, with stopping criterion based on the Newton decrement λ^2. Look for quadratic convergence. You do not have to use

an efficient method to compute the Newton step, as in exercise 9.27; you can use a general purpose dense solver, although it is better to use one that is based on a Cholesky factorization.

Hint. Use the chain rule to find expressions for $\nabla f(x)$ and $\nabla_2 f(x)$.

Solution.

(a) Gradient descent. The python code:

```
# -*- coding: utf-8 -*-

# https://docs.scipy.org/doc/scipy/reference/generated/scipy.optimize.check_grad.html

# Convex optimization. gradient method.

import numpy as np
import matplotlib.pyplot as plt
plt.close('all')

#  definition of some functions ->
def gradient_numerical(f, x0, delta = 1e-8):
"""
function calculates the numerical gradient for function f in
the point x0
"""
N = len(x0)
grad_num = np.zeros([N, 1])
for i in range(N):
xi_plus = x0.copy()
xi_plus[i] += delta

xi_minus = x0.copy()
xi_minus[i] -= delta

grad_num[i] = (f(xi_plus) - f(xi_minus)) / (2 * delta)
return grad_num


def check_grad(f, gradf, x0, delta = 1e-8, verbose = True):
grad = np.array(gradf(x0))
grad_num = gradient_numerical(f, x0, delta)
if (verbose):
print('check_grad: precise gradient = ', grad)
print('check_grad: approximate gradient = ', grad_num)
print('check_grad: gradient error = ', grad - grad_num)
```

```
return np.sqrt(np.sum((grad - grad_num) ** 2))


def f(x, a):
#  calculation of the function value
if not np.all(a.T @ x < 1):
return np.nan
if not np.all(np.abs(x) <= 1):
return np.nan
ret1 = - np.sum(np.log(1 - a.T @ x))
ret2 =  - np.sum(np.log(1 - np.square(x)))
return ret1 + ret2


def gradf(x, a):
#  calculation of the function gradient

if not np.all(a.T @ x < 1):
return np.nan
if not np.all(np.abs(x) <= 1):
return np.nan
print('x = ', x)

ret1 = a @ (1 / (1 - a.T @ x))
ret2 = 2 * x * (1 / (1 - x ** 2))
return ret1 + ret2

def L2norm(x):
return np.sqrt(np.sum(x ** 2))

def backtrack(x, grad, alpha, beta):
"""
Backtracking line search
https://stackoverflow.com/questions/52204231/implementing-backtracking-line-search-algorithm
"""
t = 1
while True:
fx = f(x - t * grad, a)
fxx = f(x, a) + alpha * t * np.dot(grad.T, grad)
if np.isnan(fx) or np.isnan(fxx):
#  print('backtrack: nan detected; multilying t: t = ', t)
t *= beta
elif fx > fxx:
t *= beta
#  print('backtrack: multilying t: t = ', t)
```

```
else:
#  print('backtrack: t found, returning: t = ', t)
return t

#  <- definition of some functions

np.random.seed(1)

m, n = 3, 2

a = np.random.random([m, n]).T

#  check the gradient calculation

x_check = np.random.random([n, 1])

grad_err = check_grad(lambda x: f(x, a), lambda x: gradf(x, a), x_check)

assert grad_err < 1e-6, 'gradient calculation incorrect'

#  parameters for gradient descent method

nu_min = 1e-8 #  tolerance

step = 0.3

x_start = np.zeros([n, 1])

x = x_start


iter_num = 0

max_iters = 1000

max_line_search_iters = 1000

alpha = 0.4

beta = 0.4

print('x_start = ', x_start)
print('x_start.shape = ', x_start.shape)


t = backtrack(x, gradf(x, a), alpha, beta)
```

```
print('t after backtrack = ', t)

# the gradient descent implementation
while True:

print('iteration number ', iter_num)
grad = gradf(x, a)
nu = L2norm(grad)
print('nu = %e' % nu)
if nu <= nu_min:
print('gradient descent: tolerance achieved, exiting...')
print('iteration number ', iter_num)
print('optimal value = %e' % f(x, a))
print('optimal x = ', x)
break
#  Backtracking line search

t = backtrack(x, grad, alpha, beta)
step = t

print('step =', step)

print('grad = ', grad)

x = x - step * grad

print('new x = ', x)
iter_num += 1
if iter_num >= max_iters:
print('gradient descent: max_iters number exeeded')
break


def gradient_descent(alpha, beta):

obj_func_arr = []
step_arr = []
iter_num = 0
x = x_start

while True:

print('iteration number ', iter_num)
obj_func_arr.append(f(x, a))
grad = gradf(x, a)
nu = L2norm(grad)
```

```
print('nu = %e' % nu)
if nu <= nu_min:
print('gradient descent: tolerance achieved, exiting...')
print('iteration number ', iter_num)
print('optimal value = %e' % f(x, a))
opt_val = f(x, a)
return np.array(obj_func_arr) - opt_val, step_arr
print('optimal x = ', x)
break
#  Backtracking line search

t = backtrack(x, grad, alpha, beta)
step = t

print('step =', step)

print('grad = ', grad)

x = x - step * grad
step_arr.append(step)

print('new x = ', x)
iter_num += 1
if iter_num >= max_iters:
print('gradient descent: max_iters number exeeded')
return None, None
break

# plot the graphs
alpha_arr = [0.2, 0.4]

beta_arr = [0.2, 0.45]

plt.figure()

for alpha in  alpha_arr:
for beta in beta_arr:
print('alpha = ', alpha)
print('beta = ', beta)
obj_func, step =  gradient_descent(alpha, beta)
x_plt = range(len(obj_func))
plt.plot(x_plt, np.log10(obj_func),
label='alpha = ' + str(alpha) + ' beta = ' + str(beta))

plt.title('logarithm of the objective function error vs iteration number')
plt.ylabel('logarithm of the objective function error')
```

```
plt.xlabel('iteration number')
plt.legend()
plt.show()

plt.savefig('9_30_a_obj_func.png', bbox_inches='tight')


plt.figure()

for alpha in  alpha_arr:
for beta in beta_arr:
print('alpha = ', alpha)
print('beta = ', beta)
obj_func, step =  gradient_descent(alpha, beta)
x_plt = range(len(step))
plt.plot(x_plt, step,
label='alpha = ' + str(alpha) + ' beta = ' + str(beta))

plt.title('step vs iteration number')
plt.ylabel('step')
plt.xlabel('iteration number')
plt.legend()
plt.show()

plt.savefig('9_30_a_step.png', bbox_inches='tight')
```

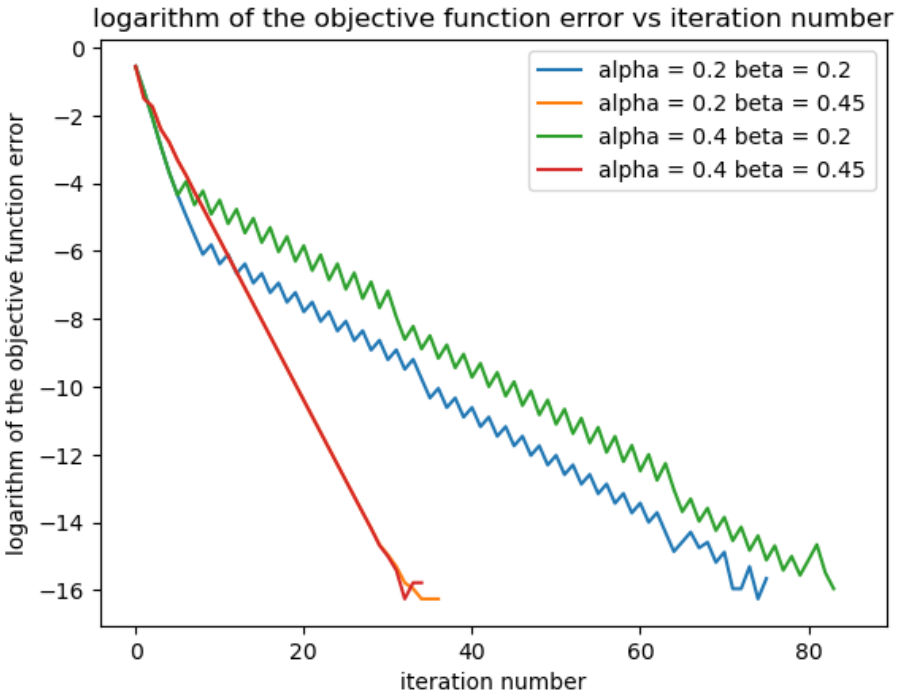

Figure 1: Gradient descent: $f - p^*$ vs. iteration number

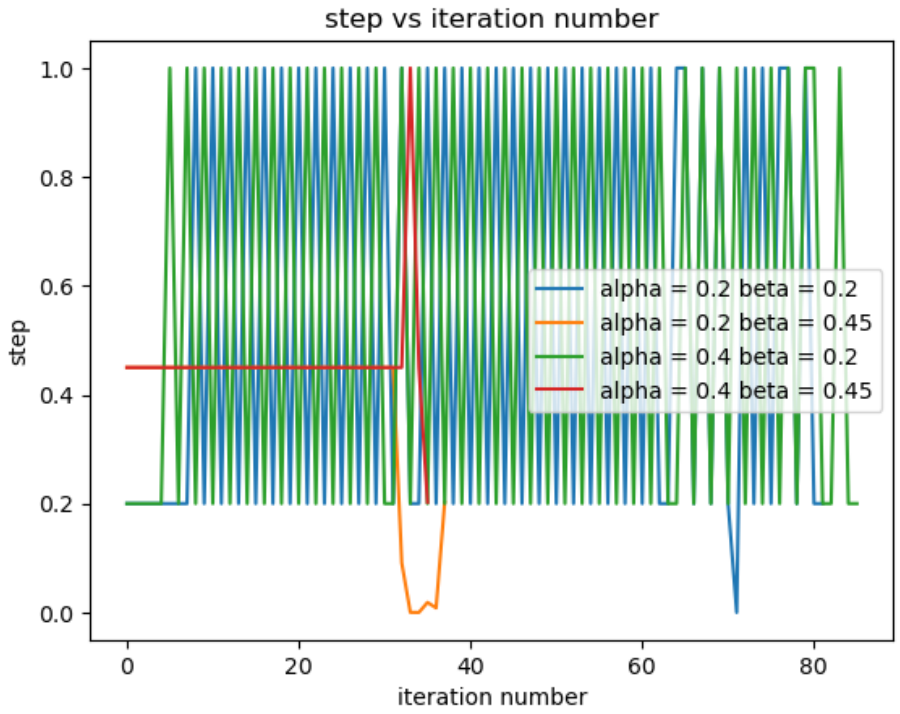

Figure 2: Gradient descent: step size vs. iteration number

(a) Newton's method. The python code:

file `ex_9_30_test_hessian.py`

```
# -*- coding: utf-8 -*-
# https://docs.scipy.org/doc/scipy/reference/generated/scipy.optimize.check_grad.html

import numpy as np

def backtrack(x, a, grad, alpha, beta):
"""
Backtracking line search
https://stackoverflow.com/questions/52204231/implementing-backtracking-line-search-algorithm
"""
t = 1
while True:
fx = f(x - t * grad, a)
fxx = f(x, a) + alpha * t * np.dot(grad.T, grad)
if np.isnan(fx) or np.isnan(fxx):
```

```
#  print('backtrack: nan detected; multilying t: t = ', t)
t *= beta
elif fx > fxx:
t *= beta
#  print('backtrack: multilying t: t = ', t)
else:
#  print('backtrack: t found, returning: t = ', t)
return t


def backtrack_2(x, a, grad, ihess, alpha, beta):
"""
Backtracking line search
https://stackoverflow.com/questions/52204231/implementing-backtracking-line-search-algorithm
"""
"""
print('grad.shape = ', grad.shape)
print('grad = ', grad)
print('ihess.shape = ', ihess.shape)
print('ihess = ', ihess)
"""
gh = grad.T @ ihess
#  print('grad.T @ ihess = ', gh)


t = 1
while True:
#  print('x = ', x)
#  print('gh = ', gh)

fx = f(x - t * gh , a)

fxx = f(x, a) + alpha * t * np.sum(gh ** 2)
#  fxx = f(x, a) + alpha * t * np.dot(gh, gh.T) #  the same as f(x, a) + alpha * t * np.sum(

if np.isnan(fx) or np.isnan(fxx):
#  print('backtrack: nan detected; multilying t: t = ', t)
t *= beta
elif fx > fxx:
t *= beta
#  print('backtrack: multilying t: t = ', t)
else:
#  print('backtrack: t found, returning: t = ', t)
return t

def L2norm(x):
```

```
return np.sqrt(np.sum(x ** 2))


def derivative_numerical(f, x0, i, delta = 1e-8):
xi_plus = x0.copy()
xi_plus[i] += delta

xi_minus = x0.copy()
xi_minus[i] -= delta
return (f(xi_plus) - f(xi_minus)) / (2 * delta)



def gradient_numerical(f, x0, delta = 1e-8):
"""
function calculates the numerical gradient for function f in
the point x0
"""
N = len(x0)
grad_num = np.zeros([N, 1])
for i in range(N):
grad_num[i] = derivative_numerical(f, x0, i, delta)
return grad_num


def check_grad(f, gradf, x0, delta = 1e-8, verbose = True):
grad = np.array(gradf(x0))
grad_num = gradient_numerical(f, x0, delta)
if (verbose):
print('check_grad: precise gradient = ', grad)
print('check_grad: approximate gradient = ', grad_num)
print('check_grad: gradient error = ', grad - grad_num)

return np.sqrt(np.sum((grad - grad_num) ** 2))

def second_derivative_numerical(f, x0, i, k, delta = 1e-5):
"""
function calculates second derivative
returns d^2f/(dx_k dx_i)
"""
xk_plus = x0.copy()
xk_plus[k] += delta

xk_minus = x0.copy()
xk_minus[k] -= delta
```

```
dfi_plus = derivative_numerical(f, xk_plus, i, delta)
dfi_minus = derivative_numerical(f, xk_minus, i, delta)

return (dfi_plus - dfi_minus) / (2 * delta)

def hessian_numerical(f, x0, delta = 1e-5):
"""
#  function calculates the hessian matrix
"""
assert x.shape[1] == 1, 'hessian_numerical: input array should have shape [N, 1]'

N = len(x)
hessian = np.zeros([N, N], dtype = np.float64)
for i in range(N):
for k in range(i, N):
hessian[i, k] = second_derivative_numerical(f, x0, i, k, delta)
if i != k:
hessian[k, i] = hessian[i, k]
return hessian

def check_hessian(f, hess_analytical, x0, delta = 1e-5, verbose = True):
"""
function checks he hessian matrix
"""
hessian_analytical = np.array(hess_analytical)
hessian_num = hessian_numerical(f, x0, delta)
if verbose:
print('check_hessian: hessian_analytical = ', hessian_analytical)
print('check_hessian: hessian_num = ', hessian_num)
print('check_hessian: hessian difference = ',
hessian_analytical - hessian_num)

return np.sqrt(np.sum((hessian_analytical - hessian_num) ** 2))

#%%
# definitions for the function, gradient and hessian

def f(x, a):
#  calculation of the function value
if not np.all(a.T @ x < 1):
return np.nan
if not np.all(np.abs(x) <= 1):
return np.nan
ret1 = 0.0
ret1 = - np.sum(np.log(1 - a.T @ x))
ret2 =  - np.sum(np.log(1 - np.square(x)))
```

```
return ret1 + ret2

#  print('f(x, a) = ', f(x, a))

def gradf(x, a):
#  calculation of the function gradient

if not np.all(a.T @ x < 1):
return np.nan
if not np.all(np.abs(x) <= 1):
return np.nan
#  print('x = ', x)

ret1 = 0.0
ret1 = a @ (1 / (1 - a.T @ x))
ret2 = 2 * x * (1 / (1 - x ** 2))
return ret1 + ret2

def hessf(x, a):
if not np.all(a.T @ x < 1):
return np.nan
if not np.all(np.abs(x) <= 1):
return np.nan
ret1 = 0
ret1 = a @ (a.T * (1 / (1 - a.T @ x) ** 2))
ret2 = 2 * (1 + x ** 2) / ((1 - x ** 2) ** 2)
ret2 = np.diagflat(ret2)

return ret1 + ret2


if __name__ == "__main__":

np.random.seed(1)

m, n = 3, 2

a = np.random.random([m, n]).T

#  a = np.array([[-1, 0], [0.5, - 0.5], [0.5, 0]], dtype = np.float64).T

x0 = 0.5 * np.ones([n, 1])
x = np.array([-0.25, 0.75], dtype = np.float64).reshape(- 1, 1)
```

```
print('a.shape = ', a.shape)


#  x = np.array([0.5, 0.75])

#x = np.array([-0.75, 0.5], dtype = np.float64).reshape(- 1, 1)

error1 = check_grad(lambda x: f(x, a), lambda x: gradf(x, a), x0)

assert error1 < 1e-6, 'error1 too big'

print('gradient error1 = ', error1)


x0 = - 0.5 * np.ones([n, 1])
fl3 = lambda x: (x[0]**2 + 3*x[1]*x[0] + 12)

def f3(x):
return (x[0]**2 + 3*x[1]*x[0] + 12)[0]

dfx1 = lambda x: (2*x[0] + 3*x[1])
dfx2 = lambda x: (3*x[0])

def gradf3(x):
return np.array([dfx1(x), dfx2(x)]).reshape([-1, 1])


error3 = check_grad(f3, gradf3, x0)

print('gradient error3 = ', error3)

assert error3 < 1e-6, 'error3 too big'

error4 = check_grad(f3, gradf3, x0)

print('gradient error4 = ', error4)

assert error4 < 1e-6, 'error4 too big'

#%%
# test function for hessian

def fh(z):
assert z.shape[0] == 2 and z.shape[1] == 1, 'fh(x): incorrect input shape'

x = z[0]
```

```
y = z[1]

return x**2 + 0.5 * y**2 + 2 * x * y + 3 * x + 4 * x**2 * y**2 + 5 * y * x**2

def fh_hessian(z):
assert z.shape[0] == 2 and z.shape[1] == 1, 'fh_hessian(x): incorrect input shape'

x = z[0]
y = z[1]

hessian = np.zeros([2, 2], dtype = np.float64)
print('')
hessian[0, 0] = 2 + 8 * y**2 + 10 * y
hessian[0, 1] = 2 + 16 * x * y + 10 * x
hessian[1, 0] = hessian[0, 1]
hessian[1, 1] = 1 + 8 * x ** 2
return hessian

#%%


#  test check_hessian function:

x0 = np.array([0.5, 8], dtype=np.float64).reshape(-1, 1)

#  x0 = np.array([0, 0], dtype=np.float64).reshape(-1, 1)

#  x0 = np.array([1, 1], dtype=np.float64).reshape(-1, 1)

hess_analytical = fh_hessian(x0)


hd = check_hessian(fh, hess_analytical, x0, delta = 1e-5, verbose = True)

print('test of check_hessian, diff = %e' % hd)

# x0 = np.array([0.5, -0.25], dtype=np.float64).reshape(-1, 1)
x0 = np.array([-0.5, -0.75], dtype=np.float64).reshape(-1, 1)

v_hess_analytical = hessf(x0, a)

print('v_hess_analytical = ', v_hess_analytical)


v_hess_num = hessian_numerical(lambda x: f(x, a), x0)
```

```
print('v_hess_num = ', v_hess_num)

hc = check_hessian(lambda x: f(x, a), v_hess_analytical, x0)
print('hc = ', hc)

assert hc < 1e-4, 'hessian seems to be incorrect'

grad = np.array([[3.26697727], [4.08950456]])
ihess = np.array([[ 0.31264018, -0.13959122],  [-0.13959122,  0.28763308]])
print('grad.shape = ', grad.shape)
print('ihess.shape = ', ihess.shape)
y = grad.T @ ihess
print('y = ', y)
print('y.shape = ', y.shape)


print(np.sum(y**2))
```

file ex_9_30_b.py

```
# -*- coding: utf-8 -*-

import numpy as np
import ex_9_30_test_hessian as h
import matplotlib.pyplot as plt
plt.close('all')


f = h.f

np.random.seed(3)

m, n = 800, 60

a = np.random.random([m, n]).T


#  parameters for gradient descent method

nu_min = 1e-8 #  tolerance

step = 0.3

x_start = np.zeros([n, 1])

x = x_start
```

```
iter_num = 0

max_iters = 1000

max_line_search_iters = 1000

alpha = 0.4

beta = 0.4

print('x_start = ', x_start)
print('x_start.shape = ', x_start.shape)


# t = h.backtrack(x, a, h.gradf(x, a), alpha, beta)
# print('t after backtrack = ', t)

# the Newton's method implementation
while True:

print('iteration number ', iter_num)
grad = h.gradf(x, a) #  gradient of f
hess = h.hessf(x, a) #  hessian of f
ihess = np.linalg. inv(hess) #  inverse hessian of f
dx = - ihess @ grad #  Newton step
lam_sq = grad.T @ (ihess @ grad) # Newton decrement

print('lam_sq = %e' % lam_sq)
if np.sqrt(lam_sq / 2) <= nu_min:
print("Newton's method: tolerance achieved, exiting...")
print('iteration number ', iter_num)
#  print('a = ', a)
print('optimal value = %e' % f(x, a))
print('optimal x = ', x)
break
#  Backtracking line search

t = h.backtrack_2(x, a, grad, ihess, alpha, beta)
step = t
# step = 1

print('step =', step)

x = x + step * dx
```

```
print('new x = ', x)
iter_num += 1
if iter_num >= max_iters:
print("Newton's method: max_iters number exeeded")
break


def newton_method(alpha, beta):

obj_func_arr = []
step_arr = []
iter_num = 0
x = x_start

while True:

print('iteration number ', iter_num)
obj_func_arr.append(f(x, a))
grad = h.gradf(x, a) #  gradient of f
hess = h.hessf(x, a) #  hessian of f
ihess = np.linalg. inv(hess) #  inverse hessian of f
dx = - ihess @ grad #  Newton step
lam_sq = grad.T @ (ihess @ grad) # Newton decrement

print('lam_sq = %e' % lam_sq)
#  if lam_sq / 2 <= nu_min:
if np.sqrt(lam_sq / 2) <= nu_min:
print("Newton's method: tolerance achieved, exiting...")
print('iteration number ', iter_num)
#  print('a = ', a)
opt_val = f(x, a)
print('optimal value = %e' % opt_val)
print('optimal x = ', x)
return np.array(obj_func_arr) - opt_val, step_arr
break
#  Backtracking line search

t = h.backtrack_2(x, a, grad, ihess, alpha, beta)
step = t
step_arr.append(step)

print('step =', step)

x = x + step * dx
```

```
print('new x = ', x)
iter_num += 1
if iter_num >= max_iters:
print("Newton's method: max_iters number exeeded")
return None, None
break

# plot the graphs
alpha_arr = [0.2, 0.4]

beta_arr = [0.2, 0.45]

plt.figure()

for alpha in  alpha_arr:
for beta in beta_arr:
print('alpha = ', alpha)
print('beta = ', beta)
obj_func, step =  newton_method(alpha, beta)
x_plt = range(len(obj_func))
plt.plot(x_plt, np.log10(obj_func),
label='alpha = ' + str(alpha) + ' beta = ' + str(beta))

plt.title('logarithm of the objective function error vs iteration number')
plt.ylabel('logarithm of the objective function error')
plt.xlabel('iteration number')
plt.legend()
plt.show()

plt.savefig('9_30_b_obj_func.png', bbox_inches='tight')


plt.figure()

for alpha in  alpha_arr:
for beta in beta_arr:
print('alpha = ', alpha)
print('beta = ', beta)
obj_func, step =  newton_method(alpha, beta)
x_plt = range(len(step))
plt.plot(x_plt, step,
label='alpha = ' + str(alpha) + ' beta = ' + str(beta))

plt.title('step vs iteration number')
plt.ylabel('step')
plt.xlabel('iteration number')
```

```
plt.legend()
plt.show()

plt.savefig('9_30_b_step.png', bbox_inches='tight')
```

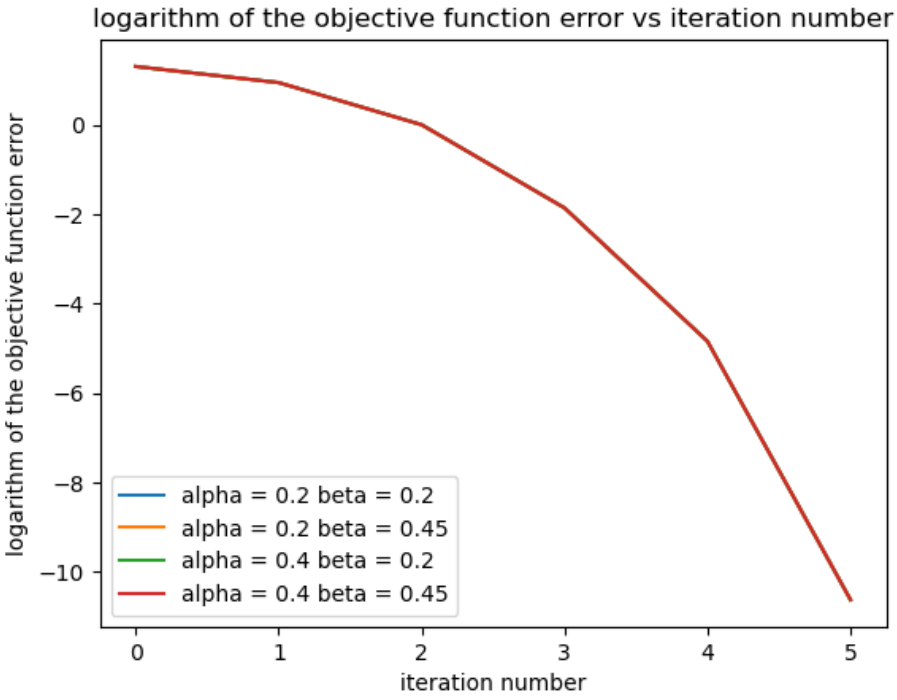

Figure 3: Newton's method: $f - p^*$ vs. iteration number

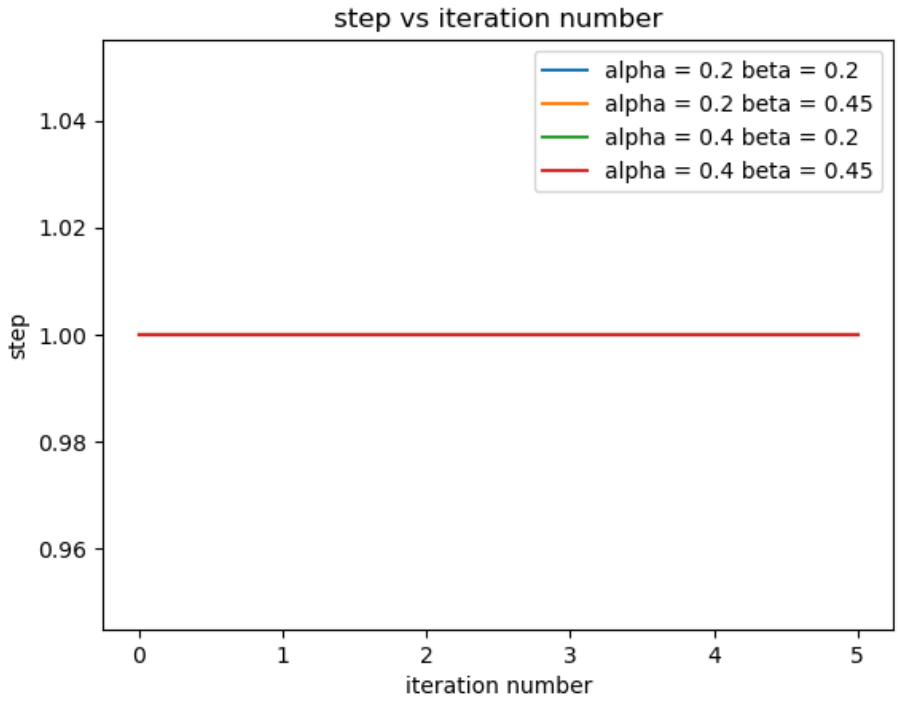

Figure 4: Newton's method: step size vs. iteration number

9.31

Some approximate Newton methods. The cost of Newton's method is dominated by the cost of evaluating the Hessian $\nabla_2 f$ and the cost of solving the Newton system. For large problems, it is sometimes useful to replace the Hessian by a positive definite approximation that makes it easier to form and solve for the search step. In this problem we explore some common examples of this idea.

For each of the approximate Newton methods described below, test the method on some instances of the analytic centering problem described in exercise 9.30, and compare the results to those obtained using the Newton method and gradient method.

(a) Re-using the Hessian. We evaluate and factor the Hessian only every N iterations, where $N > 1$, and use the search step $\Delta x = -H^{-1}\nabla f(x)$, where H is the last Hessian evaluated. (We need to evaluate and factor the Hessian once every N steps; for the other steps, we compute the search direction using back and forward substitution.)

(b) Diagonal approximation. We replace the Hessian by its diagonal, so we only have to evaluate the n second derivatives $\frac{\partial^2 f(x)}{\partial t^2}$ and computing the search

step is very easy.

Solution:
(a) The python code:

```
# -*- coding: utf-8 -*-

import numpy as np
import ex_9_30_test_hessian as h
import matplotlib.pyplot as plt
plt.close('all')


f = h.f

np.random.seed(3)

m, n = 18, 10

a = np.random.random([m, n]).T


#  parameters for gradient descent method

nu_min = 1e-8 #  tolerance

step = 0.3

x_start = np.zeros([n, 1])

x = x_start

max_iters = 1000

max_line_search_iters = 1000

alpha = 0.4

beta = 0.4

print('x_start = ', x_start)
print('x_start.shape = ', x_start.shape)

N = 3

iter_num = 0
```

```
# the Newton's method implementation

while True:

print('iteration number ', iter_num)

grad = h.gradf(x, a) #  gradient of f

"""
We need to evaluate and factor the Hessian once every N steps; for the
other steps, we compute the search direction using back and forward substitution
"""
if iter_num // N == 0:
hess = h.hessf(x, a) #  hessian of f
ihess = np.linalg. inv(hess) #  inverse hessian of f

dx = - ihess @ grad #  Newton step
lam_sq = grad.T @ (ihess @ grad) # Newton decrement

print('lam_sq = %e' % lam_sq)
if np.sqrt(lam_sq / 2) <= nu_min:
print("Newton's method: tolerance achieved, exiting...")
print('iteration number ', iter_num)
#  print('a = ', a)
print('optimal value = %e' % f(x, a))
print('optimal x = ', x)
break
#  Backtracking line search

t = h.backtrack_2(x, a, grad, ihess, alpha, beta)
step = t
# step = 1

print('step =', step)

x = x + step * dx

print('new x = ', x)
iter_num += 1
if iter_num >= max_iters:
print("Newton's method: max_iters number exeeded")
break

def newton_method(alpha, beta):
```

```
obj_func_arr = []
step_arr = []
iter_num = 0
x = x_start

while True:

print('iteration number ', iter_num)
obj_func_arr.append(f(x, a))

grad = h.gradf(x, a) #  gradient of f

"""
We need to evaluate and factor the Hessian once every N steps; for the
other steps, we compute the search direction using back and forward substitution
"""
if iter_num // N == 0:
hess = h.hessf(x, a) #  hessian of f
ihess = np.linalg. inv(hess) #  inverse hessian of f

dx = - ihess @ grad #  Newton step
lam_sq = grad.T @ (ihess @ grad) # Newton decrement

print('lam_sq = %e' % lam_sq)
if np.sqrt(lam_sq / 2) <= nu_min:
print("Newton's method: tolerance achieved, exiting...")
print('iteration number ', iter_num)
#  print('a = ', a)
print('optimal value = %e' % f(x, a))
print('optimal x = ', x)
return np.array(obj_func_arr) - f(x, a), step_arr
break
#  Backtracking line search

t = h.backtrack_2(x, a, grad, ihess, alpha, beta)
step = t
step_arr.append(step)

print('step =', step)

x = x + step * dx

print('new x = ', x)
iter_num += 1
if iter_num >= max_iters:
print("Newton's method: max_iters number exeeded")
```

```
break

# plot the graphs
alpha_arr = [0.2, 0.4]

beta_arr = [0.2, 0.45]

plt.figure()

for alpha in  alpha_arr:
for beta in beta_arr:
print('alpha = ', alpha)
print('beta = ', beta)
obj_func, step =  newton_method(alpha, beta)
x_plt = range(len(obj_func))
plt.plot(x_plt, np.log10(obj_func),
label='alpha = ' + str(alpha) + ' beta = ' + str(beta))

plt.title('logarithm of the objective function error vs iteration number')
plt.ylabel('logarithm of the objective function error')
plt.xlabel('iteration number')
plt.legend()
plt.show()

plt.savefig('9_31_a_obj_func.png', bbox_inches='tight')


plt.figure()

for alpha in  alpha_arr:
for beta in beta_arr:
print('alpha = ', alpha)
print('beta = ', beta)
obj_func, step =  newton_method(alpha, beta)
x_plt = range(len(step))
plt.plot(x_plt, step,
label='alpha = ' + str(alpha) + ' beta = ' + str(beta))

plt.title('step vs iteration number')
plt.ylabel('step')
plt.xlabel('iteration number')
plt.legend()
plt.show()

plt.savefig('9_31_a_step.png', bbox_inches='tight')
```

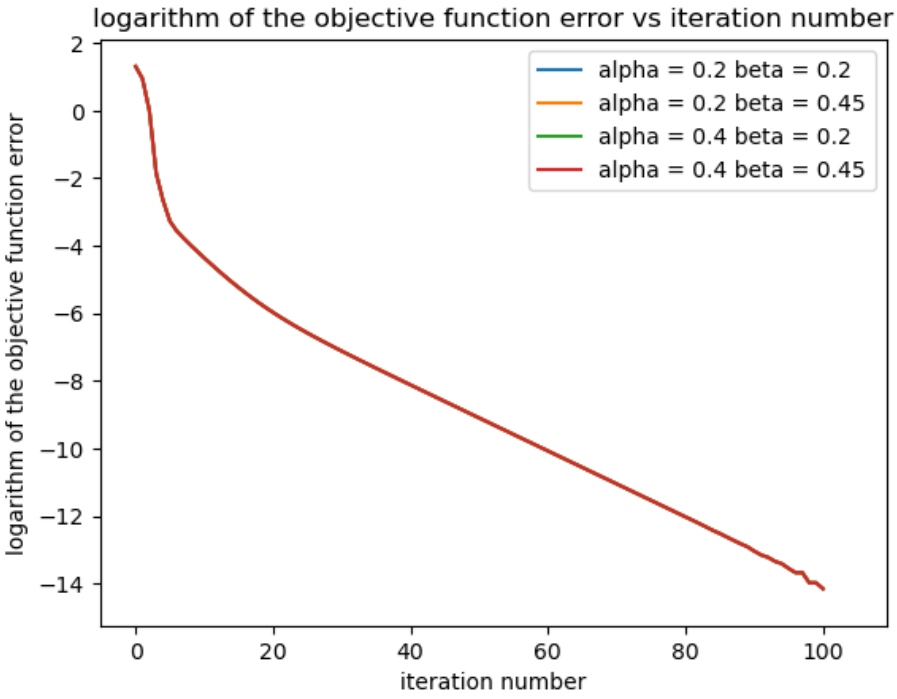

Figure 5: Newton's method: $f - p^*$ vs. iteration number

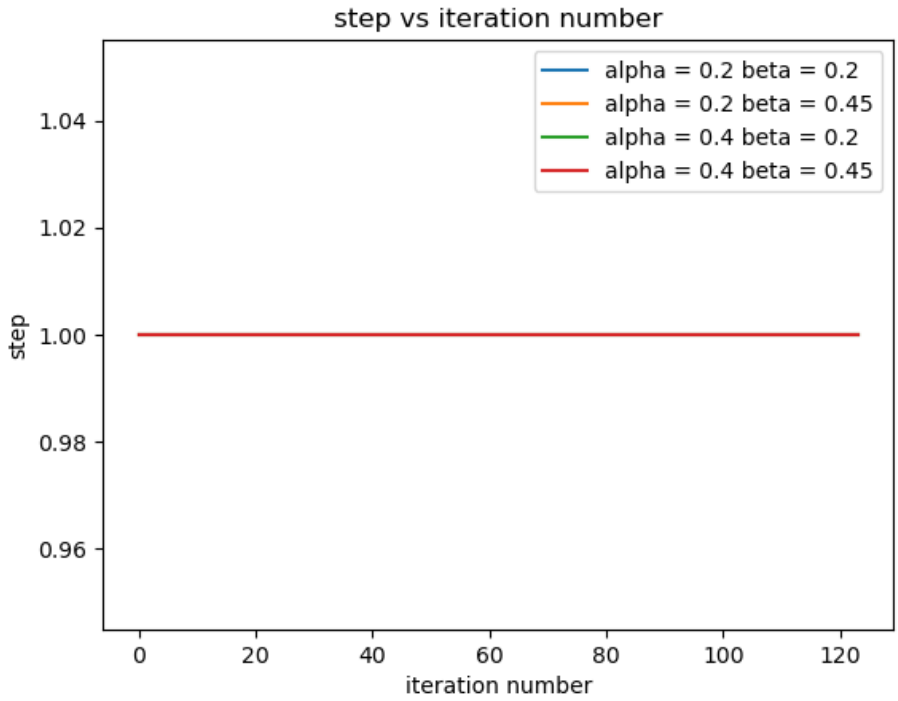

Figure 6: Newton's method: step size vs. iteration number

(b) The python code:

```
# -*- coding: utf-8 -*-

import numpy as np
import ex_9_30_test_hessian as h
import matplotlib.pyplot as plt
plt.close('all')

f = h.f

np.random.seed(3)

m, n = 18, 10

a = np.random.random([m, n]).T


#  parameters for gradient descent method
```

```
nu_min = 1e-8 #  tolerance

step = 0.3

x_start = np.zeros([n, 1])

x = x_start


iter_num = 0

max_iters = 1000

max_line_search_iters = 1000

alpha = 0.4

beta = 0.4

print('x_start = ', x_start)
print('x_start.shape = ', x_start.shape)


# the Newton's method implementation
while True:

print('iteration number ', iter_num)
grad = h.gradf(x, a) #  gradient of f
hess = h.hessf(x, a) #  hessian of f
# Diagonal approximation. We replace the Hessian by its diagonal
ihess = np.diag(1 / np.diagonal(hess))
# ihess = np.linalg. inv(hess) #  inverse hessian of f
dx = - ihess @ grad #  Newton step
lam_sq = grad.T @ (ihess @ grad) # Newton decrement

print('lam_sq = %e' % lam_sq)
if np.sqrt(lam_sq / 2) <= nu_min:
print("Newton's method: tolerance achieved, exiting...")
print('iteration number ', iter_num)
#  print('a = ', a)
print('optimal value = %e' % f(x, a))
print('optimal x = ', x)
break
#  Backtracking line search
```

```
t = h.backtrack_2(x, a, grad, ihess, alpha, beta)
step = t
# step = 1

print('step =', step)

x = x + step * dx

print('new x = ', x)
iter_num += 1
if iter_num >= max_iters:
print("Newton's method: max_iters number exeeded")
break

def newton_method(alpha, beta):

obj_func_arr = []
step_arr = []
iter_num = 0
x = x_start


while True:

print('iteration number ', iter_num)
obj_func_arr.append(f(x, a))
grad = h.gradf(x, a) #  gradient of f
hess = h.hessf(x, a) #  hessian of f
# Diagonal approximation. We replace the Hessian by its diagonal
ihess = np.diag(1 / np.diagonal(hess))

dx = - ihess @ grad #  Newton step
lam_sq = grad.T @ (ihess @ grad) # Newton decrement

print('lam_sq = %e' % lam_sq)
if np.sqrt(lam_sq / 2) <= nu_min:
print("Newton's method: tolerance achieved, exiting...")
print('iteration number ', iter_num)
#  print('a = ', a)
print('optimal value = %e' % f(x, a))
print('optimal x = ', x)
return np.array(obj_func_arr) - f(x, a), step_arr
break
#  Backtracking line search

t = h.backtrack_2(x, a, grad, ihess, alpha, beta)
```

```
step = t
step_arr.append(step)

print('step =', step)

x = x + step * dx

print('new x = ', x)
iter_num += 1
if iter_num >= max_iters:
print("Newton's method: max_iters number exeeded")
break

# plot the graphs
alpha_arr = [0.2, 0.4]

beta_arr = [0.2, 0.45]

plt.figure()

for alpha in  alpha_arr:
for beta in beta_arr:
print('alpha = ', alpha)
print('beta = ', beta)
obj_func, step =  newton_method(alpha, beta)
x_plt = range(len(obj_func))
plt.plot(x_plt, np.log10(obj_func),
label='alpha = ' + str(alpha) + ' beta = ' + str(beta))

plt.title('logarithm of the objective function error vs iteration number')
plt.ylabel('logarithm of the objective function error')
plt.xlabel('iteration number')
plt.legend()
plt.show()

plt.savefig('9_31_b_obj_func.png', bbox_inches='tight')


plt.figure()

for alpha in  alpha_arr:
for beta in beta_arr:
print('alpha = ', alpha)
print('beta = ', beta)
obj_func, step =  newton_method(alpha, beta)
x_plt = range(len(step))
```

```
plt.plot(x_plt, step,
label='alpha = ' + str(alpha) + ' beta = ' + str(beta))

plt.title('step vs iteration number')
plt.ylabel('step')
plt.xlabel('iteration number')
plt.legend()
plt.show()

plt.savefig('9_31_b_step.png', bbox_inches='tight')
```

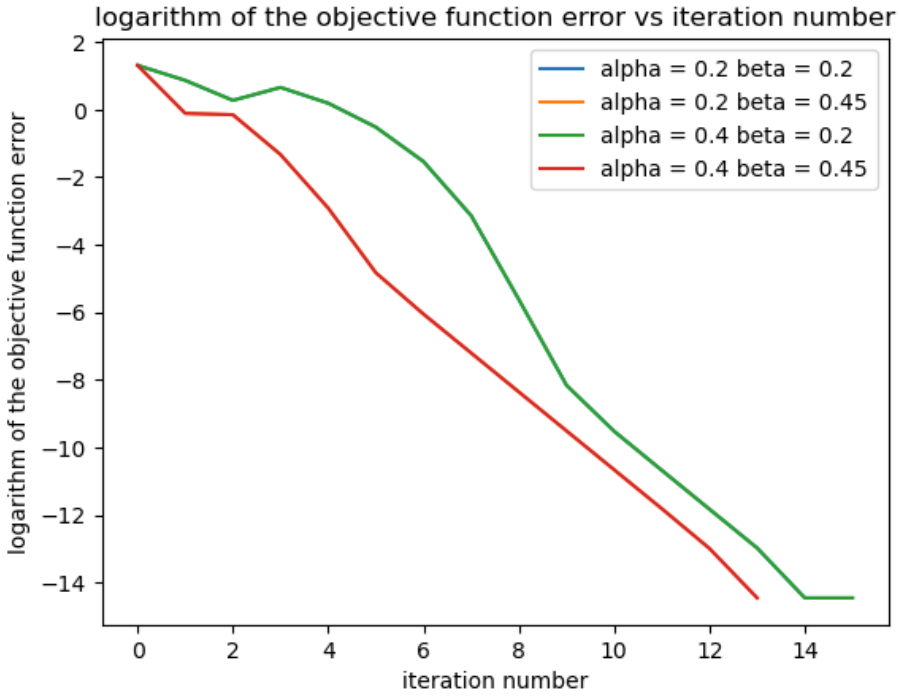

Figure 7: Newton's method: $f - p^*$ vs. iteration number

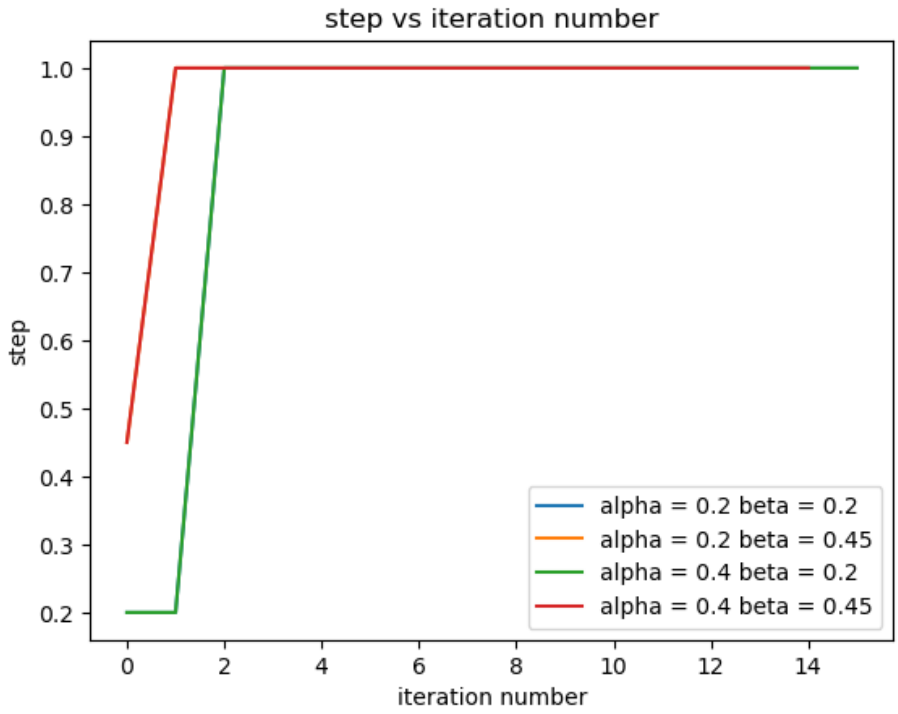

Figure 8: Newton's method: step size vs. iteration number

As it can be seen from the pictures, method (a) is not very good, but method (b) shows results somewhat comparable with the pure Newton's method ones.