In [53]: import numpy as np
import matplotlib.pyplot as plt

$$f(x_1, x_2) = (x_1 - 3)^2 + (x_2 - 1)^2$$

$$f(x_1, x_2) = 10(x_1 - 1)^2 + (x_2 - 2)^2$$

In [55]: def f2(x, \*\*kwargs):
 x1, x2 = x
 return 10\*(x1 - 1)\*\*2 + (x2 - 2)\*\*2

def grad\_f2(x, \*\*kwargs):
 x1, x2 = x
 g\_x1 = 20\*(x1-1)
 g\_x2 = 2\*(x2-2)
 return np.array([g\_x1, g\_x2])

$$f(x)=\frac{1}{2}||Ax-b||_2^2$$

In [56]: def f3(x, \*\*kwargs):
 A, b = compute\_params(\*\*kwargs)
 return 0.5 \* np.linalg.norm(A@x - b, 2)\*\*2

def grad\_f3(x, \*\*kwargs):
 A, b = compute\_params(\*\*kwargs)
 return A.T @ (A @ x - b)

def compute\_params(\*\*kwargs):
 n = kwargs['n']
 v = np.linspace(0,1,n)

 A = np.vander(v, increasing=True)

 x\_star = np.ones\_like(v)
 b = A @ x\_star

 return A, b

$$f(x) = rac{1}{2} ||Ax - b||_2^2 + rac{\lambda}{2} ||x||_2^2$$

$$f(x) = x^4 + x^3 - 2x^2 - 2x.$$

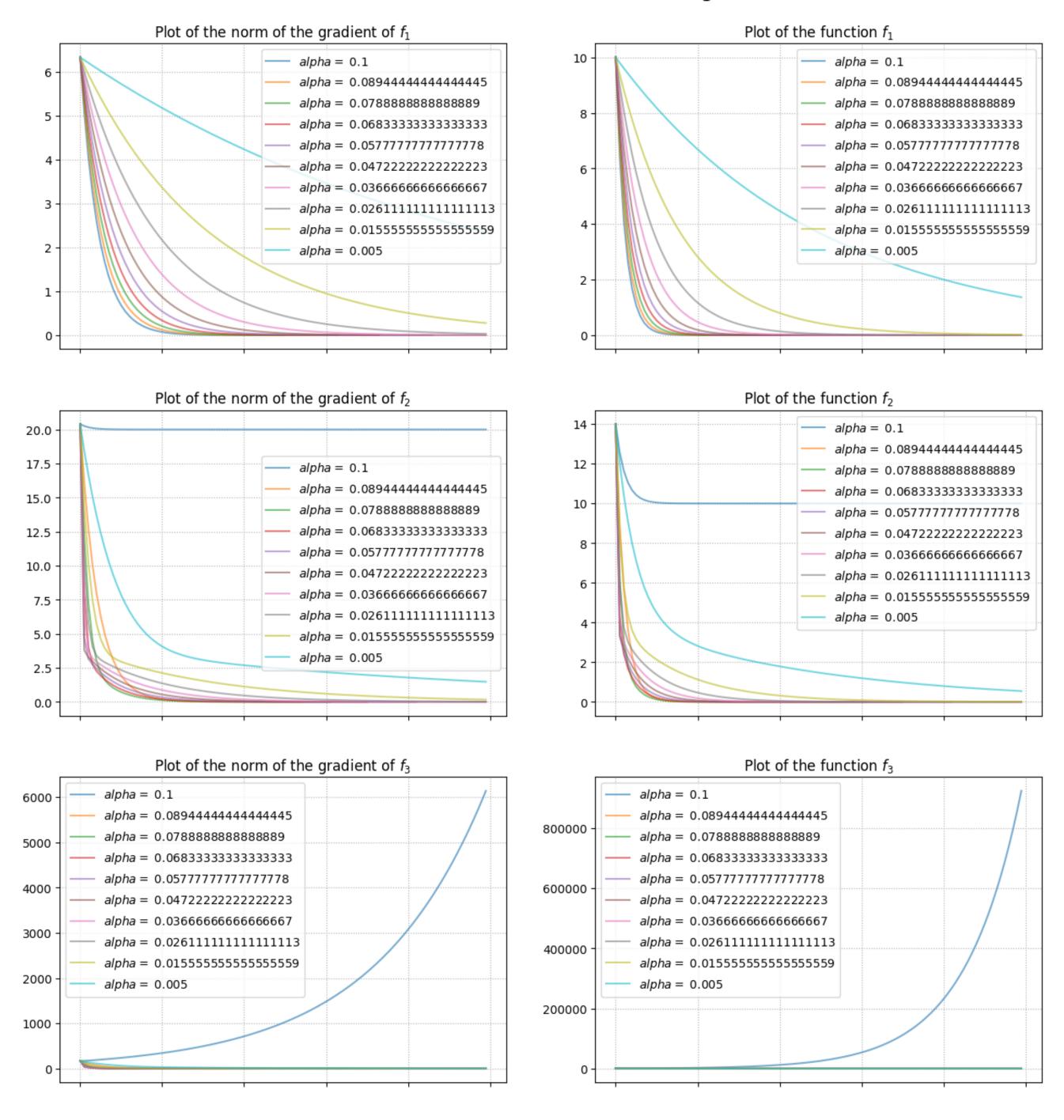
```
In [58]: def f5(x, **kwargs):
             return x**4 + x**3 - 2*x**2 - 2*x
         def grad_f5(x, **kwargs):
             return 4*x**3 + 3*x**2 - 4*x - 2
In [59]: def GD(f, grad_f, x0, alpha, maxit=100, tolf=1e-5, tolx=1e-5, **kwargs):
             f_vals = np.zeros((maxit+1,))
             grad_norms = np.zeros((maxit+1,))
             #compute f and the gradient of f in x0
             f_{vals}[0] = f(x0, **kwargs)
             grad_norms[0] = np.linalg.norm(grad_f(x0, **kwargs))
             for iter in range(maxit): # loop over maxit
                 next_x = x0 - alpha*grad_f(x0, **kwargs) # compute the next x
                 # store values for plotting
                 f_vals[iter+1] = f(next_x, **kwargs)
                 grad_norms[iter+1] = np.linalg.norm(grad_f(next_x, **kwargs))
                 #check stopping conditions
                 if (np.linalg.norm(grad_f(next_x, **kwargs))<(tolf*np.linalg.norm(x0))) or (np.linalg.norm(next_x-x0)<tolx):</pre>
                     break
                 x0 = next_x # update current x
             # truncate not used part of the array
             f_vals = f_vals[:iter+1]
             grad_norms = grad_norms[:iter+1]
             return x0, iter, f_vals, grad_norms
In [67]: alpha_vals = np.linspace(0.1, 0.005, 10)
         results = []
         n=10
         for alpha in alpha vals:
             sol1 = GD(f1, grad_f1, x0=np.array([0,0]).reshape(-1,1), alpha=alpha)
             sol2 = GD(f2, grad_f2, x0=np.array([0,0]).reshape(-1,1), alpha=alpha)
             sol3 = GD(f3, grad_f3, x0=np.zeros((n)).reshape(-1,1), alpha=alpha, n=n)
             sol4 = GD(f4, grad_f4, x0=np.zeros((n)).reshape(-1,1), alpha=alpha, n=10, lamb=0.5)
             sol5 = GD(f5, grad_f5, x0=0, alpha=alpha)
             results.append([sol1,sol2,sol3,sol4,sol5]) # shape (10, 5). On the rows we have the values of alpha, on the columns the 5 function
In [68]: fig, axs = plt.subplots(figsize=(15,25), ncols=2, nrows=5, sharex=True)
         fig.suptitle('Gradient descent no backtracking', size=18)
         fig.subplots_adjust(top=0.95)
         for i in range(5): #iterate over the functions
             fun_res = [row[i] for row in results] # array with 10 elements, contains the results obtained for the i-th function for each value of alpha
             for j in range(len(fun_res)): # iterate, for each function, over the results obtained with different values of alpha
                 x_{sol}, k, f_{vals}, grad_{norms} = fun_{res}[j] #unpack the result otained for the i-th function with the j-th value of alpha
                 # plot the norm of the gradient on the left
                 ax = axs[i][0]
                 ax.plot(range(k+1), grad_norms, label=f'$alpha = $ {alpha_vals[j]}', alpha=0.6) # plot the norm of the gradient w.r.t iterations
                 ax.legend()
                 ax.grid(linestyle=':')
                 ax.set_title(f'Plot of the norm of the gradient of $f_{i+1}$')
                 # plot the function on the right
                 ax = axs[i][1]
                 ax.plot(range(k+1), f_vals, label=f'*alpha = * {alpha_vals[j]}', alpha=0.6) # plot the value of the function w.r.t iterations
```

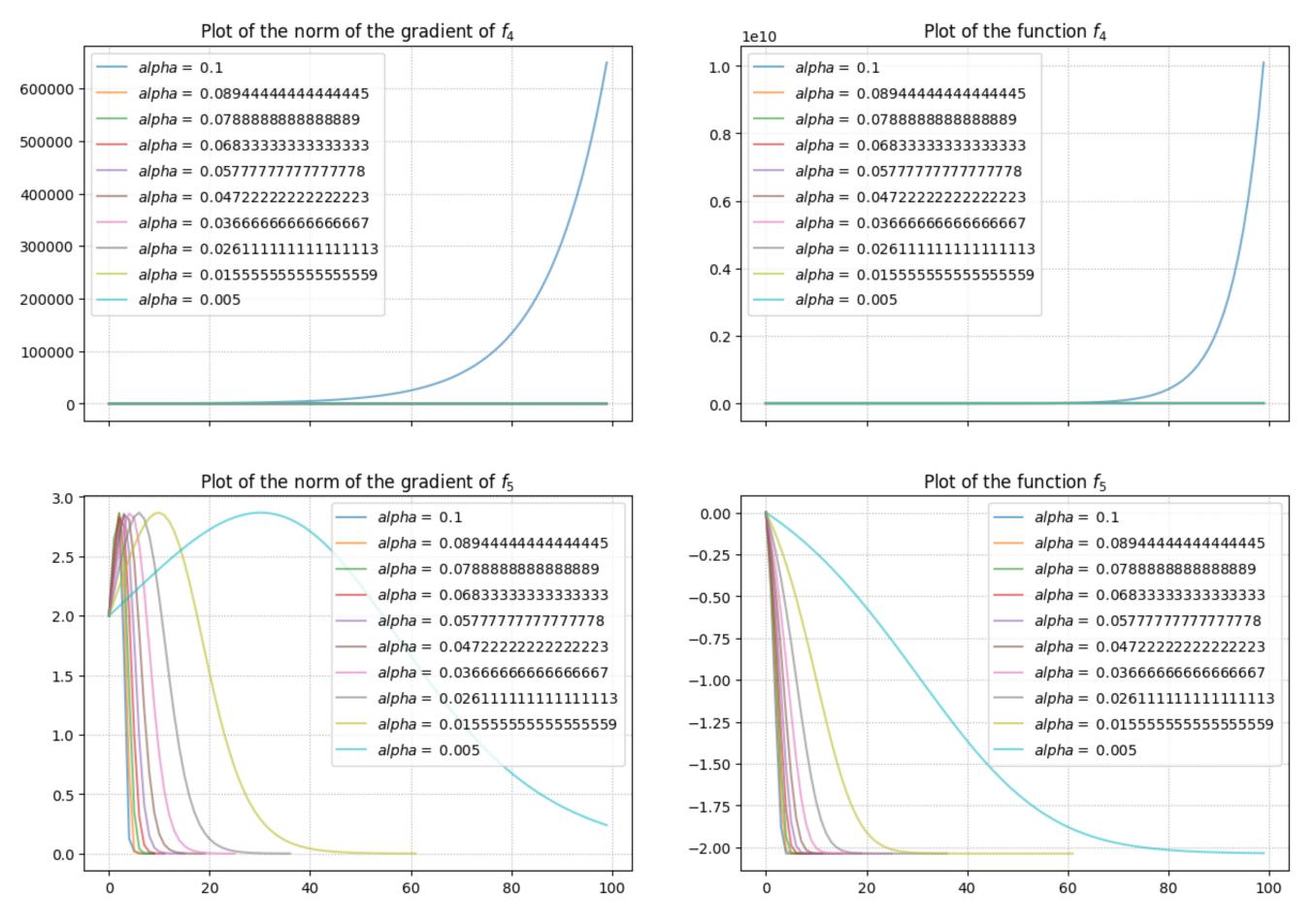
ax.legend()

ax.grid(linestyle=':')

ax.set\_title(f'Plot of the function \$f\_{i+1}\$')

## Gradient descent no backtracking



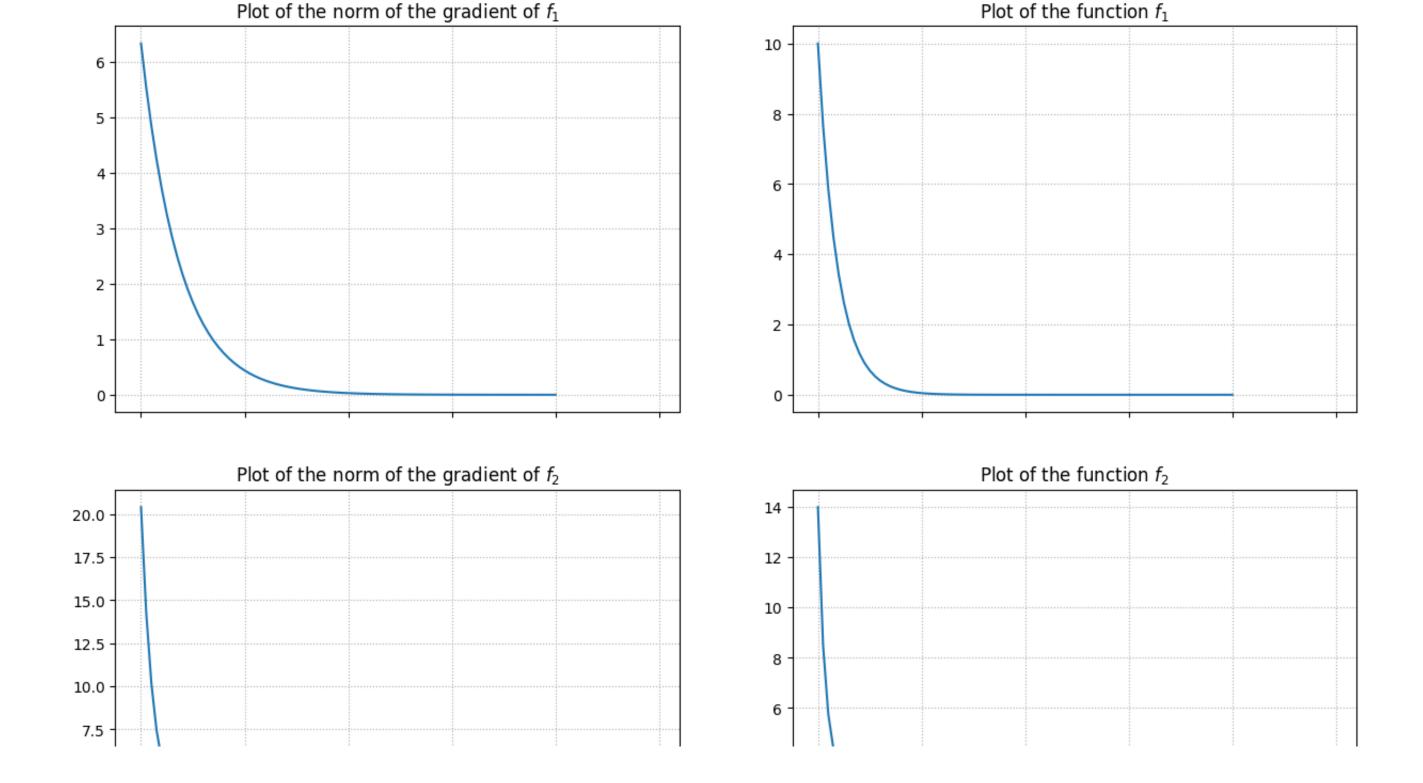


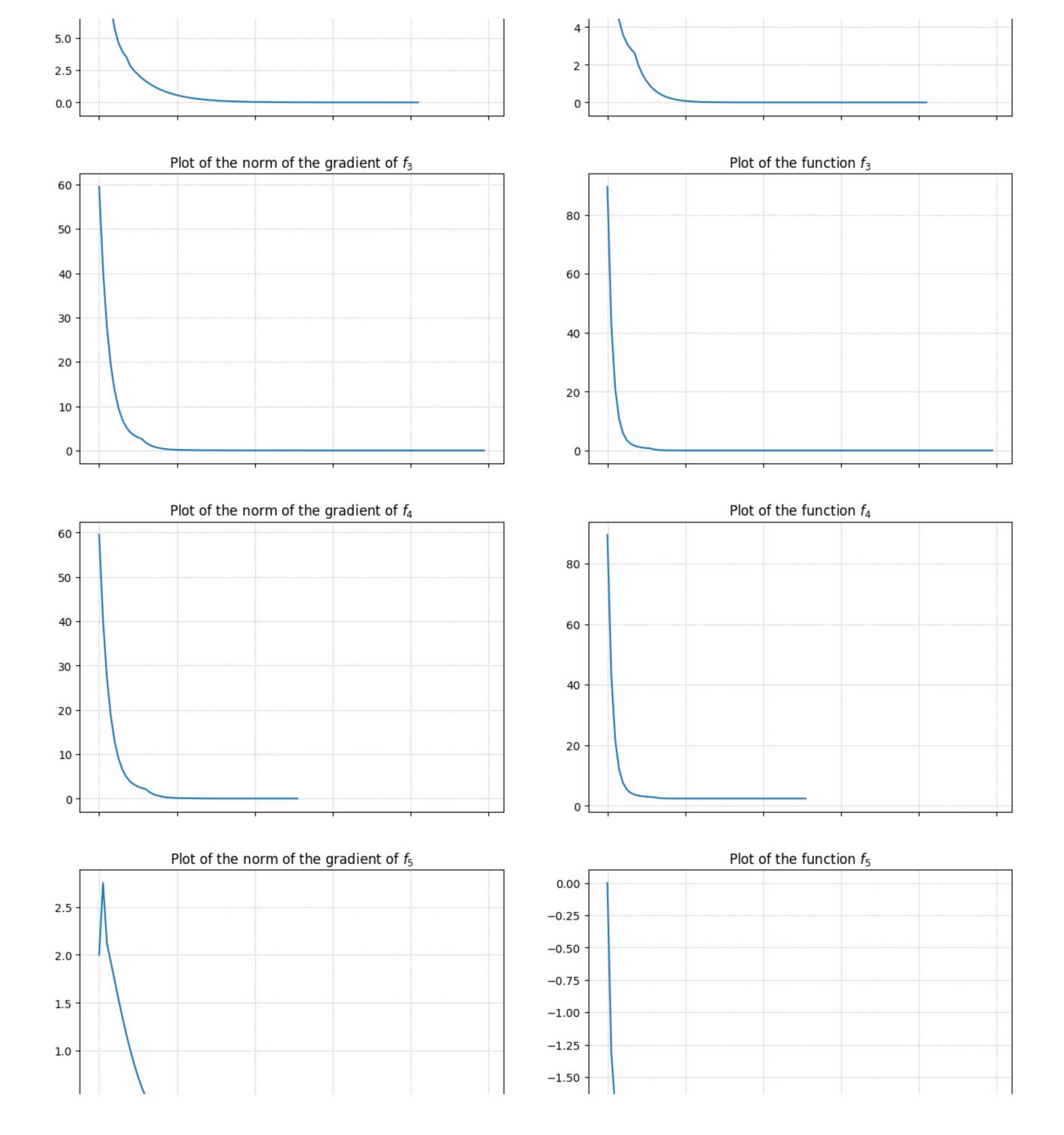
As we can see from this plot, in general, higher is alpha, faster is the convergence. However, when  $\alpha$  becomes too high ( $\alpha=0.1$  for example), the algorithm doesn't converge. This means that the step size is too long escaping from local minima. This happens for the functions  $f_2$ ,  $f_3$  and  $f_4$ . If we remove from the plot of  $f_3$  and  $f_4$  the line associated with  $\alpha=0.1$  we obtain a plot similar to those of  $f_1$  and  $f_5$ .

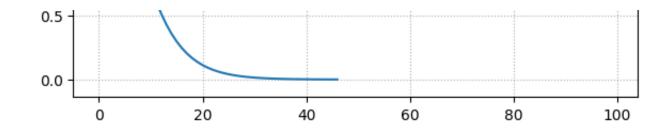
In [69]: def backtracking(f, grad\_f, x, \*\*kwargs): This function is a simple implementation of the backtracking algorithm for the GD (Gradient Descent) method. f: function. The function that we want to optimize.  $grad_f$ : function. The gradient of f(x). x: ndarray. The actual iterate x\_k. alpha = 1c = 0.8tau = 0.25 while  $f(x - alpha * grad_f(x, **kwargs), **kwargs) > f(x, **kwargs) - c * alpha * np.linalg.norm(grad_f(x, **kwargs), 2) ** 2:$ alpha = tau \* alphareturn alpha def GD\_backtracking(f, grad\_f, x0, maxit=100, tolf=1e-5, tolx=1e-5, \*\*kwargs): f\_vals = np.zeros((maxit+1,)) grad\_norms = np.zeros((maxit+1,))  $f_{vals}[0] = f(x0, **kwargs)$ grad\_norms[0] = np.linalg.norm(grad\_f(x0, \*\*kwargs)) for iter in range(maxit): alpha = backtracking(f, grad\_f, x0, \*\*kwargs)  $next_x = x0 - alpha*grad_f(x0, **kwargs)$ f\_vals[iter+1] = f(next\_x, \*\*kwargs) grad\_norms[iter+1] = np.linalg.norm(grad\_f(next\_x, \*\*kwargs)) #check stopping conditions if (np.linalg.norm(grad\_f(next\_x, \*\*kwargs))<(tolf\*np.linalg.norm(x0))) or (np.linalg.norm(next\_x-x0)<tolx):</pre> break  $x0 = next_x$ f\_vals = f\_vals[:iter+1] grad\_norms = grad\_norms[:iter+1] return x0, iter, f\_vals, grad\_norms

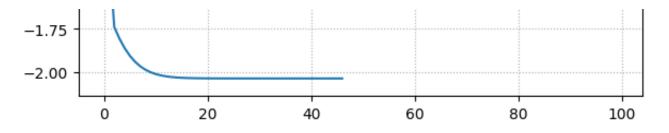
In [64]: results = []  $x_{sol} = GD_{backtracking}(f1, grad_f1, x0=np.zeros(2))$ results.append(x\_sol) x\_sol = GD\_backtracking(f2, grad\_f2,x0=np.zeros(2)) results.append(x\_sol) x\_sol = GD\_backtracking(f3, grad\_f3,x0=np.zeros(n), n=10) results.append(x\_sol)  $x_{sol} = GD_{backtracking}(f4, grad_f4, x0=np.zeros(n), n=10, lamb=0.5)$ results.append(x\_sol) x\_sol = GD\_backtracking(f5, grad\_f5,x0=np.array([0])) results.append(x\_sol) fig, axs = plt.subplots(figsize=(15,25), ncols=2, nrows=5, sharex=True) fig.suptitle('Gradient descent with backtracking', size=18) fig.subplots\_adjust(top=0.95) for i in range(5): #iterate over the functions x\_sol, k, f\_vals, grad\_norms = results[i] # plot the norm of the gradient on the left ax = axs[i][0]ax.plot(range(k+1), grad\_norms) # plot the norm of the gradient w.r.t iterations ax.grid(linestyle=':') ax.set\_title(f'Plot of the norm of the gradient of \$f\_{i+1}\$') # plot the function on the right ax = axs[i][1]ax.plot(range(k+1), f\_vals) # plot the value of the function w.r.t iterations ax.grid(linestyle=':') ax.set\_title(f'Plot of the function \$f\_{i+1}\$')

## Gradient descent with backtracking





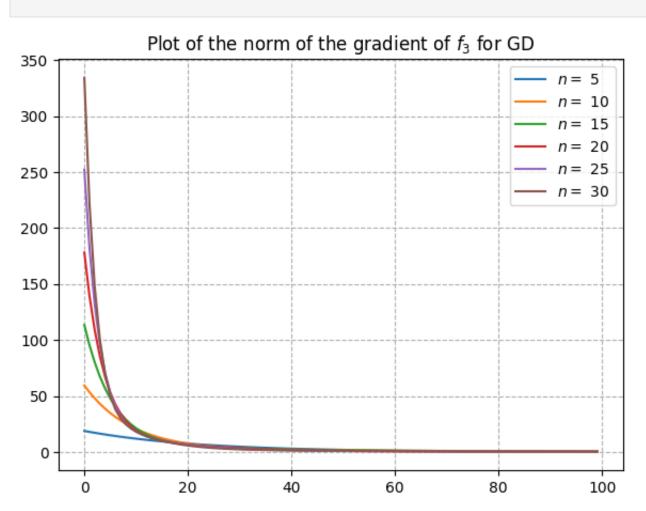


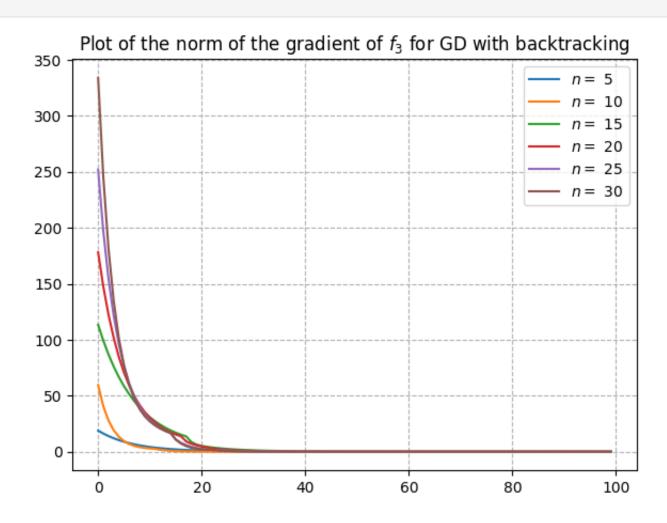


Looking at the plots of gradient descent with and without backtracking we can notice that GD with backtracking sometimes converges before sometimes converges after if compared with the best value of  $\alpha$  of GD without backtracking. This is a reasonable behavior since backtracking algorithm chooses a good value of  $\alpha$  but it is not guaranteed that the chosen value is the best.

• Try  $f_3$  for different values of n (e.g.  $n=5,10,15,\ldots$ ).

In [81]: #GD and GD\_backtracking for different values of n for f3  $n_{vals} = [5,10, 15, 20, 25,30]$ alpha = 0.005results = [] for n in n\_vals: res\_gd =  $GD(f3, grad_f3, x0=np.zeros((n)), alpha=alpha, n=n)$ res\_gdb = GD\_backtracking(f3, grad\_f3,x0=np.zeros((n)), n=n) results.append([res\_gd, res\_gdb]) fig, axs = plt.subplots(figsize=(15,5), nrows=1, ncols=2) qd results = [row[0] for row in results] gdb\_results = [row[1] for row in results] for gd, gdb, i in zip(gd\_results, gdb\_results, np.arange(n)): x\_sol\_gd, k\_gd, f\_vals\_gd, grad\_norms\_gd = gd x\_sol\_gdb, k\_gdb, f\_vals\_gdb, grad\_norms\_gdb = gdb ax = axs[0] $ax.plot(range(k_gd+1), grad_norms_gd, label=f'$n = $ {n_vals[i]}')$ ax.grid(visible=True, linestyle='--') ax.legend() ax.set\_title(f'Plot of the norm of the gradient of \$f\_3\$ for GD') ax = axs[1]ax.plot(range(k\_gdb+1), grad\_norms\_gdb, label=f'\$n = \$ {n\_vals[i]}') ax.grid(visible=True, linestyle='--') ax.set\_title(f'Plot of the norm of the gradient of \$f\_3\$ for GD with backtracking') ax.legend()

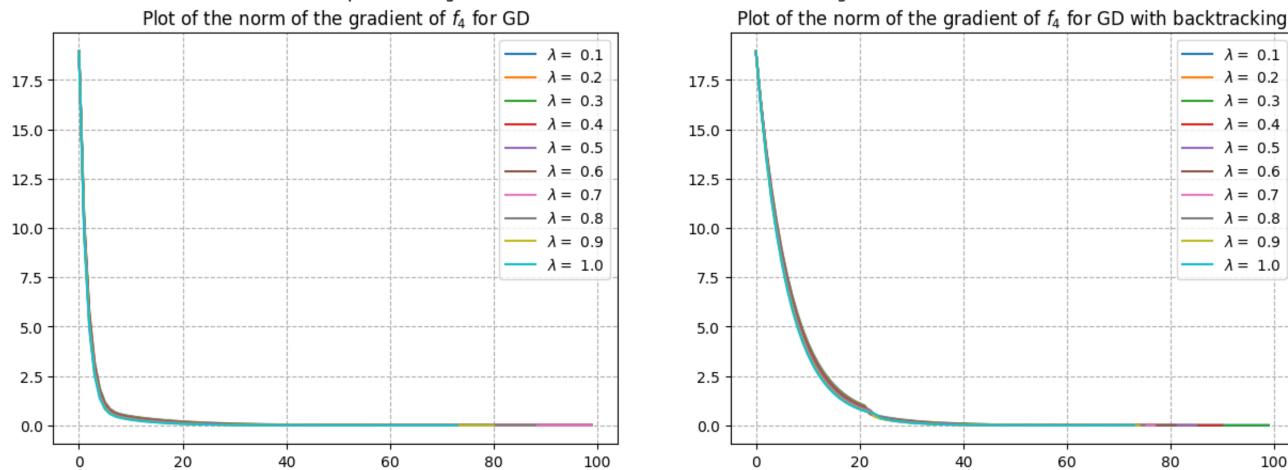




The convergence is a little bit higher for small values of n but the general behavior is the same, regardless the value of n

- Try  $f_4$  for different values of  $\lambda$  and comment the result.
- In [82]: #plot of the gradient norm of f4 for different values of lambda given n and alpha lambda\_vals = np.linspace(0.1, 1, 10) alpha = 0.05n = 5results = [] for lamb in lambda\_vals: res\_gd =  $GD(f4, grad_f4, x0=np.zeros((n)), alpha=alpha, n=n, lamb=lamb)$ res\_gdb = GD\_backtracking(f4, grad\_f4,x0=np.zeros((n)), n=n, lamb=lamb) results.append([res\_gd, res\_gdb]) fig, axs = plt.subplots(figsize=(15,5), nrows=1, ncols=2) gd\_results = [row[0] for row in results] gdb\_results = [row[1] for row in results] for i in range(len(lambda\_vals)): x\_sol\_gd, k\_gd, f\_vals\_gd, grad\_norms\_gd = gd\_results[i] x\_sol\_gdb, k\_gdb, f\_vals\_gdb, grad\_norms\_gdb = gdb\_results[i] ax = axs[0]ax.plot(range(k\_gd+1), grad\_norms\_gd, label=f'\$\lambda = \$ {lambda\_vals[i]:.1f}') ax.grid(visible=True, linestyle='--') ax.legend() ax.set\_title(f'Plot of the norm of the gradient of \$f\_4\$ for GD') ax = axs[1]ax.plot(range(k\_gdb+1), grad\_norms\_gdb, label=f'\$\lambda = \$ {lambda\_vals[i]:.1f}') ax.grid(visible=True, linestyle='--') ax.legend() ax.set\_title(f'Plot of the norm of the gradient of \$f\_4\$ for GD with backtracking')

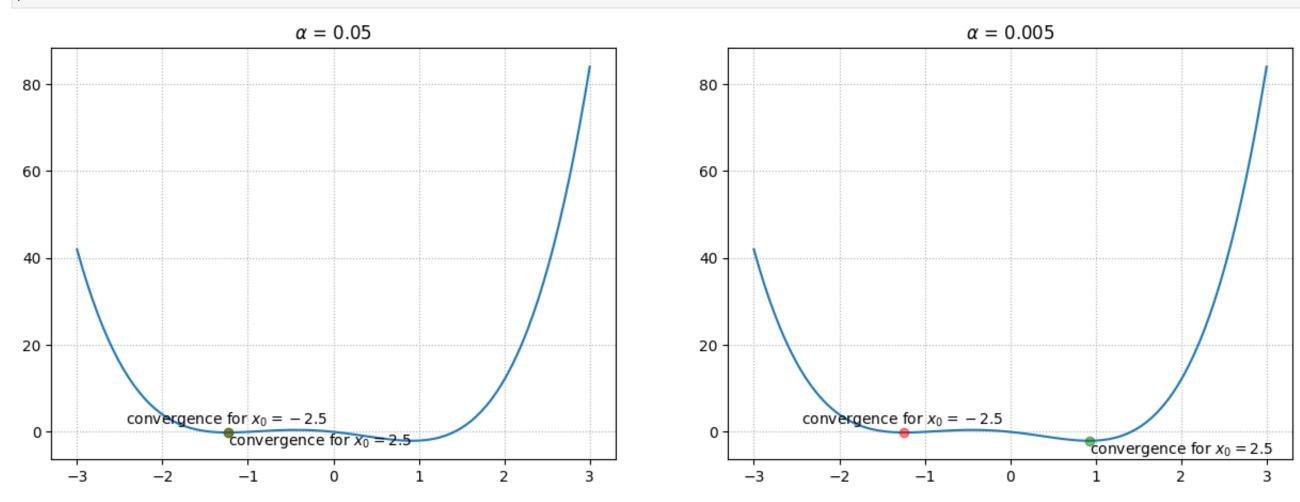
## plot of the gradient norm of f4 for different values of $\lambda$ given n and $\alpha$



As we can see from the plot, higher values  $\lambda$  result in a faster convergence.

• Plot the graph of the non-convex function 5 in the interval [-3,3], and test the convergence of GD with different values of x0 (of your choice) and different step-sizes. When is the convergence point the global minimum?

```
In [85]: import matplotlib.pyplot as plt
         alpha_vals = np.array([0.05, 0.005])
         x = np.linspace(-3,3,100)
         fig, axs = plt.subplots(figsize=(15, 5), nrows=1, ncols=2)
         for i in range(len(alpha_vals)):
             ax = axs[i]
             ax.plot(x, f5(x))
             ax.grid(linestyle=':')
             sol1 = GD(f5, grad_f5,x0=-2.5, alpha=alpha_vals[i])
             sol2 = GD(f5, grad_f5,x0=2.5, alpha=alpha_vals[i])
             ax.plot(sol1[0], f5(sol1[0]), marker='o', color='r', alpha=0.5)
             ax.text(sol1[0]-1.2, f5(sol1[0])+2, 'convergence for x_0=-2.5')
             ax.plot(sol2[0], f5(sol2[0]), marker='o', color='green', alpha=0.5)
             ax.text(sol2[0], f5(sol2[0])-3, 'convergence for x_0=2.5')
             ax.set_title(f'$\\alpha$ = {alpha_vals[i]}')
         plt.show()
```



As we can see, with an higher value of  $\alpha$  the GD with  $x_0=2.5$  escapes from the global minimum. This is not true for  $\alpha=0.005$ . This means that  $\alpha=0.05$  is too big and the global minimum is not detected.