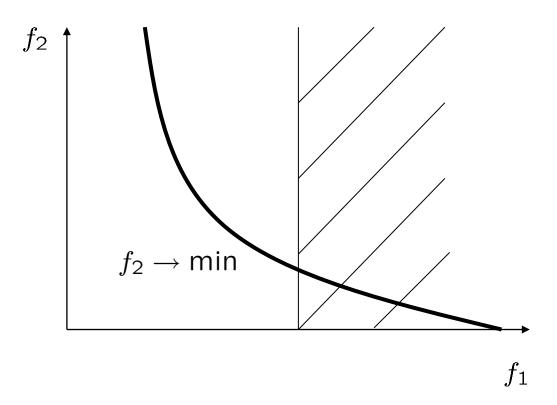
# Exercise: Hill-climbing Methods (Local Optimization)

- Exercise (based on slides and python examples)
- Pareto optimization
- Adding objective function to Desdeo



#### ε-Constraint method

 $f_2 \rightarrow min$ ,  $f_1 \leftarrow psilon$ , epsilon \in  $\{e1, e2, ....en\}$ 



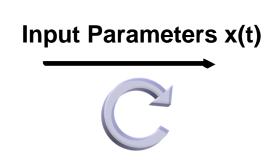
With the dimension the number of  $\epsilon$  combinations grows exponentially.

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#### Basic strategy in Black-box optimization

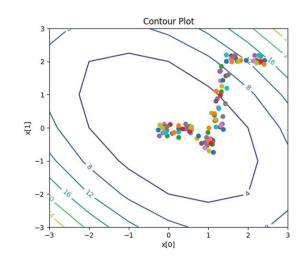
Black-box
Optimierungssoftware

- Stochastic Hillclimbing
- 2. Gradient Descent
- Newton Method
- 4. Simulated Annealing
- 5. Evolutionary Algorithm
- 6. Bayesian Optimization
- 7. Etc.



Simulator/Evaluator

Zielfunktionswerte,
Restriktionsverletzungen
f(x(t))+penalty(r(x(t))





## Hill-climbing Methods for Single-Objective Optimization

Path oriented (hill climbers) can be defined by a general iterative formula:

$$\mathbf{x}_{t+1} = \mathbf{x}_t + \sigma_t \mathbf{d}_t$$

 $x_t$ : Current search point

 $\sigma_t$ : Step size

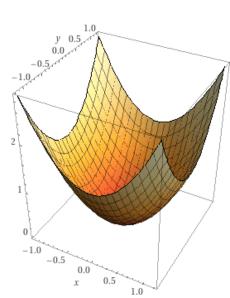
 $\mathbf{d}_t$ : Current search direction



Hill-climbers generates a sequence of points  $\{x_t\}_{t=1,2,...}$  that gradually improve the value of the objective function.



#### Simple 2-D stochastic hillclimber



```
best_eval = objective(best)
                             curr, curr_eval = best, best_eval # current working solution
                             scores = list()
                             for i in range(n_iterations):
                                 # take a step
                                 candidate = [curr[0] +rand()*step_size[0]-step_size[0]/2.0,
                                               curr[1]+rand()*step_size[1]-step_size[1]/2.0]
                                 print('>%d f(%s) = %.5f, %s' % (i, best, best_eval, candidate))
                                 #+ randn(len(bounds)) * step_size
                                 # evaluate candidate point
                                 candidate_eval = objective(candidate)
                                 # check for new best solution
                                 if candidate_eval < best_eval:</pre>
                                     # store new best point
                                     best, best_eval = candidate, candidate_eval
# objective function
                                     # keep track of scores
def objective(x):
                                      scores.append(best_eval)
    return x[0]**2+x[1]**2
                                     # report progress
                                      print('>%d f(%s) = %.5f' % (i, best, best_eval))
                                     # current best
                                      curr=candidate
                             return [best, best_eval, scores]
```

black-box optimization software

# generate an initial point

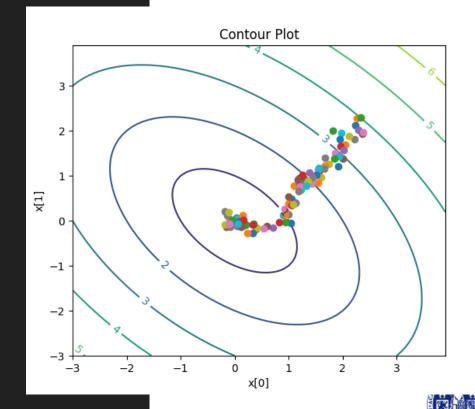
# evaluate the initial point

best = init

def local\_hillclimber(objective, bounds, n\_iterations, step\_size,init):

### Plotting the history

```
bounds=asarray([[-3.0,3.0],[-3.0,3.0]])
step_size=[0.4,0.4]
n_iterations=100
init=[2.4,2.0]
best, score, points, scores, = local_hillclimber(objective,
                                                  bounds, n_iterations,
                                                  step_size, init)
n, m = 7, 7
start = -3
x_vals = np.arange(start, start+n, 1)
y_vals = np.arange(start, start+m, 1)
X, Y = np.meshgrid(x_vals, y_vals)
print(X)
print(Y)
fig = plt.figure(figsize=(6,5))
left, bottom, width, height = 0.1, 0.1, 0.8, 0.8
ax = fig.add_axes([left, bottom, width, height])
Z = (X**2 + Y**2 + X*Y)
cp = ax.contour(X, Y, Z)
ax.clabel(cp, inline=True,
          fontsize=10)
ax.set_title('Contour Plot')
ax.set_xlabel('x[0]')
ax.set_ylabel('x[1]')
for i in range(n_iterations):
    plt.plot(points[i][0],points[i][1],"o")
plt.show()
```



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#### Penalty method for constraints

- In single objective black box optimization we can implement constraints in two ways
- Box constraints:
  - When hillclimbing method leaves the search region the point is projected back to the search region
- Implicit constraints:
  - Constraints that require the black-box function to be evaluated can be handled by a (metric) penalty value.

Metric Penalty method

$$\min f(\mathbf{x})$$

s.t. 
$$c_i(\mathbf{x}) \leq 0 \ orall i \in I$$

Replace objective f(x) by  $\Phi_k(x)$ :

$$\min \Phi_k(\mathbf{x}) = f(\mathbf{x}) + \sigma_k \ \sum_{i \in I} \ g(c_i(\mathbf{x}))$$

with:  $g(c_i(\mathbf{x})) = \max(0, c_i(\mathbf{x}))^2$   $\sigma_k$  can be constant or increasing over time (k).

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You can also add a constant C(x)>f\_max to the penal whenever constraint is violated to make sure that it is worse.

#### Simulated Annealing

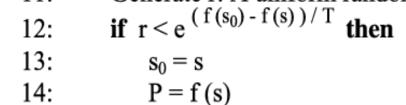
Kirkpatrick, S. Optimization by simulated annealing: Quantitative studies. *J Stat Phys* **34,** 975–986 (1984). https://doi.org/10.1007/BF01009452

```
1: Set the initial temperature (T=T0)
2: Create initial solution (s<sub>0</sub>)
3: P = Calculate f(s_0)
4: while (P > 0)
      Create Neighbor (s)
5:
      Calculate f (s)
6:
      if (f(s) \le P) then
8:
       \mathbf{s}_0 = \mathbf{s}
       P = f(s)
9:
10:
       else
11:
         Generate r: A uniform random number
```

Stochastic Hillclimbing inspired by Annealing process in crystals.

Simulated Annealing can always accepts improvements, but also worse solutions with some probability.

In order to get to global optima one might have to accept steps to get worse temporarily



15: Reduce temperature

16: **Return s<sub>0</sub>** 

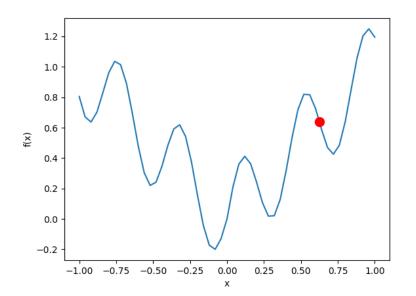




https://makeitfrommetal.com/beginners-guide-on-how-to-anneal-steel/

### Simulated Annealing

```
# objective function
def objective(x):
    return np.abs(x[0])+0.3*np.sin(x[0]*15);
```



1-D Objective Function with local optima  $f(x) = |x| + 0.3\sin(15 x), x \in [-1,1]$ 

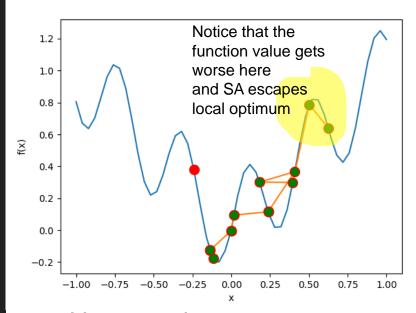
Simulated Annealing can be implemented in 2-D and N-D (homework)

```
# simulated annealing algorithm
def simulated_annealing(objective, bounds, n_iterations,
                        step_size, temp,init):
    st=[]
    c=[]
    cscore=[]
    # generate an initial point
    best=[init]
    # evaluate the initial point
    best_eval = objective(best)
    # current working solution
    curr, curr_eval = best, best_eval
    scores = list()
    # run the algorithm
    for i in range(n_iterations):
        # take a step
        candidate = curr + randn(len(bounds)) * step_size
        st.append(candidate)
        # evaluate candidate point
        candidate_eval = objective(candidate)
        # keep track of scores
        scores.append(candidate_eval)
        # check for new best solution
        if candidate_eval < best_eval:</pre>
            # store new best point
            best, best_eval = candidate, candidate_eval
            # report progress
            print('>%d f(%s) = %.5f' % (i, best, best_eval))
        # difference between candidate and current point evaluation
        diff = candidate_eval - curr_eval
        # calculate temperature for current epoch
        t = temp / float(i + 1)
        # calculate metropolis acceptance criterion
        metropolis = exp(-diff / t)
        # check if we should keep the new point
        if diff < 0 or rand() < metropolis:</pre>
            # store the new current point
            curr, curr_eval = candidate, candidate_eval
        c.append(curr)
        cscore.append(curr_eval)
    return [best, best_eval, st, scores, c, cscore]
```

```
# Random number generator initializatiom
seed(1)
# define range for input
1h = -1
ub=1
bounds = asarray([[lb, ub]])
# define the total iterations
n_{iterations} = 100
# define the maximum step size
step_size = 0.2
# initial temperature
temp = 1.0
# initial point
init=0.3
# perform the simulated annealing search
best, score, st, scores, c, cscores = \
     simulated_annealing(objective, bounds,
                         n_iterations, step_size,
                         temp, init)
def f1d(x):
    a=[]
     a.append(x)
    return objective(a)
                                # lower limit
x = np.linspace ( start = lb
                                  # upper limit
                 , num = 51
                                 # generate 51 points between 0 and 3
               # This is already vectorized, that is, y will be a vector!
y = f1d(x)
plt.plot(x, y)
plt.show()
for i in range(n_iterations):
    plt.plot(x, y)
     plt.xlabel("x")
    plt.ylabel("f(x)")
    plt.plot(c[0:i], cscores[0:i], marker="0", markersize=10,
              markeredgecolor="red", markerfacecolor="green")
    plt.plot(st[i], scores[i], '.r', ms=20)
    plt.show()
     time.sleep(1)
```

### Simulated Annealing

Plot shows the function and linked successful moves



Homework:

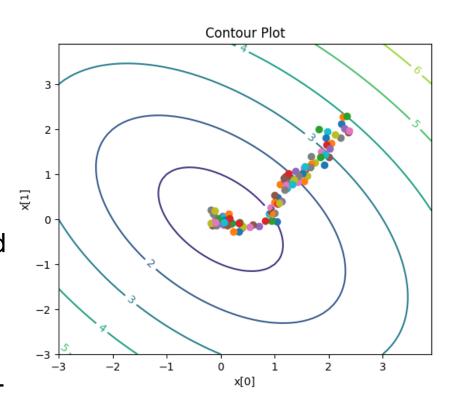
Optimize design with N-D Simulated annealing

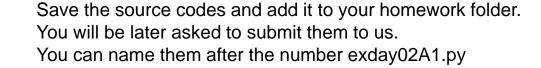
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https://trinket.io/python3/b22300f21e

# Task Local Optimization with Constraints/Penalties in Python

- A
- Modify simulated annealing example that it can be visualized in 2-D space
- Implement a method that restricts the variables to the bounds (ranges) in the random steps
- 3. Implement a penalty method for the simulated annealing
- 4. Solve and visualize the two dimensional problem:
- 5. Visualize  $|x| + 0.3\sin(15 x) + |y| + 0.3\sin(15 y)$  as a countour plot



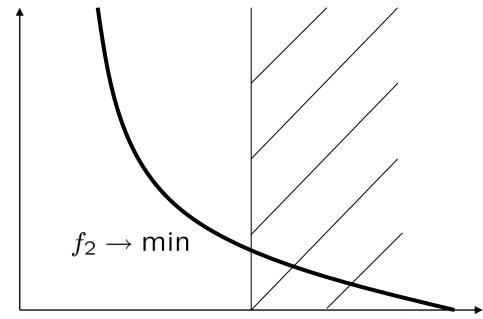


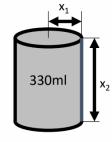


#### Exercise 2

1. Visualize the optimization  $f_2$  problem of the optimal tin as a countour plot (or cone, if you prefer this problem)

- 2. Visualize the Area objective min)
- 3. Visualize the constraint  $Volume(r,h) \ge L$ , for level 330ml
- 4. Find the efficient points by solving a series of optimization problems where the volume is a constraint  $Volume(r,h) \ge L$ , for different levels (Pareto optimization, epsilon constraint method) (scatter plot)





Problem sketch

$$f(\mathbf{x}) = 2\pi x_1 x_2 + 2\pi (x_1)^2 \to \min$$
$$h(\mathbf{x}) = 2\pi x_2 (x_1)^2 - 330 = 0$$
$$\mathbf{x} \in \begin{bmatrix} 0 \\ 0 \end{bmatrix}, \begin{pmatrix} \infty \\ \infty \end{bmatrix} \subset \mathbb{R}$$

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