

MET CS 767 - Project 1

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1 Problem Statement

Use simulated annealing to best approximate the optimal inputs, (x_1, x_2, x_3) such that $\forall x \in \{x_1, x_2, x_3\} (-10 \leq x \leq 10)$ and such that the following function, $f(x_1, x_2, x_3)$, yields the minimal output:

$$f(x_1, x_2, x_3) : x_1^2 + x_2^2 + x_3^2 \quad (1)$$

2 Approach

2.1 General Simulated Annealing

I chose to construct a general simulated annealing approach that could be used in any arbitrary problem domain, The following pseudocode defines a single sweep of a single run of the algorithm:

Require: s_{init} , Neighbor, E, P

```
 $s \leftarrow s_{init}$ 
 $e \leftarrow E(s)$ 
 $t \leftarrow 1.0$ 
 $t_{min} \leftarrow 0.0001$ 
 $\alpha \leftarrow 0.9$ 
 $k \leftarrow 1000$ 
while  $t \geq t_{min}$  do
  for all  $0 \leq i < k$  do
     $s' \leftarrow \text{Neighbor}(s)$ 
     $e' \leftarrow E(s')$ 
     $p_{accept} \leftarrow P(e, e', t)$ 
     $p_{rand} \leftarrow \text{Random}(0, 1)$ 
    if  $p_{accept} > p_{rand}$  then
       $s \leftarrow s'$ 
       $e \leftarrow e'$ 
    end if
  end for
   $t \leftarrow t * \alpha$ 
end while
return  $(s, e, t)$ 
```

Such a sweep would be conducted multiple times using the same pseudorandom generator for a single run in order to provide a better stochastic minimum to the problem. Then, multiple runs (each with multiple sweeps) would be performed. Each run uses a pseudorandom generator with a different seed.

The inputs and variables for the above formula are defined as the following:

- s_{init} : Initial state within the problem domain. This is to be chosen at random.
- $\text{Neighbor}(s)$: Function that accepts a state, s , as input and outputs a randomly chosen local neighbor, s' , within the problem domain. It should randomly select a neighbor that is considered local to s .
- $E(s)$: Function that accepts a state, s , as input and outputs the state's energy, e , with regards to the problem space. The energy should be the output of the problem function f with s as the input.

- $P(e, e', t)$ Function that accepts the energy, e , of current state, s , the potential energy, e' , of new state, s' , and the annealing temperature, t . It outputs an acceptable probability, p_{accept} , such that the new state, s' , is accepted as the current state, s , if $p_{accept} > \text{Random}(0, 1)$. This function should return $p_{accept} \leftarrow 1$ if $e' < e$. Otherwise, $p_{accept} \leftarrow \exp(e - e')/t$.
- t : The progressively cooling temperature of the annealing system. Starts at a temperature of 1 unit and eventually reduced to t_{min} . For each iteration, t is reduced in the following way: $t \leftarrow t * \alpha$.
- t_{min} : The temperature to which t should reach before the system exists
- α : The rate of temperature decrease of t for each annealing iteration.
- k : The number of steps for each temperature iteration. During each step, a new potential state, s' , and potential energy, e' , are generated and conditionally set to s and e , respectively given an acceptance probability p_{accept} .

2.2 Problem-Specific Simulated Annealing

It is the responsibility of the problem space to define s_{init} , Neighbor, E, P as defined above. For the problem at hand, a state s is defined as a tuple of (x_1, x_2, x_3) values. s_{init} is defined as randomly chosen:

$$s_{init} \leftarrow (\text{Random}(-10, 10), \text{Random}(-10, 10), \text{Random}(-10, 10)) \quad (2)$$

Furthermore, Neighbor is defined as a function that takes state, $s = (x_1, x_2, x_3)$, as input and outputs a local neighbor, s' , in which a randomly chosen x_i for $i \in \{1, 2, 3\}$ is updated with a random value, $\text{Random}(-10, 10)$:

$$\text{Neighbor}(s) : s' \leftarrow s; s'_{i \leftarrow \text{Random}(1, 2, 3)} \leftarrow \text{Random}(-10, 10) \quad (3)$$

Furthermore, E is defined as a function that takes a state, $s = (x_1, x_2, x_3)$, as input and outputs an energy, e , set to the result of problem function, f , as such:

$$E(s = (x_1, x_2, x_3)) : e \leftarrow f(x_1, x_2, x_3) \quad (4)$$

Furthermore, P is an acceptance probability function that takes as input a current energy, e , a potential energy, e' , and temperature, t . It then outputs a probability, p_{accept} such that:

$$P(e, e', t) : p_{accept} \leftarrow (1.0 \text{ If } (e' < e) \text{ Otherwise } \exp((e - e')/t)) \quad (5)$$

3 Implementation and Project Layout

The entire solution is implemented in JavaScript with a HTML frontend. Specifically, the algorithm is broken up into two files:

- `js/SimulatedAnneal.js`: General simulated annealing algorithm that takes s_{init} , Neighbor, E, P as input.
- `js/MinSumOfSquares.js`: Problem specific simulated annealing algorithm that implements s_{init} , Neighbor, E, P, and leverages `js/SimulatedAnneal.js` for the general simulated annealing heuristic.
- `js/PRG.js`: A custom pseudo random generator that accepts a seed tuple of integers, (m, c) . This was necessary because pseduorandom number generators vary from browser to browser.

In addition, the project includes a front end which allows a user to configure a seed, i , for the PRG, a number of sweeps, n , for a single run of the algorithm, as well as values: t, t_{min}, α, k , for the the simulated annealing algorithm sweep. The frontend outputs a list of the states, $s_i \leftarrow (x_1, x_2, x_3)$ and energies, e , for the t_{min} of each i . It also outputs the minimum e of all sweeps (the stochastic minimum of the problem).

The front end also outputs an animation of the problem space with each state, s_i for each annealing iteration of each sweep. The animation is a cube in which each dimation is different value, x_i for $i \in \{1, 2, 3\}$. Over time, states

s_i , should converge towards the middle of the box where $(x_1, x_2, x_3) = (0, 0, 0)$.

This functionality is implemented in the following files:

- `img/axis.png`: Represents each of the dimensions of the animation - (x_1, x_2, x_3) .
- `js/Chart.js`: Generates the animation using HTML's canvas element.
- `style/normalize.css`: Standardize the css styles across all browsers (i.e. fonts, sizes, padding, etc ...)
- `style/style.css`: Styles specific to the project frontend in `index.html`.
- `index.html`: Frontend HTML code to run algorithm and resulting animation.

3.1 Execution

Set up a simple python server to serve up `index.html` and corresponding `js/*`, `style/*`, `img/*` resources.

```
$ cd /path/to/project/  
$ python -m SimpleHTTPServer
```

This should display something like: "Serving HTTP on 0.0.0.0 port 8000 ...". At that point, go to `http://localhost:8000` to view the frontend. (Note that this is not required if viewing the frontend with Google Chrome. In such, just open `index.html` in the Chrome browser: e.g. `file:///<path-to-project>/index.html`)

4 Inputs and Solution

4.1 Inputs

The following inputs were used for a single run (multiple sweeps). Each are the default within the configuration on the frontend:

- t : 1.0
- t_{min} : 0.00001
- α : 0.95
- k : 100
- $seed$: 14759809013
- $sweeps$: 25

4.2 Solution

The resulting, best-effort, optimal solution state was: $(x_1, x_2, x_3) = (0.000875, -0.00284, -0.000318)$. The solution was generated over 225 iterations of annealing to reach a temperature of 0.000104. The resulting minimum value of f was: $\min(f(x_1, x_2, x_3)) = 0.00000895$.

The previous was repeated for multiple runs of the following seeds: $\{14759809013, 743, 113819723, 1931117\}$. The respective results were as follows:

- $seed = 14759809013$: $(x_1, x_2, x_3) = (0.000875, -0.00284, -0.000318)$, $\min(f(x_1, x_2, x_3)) = 0.00000895$
- $seed = 743$: $(x_1, x_2, x_3) = (0.0000925, -0.000734, 0.00476)$, $\min(f(x_1, x_2, x_3)) = 0.0000232$
- $seed = 113819723$: $(x_1, x_2, x_3) = (-0.00683, -0.0133, -0.00835)$, $\min(f(x_1, x_2, x_3)) = 0.000293$
- $seed = 14759809013$: $(x_1, x_2, x_3) = (0.00108, 0.00891, 0.0120)$, $\min(f(x_1, x_2, x_3)) = 0.000224$

Thus, the minimal of all runs is: $(x_1, x_2, x_3) = (0.0000925, -0.000734, 0.00476)$, $\min(f(x_1, x_2, x_3)) = 0.0000232$. These results may be reproduced by simply using the defaults on the frontend web application.

5 Resources

- https://en.wikipedia.org/wiki/Simulated_annealing