

Introduction to Programming in Python

Assignment 5 (Atomic Nature of Matter) Discussion

Part I (Warmup Problems) · Problem 1 (Sum of Integers)

Implement the function `_sumOfInts()` in `sum_of_ints.py` that takes an integer n as argument and returns the sum $S(n) = 1 + 2 + 3 + \cdots + n$, computed recursively using the recurrence equation

$$S(n) = \begin{cases} 1 & \text{if } n = 1, \\ n + S(n - 1) & \text{if } n > 1. \end{cases}$$

```
>_ ~/workspace/atomic_nature_of_matter
```

```
$ python3 sum_of_ints.py 100  
5050
```

Part I (Warmup Problems) · Problem 1 (Sum of Integers)

Base case: if $n = 1$, return 1

Recursive step: return n plus `_sumOfInts($n - 1$)`

Part I (Warmup Problems) · Problem 2 (Bit Counts)

Implement the functions `_zeros()` and `_ones()` in `bits.py` that take a bit string (ie, a string of zeros and ones) `s` as argument and return the number of zeros and ones in `s`, each computed recursively

```
>_ ~/workspace/atomic_nature_of_matter
```

```
$ python3 bits.py 1010010010011110001011111
zeros = 11, ones = 14, total = 25
```

Part I (Warmup Problems) · Problem 2 (Bit Counts)

`_zeros(s)`

- Base case: if s is empty, return 0
- Recursive step: if the first character of s is a 0, return 1 plus `_zeros(s[1 :])`; otherwise return `_zeros(s[1 :])`

`_ones(s)`

- Base case: if s is empty, return 0
- Recursive step: if the first character of s is a 1, return 1 plus `_ones(s[1 :])`; otherwise return `_ones(s[1 :])`

Part I (Warmup Problems) · Problem 3 (String Reversal)

Implement the function `_reverse()` in `reverse.py` that takes a string `s` as argument and returns the reverse of the string, computed recursively

```
>_ ~/workspace/atomic_nature_of_matter
```

```
$ python3 reverse.py bolton
notlob
$ python3 reverse.py madam
madam
```

Part I (Warmup Problems) · Problem 3 (String Reversal)

Set n to the length of s

Base case: if $n = 0$, return the empty string

Recursive step: Return a concatenation of $s[n - 1]$ and `_reverse($s[: n - 1]$)`

Part I (Warmup Problems) · Problem 4 (Palindrome)

Implement the function `_isPalindrome()` in `palindrome.py`, using recursion, such that it returns `True` if the argument `s` is a palindrome (ie, reads the same forwards and backwards), and `False` otherwise

```
>_ ~/workspace/atomic_nature_of_matter
```

```
$ python3 palindrome.py bolton
False
$ python3 palindrome.py madam
True
```


Part I (Warmup Problems) · Problem 4 (Palindrome)

Set n to the length of s

Base case: if $n = 0$, return `True`

Recursive step: return `True` if the first character in s is the same as the last character *and* `_isPalindrome(s[1 : n - 1])` is `True`; otherwise, return `False`

Part I (Warmup Problems) · Problem 5 (Password Checker)

Implement the function `_isValid()` in `password_checker.py` that returns `True` if the given password string meets the following requirements, and `False` otherwise:

- Is at least eight characters long
- Contains at least one digit (0-9)
- Contains at least one uppercase letter
- Contains at least one lowercase letter
- Contains at least one character that is neither a letter nor a number

```
>_ ~/workspace/atomic_nature_of_matter
```

```
$ python3 password_checker.py Abcde1fg
False
$ python3 password_checker.py Abcde1@g
True
```

Part I (Warmup Problems) · Problem 5 (Password Checker)

Set *check*₁, *check*₂, *check*₃, *check*₄, and *check*₅ to `False`

If length of *pwd* is at least 8, set *check*₁ to `True`

For each character *c* in *pwd*

- If *c* is a digit, set *check*₂ to `True`
- Else if *c* is upper case, set *check*₃ to `True`
- Else if *c* is lower case, set *check*₄ to `True`
- Else if *c* is *not* alphanumeric, set *check*₅ to `True`

Return the logical *and* of *check*₁, *check*₂, *check*₃, *check*₄, and *check*₅

Part I (Warmup Problems) · Problem 6 (Point Data Type)

Define a data type called `Point` in `point.py` that represents a point in 2D. The data type must support the following API:

☰ `point.Point`

<code>Point(x, y)</code>	constructs a point <code>p</code> from the given <code>x</code> and <code>y</code> values
<code>p.distanceTo(q)</code>	returns the Euclidean distance between <code>p</code> and <code>q</code>
<code>str(p)</code>	returns a string representation of <code>p</code> as <code>'(x, y)'</code>

`>_ ~/workspace/atomic_nature_of_matter`

`$ python3 point.py`

`p1 = (0, 1)`

`p2 = (1, 0)`

`d(p1, p2) = 1.4142135623730951`

Part I (Warmup Problems) · Problem 6 (Point Data Type)

Instance variables

- *x*-coordinate, `_x` (float)
- *y*-coordinate, `_y` (float)

`__init__(self, x, y)`

- Initialize the instance variables to the values of the corresponding parameters

`distanceTo(self, other`

- Return the Euclidean distance between the points *self* and *other*

`__str__(self)`

- Return a string representation of the point *self*

Part I (Warmup Problems) · Problem 7 (Interval Data Type)

Define a data type called `Interval` in `interval.py` that represents a closed 1D interval. The data type must support the following API:

```
Interval.Interval
```

<code>Interval(lbound, rbound)</code>	constructs an interval <code>i</code> given its lower and upper bounds
---------------------------------------	--

<code>i.lower()</code>	returns the lower bound of <code>i</code>
------------------------	---

<code>i.upper()</code>	returns the upper bound of <code>i</code>
------------------------	---

<code>i.contains(x)</code>	returns <code>True</code> if <code>i</code> contains the value <code>x</code> , and <code>False</code> otherwise
----------------------------	--

<code>i.intersects(j)</code>	returns <code>True</code> if <code>i</code> intersects interval <code>j</code> , and <code>False</code> otherwise
------------------------------	---

<code>str(i)</code>	returns a string representation of <code>i</code> as <code>'[lbound, rbound]'</code>
---------------------	--

```
>_ ~/workspace/atomic_nature_of_matter
```

```
$ python3 interval.py
```

```
a = [-2, 1]
```

```
b = [0, 3]
```

```
a.contains(b.lower()) = True
```

```
a.contains(b.upper()) = False
```

```
a.intersects(b) = True
```

Part I (Warmup Problems) · Problem 7 (Interval Data Type)

Instance variables

- Lower bound of the interval, `_lbound` (float)
- Upper bound of the interval, `_rbound` (float)

```
__init__(self, lbound, ubound)
```

- Initialize the instance variables to the values of the corresponding parameters

```
lower(self)
```

- Return the value of the instance variable `_lbound`

```
upper(self)
```

- Return the value of the instance variable `_ubound`

```
contains(self, x)
```

- Return `True` if the interval *self* contains *x* and `False` otherwise

```
intersects(self, other)
```


- Return `True` if the interval *self* intersects the interval *other* and `False` otherwise

```
__str__(self)
```

- Return a string representation of the interval *self*

Part I (Warmup Problems) · Problem 8 (Rectangle Data Type)

Define a data type called `Rectangle` in `rectangle.py` that represents a rectangle using 1D intervals (ie, `Interval` objects) to represent its x (width) and y (height) segments. The data type must support the following API:

 `rectangle.Rectangle`

<code>Rectangle(xint, yint)</code>	constructs a rectangle <code>r</code> given its x and y segments, each an <code>Interval</code> object
------------------------------------	--

<code>r.area()</code>	returns the area of rectangle <code>r</code>
-----------------------	--

<code>r.perimeter()</code>	returns the perimeter of rectangle <code>r</code>
----------------------------	---

<code>r.contains(x, y)</code>	returns <code>True</code> if <code>r</code> contains the point (x, y) , and <code>False</code> otherwise
-------------------------------	--

<code>r.intersects(s)</code>	returns <code>True</code> if <code>r</code> intersects rectangle <code>s</code> , and <code>False</code> otherwise
------------------------------	--

<code>str(r)</code>	returns a string representation of <code>r</code> as <code>'[x1, x2] x [y1, y2]'</code>
---------------------	---

`>_ ~/workspace/atomic_nature_of_matter`

```
$ python3 rectangle.py
a      = [-1, 1] x [-1, 1]
b      = [0, 2] x [0, 3]
area(a)      = 4
perimeter(b) = 10
a.contains(1, 5) = False
a.intersects(b) = True
```


Part I (Warmup Problems) · Problem 8 (Rectangle Data Type)

Instance variables

- x -interval of the rectangle, `_xint` (Interval)
- y -interval of the rectangle, `_yint` (Interval)

`__init__(self, xint, yint)`

- Initialize the instance variables to the values of the corresponding parameters

`area(self)`

- Return the area of this rectangle

`perimeter(self)`

- Return the perimeter of this rectangle

`contains(self, x, y)`

- Return `True` if the rectangle *self* contains the point (x, y) and `False` otherwise

`intersects(self, other)`

- Return `True` if the rectangle *self* intersects the rectangle *other* and `False` otherwise

`__str__(self)`

- Return a string representation of the rectangle *self*

Part II (Atomic Nature of Matter) · Introduction

Goal: track the motion of particles undergoing Brownian motion, fit this data to Einstein's model, and estimate Avogadro's constant

Part II (Atomic Nature of Matter) · Problem 9 (Particle Representation)

Define a data type called `Blob` in `blob.py` to represent a particle (aka blob). The data type must support the following API:

 `Blob`

<code>Blob()</code>	constructs an empty blob <i>b</i>
<code>b.add(x, y)</code>	adds a pixel <i>(x, y)</i> to <i>b</i>
<code>b.mass()</code>	returns the mass of <i>b</i> , ie, the number of pixels in it
<code>b.distanceTo(c)</code>	returns the Euclidean distance between the center of mass of <i>b</i> and the center of mass of blob <i>c</i>
<code>str(b)</code>	returns a string representation of <i>b</i>

```
>_ ~/workspace/atomic_nature_of_matter
```

```
$ python3 blob.py
1 0 0 1 -1 0 0 -1
<ctrl-d>
a          = 1 (0.0000, 0.0000)
b          = 4 (0.0000, 0.0000)
dist(a, b) = 0.0
```

Part II (Atomic Nature of Matter) · Problem 9 (Particle Representation)

Instance variables:

- x -coordinate of center of mass, `_x` (float)
- y -coordinate of center of mass, `_y` (float)
- Number of pixels, `_pixels` (int)

`Blob()`

- Initialize the instance variables appropriately

`b.add(x, y)`

- Use the idea of *running average*¹ to update the center of mass of blob `b`
- Increment the number of pixels in blob `b` by 1

`b.mass()`

- Return the number of pixels in the blob `b`

`b.distanceTo(c)`

- Return the Euclidean distance between the center of mass of blob `b` and the center of mass of blob `c`

¹If \bar{x}_{n-1} is the average value of $n - 1$ points x_1, x_2, \dots, x_{n-1} , then the average value \bar{x}_n of n points $x_1, x_2, \dots, x_{n-1}, x_n$ is $\frac{\bar{x}_{n-1} \cdot (n-1) + x_n}{n}$.

Part II (Atomic Nature of Matter) · Problem 10 (Particle Identification)

Define a data type called `BlobFinder` in `blob_finder.py` that supports the following API. Use depth-first search to efficiently identify the blobs.

BlobFinder

<code>BlobFinder(pic, tau)</code>	constructs a blob finder <code>bf</code> to find blobs in the picture <code>pic</code> using a luminance threshold <code>tau</code>
<code>bf.getBeads(pixels)</code>	returns a list of all blobs with mass \geq <code>pixels</code> , ie, a list of beads

Part II (Atomic Nature of Matter) · Problem 10 (Particle Identification)

```
>_ ~/workspace/atomic_nature_of_matter
```

```
$ python3 blob_finder.py 25 180.0 data/run_1/frame00001.jpg
```

```
13 Beads:
```

```
29 (214.7241, 82.8276)
36 (223.6111, 116.6667)
42 (260.2381, 234.8571)
35 (266.0286, 315.7143)
31 (286.5806, 355.4516)
37 (299.0541, 399.1351)
35 (310.5143, 214.6000)
31 (370.9355, 365.4194)
28 (393.5000, 144.2143)
27 (431.2593, 380.4074)
36 (477.8611, 49.3889)
38 (521.7105, 445.8421)
35 (588.5714, 402.1143)
```

```
15 Blobs:
```

```
29 (214.7241, 82.8276)
36 (223.6111, 116.6667)
1 (254.0000, 223.0000)
42 (260.2381, 234.8571)
35 (266.0286, 315.7143)
31 (286.5806, 355.4516)
37 (299.0541, 399.1351)
35 (310.5143, 214.6000)
31 (370.9355, 365.4194)
28 (393.5000, 144.2143)
27 (431.2593, 380.4074)
36 (477.8611, 49.3889)
38 (521.7105, 445.8421)
35 (588.5714, 402.1143)
13 (638.1538, 155.0000)
```

Part II (Atomic Nature of Matter) · Problem 10 (Particle Identification)

Instance variable:

- Blobs identified by this blob finder, `_blobs` (list of `Blob` objects).

`BlobFinder()`

- Initialize `blobs` to an empty list.
- Create a 2D list of booleans called `marked`, having the same dimensions as `pic`.
- Enumerate the pixels of `pic`, and for each pixel `(i, j)`:
 - create a `Blob` object called `blob`;
 - call `_findBlob()` with the appropriate arguments; and
 - add `blob` to `blobs` if it has a non-zero mass.

`bf._findBlob()`

- Base case: return if pixel `(i, j)` is out of bounds, or if it is marked, or if its luminance (use the `luminance()` method from `Color` for this) is less than `tau`.
- Mark the pixel `(i, j)`.
- Add the pixel `(i, j)` to the blob `blob`.
- Recursively call `_findBlob()` on the N, E, W, and S pixels.

`bf.getBeads(pixels)`

- Return a list of blobs from `blobs` that have a mass \geq `pixels`.

Part II (Atomic Nature of Matter) · Problem 11 (Particle Tracking)

Implement a program called `bead_tracker.py` that accepts p (int), τ (float), δ (float), and a sequence of JPEG filenames as command-line arguments; identifies the beads in each JPEG image using `BlobFinder`; and writes to standard output (one per line, formatted with 4 decimal places to the right of decimal point) the radial distance that each bead moves from one frame to the next (assuming it is no more than δ)

```
>_ ~/workspace/atomic-nature-of-matter
```

```
$ python3 bead_tracker.py 25 180.0 25.0 data/run_1/frame00000.jpg data/run_1/frame00001.jpg
7.1833
4.7932
2.1693
5.5287
5.4292
4.3962
```


Part II (Atomic Nature of Matter) · Problem 11 (Particle Tracking)

Accept command-line arguments `pixels` (int), `tau` (float), and `delta` (float)

Construct a `BlobFinder` object for the frame `sys.argv[4]` and from it get a list of beads `prevBeads` that have at least `pixels` pixels

For each frame starting at `sys.argv[5]`:

- Construct a `BlobFinder` object and from it get a list of beads `currBeads` that have at least `pixels` pixels
- For each bead `currBead` in `currBeads`, find a bead `prevBead` from `prevBeads` that is no further than `delta` and is closest to `currBead`, and if such a bead is found, write its distance (using format string `'%.4f\n'`) to `currBead`
- Write a newline character
- Set `prevBeads` to `currBeads`

Part II (Atomic Nature of Matter) · Problem 12 (Estimating Avogadro's Constant)

Implement a program called `avogadro.py` that accepts the displacements (output of `bead_tracker.py`) from standard input; computes an estimate of Avogadro's constant using the formulae described above; and writes the value to standard output

```
>_ ~/workspace/atomic_nature_of_matter
```

```
$ python3 bead_tracker.py 25 180.0 25.0 data/run_1/* | python3 avogadro.py  
6.633037e+23
```

Part II (Atomic Nature of Matter) · Problem 12 (Estimating Avogadro's Constant)

Initialize `ETA`, `RHO`, `T`, and `R` to appropriate values

Calculate `var` as the sum of the squares of the `n` displacements (each converted from pixels to meters) read from standard input

Divide `var` by `2 * n`

Estimate Boltzmann's constant as `6 * math.pi * var * ETA * RHO / T`

Estimate Avogadro's constant as `R / k`

Write to standard output the Avogadro constant in scientific notation (using the format string `"%e"`)