

Answer: D

The NumPy code creates a vector of length size of random integers that are either -1 or 1. So we want to create a list of length size of similar random integers using the random.choice function.

```
delta = np.random.choice([-1,1], steps)
walk = np.cumsum(delta)
```

Which of the following snippets are equivalent to the above NumPy code (steps is  $\geq 0$ ). NumPy vectors should be Python lists.

A. pos = 0 walk = [] for i in range(steps): cumulate pos += random.choice([-1, 1]) Walk.append(pos) Intermediate	<pre>B. walk = [] for i in range(steps):     delta=random.choice([-1,1])     walk.append(delta)</pre>
<pre>C. walk = [] for i in range(steps):     if random.randint(0,1) == 0:         walk.append(-1)         else:         walk.append(1)</pre>	<pre>D. pos = 0 for i in range(steps):     pos += random.choice([-1, 1])</pre>

Answer: A

delta is an array of randomly sample +1, -1, the cumsum function computes the cumulative sum, i.e., for [a, b, c] it computers [a, a+b, a+b+c]. Answer A implements the latter via the pos accumulator.

## scalar index = np.argmax(walk)

Which of the following snippets are equivalent to the above NumPy code?

<pre>A. index = [] for i in range(1,len(a_list)):     if a_list[i] &gt; a_list[i-1]:         index.append(i)</pre>	<pre>. index = [] for i in range(1,len(a_list)):     if a_list[i] &gt; a_list[index[-1]]:         index.append(i)</pre>
<pre>C. index = 0 for i in range(1,len(a_list)):     if a_list[i] &gt; a_list[i-1]:         index = i</pre>	<pre>D. index = 0 for i in range(1,len(a_list)):     if a_list[i] &gt; a_list[index]:         index = i</pre>

## Answer: D

We want to find the index of the largest value in the entire list. That will be a scalar, so A and B are incorrect. Answer C, would return the last index where the item was larger than its immediate predecessor (not the largest overall).

## returns = np.cumprod(np.random.laplace(mean, scale, 240))

Which of the following snippets are equivalent to the above NumPy code? Assume there is a laplace function that has the mean and scale as arguments and returns a single sample.

```
A. returns = []
                                           B. returns = []
  for i in range(240):
                                               prod = 1.0
     sample = laplace(mean, scale)
                                               for i in range(240):
                                                 sample = laplace(mean, scale)
     returns.append(sample)
                                                 prod *= sample
                                                 returns.append(prod)
C. returns = []
                                           D. returns = []
   prod = 1.0
                                               prod = 1.0
   for i in range(240):
                                               for i in range(240):
     sample = laplace(mean, scale)
                                                 sample = laplace(mean, scale)
     returns.append(sample)
                                                 returns.append(prod)
     prod *= sample
                                                 prod *= sample
```

## Answer: B

The `cumprod` function computes a \_cumulative\_ product. Answers A and C only record the samples. Answer D has a "off by one", that is starts with the initial values and doesn't record the final product.