# Case Study 1: Interfacing C++ with Python for Blahut-Arimoto Simulations

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This case study will introduce an important algorithm for data transmission and channel capacity, then demonstrating how we have used C++ compatability to optimise our computations over large dimensional (and sample size) data. Unfortunately we began work in python so this will be the main language we use, including its compatability with C++. However we hope that this will still be educational and should compliment the topic on implimenting R and C++ with the Rccp package. We will be using multiple python and c++ files during this study, the full code can be found here.

# The Blahut Arimoto Algorithm

The Blahut-Arimoto algorithm is an iterative algorithm that computes the capacity of a channel, a channel can be thought of as a box which some signal can be parsed through and causes some sort of distortion. The capacity of a channel is the maximum rate at which information can be sent through the channel with an arbitrarily small error rate, i.e it gives a lower bound on how much redundancy we must add to guarantee (asymptotically) error free communication.

If we imagine the input and output signals X and Y as catagorical random variables on the sets  $\{1, ..., m\}$  and  $\{1, ..., n\}$ , we can define the channel as a conditional probability distribution  $P_{Y|X}$ , or more concretely a matrix  $P_{ij} = P(Y = j|X = i)$ . Then we can write Y = XP. The channel coding theorem states that the capacity of a channel is given by the formula

$$C = \max_{X} I(X;Y) = \max_{X} I(X,XP)$$

where I(X;Y) is the mutual information between the input and output signals. The Blahut-Arimoto algorithm is an iterative algorithm that computes the capacity of a channel by finding the optimal input distribution  $P_X$  that maximises the mutual information. We won't dive into the specifics here but the general idea is that this single max problem can be split into a maxmax problem, where the smaller maximisations correspond to analytic substitutions of parameters into equations. This boils down to the algorithm being a for loop. Below is my original python code which performs Blahut-Arimoto.

```
We attempt to maximise I with the blahut amirito algorithm
prior: Inital quess for maximal X
channel\_matrix: is the matrix representing the transition from X to Y, i.e P
log_base: base of logarithm, typically 2 or nats
thresh: threshold to finish algorithm, when iterations are < thresh apart
max_iter: maximum number of iterations before giving up
      _____
BA is a maxmax algorithm, each maximisation has a closed form expression
We perform the maximisation as a for loop
if display_results:
    print("Arimoto: PYTHON EDITION")
#Check we have a valid dimension for the prior
assert prior.shape[0] > 1
#Checkl prior is a valid probability distribution
assert np.abs(prior.sum() - 1) < 1e-6</pre>
#Check we have a valid dimension for the channel matrix
n,m = channel_matrix.shape
m_1 = prior.shape[0]
assert m == m_1
#Initialise the prior and Phi
p = prior
p_route = np.array([p])
P = channel_matrix.transpose()
W = np.zeros((m,n))
if display_results:
    print(f"Prior for p(x): {p}")
    print(f"Channel matrix p(y|x): \n {channel_matrix}")
for iter in range(int(maxiter)):
    q = np.zeros(m)
    #Maximise I(p, W; P) over W
    for i in range(n):
       for j in range(m):
            W[j][i] = (P[i,j]*p[j])/np.dot(P[i,:],p)
    #Maximise I(p,W;P) over p
   r = np.zeros(m)
    for j in range(m):
       r[j] = np.exp(np.dot(P[:,j],np.log(W[j,:])))
    for i in range(m):
        q[i] = r[i]/np.sum(r)
    #Add to array of p values
    p_route = np.append(p_route,[q],axis=0)
    #Check if we have converged
    if np.linalg.norm(q-p) < thresh:</pre>
    p = q
#Calculate the capacity
```

As a demonstration of the algorithm, we can consider a simple binary channel with a transition matrix P given by:

$$P = \begin{bmatrix} 0.9 & 0.1 \\ 0.1 & 0.9 \end{bmatrix}$$

and a prior distribution p(x) given by: p(x=0) = 0.8 and p(x=1) = 0.2. In fact the maximisation for this type of channel (a binary symmetric channel) has an analytic solution for the capacity, which is given by:

$$C = 1 - H(0.9)$$

where H(p) is binary entropy function, and the maximising distribution is given by p(y=0) = p(y=1) = 0.5. We can then run the algorithm with the following code:

## --||Binary Symmetric Channel Simulation||--

```
## ------
## Python Capacity: 0.5310044064107188
## Python ArgMax: [0.5 0.5] :: <class 'numpy.ndarray'>
## True Capacity: 0.5310044064107188
## ------
end = time.time()
print(f"Time taken: {end-start} seconds")
```

## Time taken: 0.005226850509643555 seconds

Our search space, for an n-ary input space, is an (n-1)-simplex, and of course as the dimension of the input space increases, the number of iterations required to converge increases too. Below we run the algorithm for a 3-ary input space, with 3 different priors and a random channel matrix, and plot the path taken by the algorithm to converge.

```
import Arimito_Generator as Generator
import Arimito_Convergence as Convergence
```

## INFO: Using numpy backend

```
random.seed(124)

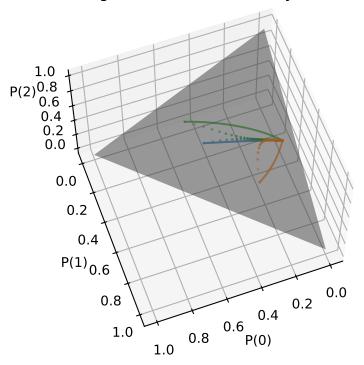
channel_Matrix = Generator.random_channel_matrix(xdim = 3,ydim = 3)

start = time.time()
Convergence.plot_path_ternary(n=3,include_geodesic = True)
```

```
## Arimoto: PYTHON EDITION
## Prior for p(x): [0.44003885 0.26384277 0.29611838]
## Channel matrix p(y|x):
## [[0.81791449 0.55734968 0.38944123]
## [0.04970083 0.38175277 0.02095369]
## [0.13238468 0.06089754 0.58960507]]
## Max Capacity: 0.3333866550177944
## ArgMax: [0.06027683 0.46752258 0.47220058]
## -----
## -----
## Arimoto: PYTHON EDITION
## Prior for p(x): [0.22511391 0.56522352 0.20966257]
## Channel matrix p(y|x):
## [[0.81791449 0.55734968 0.38944123]
## [0.04970083 0.38175277 0.02095369]
## [0.13238468 0.06089754 0.58960507]]
## Max Capacity: 0.33338665501779446
## ArgMax: [0.06027683 0.46752258 0.47220058]
## -----
## Arimoto: PYTHON EDITION
## Prior for p(x): [0.50301787 0.12901178 0.36797035]
## Channel matrix p(y|x):
## [[0.81791449 0.55734968 0.38944123]
## [0.04970083 0.38175277 0.02095369]
```

```
## [0.13238468 0.06089754 0.58960507]]
## Max Capacity: 0.3333866550177947
## ArgMax: [0.06027683 0.46752258 0.47220058]
## ------
```

# Convergence Path for a Ternary Channel



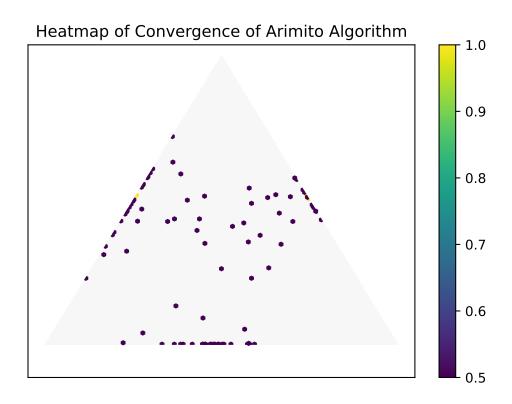
```
end = time.time()
print(f"Time taken: {end-start} seconds")
```

### ## Time taken: 0.5891118049621582 seconds

We also plotted the geodesic path (with respect to the fisher information matrix) along the simplex to study how well the algorithm converges. One thing you may notice in the above example is that the optimal solution lies on the boundary of the simplex, i.e one of the probabilities is 0. This might seem kind of weird, since we previously discussed that we have an infinite family of solutions at p = (1/3, 1/3, 1/3). The question became, what does a typical solution look like? We answer this question by simulating a collection of random channel matrices and priors, and then plotting the heatmap of solutions. Note that since the mutual information is convex, the choice of prior should not matter. This is where computational complexity begins to creep in, as the simulation took my computer  $\sim 2$  minutes, with 1000 iterations and xdim = 3. We have only plotted n=100, xdim=3 for this case to make knitting to a pdf practical, plus you get to see a better picture at the end! For the below code we use the package ternary which helps with heatmaps on simplexes.

```
import ternary
start = time.time()
```

```
#Create function which performs arimito multiple times and outputs the distribution
def arimito_multiple(prior,
                    log_base: float = 2,
                    thresh: float = 1e-12,
                    maxiter: int = 1e3,
                    n: int = 1):
    111
    Perform the arimito algorithm n times and output the distribution
    p = []
    for i in range(n):
        random_binary_channel = Generator.random_channel_matrix(xdim = 3,ydim = 3)
        p.append(arimito(prior, random_binary_channel,log_base=log_base,thresh=thresh, maxiter=maxiter, di
    return p
#Perform the arimito algorithm multiple times
p = arimito_multiple(prior = np.array([0.2,0.3,0.5]),n=100)
## <string>:50: RuntimeWarning: divide by zero encountered in log
p = np.array(p)*100
heatmap_data = {}
for i in range(p.shape[0]):
    # Round the points to 2 decimal places and convert to a tuple
    point = tuple(np.round(p[i, :], 1))
    # Add the point to the heatmap data
    if point in heatmap_data:
        heatmap_data[point] += 1
    else:
        heatmap_data[point] = 1
# Create the figure
fig, tax = ternary.figure(scale=100)
# Normalize the values in the heatmap data
max_value = max(heatmap_data.values())
heatmap_data = {key: value / max_value for key, value in heatmap_data.items()}
# Plot the heatmap
tax.heatmap(heatmap_data, style="hexagonal", use_rgba=False)
# Set labels
tax.set_title("Heatmap of Convergence of Arimito Algorithm")
tax.left_corner_label("$p_0$")
tax.right_corner_label("$p_1$")
tax.top_corner_label("$p_2$")
# Remove default Matplotlib Axes
tax.clear_matplotlib_ticks()
ternary.plt.show()
```



```
end = time.time()
print(f"Time taken: {end-start} seconds")
```

## Time taken: 0.530747652053833 seconds

As you can see, counterintuitively, the optimal solution is often on the boundary of the simplex. Which is a great observation, however as it took so long it's hard to get a good measure of the behaviour.

Since the arimito algorithm is fairly rudamental (its just a for loop where we update probabilities) we rewrote the arimito function in C instead. We also used python to interface with C to preserve our codebase.

# Python Interface to C++

Python has a slightly different method of interfacing with C++, which comes from a few different packages, we chose to use pybind11 as that's what google said was good, plus it works so that's good. All it requires is installation and inclusion of a few lines of code in your C++ code. Firstly we pip install.

```
pip install pybind11
```

The technique for mixing C++ with python is very simple, we essentially make a local package which we can import like any other python package - this is probably a side affect of python being built on C. We only need to include the following code in our preamble:

```
#include <pybind11/pybind11.h>
#include <pybind11/numpy.h>
#include <pybind11/stl.h>

namespace py = pybind11;
```

to tell the compiler that we are using pybind11. This plays a similar role to #include <Rcpp.h> in in R. Then at the bottom we included the command:

```
PYBIND11_MODULE(CArimito, m) {
    m.def("arimito", &arimito, "Arimoto algorithm");
}
```

This tells pybind to create a function called arimito in the module CArimito. With this setup, pybind does all the work for us, so long as we compile. Which we do by going to the working directory of our .cpp file and running:

```
c++ -03 -Wall -shared -std=c++11 -fPIC `python3
-m pybind11 --includes` Arimito_Quick.cpp
-o CArimito`python3-config --extension-suffix`
```

all on one line in the terminal. This creates CArimito.cpython-310-x86\_64-linux-gnu.so which can be imported into python via:

```
import CArimito
```

We can then call our function arimito as if it were a python function, so specifically we can run the following code:

```
CArimito.arimito(channel_matrix_list,prior_list,1000,1e-12)
```

Although to make the function more user friendly we created the following wrapper to match it up with or python edition:

```
max_iter: maximum number of iterations before giving up
_____
BA is a maxmax algorithm, each maximisation has a closed form expression
We perform the maximisation as a for loop
if display_results:
   print("Arimoto: C EDITION")
#Check we have a valid dimension for the prior
assert prior.shape[0] > 1
#Checkl prior is a valid probability distribution
assert np.abs(prior.sum() - 1) < 1e-6</pre>
#Check we have a valid dimension for the channel matrix
n,m = channel_matrix.shape
m_1 = prior.shape[0]
assert m == m_1
if display_results:
   print(f"Prior for p(x): {prior}")
   print(f"Channel matrix p(y|x): \n {channel_matrix}")
prior_list = prior.tolist()
channel_matrix_list = channel_matrix.tolist()
p,C,p_route = CArimito.arimito(channel_matrix_list,prior_list,1000,1e-12)
if display_results:
   print('Max Capacity: ', C)
   print('ArgMax: ', p)
   print("----")
   print("----")
return C,p,p_route
```

# Results

As a demonstration, lets begin by recreating the output from our intial symmetric examples and three random prior examples.

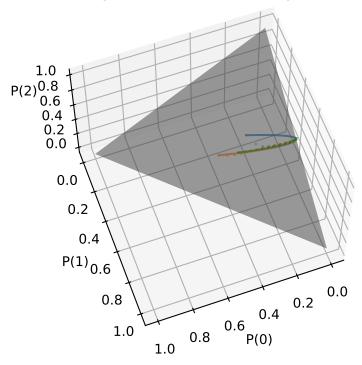
#### ## Time taken: 0.005412101745605469 seconds

## -----

Hilariously, in this simple case the difference is by a measily 0.001, lets try our 3 random priors examples and see the result:

```
random.seed(123)
start = time.time()
Convergence.plot path ternary(n=3,include geodesic = True,lang = "C++")
## Arimoto: C EDITION
## Prior for p(x): [0.23176497 0.34967348 0.41856154]
## Channel matrix p(y|x):
## [[0.81791449 0.04970083 0.13238468]
## [0.55734968 0.38175277 0.06089754]
## [0.38944123 0.02095369 0.58960507]]
## Max Capacity: 0.3465845061760716
## ArgMax: [0.
                      0.49740853 0.50259147]
## -----
## -----
## Arimoto: C EDITION
## Prior for p(x): [0.39127933 0.35114807 0.2575726 ]
## Channel matrix p(y|x):
## [[0.81791449 0.04970083 0.13238468]
## [0.55734968 0.38175277 0.06089754]
## [0.38944123 0.02095369 0.58960507]]
## Max Capacity: 0.3465845061760716
## ArgMax: [0. 0.49740853 0.50259147]
## -----
## -----
## Arimoto: C EDITION
## Prior for p(x): [0.2927475 0.3939528 0.3132997]
## Channel matrix p(y|x):
## [[0.81791449 0.04970083 0.13238468]
## [0.55734968 0.38175277 0.06089754]
## [0.38944123 0.02095369 0.58960507]]
## Max Capacity: 0.34658450617607167
## ArgMax: [0.
                      0.49740853 0.50259147]
```

# Convergence Path for a Ternary Channel

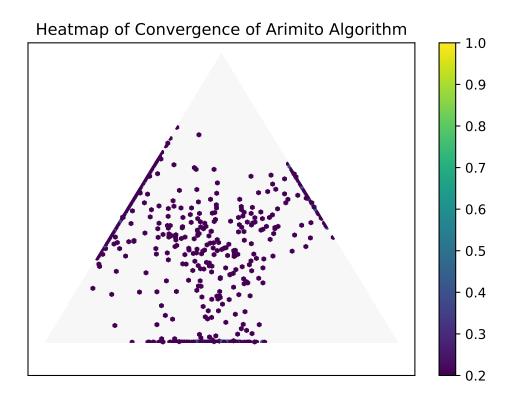


```
end = time.time()
print(f"Time taken: {end-start} seconds")
```

### ## Time taken: 0.08254814147949219 seconds

We're working at around about 4x as fast. Which should be useful with large simulations. Let's put this into practice and get a good image of how our data looks. We reconstruct our arimito\_multiple function to use the C++ version of the algorithm and plot the heatmap of the results:

```
ydim = 3)
        p.append(Arimito.Carimito(prior,
                                random_binary_channel,
                                log_base=log_base,
                                thresh=thresh,
                                maxiter=maxiter,
                                display_results = False)[1])
    return p
#Perform the arimito algorithm multiple times
p = arimito_multiple(prior = np.array([0.2,0.3,0.5]),n=1000)
p = np.array(p)*100
heatmap_data = {}
for i in range(p.shape[0]):
    # Round the points to 2 decimal places and convert to a tuple
    point = tuple(np.round(p[i, :], 1))
    # Add the point to the heatmap data
    if point in heatmap_data:
        heatmap_data[point] += 1
    else:
        heatmap_data[point] = 1
# Create the figure
fig, tax = ternary.figure(scale=100)
# Normalize the values in the heatmap data
max_value = max(heatmap_data.values())
heatmap_data = {key: value / max_value for key, value in heatmap_data.items()}
# Plot the heatmap
tax.heatmap(heatmap_data, style="hexagonal", use_rgba=False)
# Set labels
tax.set_title("Heatmap of Convergence of Arimito Algorithm")
tax.left_corner_label("$p_0$")
tax.right_corner_label("$p_1$")
tax.top_corner_label("$p_2$")
# Remove default Matplotlib Axes
tax.clear_matplotlib_ticks()
ternary.plt.show()
```



```
end = time.time()
print(f"Time taken: {end-start} seconds")
```

## ## Time taken: 0.645514965057373 seconds

For context, I didn't actually realise I had included n=1000 in the above code, because it compiled as quick as python did for n=100. Now we can increase n drastically to get a solid density plot.

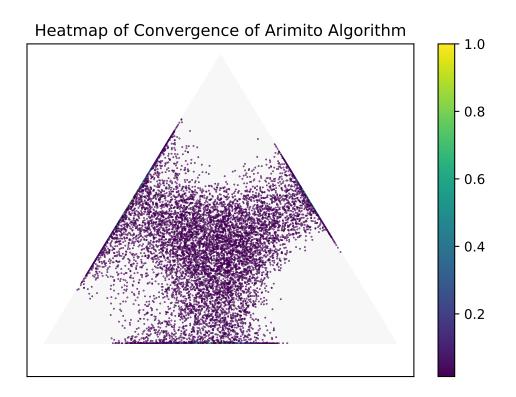
```
#Perform the arimito algorithm multiple times
p = arimito_multiple(prior = np.array([0.2,0.3,0.5]),n=30000)

p = np.array(p)*1000

heatmap_data = {}

for i in range(p.shape[0]):
    # Round the points to 2 decimal places and convert to a tuple
    point = tuple(np.round(p[i, :], 0))
    # Add the point to the heatmap data
    if point in heatmap_data:
        heatmap_data[point] += 1
    else:
        heatmap_data[point] = 1
```

```
# Create the figure
fig, tax = ternary.figure(scale=1000)
# Calculate the maximum value in the heatmap data
max_value = max(heatmap_data.values())
# Normalize the heatmap data
heatmap_data = {key: value / max_value for key, value in heatmap_data.items()}
# Plot the heatmap
tax.heatmap(heatmap_data, style="hexagonal", use_rgba=False)
# Set labels
tax.set_title("Heatmap of Convergence of Arimito Algorithm")
tax.left_corner_label("$p_0$")
tax.right_corner_label("$p_1$")
tax.top_corner_label("$p_2$")
# Remove default Matplotlib Axes
tax.clear_matplotlib_ticks()
ternary.plt.show()
```



```
end = time.time()
print(f"Time taken: {end-start} seconds")
```

#### ## Time taken: 9.373404502868652 seconds

We can see much clearer distribution of data. Additionally we can see that the accumulation of data is often in the middle of the boundaries. Which is really cool! For my own research I will be continuing to study this property, as it has been studied in other works that zero values (boundary solutions) for capacity are known for continuous channel distributions, but not much is known for discrete. Hopefully this has been a quaint adventure into information theory that compliments our work on interfacing R with C++ with Rcpp, demonstrating how the technique of segmenting code for optimisation with C carries between many languages.