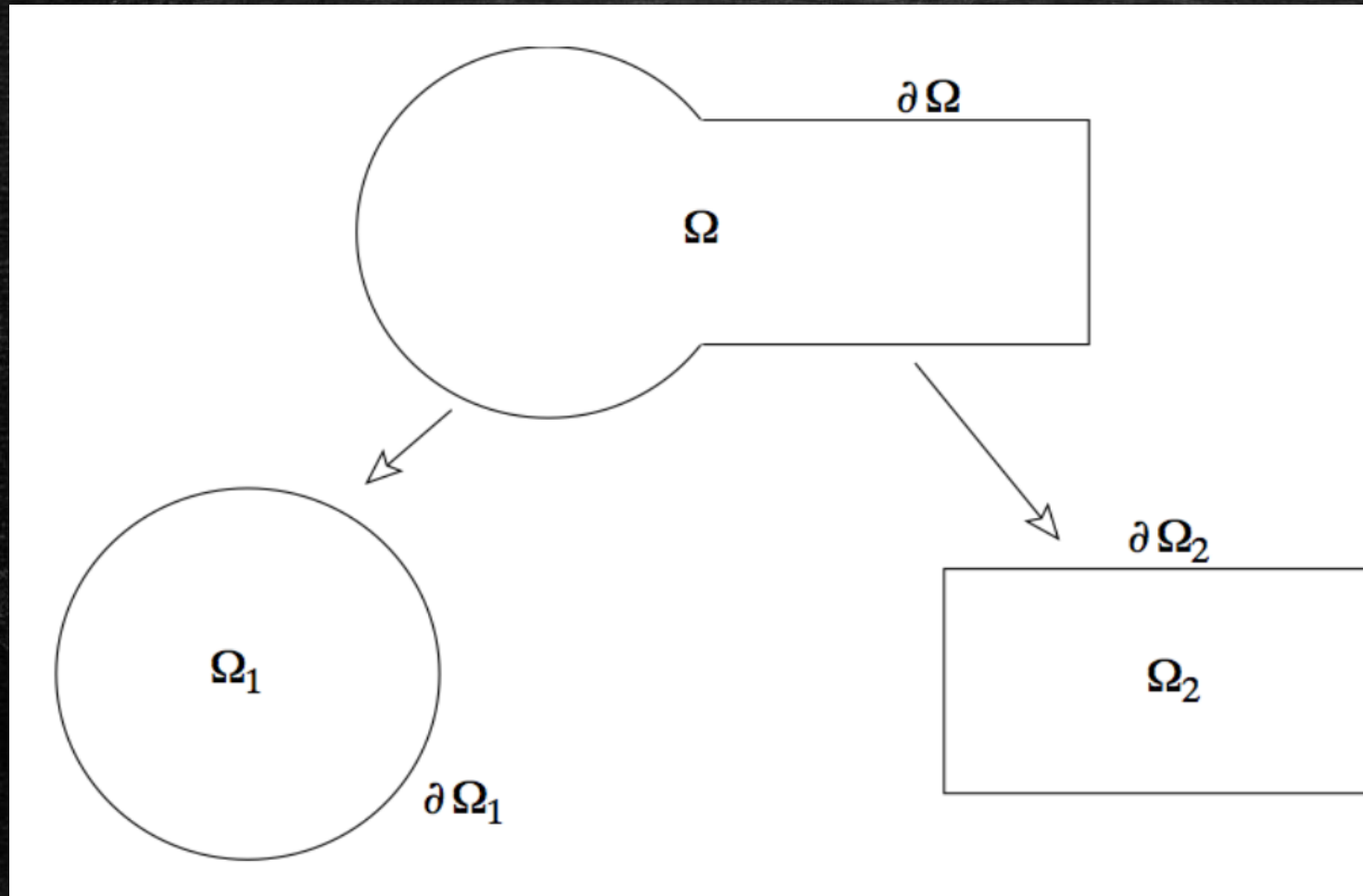


# Conducción de calor en 1D empleando el Algoritmo alternante de Schwarz

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# ¿En qué consiste el algoritmo?



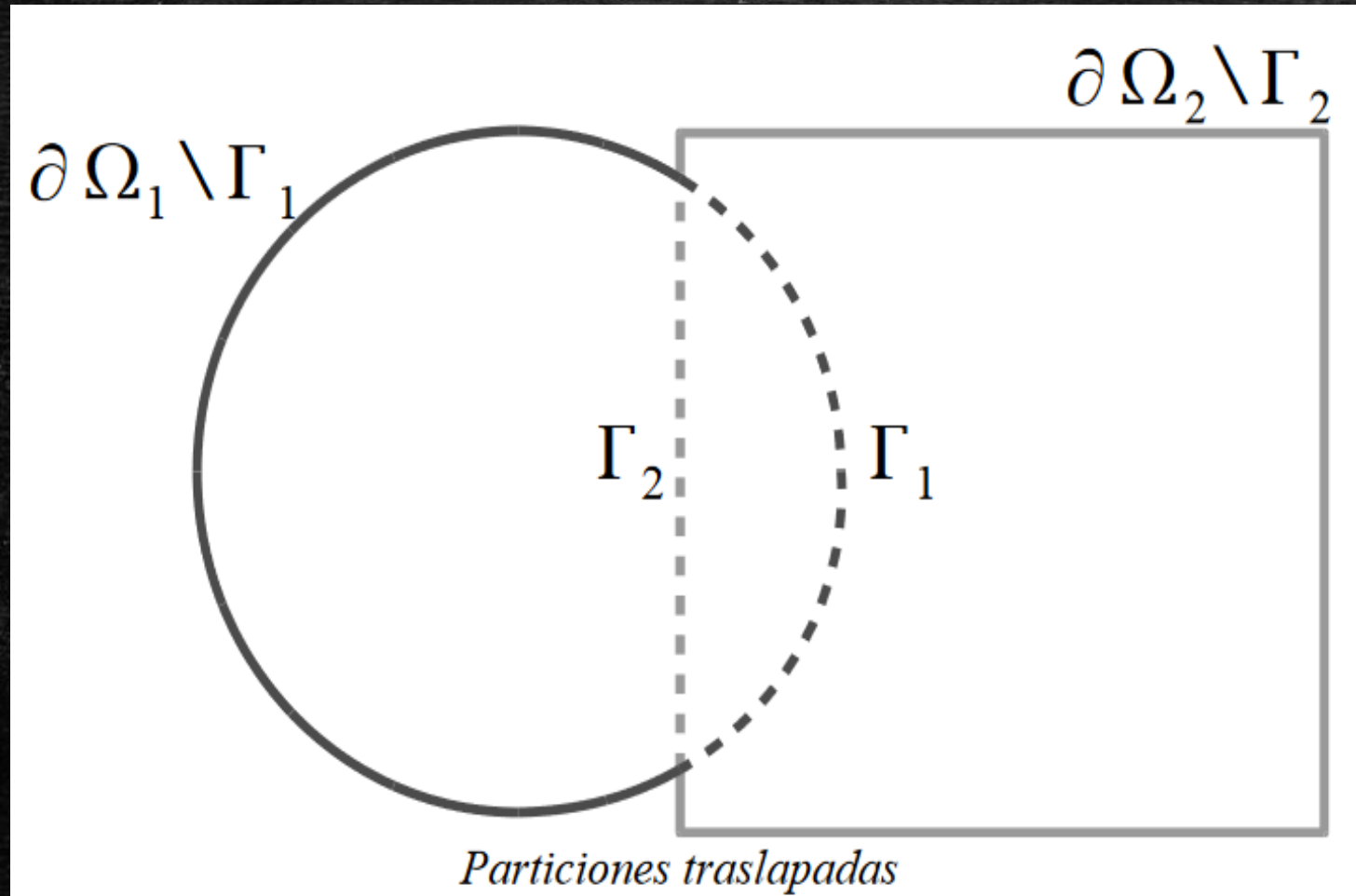
Dominio original  
 $\Omega$

Particiones  
 $\Omega_1$  y  $\Omega_2$

Fronteras  
 $\partial\Omega_1$  y  $\partial\Omega_2$



¿En qué consiste el algoritmo?



Particiones se traslapan

$$\Omega = \Omega_1 \cup \Omega_2$$

Fronteras artificiales

$\Gamma_1$  y  $\Gamma_2$

Fronteras reales

$$\partial \Omega_1 \setminus \Gamma_1 \text{ y } \partial \Omega_2 \setminus \Gamma_2$$

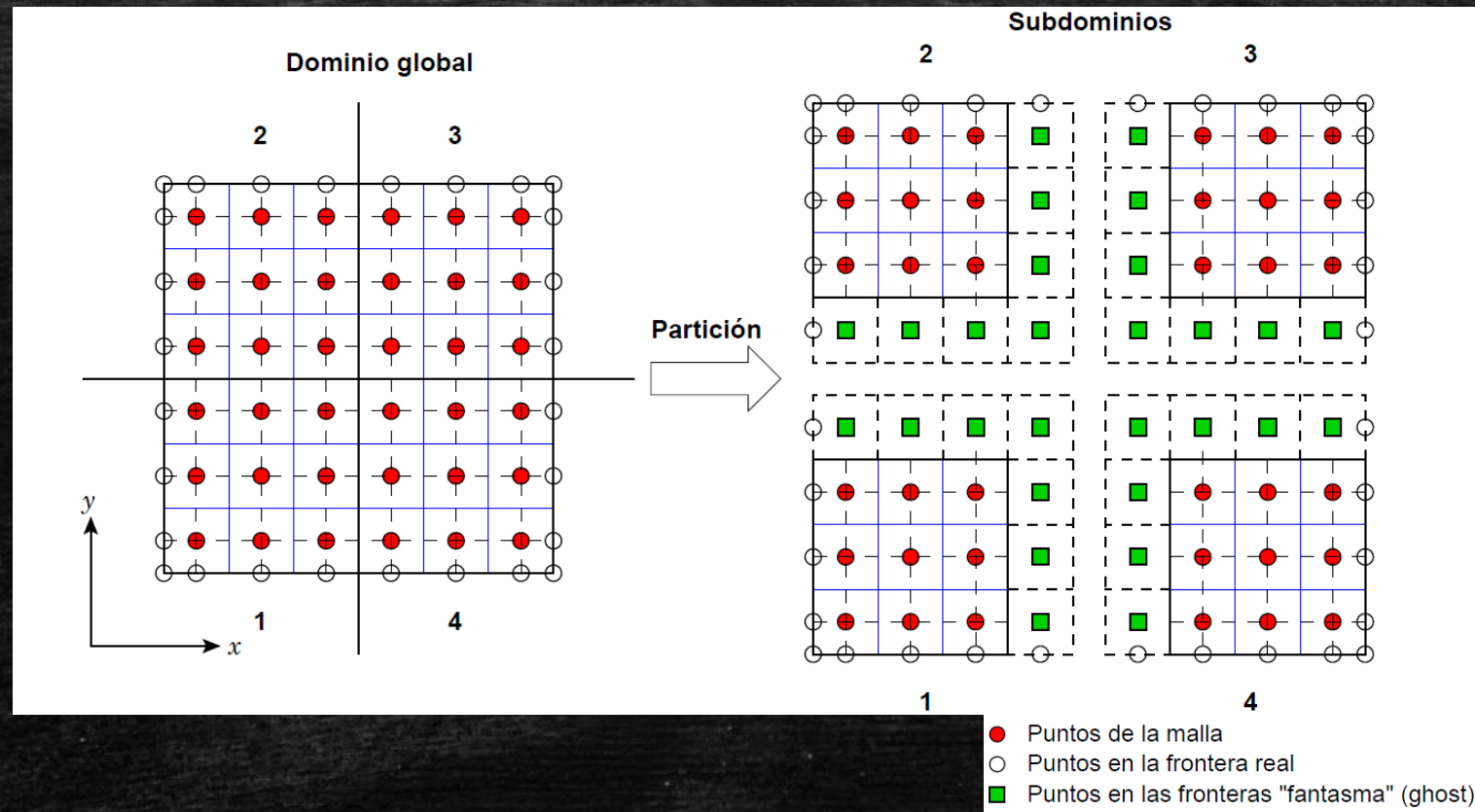
# ¿En qué consiste el algoritmo?

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- Se suponen conocidas las fronteras internas  $\Gamma_1$  y  $\Gamma_2$
- Se resuelve de manera simultanea en cada subdominio
  - Con la condición  $g$  en la frontera real  $\delta\Omega_1 \setminus \Gamma_1$  y  $\delta\Omega_2 \setminus \Gamma_2$
- Se modifican los valores de las fronteras  $\Gamma_1$  y  $\Gamma_2$  con iteración anterior
- Se resuelve de manera simultanea en cada subdominio
  - Con la condición  $g$  en la frontera real  $\delta\Omega_1 \setminus \Gamma_1$  y  $\delta\Omega_2 \setminus \Gamma_2$
  - Valores de la iteración anterior en las fronteras  $\Gamma_1$  y  $\Gamma_2$



# Partición del dominio (MPI)



`MPI.COMM_WORLD.Create_cart(dim, periods, reorder)`

# Comunicación de los subdominios (1D)

- Get\_coords()
- Shift()

```
grid_rows = int(1)
grid_cols = comm.size
```

```
Building a 1 x 1 grid topology
Process = 0          row = 0          column = 0
    neighbor_processes[LEFT] = -1
    neighbor_processes[RIGHT] = -1
    neighbor = [-1, -1]
```

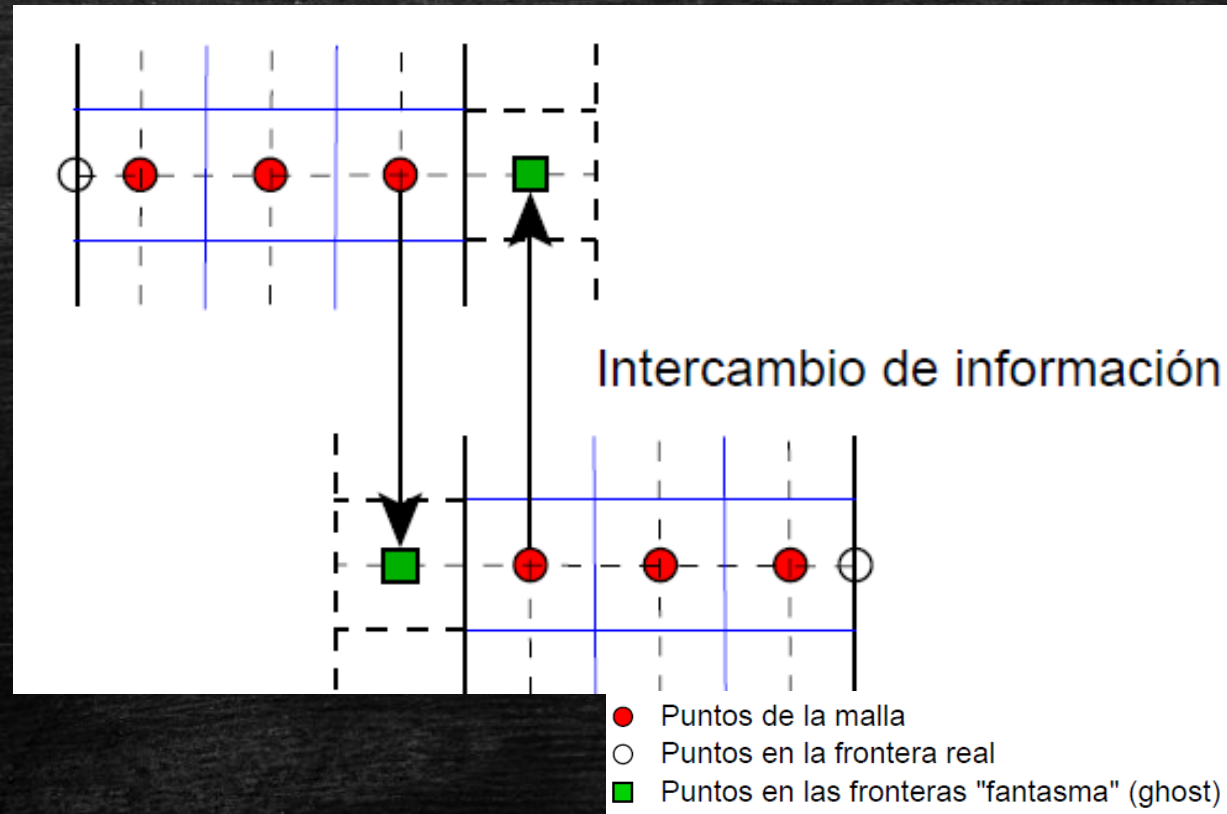
```
Building a 1 x 2 grid topology
Process = 0          row = 0          column = 0
    neighbor_processes[LEFT] = -1
    neighbor_processes[RIGHT] = 1
    neighbor = [-1, 1]
Process = 1          row = 0          column = 1
    neighbor_processes[LEFT] = 0
    neighbor_processes[RIGHT] = -1
    neighbor = [0, -1]
```

```
Building a 1 x 3 grid topology
Process = 0          row = 0          column = 0
    neighbor_processes[LEFT] = -1
    neighbor_processes[RIGHT] = 1
    neighbor = [-1, 1]
Process = 1          row = 0          column = 1
    neighbor_processes[LEFT] = 0
    neighbor_processes[RIGHT] = 2
    neighbor = [0, 2]
Process = 2          row = 0          column = 2
    neighbor_processes[LEFT] = 1
    neighbor_processes[RIGHT] = -1
    neighbor = [1, -1]
```

`mpiexec -n numproc python programa.py`

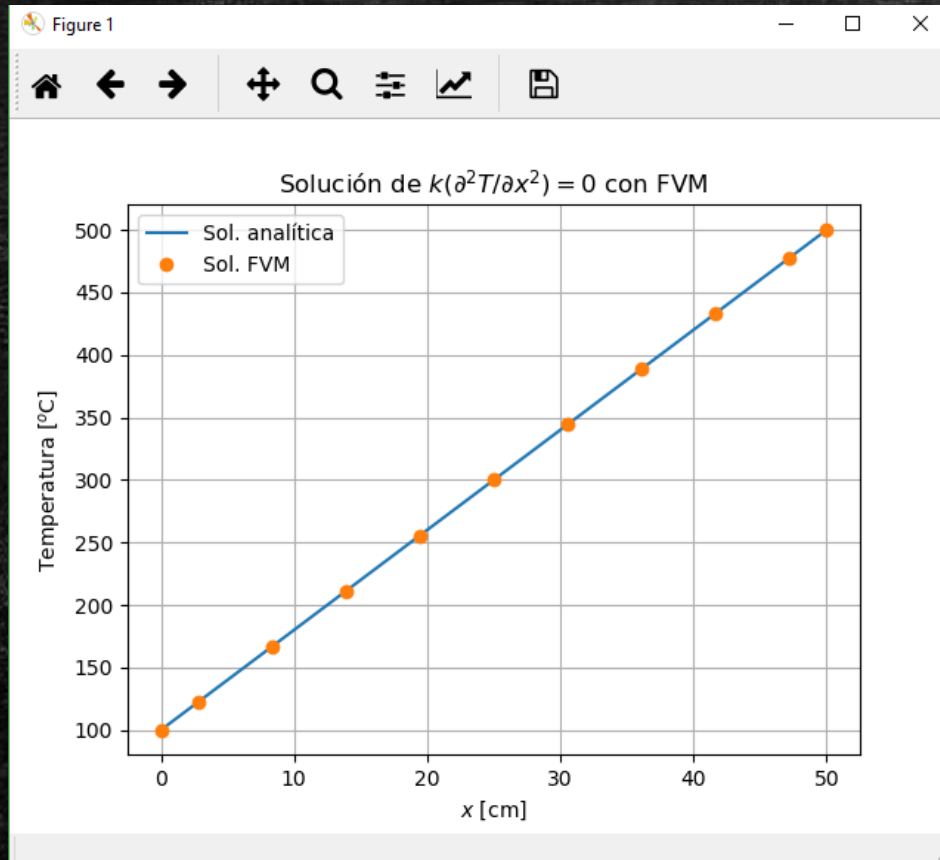


# Comunicación de los subdominios (1D)



```
comm.sendrecv(data, dataType, dest, tag, origen, tag )
```

# Resultados grid 1x1



Building a 1 x 1 grid topology

Solucionaremos la  $T = [0. \ 0. \ 0. \ 0. \ 0. \ 0. \ 0. \ 0. \ 0. \ 0.]$  en el rank = 0

ahora t vale =  $[100. \ 0. \ 0. \ 0. \ 0. \ 0. \ 0. \ 0. \ 0. \ 500.]$

Solución para  $T = [100. \ 122.22222222 \ 166.66666667 \ 211.11111111 \ 255.55555556$

$300. \ 344.44444444 \ 388.88888889 \ 433.33333333 \ 477.77777778$

$500. \ ]$

.....



# Resultados grid 1x2

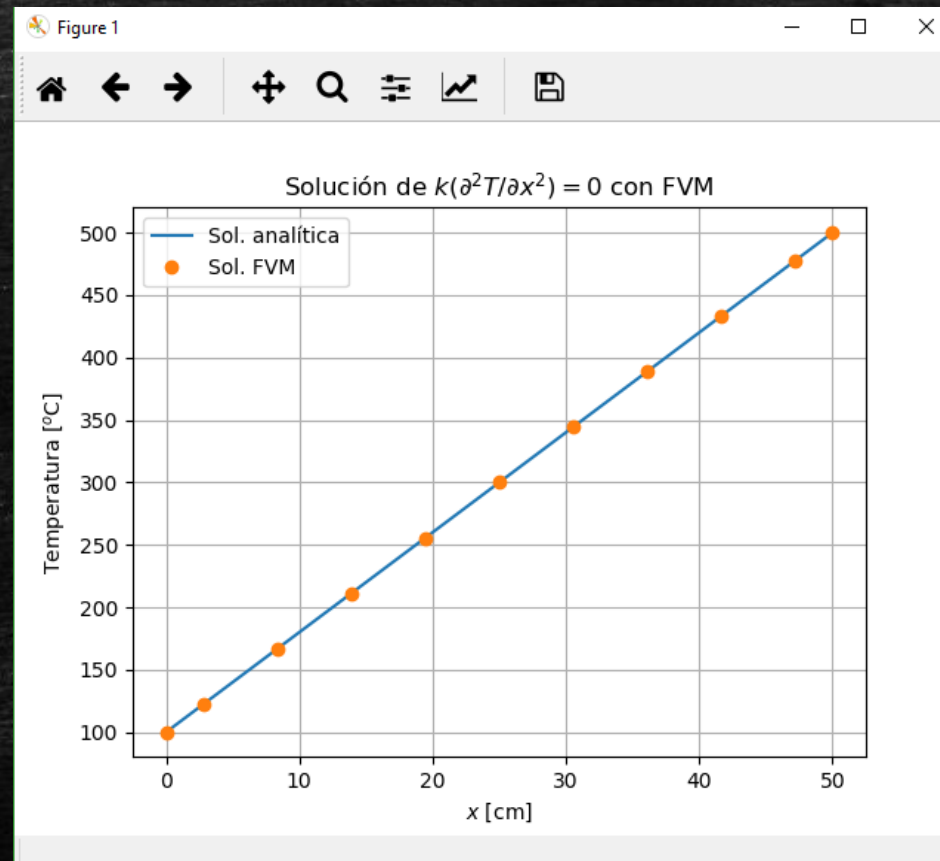
```
Building a 1 x 2 grid topology
Solucionaremos la T = [0. 0. 0. 0. 0. 0. 0.] en el rank = 0
ahora t vale = [100. 0. 0. 0. 0. 0. 0.]
Solución para T= [100. 90. 70. 50. 30. 10. 0.]
.-----
Solución para T = [100. 95. 85. 75. 65. 55. 50.]
.-----
Solución para T = [100. 95.9 87.7 79.5 71.3 63.1 59. ]
.-----
Solución para T = [100. 99.95 99.85 99.75 99.65 99.55 99.5 ]
.-----
Solución para T = [100. 100.679 102.037 103.395 104.753 106.111 106.79 ]
.-----
Solución para T = [100. 103.9595 111.8785 119.7975 127.7165 135.6355 139.595 ]
.-----
Solución para T = [100. 104.54999 113.64997 122.74995 131.84993 140.94991 145.4999 ]
.-----
Solución para T = [100. 107.207195 121.621585 136.035975 150.450365 164.864755
172.07195 ]
.-----
Solución para T = [100. 107.6854919 123.0564757 138.4274595 153.7984433 169.1694271
176.854919 ]
.-----
Solución para T = [100. 109.83782795 129.51348385 149.18913975 168.86479565
188.54045155 198.3782795 ]
.-----
Solución para T = [100. 110.22524844 130.67574532 151.12624219 171.57673907
192.02723595 202.25248439]
.-----
Solución para T = [100. 111.96864064 135.90592192 159.8432032 183.78048448
207.71776576 219.68640639]
.-----
Solución para T = [100. 112.28245124 136.84735371 161.41225618 185.97715865
```

```
Solucionaremos la T = [0. 0. 0. 0. 0. 0. 0.] en el rank = 1
ahora TT vale = [ 0. 0. 0. 0. 0. 0. 500.]
Solución para TT= [ 0. 50. 150. 250. 350. 450. 500.]
.-----
Solución para TT = [ 10. 59. 157. 255. 353. 451. 500.]
.-----
Solución para TT = [ 55. 99.5 188.5 277.5 366.5 455.5 500. ]
.-----
Solución para TT = [ 63.1 106.79 194.17 281.55 368.93 456.31 500. ]
.-----
Solución para TT = [ 99.55 139.595 219.685 299.775 379.865 459.955 500. ]
.-----
Solución para TT = [106.111 145.4999 224.2777 303.0555 381.8333 460.6111 500. ]
.-----
Solución para TT = [135.6355 172.07195 244.94485 317.81775 390.69065 463.56355 500. ]
.-----
Solución para TT = [140.94991 176.854919 248.664937 320.474955 392.284973 464.094991
500. ]
.-----
Solución para TT = [164.864755 198.3782795 265.4053285 332.4323775 399.4594265 466.4864755
500. ]
.-----
Solución para TT = [169.1694271 202.25248439 268.41859897 334.58471355 400.75082813
466.91694271 500. ]
.-----
Solución para TT = [188.54045155 219.68640639 281.97831608 344.27022577 406.56213546
468.85404515 500. ]
.-----
Solución para TT = [192.02723595 222.82451236 284.41906517 346.01361798 407.60817079
469.2027236 500. ]
.-----
Solución para TT = [207.71776576 236.94598918 295.40243603 353.85888288 412.31532973
```

# Resultados grid 1x2

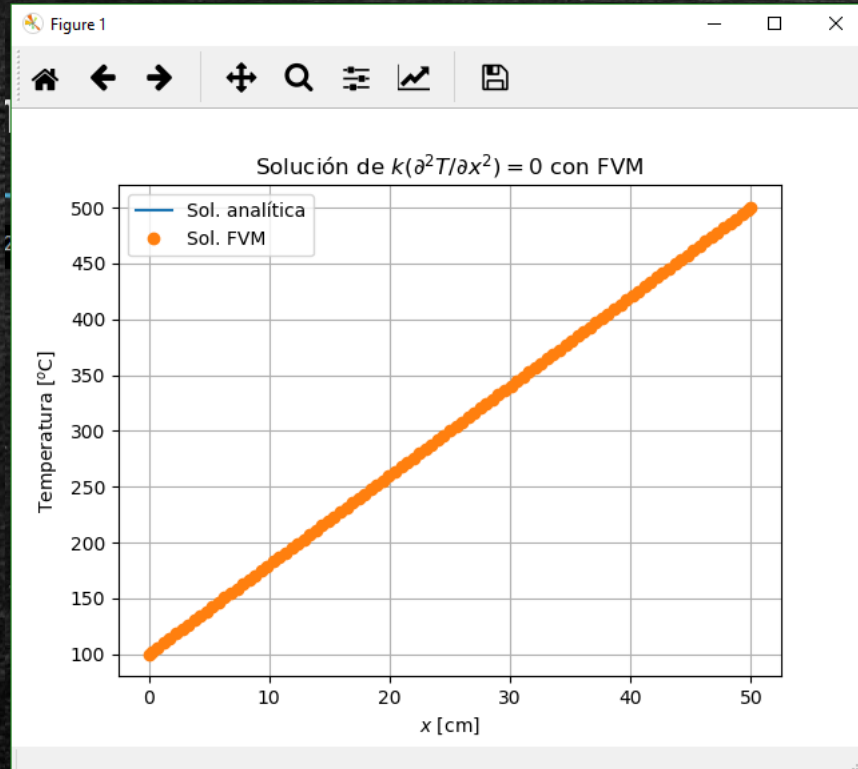
```
Solucion = [100.0, 121.05263157894733, 163.15789473684197, 205.26315789473665, 247.36842105263133,
```

```
310.52631578947336, 352.63157894736815, 394.73684210526295, 436.8421052631578, 478.9473684210526, 500.0]
```





# Comparación Secuencial vs Paralelo

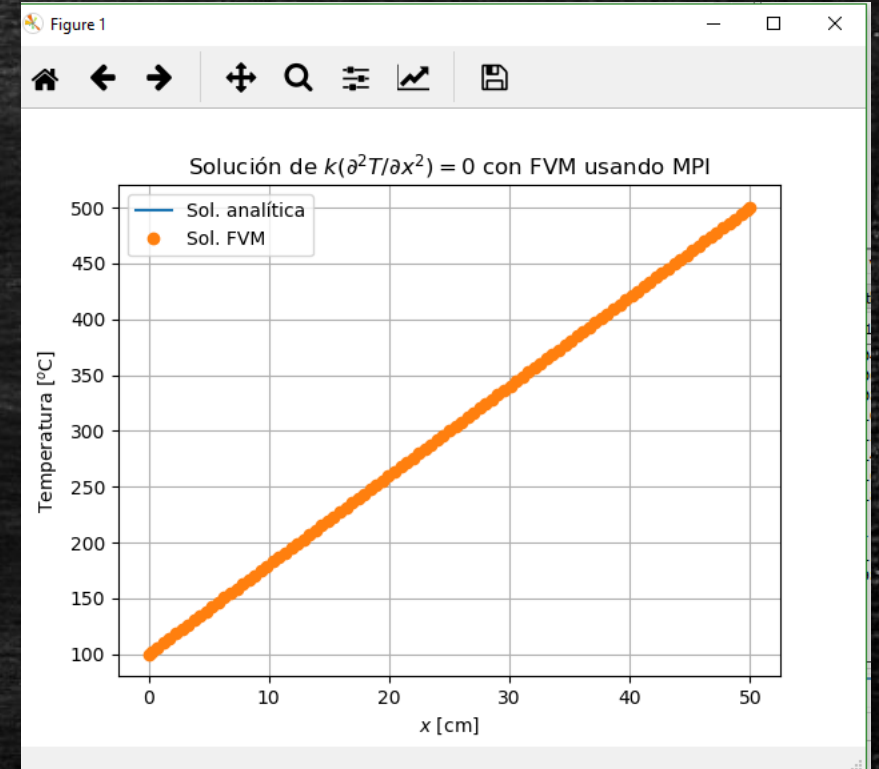


N = 100

500

Time = 0.006984

0.022998



N = 100

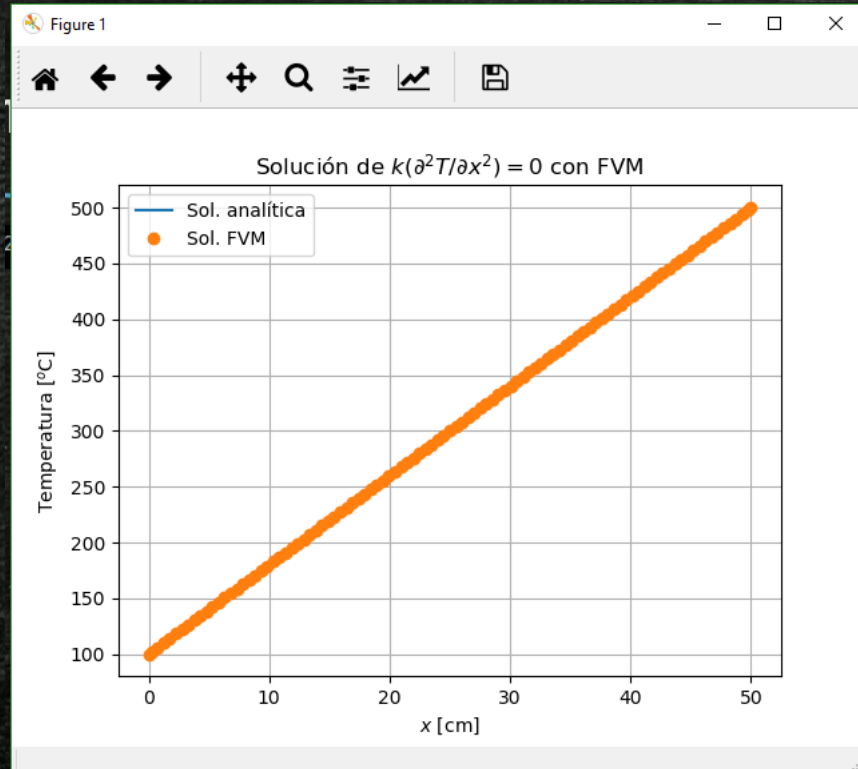
500

Time = 0.002298

0.016984

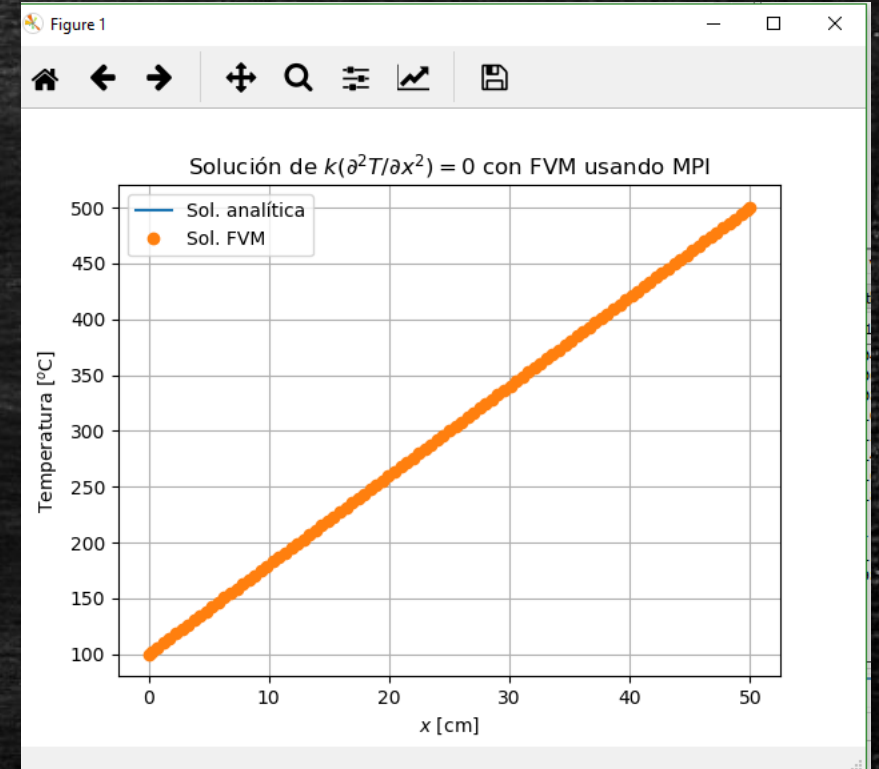
1x2 grid

# Comparación Secuencial vs Paralelo



N = 1000      10,000

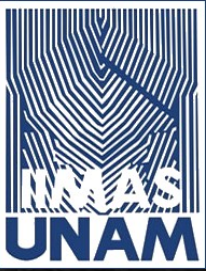
Time = 0.03298    8.43076



N = 1000      10,000

Time = 0.024974    6.808732





# Gracias

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