# Wildfire Spreading Model using a Parallel Implementation of Cellular Automata

ADVANCED APPLIED PARALLEL PROGRAMMING

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## **OUTLINE**

- Introduction
- PROBLEM
  - Formalization
  - Solution
- 3 Results
  - Evaluation
  - Validation/Discussion
  - Conclusion
  - Future Work





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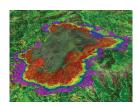




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## FORMALIZATION

#### PROBLEM

The integration of factors involved on wildfires is costly computationally.





#### SOLUTION

Construct a simulation model based on Cellular Automata which it allows decomposing the computations in order to map tasks to threads.





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- A parallel computing approach is introduced to reduce the execution time (more realistic simulations).





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- Forest fuels: All the vegetation elements, woody or herbaceous, alive or dead that may be flammable.
- **Topography**: It influences over the two above factors, the fuel and weather, modifying or altering them.



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• Cellular Automata





#### • Cellular Automata

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- Neighborhood: This CA model use the Moore Neighborhood, defined as

$$N_{(i_0,j_0)}^M = \{(i,j): |i-i_0| \le r, |j-j_0| \le r\},$$
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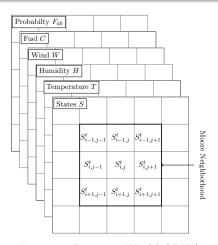
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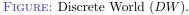
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- Transition Function:  $F: DW \to S$  where,



DW is a discretized world that contains states S, temperature T, humidity H, wind speed  $W_s$  and direction  $W_d$ , fuel C and the probability threshold of change from state a to b  $F_{ab}$ . All this variables are matrices with values between 0 to 1.







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Formally, we define the transition as

$$S_{i,j}^{t+1} = \begin{cases} 2, & \text{if } S_{i,j}^t = 1 \text{ and } f_{12} \ge F_{12} \\ 2, & \text{if } S_{i,j}^t = 4 \text{ and } f_{42} \le F_{42} \\ 3, & \text{if } S_{i,j}^t = 2 \text{ and } f_{23} \le F_{23} \\ 4, & \text{if } S_{i,j}^t = 3 \text{ and } f_{24} \le F_{24} \end{cases}$$

$$(2)$$

where functions f are defined by



- Function  $f_{12}$ :
  - The environmental factors E:

$$E_{i,j} = \frac{W_{d_{i,j}} C_{i,j} T_{i,j} W_{s_{i,j}}}{H_{i,j} P_{i,j}}.$$
 (3)

• The burning states of neighborhood  $p(N_b)$ :

$$p(N_b)_{i,j} = \frac{N_b}{8},\tag{4}$$

where  $N_b$  is the number of burning state neighboring cells. Then,  $f_{12}$  is computed by

$$f_{12} = \alpha \ E_{i,j} + \beta \ p(N_b)_{i,j}.$$
 (5)

• Functions  $f_{23}$ ,  $f_{24}$ ,  $f_{42}$  are random values between 0 and 1 and  $F_{23}$ ,  $F_{24}$ ,  $F_{42}$  are threshold parameters.

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- The model works with a world discretized by layers (DW), where each layer contains the information of the components described before.
- Model evolves in discrete times.



## ALGORITHM

### Algorithm 1 Main Algorithm

 $S^0 \leftarrow$  Initialize cell's states.

 $DW \leftarrow$  Initialize discrete world.

for t = 0 to  $T_{max}$  do  $S^{t+1} \leftarrow spreading(S^t, DW)$ 

end for



#### ALGORITHM

## Algorithm 2 Spreading Algorithm

```
procedure SPREADING(S^t, DW)
    N \leftarrow \text{number of rows in } DW
    for i = 0 to N_{threads} - 1 do
        delta \leftarrow N/N_{threads}
        start \leftarrow i \cdot delta
        if i = N_{threads} - 1 then
            end \leftarrow N
        else
            end \leftarrow (i+1) \cdot delta
        end if
        S_{start:end,\cdot} \leftarrow subSpreading(start, end, S^t, DW)
    end for
    for i = 0 to N_{threads} - 1 do
        Thread's join
    end for
    return S
end procedure
```



### ALGORITHM

```
Algorithm 3 Sub-spreading Algorithm
```

```
\begin{aligned} & \textbf{procedure} \ \text{SUBSPREADING}(start, end, S^t, DW) \\ & N \leftarrow \text{number of columns in } DW. \\ & \textbf{for } i = start \ \text{to } end \ \textbf{do} \\ & \textbf{for } j = 0 \ \text{to } N \ \textbf{do} \\ & \text{Compute } S^t_{i,j} \ \text{using equation (2) with } S^t \ \text{and } DW. \\ & \textbf{end for} \\ & \textbf{end for} \\ & \textbf{return } S^t_{start:end,\cdot} \\ & \textbf{end procedure} \end{aligned}
```



## ALGORITHM

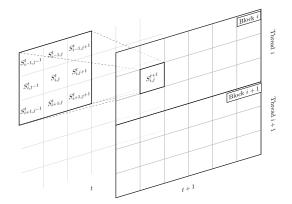


FIGURE: SubSpreading computing a matrix block.



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- If we choose  $T_{max} \ll N$  the complexity is about  $O(N^2)$ , but it may change a lot if  $T_{max} \sim N$  increasing the complexity to  $O(N^3)$ .
- This is the main motivation for including threads in the computing of the model.



The experiments were made using a fixed  $T_{max} = 50$ , for 1 to 4 number of threads, repeating the simulations 10 times per N.

Table: Summary of the average times in seconds.

Threads	N = 100	N = 500	N = 1000	N = 1500
1	0.198	8.023	32.861	72.464
2	0.135	5.147	19.695	44.407
3	0.163	4.825	18.852	41.626
4	0.220	4.650	18.055	41.458

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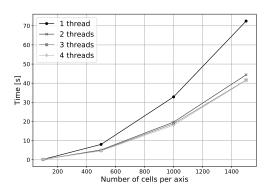


FIGURE: Threads performance comparison.



Table: Speedups results.

Threads	N = 100	N = 500	N = 1000	N = 1500
2	1.47	1.56	1.67	1.63
3	1.21	1.66	1.74	1.74
4	0.90	1.73	1.82	1.75





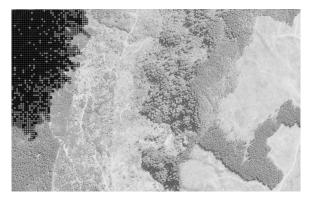


FIGURE: A portion of map simulated for  $N=1500,\,T=30^{\circ}{\rm C},\,H=50\%,\,W_s=40$  km/hr,  $W_d=90^{\circ},\,P=50$  hPa,  $F_{23}=F_{24}=F_{42}=0.1.$ 



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- The wildfire dynamics generated with the prototypes are consistent respect to the past events reported by CONAF.
- The complexity of the model exhibits a direct relationship between realism the amount fire factors.





### CONCLUSION

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- According to model features, assign tasks to threads is an effective approach for reducing the execution time.
- Respect to the EDP-based models, the Cellular Automata exhibits a favorable trade-off between the fire factors and the model complexity.



# FUTURE WORK

• Increase the resolution of the simulated area in order to improve the model approximation.





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- Increase the resolution of the simulated area in order to improve the model approximation.
- In order to take advantage of computation of several contiguous CA states, implementing a parallel approach based on pipeline.



## References I

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