

# Simplified Combustion Modeling Via Reduced-Physics Approximation In COMSOL Multiphysics®

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## Abstract

Accurately resolving combustion normally requires solving large, stiff sets of elementary reaction equations. We present a reduced-physics approximation that replaces detailed chemistry with a single, smoothed reaction-rate expression  $\{rr = kk \cdot \text{step1}(T[1/K]) \cdot c\}$ , where  $kk$  is a tunable rate constant,  $c$  is the local mass fraction of the combustible mixture, and  $\text{step1}(T)$  is COMSOL's continuously differentiable Heaviside function, calibrated to mimic Arrhenius activation. The corresponding volumetric heat source is  $\{Q = rr \cdot Hu\}$  with  $Hu$  the lower heating value. This formulation captures the essential coupling between species depletion and temperature rise while avoiding the stiffness and cost of multi-step kinetics.

Problem setup in COMSOL Multiphysics®:

The model is implemented with the Heat Transfer in Fluids and Transport of Diluted Species interfaces (Heat Transfer Module + Chemical Reaction Engineering Module). Temperature-dependent density and specific heat are imported from the Materials Library; gas thermal conductivity  $\{k\}$  is treated as a user parameter  $\text{cond0}$ . A 2D domain representing a generic gaseous layer is initialized with a narrow hot spot to trigger ignition.

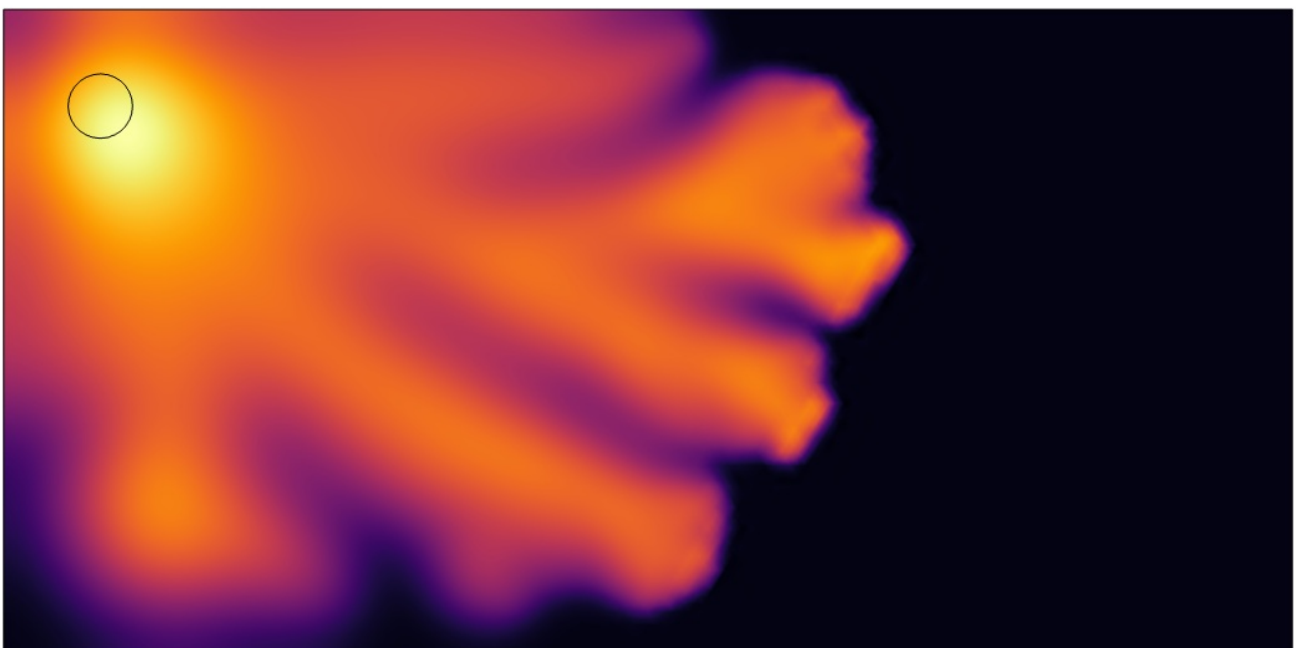
Results:

Across 18 parameter combinations, the model reproduces key combustion behaviors at a fraction ( $< 5\%$ ) of the computational cost of detailed chemistry. For high  $kk$ , high  $Hu$ , and low  $\text{cond0}$  we observe dendritic, finger-like flame propagation reminiscent of wild-fire spread. Quantitatively, the front speed increases up to 60 % relative to planar cases, and the calculated fractal dimension rises from 1.00 to 1.35, indicating significant front roughening. Conversely, higher thermal conductivities damp this instability and yield smooth, quasi-planar fronts.

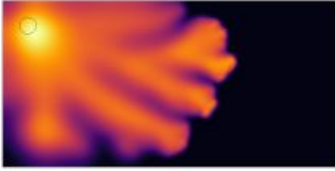
Conclusions and implications:

The reduced-physics approach retains essential thermo-chemical coupling while enabling rapid screening of material properties and operational envelopes. It is particularly useful for preliminary design of safety measures in low-conductivity gas environments (e.g., hydrogen-rich blends or forest under-canopies).

## Figures used in the abstract



**Figure 1** : Fractal Flame Front



**Figure 2** : thumbnail