

QES-Fire: a dynamically coupled fast-response wildfire model

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ABSTRACT

A microscale wildfire model, QES-Fire, that dynamically couples the fire front to microscale winds was developed using a simplified physics rate of spread (ROS) model, a kinematic plume-rise model and a mass-consistent wind solver. The model is three-dimensional and couples fire heat fluxes to the wind field while being more computationally efficient than other coupled models. The plume-rise model calculates a potential velocity field scaled by the ROS model's fire heat flux. Distinct plumes are merged using a multiscale plume-merging methodology that can efficiently represent complex fire fronts. The plume velocity is then superimposed on the ambient winds and the wind solver enforces conservation of mass on the combined field, which is then fed into the ROS model and iterated on until convergence. QES-Fire's ability to represent plume rise is evaluated by comparing its results with those from an atmospheric large-eddy simulation (LES) model. Additionally, the model is compared with data from the FireFlux II field experiment. QES-Fire agrees well with both the LES and field experiment data, with domain-integrated buoyancy fluxes differing by less than 17% between LES and QES-Fire and less than a 10% difference in the ROS between QES-Fire and FireFlux II data.

Keywords: buoyant plume, diagnostic wind solver, fire–atmosphere coupling, level set method, merging plumes, plume rise model, rate of spread, simplified fire spread physics.

Introduction

Winds are one of the primary drivers at the fire front (Albini 1982) dictating fire-spread behaviour. Currently, only models that directly include the physical coupling between combustion processes and fluid momentum and scalar transport represent this phenomenon with minimal parameterisation or empiricism. For wildfires, these include models such as FIRETEC (Linn *et al.* 2002), the Wildland–Urban Interface Fire Dynamics Simulator (WFDS, Mell *et al.* 2007) and Firestar (Morvan and Dupuy 2004; Morvan *et al.* 2009). These models resolve fire progression based on physical representation of the combustion without relying on rate of spread (ROS) parameterisations, and account for the impact of fire–atmosphere interactions on fire behaviour. These models directly resolve entrainment due to combustion, as well as buoyancy, by solving the equations for conservation of mass, momentum and energy. However, the resolution required to resolve the physics of atmosphere–fire interactions precludes simulations that cover a large enough range of spatial and temporal scales to represent large wildfires such as those that have transpired in recent years, e.g. the 2018 British Columbia wildfires, which burned 1 354 284 ha (BC Wildfire Services 2020), the 2020 Oregon wildfires, which burned more than 400 000 ha (National Interagency Fire Center (NIFC) 2020), and the August complex fire in California, which burned more than 417 000 ha (NIFC 2020).

One solution used to mitigate the high computational cost of full-physics models is to simplify and parameterise different components of the fire–atmosphere system, for example the ROS. For wildfire, this is commonly done with empirically derived parameterisations. One example is the Fire Area Simulator (FARSITE, Finney *et al.* 1995), a deterministic modelling system capable of simulating fire spread in heterogeneous terrain with heterogeneous fuel and fuel moisture conditions. It includes several semi-empirical

Received: 30 April 2021

Accepted: 21 January 2022

Published: 18 March 2022

Cite this:

Moody MJ *et al.* (2022)

International Journal of Wildland Fire

31(3), 306–325. doi:[10.1071/WF21057](https://doi.org/10.1071/WF21057)

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ROS models: the [Rothermel \(1972\)](#) surface fire spread model, the [Van Wagner \(1977\)](#) crown fire initiation model, the [Rothermel \(1991\)](#) crown fire spread model, the [Albini \(1979\)](#) spotting model, and the [Nelson \(2000\)](#) dead fuel moisture model. Another example of an empirically parameterised model is the narrow-band Eulerian Level set Method of Fire spread (ELMFire, [Lautenberger 2013](#)). ELMFire includes the [Rothermel \(1972\)](#) surface fire spread model as the forcing of a level set method ([Sethian 1999](#)). Each point of the fire front is advanced by applying Huygens' principle, which assumes that the fire front propagation can be approximated as the sum of elliptical wavelets emanating from each point of the fire front ([Baker and Copson 2003](#)). Models such as FARSITE and ELMFire represent fire behaviour with ingested meteorological data as the forcing, but are limited because fire heat sources do not affect atmospheric winds.

Models such as the Coupled Atmosphere–Wildland Fire Environment (CAWFE, [Clark et al. 2004](#); [Coen 2013](#)), MesoNH-ForeFire ([Filippi et al. 2011](#)), Weather Research and Forecasting – SFIRE (WRF-SFIRE, [Mandel et al. 2011, 2014](#)), and WRF-Fire ([Coen et al. 2013](#); [Muñoz-Esparza et al. 2018](#)) focus on the coupling between mesoscale meteorological conditions and wildfire behaviour at spatial scales of the order of hundreds of metres ($\mathcal{O}(100\text{m})$) using numerical weather prediction (NWP) models. For example, WRF-SFIRE couples the WRF model ([Skamarock and Klemp 2008](#)) with semi-empirical ROS and simplified fire heat flux parameterisations. A level set method is used to track and advance the flame front, and winds from the WRF model are interpolated to the fire front height for use in the ROS and fuel consumption models. In turn, heat fluxes from SFIRE are injected into the WRF model as a vertically distributed flux using an extinction depth. MesoNH-ForeFire ([Filippi et al. 2011, 2013](#)) uses the model of [Balbi et al. \(2009\)](#) as a forcing to a Lagrangian front tracking scheme. The ForeFire module interpolates the atmospheric winds to the midflame height and then calculates the fire-induced winds through the heat and water vapour surface fluxes. These combined winds are used to drive the ROS model. The main differences in these models are in the methodology for how the winds are interpolated to the fire front, the choice of ROS parameterisation, and how the heat and mass fluxes are injected into the atmospheric model. The feedback between meteorological variables and fire heat and mass fluxes enables these models to simulate fire-related winds and wildfire evolution. [Kochanski et al. \(2013\)](#) demonstrated that WRF-SFIRE can run coupled atmosphere–fire simulations faster than real time for large-scale wildfires. The primary limitation is that the grid resolution for such fires must be fairly modest ($\mathcal{O}(100\text{ m})$) to keep the computational time low enough for forecasting applications.

WRF-SFIRE and similar models use a full set of prognostic equations to represent the atmosphere in a coupled modelling framework, making them computationally intensive. They also typically use near-surface atmospheric and fire

front-scale parameterisations that were not developed for the small-scale three-dimensional (3-D) heterogeneity commonly found in most ecosystems. Recently, a third approach has emerged to couple atmospheric flow to fire spread that attempts to bridge the gap between the FIRETEC-WFDS approach and the coupled mesoscale atmosphere-to-fire front approach. The basic idea is to employ a simplified but still 3-D physics methodology for the microscale flow while representing fire behaviour with parameterisations similar to those used by WRF-SFIRE. QUIC-fire ([Linn et al. 2020](#)) does this by combining the diagnostic wind solver QUIC-URB ([Pardyjak and Brown 2003](#)) with a cellular automata model ([Achtemeier 2013](#)). QUIC-fire improves on the cellular automata model to include a more physics-based approach. It moves ‘energy packets’ (EP) through diffusion or advection based on the local wind. The transferred EP is used to either start new fires, intensify fires already ignited, or transfer the energy to the environment using a probability-based ROS. In addition, the impacts of fluxes originating at the fire front on the wind field are parameterised in QUIC-fire by converting the heat flux at the fire source to a buoyant plume with the plume centreline vertical velocity and lateral expansion calculated using the Briggs theory parameterisations ([Davidson 1989](#)).

Here, we present a new fast-response microscale model for wildfire spread, QES-Fire. It shares some characteristics with QUIC-fire, including the use of a similar diagnostic wind model, and its general strategy is similar while the specific submodels, methods and system flow are unique. The model is based on the QES (Quick Environmental Simulation) framework, a fast-response microscale model that can rapidly calculate wind fields in complex terrain ([Bozorgmehr et al. 2021](#)). QES-Fire uses quasi-steady-state simplified physics parameterisations to dynamically couple the ROS and fire heat fluxes with local winds. We first describe the modelling methodology used to create QES-Fire in the next section, including the fire front advancement method, the microscale wind solver, the ROS model and the parameterisations for the fire-induced winds through use of a convective plume model. The convective plume model includes single axisymmetric plumes as well as a methodology for determining the criteria for plume merging. Next, we compare the model in the ‘QES-Fire comparison’ section against an idealised plume and *in situ* measurements from an experimental burn.

QES-Fire modelling methodology

The goal of QES-Fire is to dynamically link near-surface atmospheric flow with fire front fluxes and fire front ROS in a physically consistent and computationally efficient manner. To accomplish this, a unique combination of reduced-order physical and chemical models, numerical methods and computational approaches is used. The model is composed of a set of linked submodels, each representing an aspect of the

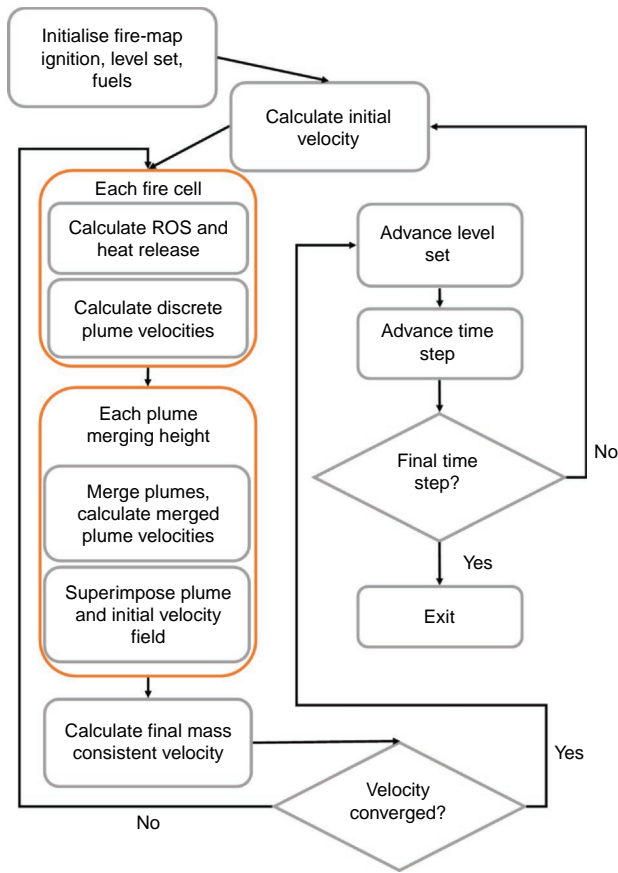


Fig. 1. Flowchart of QES-Fire depicting the dynamic coupling of the initial atmospheric winds with the fire-induced winds used to drive a ROS model to advance a fire front.

wildfire–atmosphere system (Fig. 1). Briefly, the coupling between atmospheric winds and fire heat fluxes is accomplished using a ROS model that advances the fire front through the level set method (Sethian 1999). The level set advances through QES-Fire’s discrete, two-dimensional (2-D) surface mapper that tracks fuel conditions at the domain’s surface. At the beginning of a simulation, the 2-D mapper is initialised with the fire’s ignition source and fuel characteristics. Next, an initial wind field is specified throughout the 3-D domain using QES-Winds and at each discrete cell that is on fire, a ROS model (Balbi et al. 2020) calculates the forcing used to advance the level set as well as the heat release of each discrete section of the fire front. A plume rise and merging model is then used to calculate a velocity field from the discrete fire map. The empirical plume velocity field is superimposed onto the initial velocity field and QES-Winds calculates a final velocity field. This iterates until convergence at which point the level set (fire front) and time step are advanced.

The following subsections give detailed descriptions of each component depicted in Fig. 1 and how they contribute to QES-Fire. All variables used in QES-Fire are listed in Table 1.

Table 1. Variable nomenclature.

Latin symbols		Value if constant
A	Radiation coefficient	
B	Stefan–Boltzman constant ($\text{W m}^{-2} \text{K}^{-4}$)	5.6×10^{-8}
b	Buoyancy velocity in flame (m s^{-1})	
b_n	Buoyancy velocity in flame normal to flat ground (m s^{-1})	
C	Courant number	
c_p	Specific heat of air ($\text{J kg}^{-1} \text{K}^{-1}$)	1150
d	Fuel depth (m)	
D	Level set spatial derivatives	
F	Level set forcing function	
g	Acceleration due to gravity (m s^{-2})	9.81
h	Height of flame base (m)	
H	Height of flame (m)	
I	Ignition energy (J kg^{-2})	
K_1	Drag coefficient (s m^{-1})	130
l_0	Roughness length (m)	
L_c	Characteristic plume length (m)	
m	Fuel moisture content (%)	
n_x	Normal to level set along x-axis	
n_y	Normal to level set along y-axis	
\tilde{p}	Perturbation from hydrostatic pressure	
\dot{q}'''	Rate of energy release from plume source (W m^{-3})	
Q_0	Averaged power output of fire (W)	
r	Radial distance from plume centreline (m)	
r_{00}	Balbi model coefficient	2.5×10^{-5}
R	Plume radius (m)	
ROS	Rate of spread (m s^{-1})	
s_c	Stoichiometric coefficient	17
S	Leaf area per square metre ($\text{m}^2 \text{m}^{-2}$)	
SAV	Surface area-to-volume ratio ($\text{m}^2 \text{m}^{-3}$)	
t	Time (s)	
T	Temperature (K)	
u	Velocity component along x-axis (m s^{-1})	
$U(H/2)$	Horizontal velocity at midflame height (m s^{-1})	
v	Velocity component along y-axis (m s^{-1})	
\bar{v}_r	Radial plume velocity (m s^{-1})	
\bar{v}_z	Vertical plume velocity (m s^{-1})	
\vec{V}	Velocity (m s^{-1})	
w_c	Characteristic plume velocity (m s^{-1})	

(Continued on next page)

Table 1. (Continued)

Latin symbols		Value if constant
w	Velocity component along z-axis (m s^{-1})	
x	Horizontal Cartesian axis	
x_0	Distance between two plume sources	
y	Horizontal Cartesian axis	
z	Vertical Cartesian axis	
Greek symbols		Value if constant
$\alpha_{1,2}$	Gaussian precision moduli	
β	Fuel packing ratio	
ΔH	Heat of combustion of pyrolysis gases (J kg^{-1})	
Δx	Cell size along x-axis (m)	
Δy	Cell size along y-axis (m)	
Δz	Cell size along z-axis (m)	
ϵ	Plume entrainment constant	0.09
γ	Flame tilt angle (radians)	
Γ	Plume-merging filter index	
κ	Relative strength of unequal plume sources	
λ	Lagrange multipliers for Poisson equation	
Λ	Plume potential	
μ_m	Plume mixing length (m)	
η	Ratio of thermal to momentum plume radii	
ω	Vorticity (radians s^{-1})	
Ω	Empirically derived velocity profile in plume	
ϕ	Level set function	
Ψ	Streamfunction	
ρ	Density (kg m^{-3})	
ρ_v	Fuel particle density (kg m^{-3})	
σ	Total fuel load (kg m^{-2})	
τ	Flame residence time (s)	
τ_0	Residence time parameter (s m^{-1})	75 591
θ	Terrain slope angle (radians)	
Θ	Empirically derived temperature profile in plume	
X	Flame radiative fraction	
ζ	Vertical distance above plume source (m)	
ζ_v	Virtual origin below sources for merged plume (m)	
Superscripts		
0	Original value	
*	Non-dimensional value	
+	Forward in space	
-	Backward in space	

Subscripts

a	Air
c	Characteristic value
cl	Plume centre-line value
f	Flame
h	With respect to horizontal
i	x-axis cell index
j	y-axis cell index
k	z-axis cell index
r	Along radial direction
x	Along x-axis
y	Along y-axis
z	Along z-axis

Fire map

QES-Fire uses a cellular 2-D surface map to track fuels, moisture, local terrain slope and the fire front. All values are calculated at cell centres. The 2-D mapper (Fig. 2) defines four types of cells: not burning, partially burning, fully burning and burned. Additionally, the 2-D mapper tracks the fire front using the level set method (see 'Fire front' section). While the fire front is tracked as a continuous function (or set of continuous functions for multiple fires), each cell of the 2-D mapper is defined independently. Using a cellular approach gives QES-Fire the ability to build a fire front for heterogeneous conditions by varying the fuel loading, moisture and local slope for each distinct ground cell.

Fuel properties for each ground cell are specified according to the Rothermel (1972) and Anderson (1982) fuel loading models. These models include all the inputs necessary to compute the Rothermel ROS model. Each cell is considered homogeneous throughout; this is required because many ROS models assume a uniform fuel bed (e.g. Rothermel 1972; Pagni and Peterson 1973; Balbi *et al.* 2009). QES-Fire uses two ROS models concurrently to advance the fire front, Rothermel (1972) and Balbi *et al.* (2020) (see 'Rate of spread model' section below).

QES-Fire's 2-D mapper uses the cell type to track where the fire front is in time and space. The ROS model calculates the ROS in all cells that are fully and partially burning and the fire front is advanced using a dynamic timestep (Δt) within all partially burning cells. Δt is calculated using a modified Courant number, C (Ferziger and Peric 2002), with the maximum ROS for the 2-D mapper replacing the velocity magnitude

$$C = \max(\text{ROS}) \left(\frac{\Delta t}{\Delta x} \right), \quad (1)$$

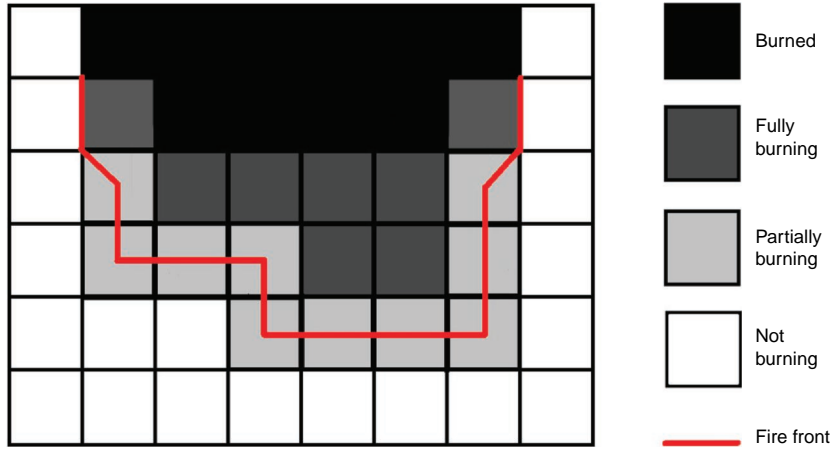


Fig. 2. Plan-view schematic illustrating how QES-Fire's 2-D mapper uses a level set function to track the fire front through discrete cells. Cells are either not burning, partially burning, fully burning, or burned. Cells are updated according to the fire front progression at each timestep.

where $C \leq 1$. The ROS-dependent timestep ensures that the flame front does not skip cells, facilitating an accurate calculation of the time on fire (t) through a simple addition of Δt to the current time on fire. At the point in time when $t \geq \tau$, where τ is the residence time calculated based on the fuel type and moisture following [Anderson \(1969\)](#), the cell is considered burned out and the cell type is set to burned so that fire cannot re-spread to that cell.

Fire front

The fire front is advanced through the level set method described by [Sethian \(1999\)](#). The level set method was chosen for its ability to track a fast-evolving interface in complex geometries, and because it has been shown to be adept at fire front tracking (e.g. [Mallet et al. 2009](#); [Rehm and McDermott 2009](#); [Mandel et al. 2011](#); [Muñoz-Esparza et al. 2018](#)). The initial fire front is prescribed in the initialisation of QES-Fire as the level set function ϕ so that $\phi_{i,j}(t = 0) = 0$, where $\phi_{i,j} = 0$ are burning cells and the subscripts i and j indicate the cell index in the x and y directions, respectively. A signed distance function (the minimum distance from a point to the level set) is then applied to all cells to determine the initial values for ϕ . Ghost cells are used by copying the values of the level set at the fire map boundaries to a buffer region around the 2-D mapper. This ensures spatial derivatives are non-complex at domain boundaries. As the progression is always forwards in a fire front (a fire cannot burn where it has already burned), it is appropriate to use an upwind numerical scheme; therefore, a first-order space convex scheme is utilised where

$$\phi_{i,j} = \phi_{i,j}^0 - \Delta t [\max(F_{i,j}, 0) \nabla^+ + \min(F_{i,j}, 0) \nabla^-], \quad (2)$$

with Δt calculated using the modified Courant number ([Eqn 1](#)). $F_{i,j}$ is the level set forcing function given by the ROS model (see 'Rate of spread model' section) and

∇^+ and ∇^- are the forward and backward gradients (in space) of the level set calculated by

$$\begin{aligned} \nabla^+ = & (\max[D_{i,j}^{-x}, 0])^2 + (\min[D_{i,j}^{+x}, 0])^2 \\ & + (\max[D_{i,j}^{-y}, 0])^2 + (\min[D_{i,j}^{+y}, 0])^2, \end{aligned} \quad (3)$$

and

$$\begin{aligned} \nabla^- = & (\max[D_{i,j}^{+x}, 0])^2 + (\min[D_{i,j}^{-x}, 0])^2 \\ & + (\max[D_{i,j}^{+y}, 0])^2 + (\min[D_{i,j}^{-y}, 0])^2, \end{aligned} \quad (4)$$

respectively. The level set spatial derivatives $D_{i,j}$ are given by

$$\begin{aligned} D_{i,j}^{+x} &= \frac{\phi(x + \Delta x, t) - \phi(x, t)}{\Delta x}, \\ D_{i,j}^{-x} &= \frac{\phi(x, t) - \phi(x - \Delta x, t)}{\Delta x}, \end{aligned} \quad (5)$$

$$\begin{aligned} D_{i,j}^{+y} &= \frac{\phi(y + \Delta y, t) - \phi(y, t)}{\Delta y}, \text{ and} \\ D_{i,j}^{-y} &= \frac{\phi(y, t) - \phi(y - \Delta y, t)}{\Delta y}. \end{aligned} \quad (6)$$

The level set forcing function $F_{i,j}$ must be applied normal to the fire front, both in calculation of the ROS and for the propagation of the level set itself. The level set normal is calculated as

$$n_x = \frac{n_{x'}}{\sqrt{n_{x'}^2 + n_{y'}^2}} \text{ and } n_y = \frac{n_{y'}}{\sqrt{n_{x'}^2 + n_{y'}^2}}, \quad (7)$$

with

$$\begin{aligned} n_{x'} &= \frac{D_{i,j}^{+x}}{\sqrt{(D_{i,j}^{+x})^2 + (D_{i,j}^{+y})^2}} + \frac{D_{i,j}^{-x}}{\sqrt{(D_{i,j}^{-x})^2 + (D_{i,j}^{-y})^2}} \\ &+ \frac{D_{i,j}^{+x}}{\sqrt{(D_{i,j}^{+x})^2 + (D_{i,j}^{-y})^2}} + \frac{D_{i,j}^{-x}}{\sqrt{(D_{i,j}^{-x})^2 + (D_{i,j}^{+y})^2}}, \end{aligned} \quad (8)$$

and

$$n_{y'} = \frac{D_{ij}^{+y}}{\sqrt{(D_{ij}^{+x})^2 + (D_{ij}^{+y})^2}} + \frac{D_{ij}^{-y}}{\sqrt{(D_{ij}^{+x})^2 + (D_{ij}^{+y})^2}} + \frac{D_{ij}^{+y}}{\sqrt{(D_{ij}^{+x})^2 + (D_{ij}^{-y})^2}} + \frac{D_{ij}^{-y}}{\sqrt{(D_{ij}^{+x})^2 + (D_{ij}^{-y})^2}}, \quad (9)$$

where n_x and n_y are the x and y components of the level set normal, respectively.

Microscale winds

The QES modelling framework is used to dynamically couple the fire front ROS and the velocity field using QES's 3-D diagnostic wind solver QES-Winds (Bozorgmehr *et al.* 2021). QES-Winds is a fast-response mass-conserving wind model based on Röckle (1990) and refined by Kaplan and Dinar (1996). The model prescribes a background velocity field and then uses empirical parameterisations to account for flow around buildings and trees in urban environments. Conservation of mass is then enforced using a variational analysis technique (Singh *et al.* 2008; Brown *et al.* 2013). The method solves the Poisson equation for Lagrange multipliers, λ :

$$\frac{\partial^2 \lambda}{\partial x^2} + \frac{\partial^2 \lambda}{\partial y^2} + \left(\frac{\alpha_1}{\alpha_2} \right)^2 \frac{\partial^2 \lambda}{\partial z^2} = \nabla \cdot \vec{V}^0, \quad (10)$$

where $\nabla \cdot \vec{V}^0$ is the divergence of the initial prescribed velocity field with $\vec{V}^0 = u_{ij,k}^0 + v_{ij,k}^0 + w_{ij,k}^0$ and u , v , and w are the velocity components in the x , y , and z directions, respectively, the superscript 0 indicates an initial value from a model parameterisation, and α_1 and α_2 are the Gaussian precision moduli that define the relative importance of horizontal versus vertical air motions, respectively.

QES-Winds solves Eqn 10 for the Lagrange multipliers using graphic processing unit (GPU) technology with a red-black successive over relaxation (SOR) scheme to achieve a speedup of two orders of magnitude compared with a serial SOR solver (Bozorgmehr *et al.* 2021). The velocity field is then updated using the Euler–Lagrange equations:

$$u_{i,j,k} = u_{i,j,k}^0 + \frac{1}{2\alpha_1^2 \Delta x} [\lambda_{i+1,j,k} - \lambda_{i,j,k}], \quad (11)$$

$$v_{i,j,k} = v_{i,j,k}^0 + \frac{1}{2\alpha_1^2 \Delta y} [\lambda_{i,j+1,k} - \lambda_{i,j,k}], \quad \text{and} \quad (12)$$

$$w_{i,j,k} = w_{i,j,k}^0 + \frac{1}{2\alpha_2^2 \Delta z} [\lambda_{i,j,k+1} - \lambda_{i,j,k}], \quad (13)$$

where Δx , Δy , and Δz are the cell sizes corresponding to the x , y , and z directions, respectively.

The modelling strategy used by QES-Winds has been validated for several urban configurations with and without

trees and gives quasi-steady-state velocity fields that are in good agreement with experimental data and full physics simulations (Hayati *et al.* 2017, 2019; Barbano *et al.* 2020). To adapt QES-Winds for wildfires, a parameterisation for the velocity field in and around fires is needed ('QES-Fire plume parameterisation' section). The initial velocity profile used in QES-Winds is an idealised rough-wall boundary layer wind profile (i.e. the stability corrected log-law). After a ground cell has been burned according to the 2-D fire map, the aerodynamic roughness for the cell is reduced owing to the lack of vegetation (fuels), which may provide a more realistic velocity profile upwind of the fire front that accounts for burned ground cover, as speculated by Sun *et al.* (2009).

Rate of spread model

The ROS model is used as the forcing function F , for the level set equation (Eqn 2). ROS models may be categorised into four separate regimes according to Sullivan (2009a, 2009b, 2009c); in order of descending complexity these are: physical, quasi-physical, quasi-empirical, and empirical models. Physical models include combustion chemistry and physics as well as momentum transfer. The main difference between physical models and quasi-physical models is the lack of combustion physics. Quasi-physical or simplified physics models may still include momentum transfer (using simplifications) and differentiate between the dominant modes of heat transfer in wildfires: radiation and convection. Quasi-empirical models lack this differentiation of heat transfer modes. Finally, empirical models simply match experimental ROS data to mathematical model parameters.

QES-Fire utilises two different quasi-steady-state approaches sequentially for its ROS model. First, the well-known quasi-empirical Rothermel (1972) model is used to provide an estimate for the ROS based on the winds, local terrain slope and fuel characteristics per grid cell. Next, the iterative quasi-physical model developed by Balbi *et al.* (2009, 2010, 2020; Chatelon *et al.* 2017) uses the ROS from the Rothermel model as an initial value for the calculation of the final ROS. Both the Rothermel and Balbi ROS models were formulated for surface fire spread, constraining QES-Fire to that mode of fire propagation.

Rothermel model

The Rothermel model is employed in the United States National Fire Danger Rating System and has the advantage that most fuel models in the United States were created for use with the Rothermel model (Deeming 1977). Additionally, it is a simple algebraic model that can be run in a single step for each burning grid cell.

Balbi model

The Balbi *et al.* (2020) model accounts for three dominant factors that transfer heat from the flame front to the

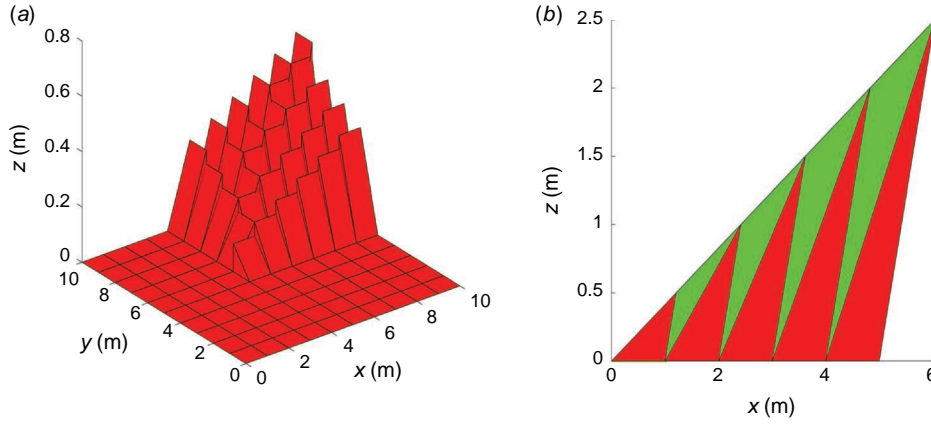


Fig. 3. (a) Flame geometry built in 3-D space using the Balbi model, with aggregated flames used to build the fire front. (b) 2-D geometric representation of Balbi model (green) using discrete prisms (red).

fuel, and the subsequent ignition of the fuel in its calculation of the ROS. These include radiation from the base of the fire (assumed to be a black body), radiation from the flame (simplified to a triangular grey body) and convection from the flame base.

Inputs for both the Rothermel and Balbi models include fuel parameters (depth of fuel, size, moisture, etc.), local slope, air temperature and the mid-flame height wind velocity calculated by QES-Winds. The Balbi model calculates the ROS (Eqn 14), flame height (Eqn 15), flame temperature (Eqn 16), buoyancy from the fire front (Eqn 17), heat release from the fire front (Eqn 18) and flame tilt angle (Eqn 19) as

$$\text{ROS} = \min\left(\frac{S}{\pi}, 1\right) \frac{BT_f^4}{\beta \rho_v I} + \frac{\text{SAV} \Delta H}{I \tau_0} \min\left(h, \frac{2\pi}{\text{SAV} \beta}\right) \left(\frac{h}{2h + H} \tan \theta + \frac{U(L) \exp(-K_1 \beta^{1/2} \text{ROS})}{b_h}\right) + A \times \text{ROS} \frac{1 + \sin \gamma - \cos \gamma}{1 + \frac{\text{ROS} \times \cos \gamma}{s \times r_{00}}}, \quad (14)$$

$$H = \frac{b_n^2}{g(T_f/T_a - 1)}, \quad (15)$$

$$T_f = T_a + \frac{\Delta H(1 - \chi)}{(s_t + 1)c_p}, \quad (16)$$

$$b = \left(\frac{h}{h + H/2}\right) \left(\frac{b_n}{2}\right), \quad (17)$$

$$\dot{Q}_0 = (1 - \chi) \left(\frac{\sigma \Delta H}{\tau}\right), \quad \text{and} \quad (18)$$

$$\tan \gamma = \tan \theta + \frac{U(L)}{b}. \quad (19)$$

In Eqns 14–19, S is the leaf area per square metre, B is the Stefan–Boltzman constant, T_f is the absolute flame temperature, β is the fuel packing ratio, ρ_v is the fuel particle density, I is the ignition energy, SAV is the surface area-to-

volume ratio of fine fuels, h is the height of the flame base, H is the flame height, b_n is the vertical gas velocity on flat terrain, T_a is the absolute ambient temperature, g is gravity, ΔH is the heat of combustion of pyrolysis gases, χ is the flame radiative fraction, c_p is the specific heat of air, \dot{Q}_0 is the heat release, σ is the total fuel load, γ is the flame tilt angle, θ is the local terrain slope angle, $U(L)$ is the midflame-height horizontal velocity calculated by QES-Winds, b is the vertical velocity inside the flame, and A is the flame radiation coefficient. The model includes three universal parameters: a flame drag coefficient K_1 , the stoichiometric coefficient for combustion s_t , and a parameter to account for flame thickness vs radiative fraction in the flame radiation model r_{00} . Heat transfer from the flames to the fuels is attenuated through the flame length in the Balbi model; this is preserved in the formulation by QES-Fire through the parameters in the Balbi model governing the convective and radiative heat transfer, K_1 and r_{00} .

The flame tilt angle for the Balbi model returns tilt normal to the fire front. For our 3-D implementation, flame tilt is calculated separately for the x and y directions for each 2-D mapper cell. The tilt angle along the direction with the strongest horizontal wind component (e.g. $U(L)$ in Eqn 19) is then chosen as the dominant direction and the flame over that cell is modelled as a triangular prism (Fig. 3). Calculating the tilt angle and flame geometry at each fully burning mapper cell in this manner enables a discrete approximation of flame structures that can represent individual point fires, complete fire fronts and fires burning in non-homogeneous terrain. While the prisms are modelled with the dominant wind direction, the ROS forcing for the level set method is calculated using the wind vector multiplied by the cosine of the angle between the wind and the normal to the fire front, thus preserving the formulation for ROS in the Balbi model.

To account for heterogeneous terrain, fuels, and winds, the flame geometry must be adapted to the QES-Winds scale ($\mathcal{O}(1 \text{ m})$) from the fire front scale ($\mathcal{O}(10 \text{ m})$) assumed in the original model (Balbi et al. 2009). Using the assumption that within a flame vertical motion is uniformly accelerated

owing to buoyancy and that flame outlines are streamlines (Chatelon *et al.* 2017), the horizontal velocity into the flame is approximately equal to the horizontal velocity out of the flame as the flame base length approaches zero. Additionally, assuming the flame residence time (τ) is inversely proportional to the surface area-to-volume ratio ($\tau = \tau_0/\text{SAV}$, where τ_0 is the residence time parameter; Anderson 1969), the discrete flame height can be decreased for a given Δx , where Δx = the flame base length for our discrete representation, linearly as a function of the time on fire (t) as

$$H = H^0 \times \left(1 - \frac{t}{\tau}\right) \quad (20)$$

where H is the modified flame height and H^0 is the original flame height. Thus, the Balbi geometry is recovered in 2-D using discrete prisms (Fig. 3).

Coupling the residence time dependence to the flame height with triangular prisms allows the construction of a fully 3-D fire front by aggregating the individual discrete flame models into a single fire front (Fig. 3). The ability of QES-Fire to represent surface conditions and fire front segments in a discrete manner enables heterogeneous fire fronts with varying fuel types, moisture contents and slope conditions to be modelled.

QES-Fire plume parameterisation

A major goal in the development of QES-Fire is to couple the ROS and local winds in the fire front vicinity, which enables future integration of QES-Fire with other fire-related phenomena (e.g. smoke dispersion, firebrand production and transport). This is accomplished using the axisymmetric plume model proposed by Baum and McCaffrey (1989) superimposed on the background QES-Winds velocity field. QES-Fire's cellular nature utilises merging plumes to account for adjacent and/or dispersed heat sources.

Velocity field from axisymmetric plume

The Baum and McCaffrey (1989) model for the flow field induced by fires provides a time-averaged quasi-steady-state velocity field for an axisymmetric plume that is scaled by the surface heat flux.

McCaffrey (1983) proposed a plume model based on the conservation of mass,

$$\frac{\partial \rho}{\partial t} + \vec{V} \cdot (\nabla \rho) + \rho (\nabla \cdot \vec{V}) = 0, \quad (21)$$

conservation of momentum,

$$\rho \frac{\partial \vec{V}}{\partial t} + \rho \vec{V} \cdot (\nabla \vec{V}) = -\nabla \tilde{p} + (\rho - \rho_0)g, \quad (22)$$

and conservation of energy equations,

$$\rho c_p \left(\frac{\partial T}{\partial t} + \vec{V} \cdot (\nabla T) \right) = \dot{q}''' \quad (23)$$

where ρ is the density of air, \tilde{p} is the perturbation from the hydrostatic pressure, and \dot{q}''' is the rate of energy release from a heat source per unit volume. For a large semi-infinite plume, the length scale for non-dimensionalisation is not readily apparent. By defining generic non-dimensional variables:

$$\begin{aligned} t^* &= t/t_c; & \rho^* &= \rho/\rho_0; & T^* &= \frac{T-T_0}{T_0} \\ \nabla^* &= \nabla \times L_c; & \tilde{p}^* &= \tilde{p}/p_c; & \vec{V}^* &= \vec{V}/\vec{V}_c; \\ \dot{q}'''^* &= \dot{q}''' L_c^3 / \dot{Q}_0, \end{aligned} \quad (24)$$

substituting them into Eqns 21–23, and then performing a scale analysis, it is possible to deduce the characteristic length

$$L_c \equiv \left(\frac{\dot{Q}_0}{\rho_0 c_p T_0 \sqrt{g}} \right)^{2/5}, \quad (25)$$

and characteristic velocity

$$w_c \equiv \left(\frac{g^2 \dot{Q}_0}{\rho_0 c_p T_0} \right)^{1/5} \quad (26)$$

scales for an axisymmetric plume, where \dot{Q}_0 is the total, volume-integrated power output of the distinct fire (Trelles 1995).

Inside the plume, the McCaffrey (1983) model assumes a non-dimensionalised Gaussian profile for both temperature,

$$\frac{T - T_0}{T_0}(r^*, \zeta^*) = \Theta_{cl}^*(\zeta^*) \exp \left[- \left(\frac{r^*}{\eta R^*(\zeta^*)} \right)^2 \right], \quad (27)$$

and vertical velocity,

$$\Omega^*(r^*, \zeta^*) = \Omega_{cl}^*(\zeta^*) \exp \left[- \left(\frac{r^*}{R^*(\zeta^*)} \right)^2 \right] \quad (28)$$

where Θ_{cl}^* is an empirically derived non-dimensionalised centreline temperature, Ω_{cl}^* is the empirically derived non-dimensional vertical velocity scaled by the characteristic velocity w_c (Eqn 26) (Baum and McCaffrey 1989), $\eta = 0.866$ is the empirically determined ratio of the thermal to momentum radii of the plume (Baum and McCaffrey 1989), $R^* = R/L_c$ is the empirically fitted non-dimensionalised plume radius, $r^* = r/L_c$ is the non-dimensional radial distance from the plume centreline, and $\zeta^* = z/L_c$ is the non-dimensional vertical distance from the plume source. For a single axisymmetric plume, the origin is at ground level; however, this value is adjusted to a virtual origin for merged plumes (see 'Merging plumes' section below).

To calculate the external flow field due to the plume, Baum and McCaffrey decompose the velocity field into potential and solenoidal flow components. Physically, the potential flow is due to the volumetric expansion of hot combustion gases while the solenoidal flow accounts for the flow induced by vorticity. Using the characteristic length and velocity, the centre-line potential can be expressed as

$$\Lambda_{cl} \equiv w_c L_c = \left(\frac{g^{1/3} \dot{Q}_0}{\rho_0 c_p T_0} \right)^{3/5}, \quad (29)$$

the centre-line vorticity as

$$\omega_{cl} \equiv \frac{w_c}{L_c} = \left(\frac{g^3 \rho_0 c_p T_0}{\dot{Q}_0} \right)^{1/5}, \quad (30)$$

and a centre-line pseudo streamfunction as

$$\Psi_{cl} \equiv w_c L_c^2 = \frac{\dot{Q}_0}{\rho_0 c_p T_0}. \quad (31)$$

In cylindrical coordinates with the assumptions of an axisymmetric plume and a Gaussian distribution for the temperature (Eqn 27) and vertical velocity (Eqn 28), the non-dimensional radial (\bar{v}_r^*) and vertical (\bar{v}_z^*) velocities can be determined from

$$\bar{v}_r^*(r^*, \zeta^*) = \frac{\partial \Lambda^*}{\partial r^*} - \frac{1}{r^*} \frac{\partial \Psi^*}{\partial r^*} \text{ and} \quad (32)$$

$$\bar{v}_z^*(r^*, \zeta^*) = \frac{\partial \Lambda^*}{\partial \zeta^*} + \frac{1}{r^*} \frac{\partial \Psi^*}{\partial r^*} \quad (33)$$

subject to the boundary conditions for Λ^* given by

$$\begin{aligned} \frac{\partial \Lambda^*(0, \zeta^*)}{\partial r^*} &= 0; \quad \Lambda^*(30, \zeta^*) = \Lambda_{\text{asymptote}}^*(30, \zeta^*); \\ \frac{\partial \Lambda^*(r^*, 0)}{\partial \zeta^*} &= 0; \quad \Lambda^*(r^*, 60) = \Lambda_{\text{asymptote}}^*(r^*, 60); \end{aligned} \quad (34)$$

that enforce that flow cannot cross the centreline axis ($r^* = 0$) or the ground ($\zeta^* = 0$). For the far field potential, the expansion velocity is governed by the asymptotic value $\Lambda_{\text{asymptote}}^*$. This is Green's function for potential due to a point source,

$$\Lambda_{\text{asymptote}}^*(r^*, \zeta^*) = \frac{-1}{2\pi\sqrt{r^{*2} + \zeta^{*2}}}, \quad (35)$$

evaluated at $r^* = 30$ and $\zeta^* = 60$. The values of r^* and ζ^* for the asymptotic solution correspond to a smooth transition from Eqns 32, 33 to the solution of Eqn 35. The streamfunction component of Eqns 32, 33 is calculated using the definition of vorticity,

$$r^* \frac{\partial}{\partial r^*} \left(\frac{1}{r^*} \frac{\partial \Psi^*}{\partial r^*} \right) + \frac{\partial^2 \Psi^*}{\partial \zeta^{*2}} = -r^* \omega_\phi^*(r^*, \zeta^*), \quad (36)$$

where ω_ϕ^* is found by taking the curl of the empirically derived vertical velocity (Eqn 28),

$$\omega_\phi^*(r^*, \zeta^*) = \frac{2r^* \vec{V}_{cl}^*(\zeta^*)}{R^{*2}(\zeta^*)} \exp\left(-\left[\frac{r^*}{R^*(\zeta^*)}\right]^2\right). \quad (37)$$

The boundary conditions for Ψ^* are

$$\begin{aligned} \Psi^*(0, \zeta^*) &= 0; \quad \Psi^*(30, \zeta^*) = \Psi_{\text{asymptote}}^*(30, \zeta^*); \\ \Psi^*(r^*, 0) &= 0; \quad \Psi^*(r^*, 60) = \Psi_{\text{asymptote}}^*(r^*, 60); \end{aligned} \quad (38)$$

enforcing the fact that the plume centreline and ground form streamlines that mass may not cross. The form of $\Psi_{\text{asymptote}}^*$ is fairly complex and described fully in Baum and McCaffrey (1989) and Trelles (1995).

The non-dimensional velocity field due to a single fire is computed as a superposition of the radial (Eqn 32) and vertical flow (Eqn 33). The result of this superposition is then scaled by the surface heat flux (Eqn 26) provided by the Balbi model (Eqn 18) for a fully burning cell, or scaled by the percentage area for a partially burning cell. This is subsequently used as a parameterisation in QES-Winds (i.e. the $u_{i,j,k}^0$, $v_{i,j,k}^0$, and $w_{i,j,k}^0$ fields) to calculate the final total velocity field as described in the 'Microscale winds' section above. For multiple distinct fires, Trelles (1995) shows that the velocity fields for individual fires can be superimposed. In QES-Fire, the superposition of multiple unconnected plumes is performed prior to the final mass conservation step (see 'Microscale winds'). When multiple distinct fires are spatially close, Trelles (1995) observed that the velocities may only be superimposed up to a height at which the plumes width begin to overlap. Above this height, the superimposed vertical velocities will overestimate the vertical velocity. Trelles (1995) gives an overview of how this may be accomplished at the asymptotic extent of the plumes; however, a new methodology is needed for merging plumes below the far field behaviour. QES-Fire's methodology for this is described in the next section.

Merging plumes

Many studies have examined the dependence of plume merging on source buoyancy and momentum fluxes (e.g. Gebhart and Pera 1971; Davidson and Slawson 1982; Davidson 1989; Macdonald et al. 2002). Kaye and Linden (2004) is the first detailed one focused on merging turbulent axisymmetric plumes. Of particular interest for QES-Fire is the development of a criterion for when turbulent plumes merge. Prior models use the methodology that when two plumes have grown laterally enough to overlap, the height where the overlap occurs is defined as the merging height (e.g. Bjørn and Nielsen 1995). This simplistic view does not take into account mutual entrainment (Kaye and Linden 2004). Finney and McAllister (2011) show that mutual entrainment from spot fires intensifies the burning and heat release rates of fires, and therefore should be included.

QES-Fire includes the impact of mutual plume entrainment using the assumption that in the absence of mean wind, individual plumes are passively advected owing to the entrainment field of other plumes in the domain, and that for equal-strength plumes, plumes will be deflected towards each other equally (Kaye and Linden 2004). In the QES-Fire formulation, the plume-merging height is calculated based on a non-dimensional mixing height μ_m defined by

$$\mu_m = \frac{\zeta}{x_0}, \quad (39)$$

where x_0 is the separation distance between plume sources. μ_m is then calculated using

$$\mu_m = \frac{1}{\epsilon} \sqrt{\frac{25}{132}}, \quad (40)$$

where $\epsilon = 0.09$ is an entrainment constant and $\sqrt{\frac{25}{132}}$ is the upper bound for the height when two non-interacting Gaussian plumes merge together (Kaye and Linden 2004). After the merging height is calculated, a new virtual origin for the merged plume is determined by (Kaye and Linden 2004)

$$\zeta_v = 0.91 \sqrt{2} \mu_m x_0 - \mu_m x_0, \quad (41)$$

where ζ_v is the distance below the merged plume sources.

For the case of unequal plumes, Eqn 40 can be modified to include the effects of the relative strengths of the plume sources as

$$\mu_m = \frac{1}{\epsilon} \sqrt{\frac{5}{6}} \{6/[5(R/x)^2] + \kappa^{1/3} + \kappa^{-1/3}\}^{-1/2}, \quad (42)$$

where $\kappa \leq 1$ is the ratio of the plume buoyancy fluxes. For $\kappa = 1$, Eqn 42 reduces to Eqn 40. μ_m has minimal variation for a wide range of κ values, approximately 25% for $0.15 \leq \kappa \leq 1$ (Kaye and Linden 2004) and, therefore, QES-fire assumes the equal plume strength value for μ_m for all plume merging. This is justified based on the small grid spacing typically used in QES-Fire, which limits the variation in heat release per grid cell (Eqn 18) to approximately 30% using Anderson's (1982) fuel model.

The discrete nature of the fire map (see 'Fire map' section above) means that nearly all simulations will require the merging of more than two plumes. This is accomplished through a multiscale aggregation operation applied to the surface heat fluxes and subsequently through the plume parameterisation (see 'Velocity field from axisymmetric plume') to the plume velocities. The goal is to merge adjacent plumes at the appropriate vertical position. This height is calculated by modifying Eqn 39 to yield

$$\zeta_r = \mu_m \times (\Gamma \times x_0), \quad (43)$$

where $x_0 = \Delta x$ for the fire map and Γ is the filter index making the product $\Gamma \times x_0$ the plume separation distance for filter level Γ . For the ground level to the first mixing height, $\Gamma = 1$ and each grid cell in the fire map is considered a distinct plume source. The velocity field from each grid cell is then calculated and superimposed following the 'Velocity field from axisymmetric plume' section starting at the ground level and up to the mixing-height level ζ_1 . The filter is then enlarged to merge two adjacent plume sources whose separation distance is $\Gamma \times x_0 = 2\Delta x$. The next merging height ζ_2 and new virtual origin $\zeta_{v,2}$ are then computed and the velocity field from the two merged plumes is then calculated as a single source. The new source is scaled by the total heat release of the merged plumes found by summing adjacent fire map cells. The velocity field is then superimposed for each merged plume in the vertical from ζ_1 to ζ_2 . This continues until the top of QES-Fire's domain is reached or until all plumes are merged into a single source. A schematic view of this plume-merging process is given in Fig. 4.

The plume-merging methodology serves two critical purposes in QES-Fire. First, it allows for the superposition of distinct plumes parameterised with the axisymmetric plume model described in the 'Velocity field from axisymmetric plume' section in a physically realistic manner and second, it

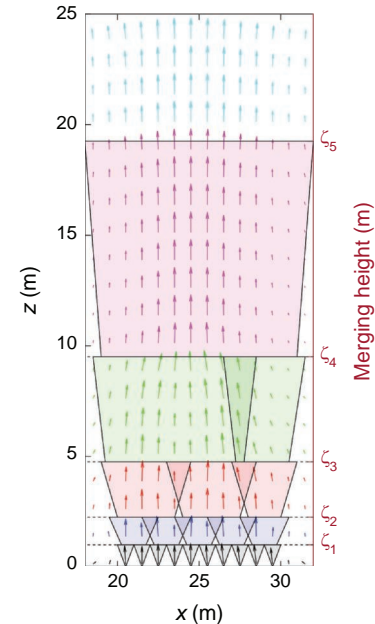


Fig. 4. Schematic representation of plume merging from distributed sources in QES-Fire. Vectors indicate the velocity field calculated by the plume parameterisation. From ground level to ζ_1 , each cell is considered a single plume source. At ζ_1 , QES-Fire merges adjacent plumes into a single plume with a new virtual origin. The velocity field is calculated for the new merged plumes from ζ_1 to ζ_2 , at which point two adjacent merged plumes are merged again. For each merging height, a new velocity field is calculated until the top of the domain is reached or all plumes have been merged. Dark shaded regions are areas of plume overlap.

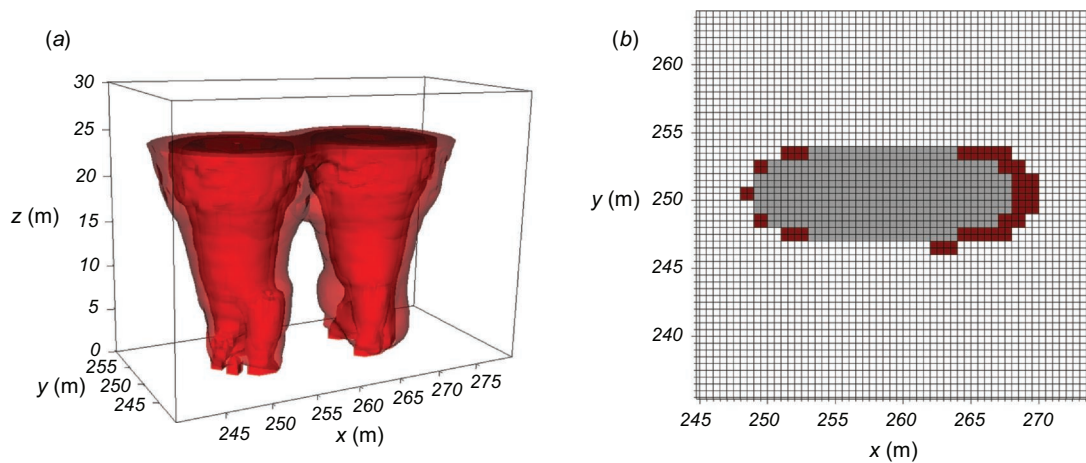


Fig. 5. (a) Vertical velocity isosurfaces from QES-Fire of merged plumes from distinct heat flux sources. (b) Plan view of QES-Fire 2-D map, red represents fully burning cells, dark grey are burned cells. Heat release per fully burning grid cell is 2.1 kW m^{-3} with a maximum vertical velocity of 5.4 m s^{-1} . Cell dimensions are $1 \times 1 \times 0.25 \text{ m}$ and background winds are 5 m s^{-1} along the positive x-axis.

gives QES-Fire the ability to have a cell-based fire map that is computationally efficient. The total merged plume from a fire front made of several sources is shown in Fig. 5.

QES-Fire comparison

Examination of QES-Fire's performance is accomplished in two phases. First, the plume structure parameterisations (independent of the ROS) are compared with high-resolution unsteady plume resolving large-eddy simulations (LESs) using WRF-SFIRE (Mandel et al. 2011, 2014). Next, the FireFlux II field campaign (Clements et al. 2019) is used to validate QES-Fire's ROS and compare its calculated wind velocities with *in situ* velocity measurements taken during an experimental burn.

Idealised plume validation

High-resolution LES runs of WRF-SFIRE were performed to examine the quality of QES-Fire's plume parameterisations. Two WRF-SFIRE simulations were performed, a 'large' burn area case and a 'small' burn area case. Both simulations used a $3 \text{ km} \times 3 \text{ km} \times 3 \text{ km}$ atmospheric domain discretised with $120 \times 120 \times 120$ grid points, resulting in a uniform grid spacing of $\Delta x = \Delta y = \Delta z = 25 \text{ m}$. The burn area was modelled using the SFIRE land surface module run on a grid one fifth the size of the atmospheric grid with a grid spacing of $\Delta x = \Delta y = \Delta z = 5 \text{ m}$. The small burn area was a $25 \text{ m} \times 25 \text{ m}$ square and the large burn area was $50 \text{ m} \times 50 \text{ m}$, both with constant heat fluxes of 72.14 kW m^{-2} . The heat fluxes were injected into the atmospheric model over an extinction depth (Mandel et al. 2011). For the small burn area, the depth was specified as 40 m , and for the large burn area, the depth was specified as 55 m . The extinction depth was

chosen based on the McCaffrey (1983) temperature profile for a flame with a comparable heat flux. The simulations were run for 20 min each with zero ambient winds and neutral background stratification. Prognostic variables were output every 20 s. After a 15-min spin-up, output data from the final 5 min of the WRF-SFIRE simulations were time-averaged to calculate plume statistics for comparison with QES-Fire's steady-state fields. Tests with different spin-up times and grid resolutions resulted in negligible differences in presented statistics. The computation time for each of the WRF-SFIRE simulations was 3300 s on 32 Intel XeonSP Skylake cores. QES-Fire was run for the same two test cases with the same grid resolution ($\Delta x = \Delta y = \Delta z = 25 \text{ m}$) and took 12 s on a single Intel XeonSP Skylake core with an NVIDIA® Titan V GPU.

Fig. 6 depicts the time-averaged vertical velocity through the axisymmetric plume centre for both WRF-SFIRE and QES-Fire. Near the ground, QES-Fire has a broader area of strong upward vertical velocity compared with WRF-SFIRE. This is due to QES-Fire's use of individual plumes at each surface grid cell, all of which contribute to the near-ground vertical velocity. The wider plume base in QES-Fire results in a wider plume at all vertical levels and translates into a shorter core-region of high vertical velocity (e.g. velocities greater than 5 m s^{-1}).

Although the core region of high vertical velocity is shorter, QES-Fire's plume extends to a higher level in the vertical direction when all positive vertical velocity values are considered. QES-Fire does not account for atmospheric stability. As a result, the buoyant plume in QES-Fire is only weakly confined in the vertical direction. LES models such as WRF-SFIRE that solve the momentum and energy equations limit the extent of vertical transport through the temperature field. The only limitation on the vertical extent

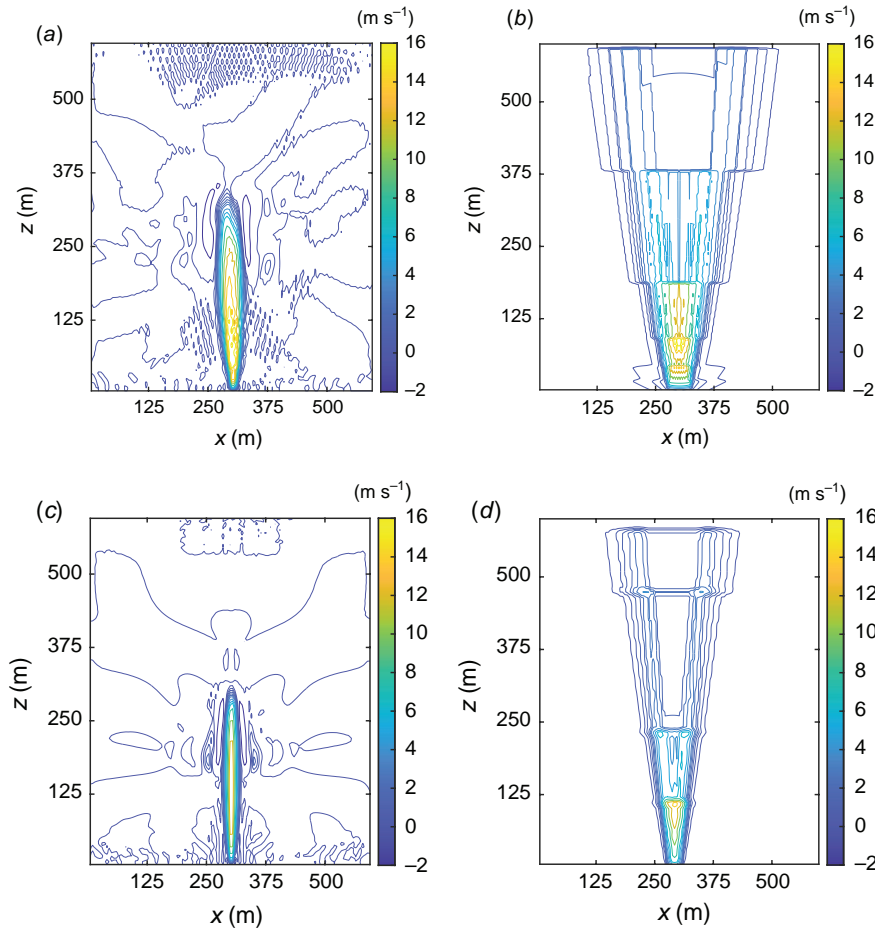


Fig. 6. Mid-plume vertical velocity contours of idealised plume. The source is a square burn area with neutral background stratification, zero winds and a constant heat flux of 72.14 kW m^{-2} specified. Large burn area = $50 \text{ m} \times 50 \text{ m}$, (a) WRF-SFIRE and (b) QES-Fire. Small burn area = $25 \text{ m} \times 25 \text{ m}$, (c) WRF-SFIRE and (d) QES-Fire.

of the plume rise in QES-Fire is from the Baum and McCaffrey (1989) plume model, particularly the asymptotic behaviour of the model at $\zeta^* = 60$ ($\zeta = 460 \text{ m}$ for the large and $\zeta = 262 \text{ m}$ for the small heat flux plots).

In addition to the structure of the vertical plume winds, it is of interest to understand the performance of QES-Fire's horizontal winds compared with WRF-SFIRE's owing to their key role in coupling the fire front–atmospheric winds to the ROS model in QES-Fire. The horizontal winds were examined at heights for which the Baum and McCaffrey (1989) model predicts the equivalent for a mid-flame region for the calculated heat flux: 10 m for the large burn area and 7 m for the small burn area (Fig. 7). Vectors in the figure have nearly identical patterns and a similar entrainment velocity at the heat flux perimeter for both models. For the large burn area, the average horizontal velocity across the heat flux perimeter for the WRF-SFIRE idealised plume is 4.75 m s^{-1} , vs 5.34 m s^{-1} for QES-Fire. For the small burn area, the average horizontal velocity across the heat flux perimeter for the WRF-SFIRE idealised plume is 1.92 m s^{-1} vs 2.11 m s^{-1} for QES-Fire.

A clear difference between QES-Fire and WRF-SFIRE plumes is the vertical and horizontal extent of the plumes they create. Still, the centreline vertical velocity values of

the two models agree well near the ground (Fig. 8). The magnitudes of the centreline vertical velocities agree but the profile shapes have two clear differences. The first is the distinct steps in the QES-Fire profile. These steps correspond to the plume mixing height levels and are a direct result of the merging plumes modelling methodology (see 'Merging plumes' above). The second difference is also evident from the vertical velocity contour plots (Fig. 6). The QES-Fire centreline vertical velocity decreases significantly at a lower vertical level than in WRF-SFIRE (200 m vs 250 m for the large burn area, and 110 m vs 250 m for the small burn area) but then continues with a small value to a much greater height.

A direct result of QES-Fire's broader plume but similar vertical velocity values is an increase in the integrated buoyancy flux compared with WRF-SFIRE at the surface (Fig. 8). However, as we move away from the ground, the total buoyancy fluxes through the x – y plane at a given height are comparable between the two models. Overall, the domain-integrated buoyancy flux for the plume from the large burn area is $1.36 \times 10^6 \text{ m}^4 \text{ s}^{-3}$ in WRF-SFIRE whereas in QES-Fire it is $1.61 \times 10^6 \text{ m}^4 \text{ s}^{-3}$, a relative difference of less than 15%. For the small burn area, the domain integrated buoyancy flux is $1.58 \times 10^5 \text{ m}^4 \text{ s}^{-3}$ in

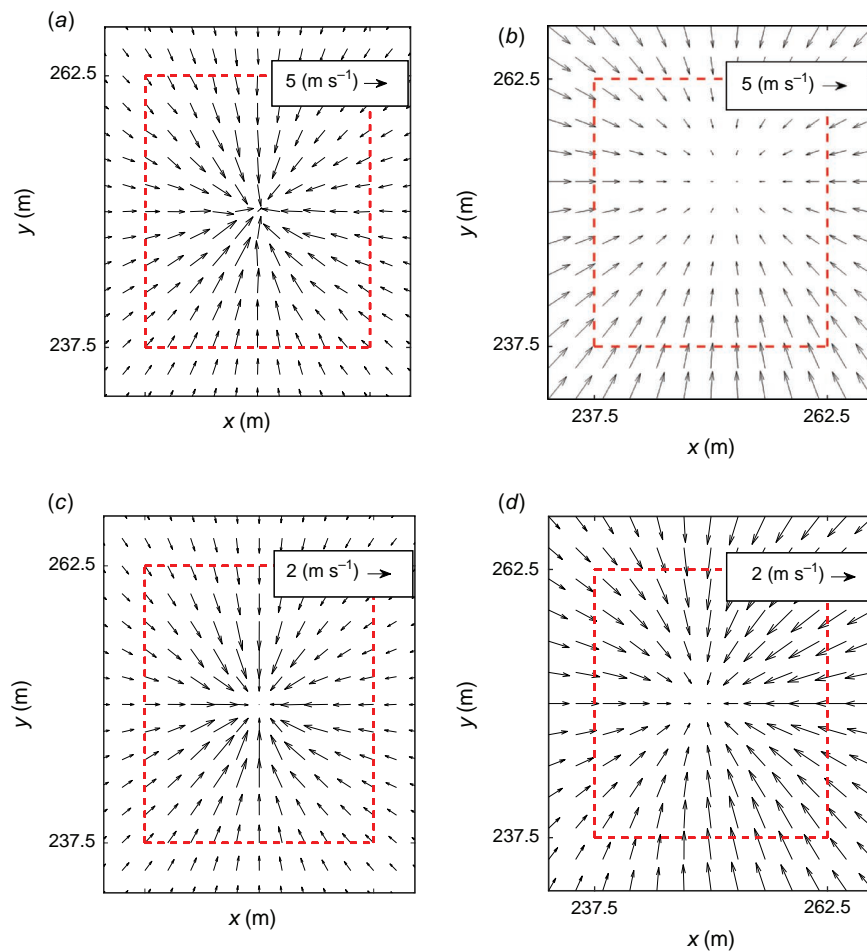


Fig. 7. Horizontal winds from idealised plume simulations. The source is a square burn area (noted by --), with neutral background stratification, zero winds, and a constant heat flux of 72.14 kW m^{-2} specified: large burn area = $50 \text{ m} \times 50 \text{ m}$ at 10-m height, (a) WRF-SFIRE and (b) QES-Fire. Small burn area = $25 \text{ m} \times 25 \text{ m}$ at 7-m height, (c) WRF-SFIRE and (d) QES-Fire. Legends show length of a reference value velocity vector.

WRF-SFIRE whereas in QES-Fire it is $1.89 \times 10^5 \text{ m}^4 \text{ s}^{-3}$, a relative difference of less than 17%.

To examine the grid resolution dependence of the QES plume-merging methodology, the large burn area was simulated with QES-Fire for a range of grid cell sizes. For a cell the size of the burn area, there is no merging and the Baum and McCaffrey (1989) plume model (see ‘Velocity field from axisymmetric plume’) holds exactly. As the burn area is broken into multiple QES cells, the plumes should merge in the vertical direction to create a single plume with a maximum velocity that is the same as the single plume case. For a $50 \text{ m} \times 50 \text{ m}$ burn area, the maximum vertical velocity from a single plume was calculated as 19.53 m s^{-1} . The maximum velocities for $\Delta x = \Delta y = 2\text{--}50 \text{ m}$ varied between 19.53 and 20.72 m s^{-1} , a maximum relative difference of 7%, demonstrating QES-Fire’s ability to represent the vertical velocity structure for a single plume using its discrete surface representation (see ‘Merging plumes’).

Experimental burn validation

The FireFlux II (FFII) experiment provides a valuable dataset of *in situ* measurements during a prescribed burn. FFII was

conducted in a Texas grass plot over flat land where topography plays a minimal role in the fire spread. The site is described in detail in Clements *et al.* (2007), and the experimental setup for FFII in Clements *et al.* (2019) (see Fig. 9 for a schematic). Here, we briefly review the most salient measurements used to validate QES-Fire.

The ROS calculations for the field data are made from an array of 24 HOBO[®] temperature loggers recording at 1 Hz. Each HOBO logger was equipped with an Omega 20-gauge K-type thermocouple to measure the temperature just above the soil. The HOBO loggers were arranged in a grid with 50-m spacing in the burn plot to track the fire front progression. The arrival time of the fire front was taken to be when the temperature first exceeded 400°C .

Ambient wind forcing data required to run QES-Fire are provided by the University of Houston Atmospheric Laboratory (HAL) tower located on the western edge of the burn plot. Additionally, a 43-m micrometeorology tower located in the burn plot fitted with 3-D sonic anemometers and type T thermocouples at heights of 5.8-m, 10-m and 20-m provides wind and temperature data for validation purposes. A planar fit tilt correction was applied to the sonic anemometer velocity components (Wilczak *et al.* 2001)

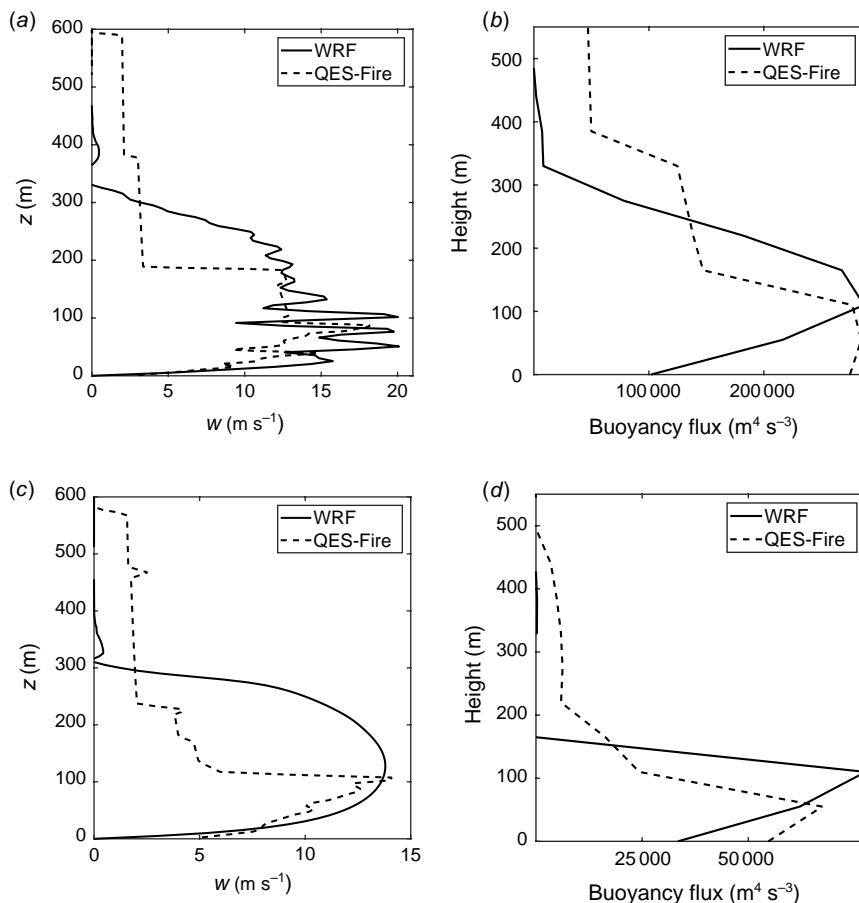


Fig. 8. Comparison of centreline vertical velocity and integrated buoyancy flux calculated by WRF-SFIRE and QES-Fire in idealised plume simulations. (a) Large burn area centreline vertical velocity; (b) large burn area x-y plane integrated buoyancy flux; (c) small burn area centreline vertical velocity; and (d) small burn area x-y plane integrated buoyancy flux.

using the 24-h period prior to the experimental burn. Wind forcing for QES-Fire was specified using the average wind-speed (10 m s^{-1}) and direction (310°) over the 30-min time period prior to the start of the experiment taken from the HAL tower sonic anemometer at a height of 31 m. Winds were then interpolated using a log-law profile with an aerodynamic roughness length (l_0) of 0.15 m corresponding to a field of uncut tall grass (Stull 2003). The horizontal turbulence strength from the HAL tower was calculated as the root mean square of the turbulent fluctuations (2.09 m s^{-1}) beginning 30 min prior to ignition through the end of the burn. The turbulence intensity was then determined to be 0.21, justifying the use of a quasi-steady-state model with a single wind forcing.

Fuels were sampled throughout the burn plot 30 min prior to the burn. Based on this sampling, the appropriate fuel type was determined to be tall grass. The fuel load was specified as 0.64 kg m^{-2} with a moisture content of 18% (Clements *et al.* 2019). The Balbi model requires the fuel load for both live and dead fuels. Because the composition of live vs dead fuel load was unknown for the FFII burn, a conservative estimate that the dead fuel load accounted for 10% of the total fuel load was specified. Sensitivity testing found a 3% relative error in the ROS for a 25% change in dead fuel load vs total fuel load. The ambient temperature

taken via a radiosonde sounding just before ignition was 16°C . The ignition line for FFII was tracked and timed based on video taken from a GoPro[®] camera mounted at 40-m on an observation tower in the northwest corner of the burn plot. This ignition sequence was read into QES-Fire, and at the specified times and locations in the burn plot, a new ignition was created to accurately describe the ignition procedure. Once the fire front is initiated and the background winds are specified, QES-Fire is used to propagate the fire front following the procedure detailed in 'QES-Fire modelling methodology' above (see Table 2 for a summary of input parameters).

One of the most critical aspects to evaluate for any wild-fire spread model is its ability to create realistic fire fronts that move at realistic ROSs. QES-Fire's calculated fire front progression for the FFII controlled burn is shown in Fig. 10. Based on qualitative comparisons with video taken from the observation tower GoPro camera, QES-Fire captures the general shape and progression pattern (i.e. isochrones) of the experiment. It is noteworthy that while the forward progression of the fire front matches field data quite well, the backing fire front ROS is overestimated. This is likely a result of the model's induced winds because the ROS model predicts negligible forcing for the level set when winds are opposed to the fire front.



Fig. 9. Map of the experimental design with instrument placement and ignition sequence for the FFII prescribed burn (Google Earth).

A more quantitative examination of the ability of QES-Fire to accurately represent the FFII fire front progression can be preformed by comparing the QES-Fire ROS with the ROS calculated from the HOBOS data loggers. This is facilitated by QES-Fire's calculation of the ROS vector at each grid cell that is on fire (see 'Fire map' section) at each model time step, allowing for a direct comparison at points matching the HOBOS data (Fig. 11). Additionally, isochrones from the QES-Fire fire front and the experimental data fire front have been calculated to enable direct comparisons of the fire front shape as it progresses through the HOBOS array.

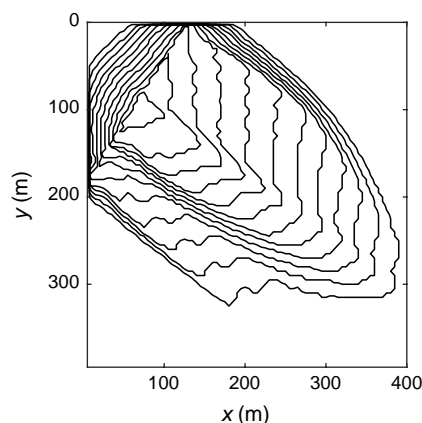
The isochrones calculated by QES-Fire have the same general shape as those calculated for FFII. The primary differences in the ROS and isochrones between the model and the experiment occurs at two HOBOS points in a region near 50–125 m east and 75–150 m north where QES-Fire predicts a rapid acceleration of the fire front that is not

observed in the data. Comparison of the ROS from FFII and QES-Fire through the HOBOS array shows a root mean squared error (RMSE) of 0.22 m s^{-1} with a mean ROS of 0.53 m s^{-1} for QES-Fire compared with 0.43 m s^{-1} from the field data. There is an overestimation of the ROS for the western flank of the fire front. This is most likely a result of the acceleration of the fire front due to the level set combined with the fact that QES-Fire does not account for changes in ambient winds present during the experiment. On the leading edge of the fire front, there is a slight underestimation of the ROS, 0.92 m s^{-1} for QES-Fire vs 0.95 m s^{-1} observed in the experiment. If the Balbi model is used without flame mid-height surface wind feedback from QES-Fire, the average ROS drops below the experimental data to a value of 0.32 m s^{-1} .

Of particular interest is the accuracy of QES-Fire's horizontal wind predictions in the vicinity of the fire front.

Table 2. QES-Fire model parameters for FFII case study.

Domain size	450 m × 800 m × 300 m
Horizontal grid spacing	$\Delta x = \Delta y = 2$ m
Vertical grid spacing	$\Delta z = 0.25$ m
Simulation time	20 min
Ignition sequence	Ignition points and times matched to observed
Inflow winds	10 m s^{-1} at 31-m height, log-law profile, $l_0 = 0.15$ m
Inflow wind direction	310°
Courant number	0.9
Fuel type	Tall grass (Rothermel 1972 Type 3)
Fuel load (total)	$\sigma = 0.64 \text{ kg m}^{-2}$
Fuel load (dead)	$\sigma_{\text{dead}} = 0.064 \text{ kg m}^{-2}$
Fuel moisture content	$m = 18\%$
Fuel depth	$d = 1.35$ m
Fuel density	$\rho_v = 1500 \text{ kg m}^{-3}$ (Rothermel 1972)
Fuel packing ratio	$\beta = \sigma/(d \times \rho_v) = 3.16 \times 10^4$
Air temperature	289 K
Slope	$\theta = 0$

**Fig. 10.** Fire front of FireFluxII experiment modelled by QES-Fire, isochrones calculated at 30 s intervals.

These winds drive the ROS and serve as one of the primary coupling mechanisms between the fire front and the atmospheric flow. To validate QES-Fire's fire-induced winds, comparisons are made with the anemometers on the 43-m tower in the FFII burn plot (Fig. 12). Because QES-Fire is a quasi-steady-state model that represents the ensemble mean value of the velocity vector and to facilitate visual comparisons, the QES-Fire horizontal winds are examined relative to the experimental horizontal winds averaged with a 3-s moving average, corresponding to the fire front residence time. The most notable feature in the horizontal winds is a

decrease as the fire front approaches the 43-m tower. The decrease is linked to the fire-induced winds causing entrainment into the flame plume. Once the fire front passes the tower, there is an increase in the velocity as the fire-induced winds add to the ambient winds. The effect is most prominently exhibited at the 6-m measurement height but persists up to the 20-m level. Both QES-Fire and the field data exhibit this behaviour, with the trend more pronounced in QES-Fire. QES-Fire's behaviour is a result of the fire line and plume parameterisation models. Additionally, the aerodynamic roughness length (l_0) is decreased as the fuel is consumed by the fire front. This decrease in l_0 corresponds to an increase in the near-surface horizontal winds calculated by QES-Winds supporting the conclusion of Sun *et al.* (2009) that a decrease in l_0 may have a significant impact on ROS. Overall, the general ability of QES-Fire to predict horizontal winds is considered to be acceptable. RMSEs of horizontal velocity for QES-Fire vs field data are 1.48 m s^{-1} at the 6-m height measurement position, 1.90 m s^{-1} at the 10-m height, and 0.69 m s^{-1} at the 20-m height.

To examine how well QES-Fire captures key features of the vertical velocity near the fire line, it is compared against the sonic anemometer data smoothed with a 3-s moving average from the 43-m tower (Fig. 13).

The clearest signature in the vertical velocity time series is an increase in upward magnitude. The increase is evident in the field data and the QES-Fire modelled winds. Away from the ground, at the 10-m and 20-m heights, the magnitude of QES-Fire's peak vertical velocity agrees reasonably well with the values in the field data. Near the ground at the 6-m measurement height, QES-Fire clearly overestimates the upward vertical velocity. The increased overestimation near the ground is consistent with our examination of QES-Fire's plume-merging parameterisation (see 'Merging plumes' above) where the distributed nature of the model near the ground results in enhanced vertical velocity. During the experimental burn, a large-scale subsidence of approximately -0.5 m s^{-1} was observed, leading to a mean negative value in the vertical velocity prior to, and after, the fire front passage at the 43-m tower. Because the subsidence is presumably not local to the burn plot and not explicitly included in the QES-Fire forcing, it is not captured by QES-Fire and explains some of the discrepancy in the comparison of vertical velocities. The arrival time of the peak vertical velocity in the QES-Fire predictions is also observed to lag the field observation's arrival at all heights. The delay agrees with the slightly delayed fire line progression (Fig. 11) through the HOB0 array. To a lesser extent, a second likely explanation for the delayed timing is the lack of direct inclusion of cross-wind advection in the QES-Fire plume parameterisation. The FFII burn occurred under strong winds, which tilt the plume in the downwind direction resulting in earlier plume arrival times for measurement points that are higher off the ground. Although not strong, this phenomenon can be observed in the tower wind

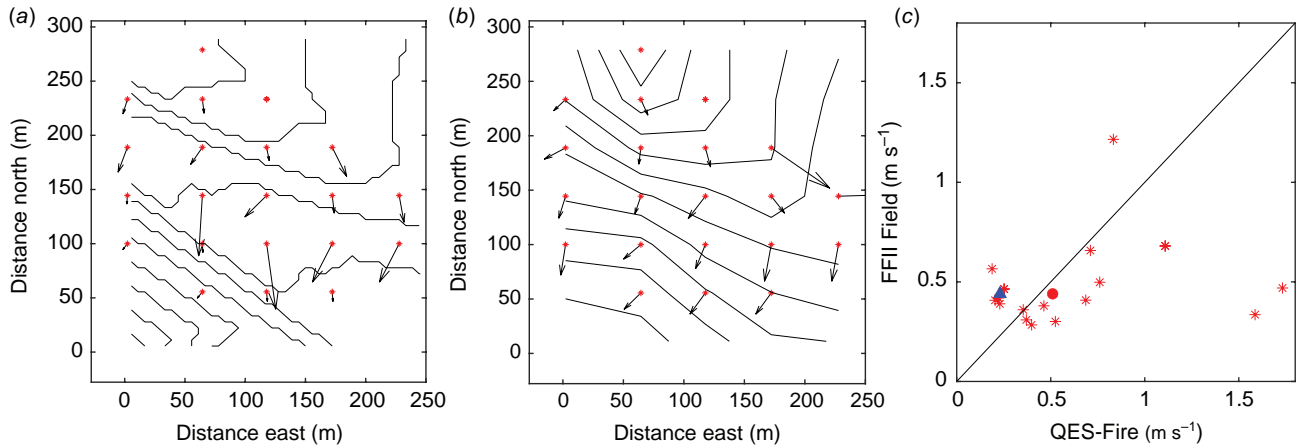


Fig. 11. Comparison of isochrones and ROS vectors through HOB0 array in FFII burn plot; distances measured from southwest corner of HOB0 array; (*) are HOB0 locations in FFII burn plot. (a) Modelled by QES-Fire, and (b) calculated from FFII field data. (c) ROS comparison for QES-Fire calculated vs FFII field data for HOB0 locations (*). Mean ROS: QES-Fire vs FFII (●). Mean ROS: Balbi model with no winds feedback vs FFII (▲).

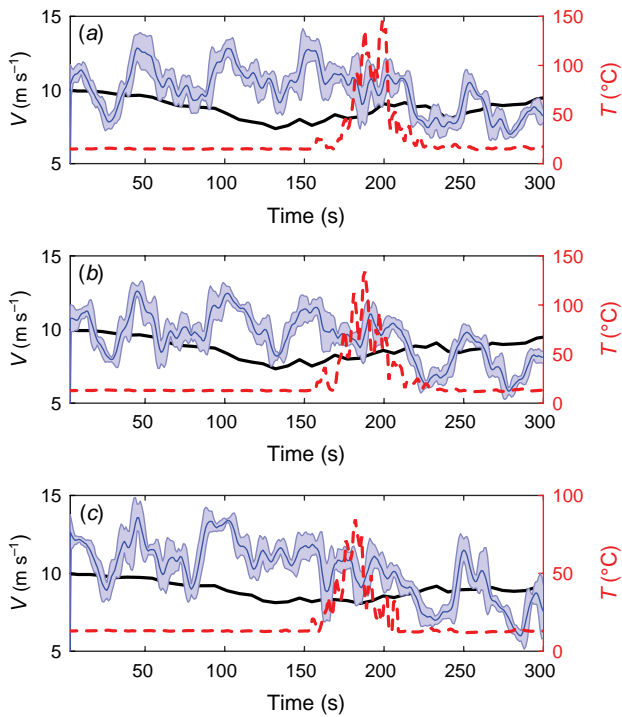


Fig. 12. Horizontal velocity (V) at the 43-m tower location during the FireFlux II controlled burn. The blue line (—) indicates the 3-s filtered sonic anemometer data, the solid (—) black line the QES-Fire horizontal velocity, and the shaded region denotes ±1 s.d. from the field data over the 3-s moving average. The dashed red (---) line is the temperature measured with a type T thermocouple, shown to indicate the fire arrival time: (a) 6-m height; (b) 10-m height; (c) 20-m height.

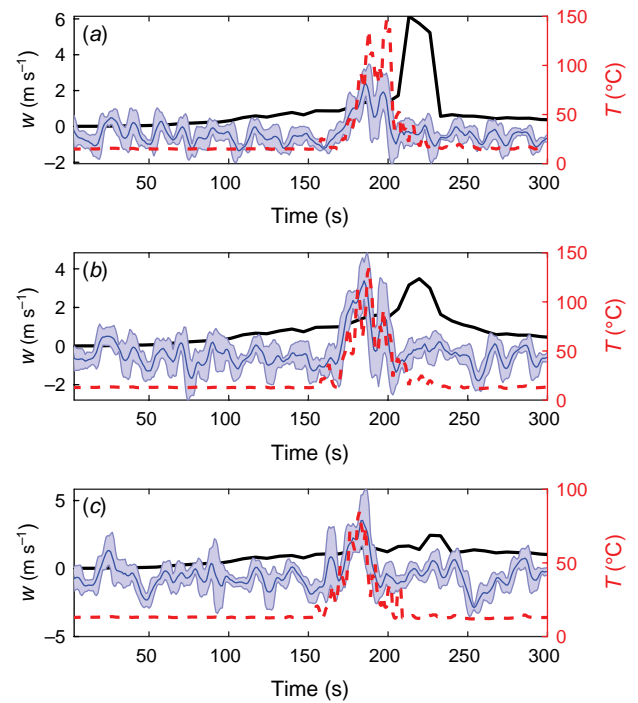


Fig. 13. Vertical velocity (w) at the 43-m tower location during the FireFlux II controlled burn. The blue line (—) indicates the 3-s filtered sonic anemometer data, the solid (—) black line the QES-Fire vertical velocity, and the shaded region denotes ±1 s.d. from the field data over the 3-s moving average. The dashed red (---) line is the temperature measured with a type T thermocouple, shown to indicate the fire arrival time: (a) 6-m height, (b) 10-m height, (c) 2-m height.

and temperature fields (Figs 12, 13). In addition to altering the plume arrival time, the lack of a specific cross-wind parameterisation in QES-Fire likely contributes to the overestimation of the maximum vertical winds in the plume,

especially for the higher measurement heights. While not as good as the horizontal velocity predictions, the vertical velocity predicted by QES-Fire is also acceptable considering the complexity of the simulated process. Vertical velocity

RMSEs for QES-Fire vs field data are 1.15 m s^{-1} at the 6-m height measurement position, 1.00 m s^{-1} at the 10-m height, and 1.22 m s^{-1} at the 20-m height.

Summary

This work presents a fast-response microscale wildfire progression model, QES-Fire. It incorporates three simplified physics quasi-steady-state models: a diagnostic wind solver (QES-Winds), a surface fire ROS model (the Balbi model), and a buoyant plume rise model that includes merging plumes. QES-Fire dynamically couples the background wind velocities to the fire-induced winds through the plume rise model and wind solver, and incorporates this new velocity field as a direct driver to the ROS model, providing high-resolution inputs for calculation of ROS and a highly realistic wind field in the vicinity of the fire front.

QES-Fire's fire-induced winds are compared against high-resolution LES of idealised buoyant plumes, yielding results with relative errors for the maximum centreline vertical velocity of 10% for a large burn area and 6% for a small burn area at a fraction of the computational cost. In addition, the buoyant fluxes have a relative error of 15% for the large burn plot and 17% for the small burn plot. To challenge QES-Fire to a more realistic scenario, ROS and wind velocities, both vertical and horizontal, are compared against the FireFlux II field experiment data and are in good agreement with observed results for flat terrain with homogeneous fuel. Backing fire front ROS is overestimated and requires further investigation.

Additional model development for QES-Fire is needed to extend its use to more general cases. This includes the effects of terrain on the coupled feedback with the winds, both the effects the terrain has on the local winds and the impact of terrain on the plume parameterisations. The general impact of ambient winds on the fire plume should also be explored, especially how coupling with terrain-induced winds can lead to extreme fire behaviours (Sharples *et al.* 2012). The overall goal of QES-Fire is to provide rapid fire evolution forecasts that account for fire-atmosphere coupling at the microscale. These forecasts could be used to support fire operations and prescribed burns or used as a research tool to explore coupling processes. QES-Fire's faster-than-real-time calculations using quite modest computational resources allows examination of how changes in variables (e.g. wind speed, wind direction, fuel moisture) affect the ROS. Ensemble runs of QES-Fire can help in assessing the risk fire fronts pose due to changes in coupled fire-atmospheric conditions.

Future work will also take advantage of other aspects of the QES framework including its Lagrangian particle dispersion model (Singh *et al.* 2011), which may be used for firebrand and smoke transport. In addition, QES-Winds

was primarily developed for and validated in urban environments with resolved buildings (Bozorgmehr *et al.* 2021). Hence QES-Fire is well positioned to be adapted for fires at the wildland-urban interface.

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Data availability. Data sharing is not applicable as no new data were generated or analysed during this study.

Conflicts of interest. The authors declare no conflicts of interest.

Declaration of funding. This research was supported by the National Science Foundation under grant PREEVENTS I664175, the University of California Office of the President award LFR-20-651032, USDA NIFA Hatch project no. 1013396, NASA grant 80NSSC19K1091, and NSF grants DEB-2039552.

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