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## An Implementation of the Rothermel Fire Spread Model in the R Programming Language

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*Published version:*

DOI:10.1007/s10694-014-0405-6

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Article Sub-Title		
Article CopyRight	Springer Science+Business Media New York (This will be the copyright line in the final PDF)	
Journal Name	Fire Technology	
Corresponding Author	Family Name	<b>Vacchiano</b>
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	Division	DISAFA
	Organization	Universita degli Studi di Torino
	Address	Via Da Vinci 44, 10095, Grugliasco, TO, Italy
	Email	d.ascoli@unito.it
Schedule	Received	11 February 2014
	Revised	
	Accepted	4 April 2014
Abstract	<p>This note describes an implementation of the Rothermel fire spread model in the R programming language. The main function provided, <code>ros()</code>, computes the forward rate of spread at the head of a surface fire according to Rothermel fire behavior model. Additional functions are described to illustrate the potential use and expansions of the package. The function <code>rosunc()</code> carries out uncertainty analysis of fire behavior, that has the ability of generating information-rich, probabilistic predictions, and can be coupled to spatially-explicit fire growth models using an ensemble forecasting technique. The function <code>bestFM()</code> estimates the fit of Standard Fuel Models to observed fire rate of spread, based on absolute bias and root mean square error. Advantages of the R implementation of Rothermel model include: open-source coding, cross-platform availability, high computational efficiency, and linking to other R packages to perform complex analyses on Rothermel fire predictions.</p>	
Keywords (separated by '-')	Fire behaviour - Fuel models - Fire spread - Prescribed fire - Wildfire	
Footnote Information		

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Journal: 10694  
Article: 405



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# An Implementation of the Rothermel Fire Spread Model in the R Programming Language

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Received: 11 February 2014/Accepted: 4 April 2014

**Abstract.** This note describes an implementation of the `Rothermel` fire spread model in the R programming language. The main function provided, `ros()`, computes the forward rate of spread at the head of a surface fire according to Rothermel fire behavior model. Additional functions are described to illustrate the potential use and expansions of the package. The function `rosunc()` carries out uncertainty analysis of fire behavior, that has the ability of generating information-rich, probabilistic predictions, and can be coupled to spatially-explicit fire growth models using an ensemble forecasting technique. The function `bestFM()` estimates the fit of Standard Fuel Models to observed fire rate of spread, based on absolute bias and root mean square error. Advantages of the R implementation of Rothermel model include: open-source coding, cross-platform availability, high computational efficiency, and linking to other R packages to perform complex analyses on Rothermel fire predictions.

**Keywords:** Fire behaviour, Fuel models, Fire spread, Prescribed fire, Wildfire

## 1. Introduction

Mathematical models of wildland fire behaviour have been of great importance in both fire ecology research and fire management (e.g., [6, 26, 27, 38]). Rothermel model for forward fire rate of spread (hereafter ROS) in surface fuels is one of the most widely used fire models [29].

Rothermel model has been programmed into computer code-based versions [2], and included as a fundamental part of several fire modeling software. Examples of simulators operating at the stand scale are `Behave/BehavePlus` [4, 5], and the `Fire and Fuel Extension to the Forest Vegetation Simulator` [28], both programmed in Fortran. Furthermore, Rothermel model has been included in spatially-explicit fire

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37 simulators (e.g., [1, 17, 19, 24, 25]), or as extension to proprietary (e.g. [18]) or  
 38 open-source Geographical Information Systems (e.g., the r.ros module for GRASS  
 39 GIS [42]).

40 However, these packages often operate as a black-box, i.e., are opaque to cus-  
 41 tomization of input parameters (except for those allowed by the Graphical User  
 42 Interface), model form, and cross-format analysis of model output. We identified  
 43 a need for scientists and managers to run surface fire simulations based on Roth-  
 44 ermel model within a larger, seamless workflow of pre- and post- wildfire model-  
 45 ing analyses, such as input data preparation, iterative model runs, or plotting and  
 46 statistically manipulating model results (e.g., [7, 10, 16]).

47 The aim of this Research Note is to present the Rothermel package for the R  
 48 programming language (R Core Team, 2013). The package currently resides on  
 49 the CRAN repository (URL: [cran.r-project.org/web/packages/rothermel](http://cran.r-project.org/web/packages/rothermel)). R is an  
 50 open-source programming language and statistical analysis framework that is  
 51 rapidly becoming standard in scientific research. It allows data handling (Appen-  
 52 dix 1), statistical analysis, and graphical representations, thanks to a suite of pre-  
 53 installed statistical methods, and more than 4,000 add-on packages. It functions  
 54 under all operating systems, including Windows, Linux and OSX. To date, some  
 55 fire-related packages have been developed for R (e.g., `paleofire` [21],  
 56 `fume` [34], and `fwi.fbp` [41]), but the Rothermel fire spread model has not  
 57 been ported yet.

## 58 **2. The `ros()` Function**

### 59 **2.1. Description**

60 The `ros()` function computes ROS [ $\text{m min}^{-1}$ ] and other output variables from  
 61 Rothermel model (Table 1). Rothermel model has been subject to several correc-  
 62 tions. The model implemented here includes the following changes to the original  
 63 system of equations: an updated weighting factor for reaction intensity by fuel  
 64 category [20], updated equations for mineral content, damping coefficient, reaction  
 65 velocity, weighting factor for fuel loadings, and live fuel moisture of extinction [2],  
 66 and removing the maximum wind factor limit [7].

67 Inputs required by the fire spread model are specified by the fire behavior fuel  
 68 model (hereafter: fuel model). Other inputs are related to environmental variables  
 69 such as slope steepness, midflame wind speed, and the moisture content of each  
 70 fuel category and size class (Table 1). Rothermel model is static, therefore it  
 71 assumes constant weather variables for each simulation [29].

72 The inputs and outputs of `ros()` are in metric units, but the function con-  
 73 verts all inputs to imperial units in order to apply the original coefficients of  
 74 Rothermel model. The function accepts both single values, and `data.-`  
 75 `frames` with multiple observations. If `modeltype` is set to D, a dynamic



**Table 1**  
**Input and Output Variables for the ROS ( ) Function**

Input	Units	Description
modeltype	-	S(static), D(dynamic)
w	t ha <sup>-1</sup>	A vector or data frame of fuel load for fuel classes 1-h, 10-h, 100-h, live herbs and live woody, respectively (5 values or columns; 0 if fuel class is absent)
S	m <sup>2</sup> m <sup>-3</sup>	A vector or data frame of surface-to-volume ratio for fuel classes 1-h, 10-h, 100-h, live herbs and live woody, respectively (5 values or columns; 0 if fuel class is absent)
delta	cm	A value or vector of fuel bed depth
mx.dead	%	A value or vector of dead fuel moisture of extinction
h	kJ kg <sup>-1</sup>	A vector or data frame of heat content for fuel classes 1-h, 10-h, 100-h, live herbs and live woody, respectively (5 values or columns; 0 if fuel class is absent)
m	%	A vector or data frame of percent moisture on a dry weight basis for fuel classes 1-h, 10-h, 100-h, live herbs and live woody, respectively (5 values or columns; 0 if fuel class is absent)
u	km h <sup>-1</sup>	A value or vector of midflame windspeed
slope	%	A value or vector of site slope
Output	Units	Description
m.live	%	Characteristic dead fuel moisture
m.dead	%	Characteristic live fuel moisture
mx.live	%	Live fuel moisture of extinction
mx.live	m <sup>2</sup> m <sup>-3</sup>	Characteristic (weighted) SA/V
cSAV	kg m <sup>-3</sup>	Fuel bulk density
rho	-	Packing ratio
beta	-	Relative packing ratio
rpr	kW m <sup>-2</sup>	Dead fuel reaction intensity
IR dead	kW m <sup>-2</sup>	Live fuel reaction intensity
IR live	kW m <sup>-2</sup>	Reaction intensity
IR	0-100	Wind correction factor
fw	0-1	Slope correction factor
fs	kW m <sup>-2</sup>	Numerator of Rothermel model
Heat source	kJ m <sup>-3</sup>	Denominator of Rothermel model
Heat sink	m min <sup>-1</sup>	Rate of spread
ROS		



76 fuel model will be invoked, where part of the cured herbaceous fuel is transferred  
 77 to the 1-h fuel size class, as a function of herb fuel moisture [35]. If characteristic  
 78 fuel moisture is higher than the fuel moisture of extinction, both for live and dead  
 79 fuels, the respective reaction intensity is set to zero [5]. The following two exam-  
 80 ples demonstrate the usage of `ROS()`.

## 81 2.2. Example 1

82 This example computes Rothermel equations by using a single fuel model, mois-  
 83 ture scenario, and unique slope and wind values.

```
> library(Rothermel)
> modeltype <- "D"
> w <-c (2, 1, 0.5, 3, 8)
> s <- c (5600, 358, 98, 6200, 8000)
> delta <- 50
> mx.dead <- 30
> h <- c (18622, 18622, 18622, 19500, 20000)
> m <- c (7, 8, 9, 40, 60)
> u <- 5
> slope <- 10
> ros (modeltype, w, s, delta, mx.dead, h, m, u, slope)
```

The result is a list of the following values:

- [1] Characteristic dead fuel moisture [%] 7.02
- [2] Characteristic live fuel moisture [%] 59.37
- [3] Live fuel moisture of extinction [%] 128.40
- [4] Characteristic SA/V [m<sup>2</sup> m<sup>-3</sup>] 7325.13
- [5] Bulk density [kg m<sup>-3</sup>] 2.90
- [6] Packing ratio [dimensionless] 0.01
- [7] Relative packing ratio [dimensionless] 0.93
- [8] Dead fuel Reaction intensity [kW m<sup>-2</sup>] 553.34
- [9] Live fuel Reaction intensity [kW m<sup>-2</sup>] 933.21
- [10] Reaction intensity [kW m<sup>-2</sup>] 1486.55
- [11] Wind factor [0-100] 6.75
- [12] Slope factor [0-1] 0.25
- [13] Heat source [kW m<sup>-2</sup>] 501.85
- [14] Heat sink [kJ m<sup>-3</sup>] 4682.05
- [15] ROS [m min<sup>-1</sup>] 6.43





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88 [4] Characteristic SA/V [ $\text{m}^2 \text{m}^{-3}$ ] 7325.13  
89 [5] Bulk density [ $\text{kg m}^{-3}$ ] 2.90  
90 [6] Packing ratio [dimensionless] 0.01  
91 [7] Relative packing ratio [dimensionless] 0.93  
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93 [9] Live fuel Reaction intensity [ $\text{kW m}^{-2}$ ] 933.21  
94 [10] Reaction intensity [ $\text{kW m}^{-2}$ ] 1486.55  
95 [11] Wind factor [0–100] 6.75  
96 [12] Slope factor [0–1] 0.25  
97 [13] Heat source [ $\text{kW m}^{-2}$ ] 501.85  
98 [14] Heat sink [ $\text{kJ m}^{-3}$ ] 4682.05  
99 [15] ROS [ $\text{m min}^{-1}$ ] 6.43

### 100 2.3. Example 2

101 Here we illustrate how to compute ROS using data from fire field experiments,  
102 and validate Rothermel predictions against observed rate of spread. This example  
103 uses the dataset `firexp` of the `Rothermel` R package. The dataset includes  
104 ROS measured using a microplot scale approach [36] during field fire experiments  
105 in heathland fuels (mixed grass-shrub). The experiments were carried out on flat  
106 terrain under variable fire weather [8, 39]. For each observed ROS, environmental  
107 and fuel parameters were measured before and during the fire. Some ranges in the  
108 dataset are: ROS 0.9–26.3  $\text{m min}^{-1}$ ; wind speed 0.4–7.9  $\text{km h}^{-1}$ ; 1-h fuel mois-  
109 ture 10–27%. We predict ROS using data from three Standard Fuel Models ([35])  
110 and environmental variables measured in the field, and validate it against  
111 observed values.



```

> library (Rothermel); data (firexp); data (SFM_metric)
> # Observed variables
> m <- firexp [, 18:22]
> u <- firexp [, "u"]
> slope <- firexp [, "slope"]
> obs <- firexp[, "ros"]
> # Predict ROS using Standard Fuel Models GR5, GS3 and SH7
> a = list ( )
> models = which (rownames (SFM_metric) == "GR5" |
  rownames (SFM_metric) == "GS3" |
  rownames (SFM_metric) == "SH7")

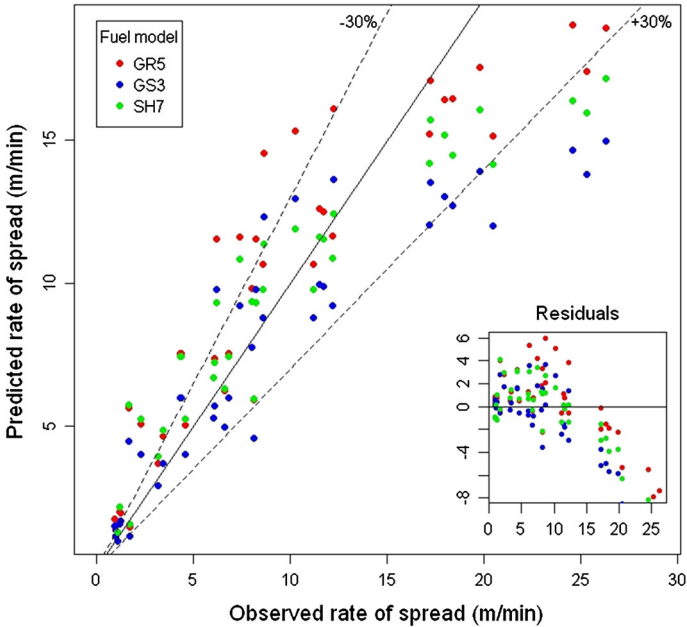
> for (i in 1 : length (models) ) {
  modeltype <- SFM_metric [models [i], 1]
  w <- SFM_metric [models [i], 2:6]
  s <- SFM_metric [models [i], 7:11]
  delta <- SFM_metric [models [i], "Fuel_Bed_Depth"]
  mx.dead <- SFM_metric [models [i], "Mx_dead"]
  h <- SFM_metric [models [i], 14:18]
  a [i] <- ros (modeltype, w, s, delta, mx.dead, h,
    m, u, slope)[15]}

> # Plot
> plot (obs, a [[1]], xlab = "Observed rate of spread (m/min)",
  ylab = "Predicted rate of spread (m/min)", col = "red",
  pch =19, xlim = c (0, 30), cex.lab = 1.1)
> points (obs, a [[2]], pch = 19, col = "green2")
> points (obs, a [[3]], pch = 19, col = "blue2")
> abline (coef = c(0, 1))
> abline (coef = c(0, 0.7), lty = 2); text (13.6, 19.2, "-30%")
> abline (coef = c(0, 1.3), lty = 2); text (28.7, 19.2, "+30%")
> legend (0, 19.2, c("GR5", "GS3", "SH7"), pch = 19,
  col = c("red", "green2", "blue2"), title = "Fuel model")

> # Inset Residual plot
> par (fig = c (.57, .98, .07, .55), new = T)
> plot (obs, a [[1]] - obs, xlab= "", ylab= "", col = "red",
  main= "Residuals", font.main = 1, pch=19, cex=.7)
> points (obs, a [[2]] - obs, pch = 19, cex = .7, col = "green2")
> points (obs, a [[3]] - obs, pch = 19, cex = .7, col = "blue2")
> abline (h = 0)
> par (fig = c (0, 1, 0, 1))

```





**Figure 1. Observed vs. Predicted ROS for the fireexp dataset using Standard Fuel Models GR5, GS3 and SH7.**

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### 3. Potential Expansion of the Package: Example of Functions

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The `ros()` function can be implemented in more complex analyses of fire behavior and effects. We illustrate below two cases of the potential development of new functions based on `ros()`. The first case is a function for uncertainty analysis of rate of spread, that implements methods already explored by the literature [9, 14, 23, 37]. The second example is a newly developed function to evaluate the fit of preset fire behavior fuel models to observed ROS (Figure 1).

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#### 3.1. The `rosunc()` Function

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Several authors have stressed the importance of introducing stochasticity in fire behavior prediction [9, 14, 23, 37]. The advantage of stochastic fire models is to obtain error bounds and probability-based outcomes for the main fire behavior parameters. Although Rothermel model is essentially deterministic, a probability density function of ROS or other model outputs can be obtained by perturbing one or more input variables (usually environmental ones). The probability associated to each output value is represented by the relative frequency of such output among all model realizations. Manually perturbing model inputs is a tedious task. The `rosunc()` function of the `Rothermel` package automatically perturbs inputs by randomly sampling from gaussian distributions, where the mean is the observed value and the standard deviation is specified by the user (in the form of

132 coefficient of variation, 0–1). The output is a vector of ROS. The function accepts  
 133 the same arguments as in `ros()`, plus the desired coefficients of variations for  
 134 wind speed, fuel moisture, slope, fuel load, and fuel bed depth, and the number of  
 135 simulations desired to produce a Monte-Carlo based probability density function  
 136 for ROS [14, 23]. Consequently, the function runs on one fuel set at a time (i.e.,  
 137 no `data.frames` allowed as input).

### 138 3.2. Example 3

139 Here, one observation (row) is selected from the `firexp` dataset. Input values  
 140 are selected similarly to `ros()`, and a coefficient of variation of 0.3 is specified  
 141 to generate a gaussian distribution of fuel moisture values. The probability distri-  
 142 bution function of ROS is generated by 1000 Monte Carlo simulations and graphi-  
 143 cally compared with the observed value. This example's output may differ from  
 144 actual results due to the stochastic simulation of moisture values.

```
> data("firexp"); varnames <- names(firexp)
> firexp <- as.numeric(firexp[5,]); names(firexp) <- varnames

> pred <- rosunc(modeltype = "D",
  w = firexp[1:5],
  s = firexp[6:10],
  delta = firexp["Fuel_Bed_Depth"],
  mx.dead = firexp["Mx_dead"],
  h = firexp[13:17],
  m = firexp[18:22],
  u = firexp["u"],
  slope = firexp["slope"],
  sdm = 0.3,
  nsim = 1000)

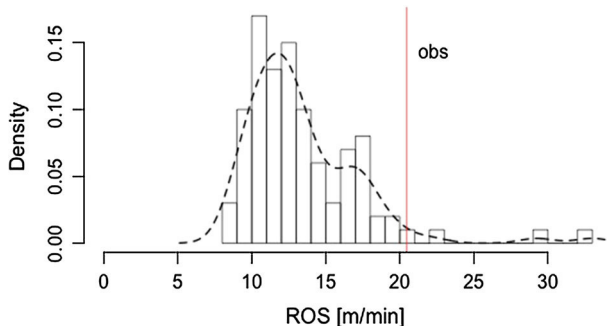
> summary(pred)
  Min. 1st Qu.  Median    Mean 3rd Qu.    Max.
  6.11  11.06   12.19   13.34   14.56   28.98
```

### 145 3.3. The `bestFM()` Function

146 A set of Standard Fuel Models (SFM) was developed to parameterize fuel proper-  
 147 ties of different fuel complexes [3, 35]. In the process of testing the predictions of  
 148 Rothermel model vs. observed ROS in a given vegetation, one of the first steps is  
 149 to verify whether any of the SFM yields a satisfactory prediction [22, 30, 35]. This  
 150 is a crucial step before undertaking the calibration of a custom fuel model [11].

151 The function `bestFM()` estimates the fit of the 53 SFM to a vector of  
 152 observed ROS, based on absolute bias (predicted - observed ROS), and root mean  
 153 square error (RMSE). Arguments of the function include environmental variables,  
 154 which are not a part of SFM, and the observed value or vector of ROS. The  
 155 function calls a dataset of SFM that has been embedded in the `Rothermel`  
 156 package (dataset `SFM_metric`), simulates ROS using SFM data and environ-





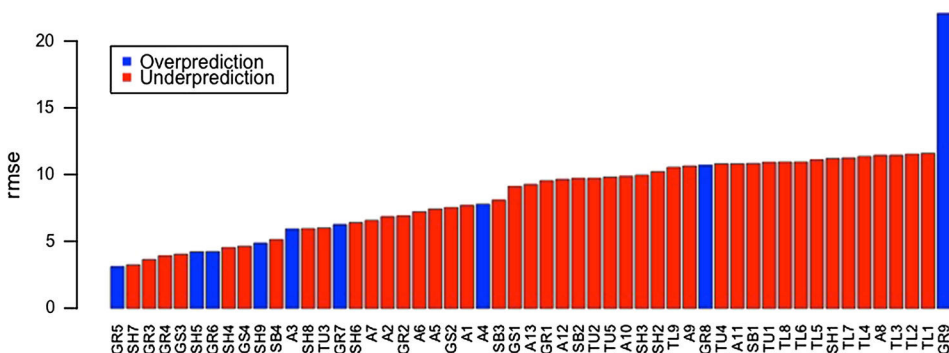
**Figure 2. Probability density function of ROS and the observed value.**

157 mental variables, and outputs a `data.frame` of RMSE and/or absolute bias.  
 158 Simulations can also be run under predefined fuel moisture scenarios [35] by call-  
 159 ing the dataset `scenarios` (Figure 2).

### 160 3.4. Example 4

161 This example loads a vector of observed ROS and environmental parameters from  
 162 the `firexp` dataset, and compares them with `ros()` predictions from a data-  
 163 set of 53 Standard Fuel Models. A sorted barplot of increasing RMSE is pro-  
 164 duced to illustrate the output of the function. The sign of prediction bias is  
 165 indicated by the bar color (Figure 3).

```
> data("firexp")
> a <- bestFM(obs = firexp[, "ros"],
             m = firexp[, 18:22],
             u = firexp[, "u"],
             slope = firexp[, "slope"])
```



**Figure 3. RMSE of 53 SFM against a dataset of observed ROS in heathland mixed grass-shrub fuels.**

## 166 4. Discussion and Practical Implications

167 The main function of the `Rothermel` package computes ROS from user-  
 168 defined (or standard) fuel and environmental parameters. The `ros()` function  
 169 computes parameters of the Rothermel model with its most common modifica-  
 170 tions [2, 7, 20]; however, the code is open to host additional formulations, such as  
 171 those by the Fuel Characteristic Classification System (FCCS) [33], or alternative  
 172 fire spread models [15].

173 The `ros()` function is functionally similar to the US Forest Service software  
 174 `BehavePlus` [5], and ROS predictions for aligned head fires are equivalent between  
 175 the two softwares. Compared to `BehavePlus`, R provides an open-source platform  
 176 that runs on multiple operating systems (Windows, OSX, Linux). However,  
 177 `ros()` lacks the additional functionality of the latter, i.e., supplementary fire  
 178 behavior and spread models, together with the user friendly interface that made  
 179 `BehavePlus` so popular among fire managers. The `ros()` function is not inten-  
 180 ded as a decision support system for fire management alternative to `BehavePlus`.  
 181 Rather, it is a new tool for fire scientists who need to carry out complex analyses  
 182 using the Rothermel model. To this regard, its objective is similar to the `Firelib C`  
 183 function library [10], that was written to give fire simulation modellers a common  
 184 programming interface to use in building fire growth applications models.

185 However, compared to existing software, the R implementation of Rothermel  
 186 model allows to perform many simulations at the same time (Example 2), plot  
 187 and export the results, and nest the computation of ROS (and of all intermediate  
 188 outputs of Rothermel model) within more complex analyses, such as `if()` state-  
 189 ments or `for()` loops, or sensitivity analysis of model output [32]. Additionally,  
 190 the R framework can generate web-based user interfaces (package `shiny` [31]),  
 191 and complex plots such as fire characteristic charts [11].

192 Much potential is associated to the newly programmed function `rosunc()`  
 193 that carries out uncertainty analysis of ROS. This method has recently been  
 194 praised for its ability to generate more information-rich, probabilistic predictions,  
 195 as compared to traditional deterministic models [23]. Furthermore, by dynamically  
 196 linking to spatially-explicit fire growth models and forest dynamics simulators at  
 197 the stand or landscape scale [13], the `rosunc()` function enables modellers to  
 198 generate probabilistic predictions of fire growth and ensemble forecasts resulting  
 199 from variable weather or fuel inputs [19].

200 Finally, the function `bestFM()` is intended as an exploratory analysis of  
 201 observed ROS in a fuel complex. RMSE from Standard Fuel Models can show  
 202 which group of models (i.e., GR, GS, SH, TU, TL, SB) have a similar fit to the  
 203 data. In Example 4, observed ROS in mixed grass-shrub heath fuels from `fi-  
 204 rep` showed increasing RMSE starting from GR, SH, GS up to TL models,  
 205 excluding GR9. Within the first 10 best fuel models, the GR group performed  
 206 slightly better than SH and GS. Our interpretation is that the herbaceous compo-  
 207 nent in heath fuels is driving the rate of successive ignitions. Consequently, when  
 208 building a custom fuel model [12] for dry heaths, particular attention should be  
 209 focused on setting the parameters of the herbaceous fuel category.



Author Proof

210 The Rothermel package is one of the first tools to support fire science in the R  
211 programming language. A wealth of packages exists for other research fields in  
212 ecology and environmental science, such as climate modelling, biodiversity, natu-  
213 ral hazard modelling, or genetics. Similarly, R has the potential to become a privi-  
214 leged platform to carry out data analysis and modelling in fire science. In fact, the  
215 R architecture is much suitable to develop tools such as decision support systems  
216 and cross-scale hierarchical models, i.e., systems of interacting simulators that  
217 take advantage of different modelling approaches (e.g., spatially-explicit fire  
218 spread, coupled physical fire models, stochastic weather generation, treatment of  
219 remotely sensed imagery...), and may effectively interact with local or remote data  
220 repositories.

221 We believe that the present package nicely fits in what a recent overview of the  
222 most up-to-date fire simulator pointed out [5]: 'Care must be taken to avoid black  
223 box modelling and to avoid use of default values. (...) A rebuild of the code from  
224 the bottom up [is desired] to facilitate integration of fire behaviour, fire effects and  
225 fire danger rating systems, as well as point and spatial systems'. Additional contri-  
226 butions to the package are welcome, and will implement complementary functions  
227 to enrich the range of fire modeling tools able to exploit the potential of the  
228 Rothermel model within the R statistical framework.

## 229 **Acknowledgments**

230 We would like to thank the CRAN staff for useful support and testing of the  
231 package.  
232

## 233 **Appendix 1: A Primer on the R Language**

234 A complete introduction to the R language goes beyond the scope of this paper.  
235 We will briefly illustrate the meaning of some key terms in order for the reader to  
236 understand the examples and data structures referenced in this paper. For an  
237 introduction to the R language, tutorials and working examples, refer e.g. to 'An  
238 introduction to R' [40], from which this section is borrowed, and to the documen-  
239 tation available on the CRAN website (URL: <http://cran.r-project.org>).

240 The user operates R via commands entered at the prompt '>'. Elementary  
241 commands consist of either expressions or assignments. Expressions are evaluated,  
242 printed (unless specifically made invisible), and the value is lost. An assignment  
243 evaluates an expression and passes the value to an object stored in a 'workspace'  
244 for future retrieval. The assignment operator is '<-'. R commands are case sensi-  
245 tive; comments can be put almost anywhere, starting with a hashmark ('#').

246 R operates on named data structures. The simplest such structure is the vector,  
247 which is a one-dimensional entity consisting of an ordered collection of numeric  
248 or string elements. To set up a vector named `x`, say, consisting of five numbers,  
249 namely 10.4, 5.6, 3.1, 6.4 and 21.7, use the R command `x <- c(10.4, 5.6, 3.1, 6.4, 21.7)`. An R data frame is a two-dimensional entity



consisting of rows (i.e., observational units) and columns (i.e., observed variables). Vectors of the same length, for example  $x$  and  $y$ , can be concatenated to form columns in a data frame named  $df$  using the R command  $df <- cbind(x, y)$ . An R list is an object consisting of an ordered collection of other objects, be them vectors, data frames, or other R data structures. List elements are numbered and may be referred to by the subsetting operator  $[ [ ] ]$ .

Finally, functions are R objects that evaluate the result of an expression using user-defined arguments. A call to the function usually takes the form `function.name(argument1, argument2)`. The Rothermel package for R operates mainly by some newly programmed functions.

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Author Proof

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