



Discrete event heat transfer simulation of a room



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ABSTRACT

The paper presents a new proposal of a discrete event simulation (DEVS) model for heat transfer in an enclosure. To our knowledge this is the first proposal in this direction, therefore the model is not comprehensive (has no windows, no ground coupling, etcetera). It is a first step of DEVS into the heat transfer in buildings. The model presented is just one possible implementation for a single thermal zone, although it has been designed to be used in multi-zone models. Common methods like CTF (conduction transfer functions) or RTF (response transfer functions) cannot be used. Instead it employs the successive transition state formulation for the 1D conduction heat transfer through multi-layered walls. An example test room has been calculated and the results compared with software that uses well known methods (CTF, RTF). Finally the DEVS model has been tested with a random convective internal-load signal.

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1. Introduction

The origin of this paper was to study the possibility of using a discrete event simulation (DEVS) model for the calculation of the annual thermal energy needs; demand, consumption, and other related magnitudes of an enclosure. Although there are many well known and powerful software tools as; TRNSYS [1], EnergyPlus [2], Esp-r [3] to name a few, which already perform such computations, we want to start exploring DEVS. The DEVS formalism has many advantages which can be exploited within this field. In the engineering practice it is desirable to compute quickly the heat transfer in buildings to size the equipment (heat exchangers and so on). But at the same time, the heat transfer model employed previously, should be able to cope with more complex calculations, including the thermal dynamical response of the equipment in order to increase the overall efficiency. Unfortunately, the heat transfer problem becomes more difficult because the characteristic response time of the walls and the equipment is very different. In contrast with the aforementioned solutions, DEVS deals with it naturally. Moreover since DEVS calculation is driven by events the model could be used afterwards in optimal real-time control of the heat transfer processes.

Since DEVS method is not very widespread, readers not familiarized with the DEVS formalism are advised to read [4] or other books on the subject.

1.1. Motivation

In general the software tools face the simulation by making an integration which focuses on the time-axis as the leading dimension. The simulation manager keeps track of the components by looking at them from the time axis, just making a picture of everything at a time. EnergyPlus [5] and Esp-r [6] use an adaptive time step. Roughly, the first uses a fixed zone time step and chops it into smaller time steps if the quickest temperature evolution of a zone crosses a tolerance value (0.3 °C/per zone time step) then all the evolutions of all components proceed with this shorter time step inside the current zone time step. The second splits the overall matrix of the simulation into sub-matrix. They are grouped by their membership (or domains) into; the building fabric, air flow network, the HVAC system, and so on. Each domain is evaluated at its own frequency. An analogy, for the EnergyPlus case, would be like using a single stroboscope and adjusting its frequency to see everything at the speed of the quickest component, while Esp-r uses several stroboscopes tuned to keep track of each domain.

This way of thinking of integration comes from classical integration methods of differential equations which in turn were designed in times where no high computing devices were available. DEVS formulation was devised to solve the time axis view constraint. Initially it dealt with discrete state computer systems

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Nomenclature			
N_i	number of boundary conduction elements of the i -zone, equal as well to the number of its boundary surfaces	E_i	internal energy of the air in the zone i [J]
N	number of volumes or thermal zones	m_{da}	mass of dry air [kg]
ΔU_{lg}	change of internal energy of H ₂ O from liquid to vapour [J kg ⁻¹]	t	time [s]
ϵ_k	emissivity of surface k	T_{da}	dry bulb air temperature [°C]
$\sigma_{\text{Stefan-Boltzmann}}$	Stefan–Boltzmann constant 5.67×10^{-8} [W m ⁻² K ⁻⁴]	W	absolute humidity [kg _{H₂O} kg _{da} ⁻¹]
F_{kj}	view factor between surface k and j	<i>Discrete event simulation</i>	
$q_{\text{conv},j}$	convective heat flux into the zone air from surface j [W m ⁻²]	$\delta_{\text{ext}}()$	external function
$Q_{\text{conv},\text{side}}$	convective heat power at either side of a conduction element [W]	$\delta_{\text{int}}()$	internal transition function
$q_{\text{conv},\text{src}}$	convective heat flux into the zone air from sources [W m ⁻²]	$\lambda()$	output function
$q_{\text{rad},\text{src},\text{side}}$	radiation heat input to <i>side</i> of any wavelength, due to known sources [W m ⁻²]	σ	scheduled time for the next call to a certain component [s]
$q_{\text{rad}-lw,j}$	longwave radiation heat exchange which depends on superficial temperatures, at surface j [W m ⁻²]	dQ_{abs}	the minimum absolute quantum of the QSS integrator
$q_{\text{rad}-lw,\text{side},m}$	longwave radiation heat exchange at side of conduction element m [W m ⁻²]	dQ_{rel}	relative quantum of the QSS integrator
$q_{\text{rad}-lw}$	vector of longwave radiation heat exchanges within a zone [W m ⁻²]	$ta()$	time advance function
T	vector of superficial temperatures of the boundary surfaces of the zone	e	elapsed time from the last call of a certain component [s]
T_0	superficial temperature at one side of a conduction element [°C]	S	set of internal states of a model
T_1	superficial temperature at the other side of a conduction element [°C]	s	internal state of a model
T_{zone}	temperature of the zone air [°C]	X	set of input events
C_{da}	specific heat capacity of dry air [J kg ⁻¹ K ⁻¹]	x	input event
$C_{\text{H}_2\text{O}}$	specific heat capacity of water vapour [J kg ⁻¹ K ⁻¹]	Y	set of output events
		y	output event
		<i>Subscripts</i>	
		01	refers to the set formed by the magnitude at both sides
		cond	conduction
		conv	convection
		i	index for zone or volume i
		j	index for surface j
		m	index for conduction element m
		side	takes values 0 or 1 to represent each side of a 1D conduction element
		src	sources

but recently it was extended to continuous systems, like the differential algebraic systems DAEs [7].

When trying to simulate real buildings and HVAC systems, one set of problems comes from the fact that controls like thermostats or forcing functions like lighting, occupancy and equipment (i.e. the internal gains) act suddenly in an ON/OFF manner at any time. This and other types of highly non-linear behaviour causes difficulties to the smooth and “stroboscopic” classical methods in many fields. Even the structure of the differential equations can change due to the action of controls. The other set of problems comes from the fact that the time response of different parts of the model may be quite different. An evident example is the change of the conduction heat flux through walls in common constructions, compared with the response time of the heat power output from the HVAC equipment, as was pointed out before. The first may take hours while the second may take minutes. Moreover the spread of the response times of different parts of the system may lead to stiffness problems. Adaptive methods try to adapt the time step to the most rapidly changing subsystem or even try to discover when a sharp change (ON/OFF type) event is going to occur. Esp-r simulation methods resemble the most, the DEVS proposal of this article. The subdivision is done based on some features shared by the elements of that part: time response, spatial resolution level, the type of problem to be solved (algebraic linear/non-linear equations, differential equations). Upon this subdivision a coordinator algorithm tries to solve the coupling among the parts or modules (which they call onion strategy, while for the decoupled components they use

the term “ping-pong”) [8]. Anyway it drives the simulation by tracking the time axis and uses several strategies to adapt the time step; “...boundary condition look ahead (monitors user specified control variable(s) and reduces time-step value if rate of change greater than user specified value), time-step reduction by iteration (reduces time-step value until difference in control variable for current time-step and previous time-step is within user specified limit), user specified time-step value, iteration without time-step reduction, simulation rewind (rewind simulation clock to user specified start period if user specified control variable is outside user specified limit)...” [9]. In article [10], Clarke makes a resume of the iterative solution of nested domains. In the case of EnergyPlus as the Engineering Manual points out [5], the simulation is driven by the integration of a set of ODEs (ordinary differential equations). (To see a discussion about the coordination of the coupling among these domains see Refs. [11,12].)

2. General philosophy and implementation

A DEVS model is formed by coupling atomic models, which send messages to each other through ports. Any DEVS model is defined by the following set: its internal state set S , the time advance function $ta()$ (used to schedule the time elapsed until its next call, named σ), the external function $\delta_{\text{ext}}()$ which deals with the arriving external input events (note that the time elapsed e from the last call is always, $e < \sigma$) and the *internal transition* which in turn is composed of $\lambda()$ and $\delta_{\text{int}}()$ functions (the first issues an output

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