

# Virtual Reality: the Need for Bond Graphs

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## KEYWORDS

Virtual reality, bond graphs, modelling, simulation

## ABSTRACT

In this paper we endeavour to show how the need for modelling and simulation in virtual reality is connected with bond graphs (BG). In effect, both try to immerse their users in realistic virtual worlds, as in flight simulators.

## INTRODUCTION

Virtual reality is a branch of computer science that tries to represent dynamic realities as virtual worlds in which the user feels immersed [MOCELLIN 2001].

We show how this science could benefit from bond graphs and how the concepts of modelling and simulation, which are the milestones of bondgraphing, are connected with it.

## BOND GRAPHS

Bond graphs are based on the fact that the power or energy derivative of a dynamic system is composed of a number of terms. This number is equal to the number of interactions the system is capable of.

In full generality, the power  $P$  is a sum of products of intensities and extensity flows, the so called efforts and flows of BG. This sum characterizes the total power interactions in the system:

$$P = \sum_{i=1}^n e_i f_i \quad (1)$$

where  $e_i$  stands for the efforts and  $f_i$  stands for the flows.

These considerations are related to the derivation of Tellegen's Theorem where all the junctions are placed within a conceptual boundary; they are valid in all domains of science [ATLAN 1973].

Thermodynamics, for example, takes into account the fact that matter can be converted into different forms by chemical reaction [THOMA 2006-1]. It describes the energy of a region of matter having  $n$  components with equations of the form:

$$\dot{E} = T\dot{S} - p\dot{V} + \sum_{i=1}^n \mu_i \dot{m}_i \quad (2)$$

Which can be visualized as in Fig. 1 :

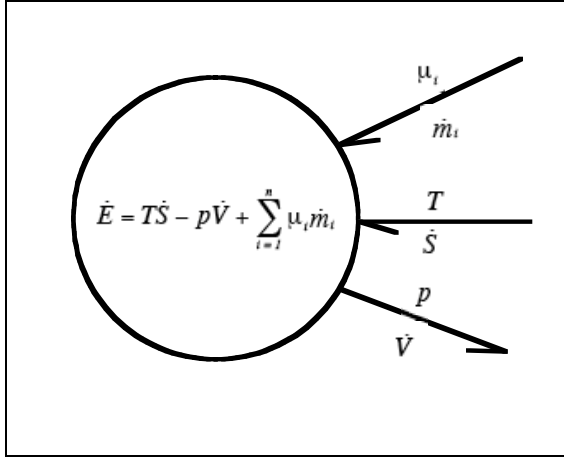


Fig 1 Equation 2 illustrated for a thermodynamic system. The power orientations correspond to the signs in equation 2.

This power equation shows the increase of energy in thermodynamic and chemical systems. It is an expression of the conservation of power, mainly chemical power.

The continuity equation for the  $i$ th extensity  $f_i$  within a control volume  $V$  under the effect of Sources and Sinks  $So_i$ , as well as Transfers  $Tr_i$ , takes the form:

$$\frac{-f_i}{-t} = So_i - Tr_i \quad (2)$$

The transfer  $Tr_i$  of an extensity  $f_i$  is due to the product of the Conductance  $Co_i$  (the so called Siemens  $S$  of German speaking countries) times the difference of intensities or efforts  $e_i$ :

$$Tr_i = Co_i (e_i - e_{i0}) \quad (3)$$

In full generality, the transfer of extensities is controlled by  $k$  intensities to take into account the cross effects among energy domains, the so called multiport  $R$  of BG:

$$Tr_i = \sum_{k=1}^{n-1} Co_{ik} (e_k - e_{k0}) \quad (4)$$

Combining the continuity equation (2) with the transfer equation (4), we obtain the integral balance equation:

$$\frac{\sum f_i}{\sum t} = So_i \sum \sum_{k=1}^{n-1} Co_{ik} (e_k - e_{k0}) \quad (5)$$

Being founded in the above transport theory, bond graphs allow the formulation of equations that describe any dynamic system. They state that the local value of the  $i$ th extensity is equal to its local density  $\sigma_i$ . By definition, the amount of a flowing extensity  $f_i$  contained in a volume equals the integral over the volume of the density  $\sigma_i$  of the extensity:

$$f_i = \int_V \sigma_i dv \quad (6)$$

The total divergence  $So_i$  generated by all the sources and sinks within a volume can similarly be characterized in terms of densities  $\sigma_i$ :

$$So_i = \int_V \sigma_i dv \quad (7)$$

The transfer  $Tr_i$  of extensive  $i$  through a surface  $F$  bounding a control volume  $V$  can be specified in terms of surface transfer densities  $t_i$ :

$$Tr_i = \oint_F t_i df \quad (8)$$

If the intensities are scalar, the nabla operator supplies the gradient. The transfer density  $\mathbf{t}_i$  of the  $i$ th extensity  $f_i$  is then stated as a sum of the products of the conductance  $Co_{ik}$  with the gradients of the individual intensities  $e_k$ :

$$\mathbf{t}_i = \sum_{k=1}^{n-1} Co_{ik} \text{grad } e_k \quad (9)$$

Introducing densities into the integral continuity equation (5) we obtain:

$$\frac{d}{dt} \int_V \rho_i dv = \int_V \rho_i dv - \int_F \mathbf{t}_i d\mathbf{f} \quad (10)$$

This balance equation of the  $i$ th extensity refers to the entire control volume  $V$ .

The second term on the right hand side of equation (10) is a surface integral which can be converted into a volume integral. The surface integral of a vector over a closed surface  $F$  equals the volume integral of the vector's divergence over the volume enclosed by the surface (sometimes called the Stokes theorem):

$$\oint_F \mathbf{t}_i d\mathbf{f} = \int_V \text{div } \mathbf{t}_i dv \quad (11)$$

The variation of an extensity contained in a given volume is equal to the volume integral of the time variation of its density. We can add it to the volume integral of the volumetric transfer  $\text{div} \mathbf{t}_i$  and subtract the volume integral of sources and sinks densities.

Interchanging differentiation and integration with respect to time, we find the volume integral:

$$\int_V \left\{ \frac{d}{dt} \int_V \rho_i + \text{div } \mathbf{t}_i - \int_V \rho_i \right\} dv = 0 \quad (12)$$

For the integral to be identically zero irrespective of the limits of integration, the integrand itself must be equal to zero, which gives the general continuity equation upon which Paynter founded bond graphs [PAYNTER 1960], that is:

$$\frac{\sigma \sigma_i}{\sigma} + \text{div } \mathbf{t}_i = \sigma_i \quad (13)$$

Replacing transfer densities  $\mathbf{t}_i$  by the general transfer equation (9), we get the equation of any dynamic system, taking into consideration the relevant extensities and intensities of the reality under consideration and the source and sink densities due to the relation to the environment:

$$\frac{\sigma \sigma_i}{dt} + \text{div} \left( \sum_{k=1}^{n-1} Co_{ik} \text{grad } e_k \right) \frac{\sigma}{\sigma} = \sigma_i \quad (14)$$

If the considered reality is itself in motion, this motion carries along the extensities because they are properties of the bulk of reality: a global convective transfer must be added.

Convective transfer densities are expressed by the product of the extensity densities  $\sigma_i$  with the velocity of displacement  $\mathbf{v}$ , giving the convective transfer density  $\sigma_i \mathbf{v}$  of the  $i$ th extensity:

$$\frac{\sigma \sigma_i}{dt} + \text{div} \left( \sum_{k=1}^{n-1} Co_{ik} \text{grad } e_k \right) \frac{\sigma}{\sigma} + \sigma_i \mathbf{v} = \sigma_i \quad (15)$$

This kind of mathematical relation is represented by bond graphs such as in Fig 2:

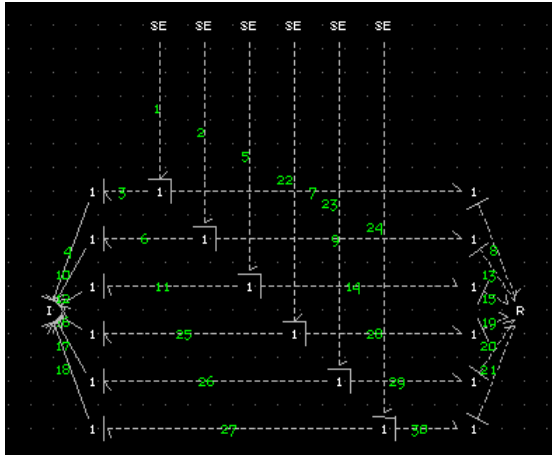


Fig 2 Representation of a dynamic reality with a bond graph (using the Camp-G software).

which give a complete knowledge of the dynamic reality at hand, including its dimensional analysis.

## SIMILITUDE

Two dynamic realities are similar if the input vector  $x$  and the output vector  $y$  of both realities are related by the same function.

Coupled with dimensional analysis [SZIRTES 1997], which allows, for example, to understand why pressure can be conceived as well as force per surface or energy per volume, bond graphs give direct access to the concept of similitude between two dynamic realities [THOMA 2006-2].

For example, a thermal diffusion phenomenon is represented by the Fourier equation of heat conduction:

$$\frac{\partial T}{\partial t} = a \frac{\partial^2 T}{\partial x^2} \quad (16)$$

where  $T$  stands for temperature of the heat source and  $a$  stands for the temperature conduction coefficient.

And a matter diffusion phenomenon is described by the Fick equation of diffusion:

$$\frac{\partial c}{\partial t} = D \frac{\partial^2 c}{\partial x^2} \quad (17)$$

where  $c$  stands for concentration and  $D$  stands for the coefficient of diffusion.

Both phenomena are similar realities because they share the same mathematical model, a fact that can be used by people working in virtual reality.

This kind of comparison is very simple, with the state space equations automatically generated by bond graph software, as in Fig. 3:

```

Inputs vector
u=[ SE1 SE2 SE5 ]
State variables vector
p,q=[P10;P12;P41;
A MATRIX
[ R13x8 R13 R13x15 R13x8 R13 R13x15
[ 14x10 110 112x10 14x12 110x12 112
R13x8 R13 R13x15
14 110x4 112x4
[ R15x8 R15x13 R15 R15x8 R15x13 R15
[ 14x10 110 112x10 14x12 110x12 112
R15x8 R15x13 R15
14 110x4 112x4
[ R8 R8x13 R8x15 R8 R8x13 R8x15
[ 14x10 110 112x10 14x12 110x12 112
R8 R8x13 R8x15
14 110x4 112x4
B MATRIX
[0 1 0]
[0 0 1]
[1 0 0]
C MATRIX
[sb sb sb]
D MATRIX
[sb sb sb]
System Order = 3

```

Fig 3 Example of a state space equation automatically generated by the Camp-G software.

Whatever the complexity of two considered realities, the necessary and sufficient condition of similitude between them is that the mathematical model of one be related by a bi-univoque relation to that of the other. In

other words that they have the same bond graph.

Modern computer programs allow the generation of simulators of the system from BGs, as in Fig.4 below. Moreover, programs like Systar (<http://www.systar.ch>) also allow the dimensional analysis of the variables describing the dynamic reality to be checked.

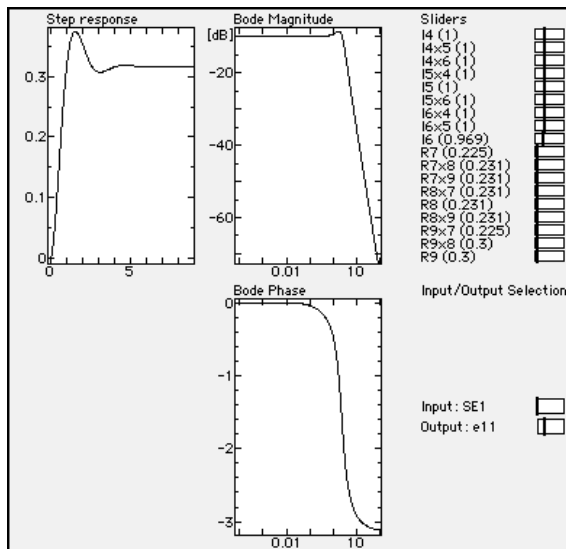


Fig 4 Front panel of an interactive simulator for a dynamic reality similar to the one in Fig. 2.

## CONCLUSION

Bond graphs are graphical representations of dynamic realities. They consequently give a mathematical and computable formulation of the behaviour of the considered realities. They provide us with a tool that allows us to represent their dynamic conditions.

Moreover, this method of modelling and simulation can draw heavily on similitude between energy or power domains and experimentally established functional relationships among their invariants.

Finally, coupled with dimensional analysis, they allow criteria of similitude to be derived

by comparison, providing solutions to many computer science virtual reality problems, within the domain of validity of the mathematical model.

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