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The multi-physics improved quasi-static method - Application to a neutronics-thermomechanics coupling

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ABSTRACT

The quasi-static method is widely used for space- and time-dependent neutron transport problems. It is based on the factorization of the flux into the product of two functions, an "amplitude" depending only on time and a "shape" which depends on all variables. Thanks to this factorization, long time-steps can be used for the computation of the shape, leading to a substantial reduction of the calculation time. Two algorithms, based on the quasi-static factorization, can be found in the literature: the "Improved Quasistatic Method" (IQM), and the "Predictor-Corrector Quasi-static Method" (PCQM).

In this paper we show, on the example of the Godiva experiment, that the IQM algorithm can be easily adapted to multi-physics simulations. Moreover, most of the common coupling or time-step control strategies are compatible with this algorithm and we test some of them here. In particular, a technique taken from existing codes with point-kinetic modules and based on feedback coefficients is found, in our case, to be especially efficient and gives precise and fast results. This shows that the multi-physics IQM presented in this paper is compatible with these existing codes, and may be a way to couple them with neutron transport solvers.

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1. Introduction

The quasi-static approach (introduced more than fifty years ago by Allan Henry ([Henry, 1958; Henry and Curlee, 1958\)](#page--1-0)) is widely used for solving space- and time-dependent neutron transport problems. It is based on the factorization of the neutron flux into an amplitude and a shape functions:

$$
\phi(\mathbf{r}, \mathbf{\Omega}, E, t) = N(t)\psi(\mathbf{r}, \mathbf{\Omega}, E, t).
$$
\n(1)

We introduce the diagonal "inertial operator" $I_{1/V}$ multiplying the flux by the inverse of the neutron speed. $\langle \cdot, \cdot \rangle$ denotes the classic scalar product. With these notations, the factorization (1) is defined by a constraint on the shape:

$$
\frac{\partial}{\partial t}\left\langle I_{1/V}\psi,f\right\rangle = 0, \tag{2}
$$

where f is an arbitrary weighting function. The steady-state adjoint flux ϕ_0^* is usually chosen as weighting function, and is used in this work.

There are two well-known ways of implementing the

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factorization (1). The historical approach [\(Ott and Meneley, 1969;](#page--1-0) [Dahmani, 1999](#page--1-0)) is called "Improved Quasi-static Method" (IQM) and is based on an amplitude (N in (1)) computation followed by a shape (ψ in (1)) one. This procedure is repeated until a convergence criterion on the constraint (2) is fulfilled. A more recent approach ([Dulla et al., 2008\)](#page--1-0), called "Predictor-Corrector Quasi-static Method" (PCQM) is based on a flux (ϕ in (1)) computation followed by an amplitude one, used to correct the flux prediction. No iterations are made. Both methods use two different time-scales: a micro one (dt) for the amplitude calculation and a macro one (Δt) for the shape or flux calculation.

The structure of the IQM algorithm seems better suited for solving multi-physics problems, and could lead to efficient coupling strategies. The purpose of this paper is to highlight the potential of IQM for multi-physics calculations, and to discuss some aspects of its use in this frame. We illustrate this idea with an original neutron transport-thermomechanics coupling example: a prompt critical burst of the Godiva experiment ([Wimett, 1956](#page--1-0)), which provides a strong and non-monotonic time-dependent coupling. CAST3M ("[CAST3M](#page--1-0)", 2015), a finite element code dedicated to structural mechanics, is used to compute the transient. In particular, a neutron transport solver, presented in ([Patricot et al., 2016](#page--1-0)) and able to deal * Corresponding author. with irregular meshes, is used.

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Section 2 recalls the quasi-static method and its two approaches, IQM and PCQM. Section [3](#page--1-0) presents the Godiva experiment and the models used. Sections [4 and 5](#page--1-0) discuss the transient computation and compare some coupling strategies with, respectively, fixed and variable time-steps. Finally, section [6](#page--1-0) concludes the paper.

2. The quasi-static method

2.1. The quasi-static equations

Consider the time-dependent neutron transport and delayed neutron precursor equations:

$$
\begin{cases}\nI_{1/V} \frac{\partial \phi}{\partial t} = \mathcal{L}\phi + F_p \phi + \sum_l \chi_l \lambda_l C_l \\
\frac{\partial C_l}{\partial t} = -\lambda_l C_l + \frac{1}{\chi_l} F_l \phi,\n\end{cases}
$$
\n(3)

where C_l denotes the normalized concentration of delayed neutron precursors of type l, \mathcal{L} the transport operator including streaming, absorption and scattering, F_p the prompt neutron production rate operator, F_l the delayed neutron production rate operator and $I_{1/V}$ the inertial operator. χ_l and λ_l are the delayed fission spectrum and the decay constant of the precursor of type l.

Taking the scalar product of (3) with the weighting function, and using [\(2\),](#page-0-0) we obtain the point-kinetic equations satisfied by the amplitude function:

$$
\begin{cases}\n\frac{\partial N}{\partial t} = \frac{\rho - \overline{\beta}}{\Lambda} N + \sum_{l} \lambda_{l} c_{l} \\
\frac{\partial c_{l}}{\partial t} = -\lambda_{l} c_{l} + \frac{\overline{\beta}_{l}}{\Lambda} N.\n\end{cases} \tag{4}
$$

The point-kinetic parameters are $(F = F_p + \sum_l F_l$ is the total neutron production rate operator):

$$
\rho = \frac{\langle (\mathcal{L} + F)\psi, \phi_0^* \rangle}{\langle F\psi, \phi_0^* \rangle},\tag{5}
$$

$$
\Lambda = \frac{\langle I_{1/V}\psi, \phi_0^* \rangle}{\langle F\psi, \phi_0^* \rangle},\tag{6}
$$

$$
\overline{\beta}_l = \frac{\langle F_l \psi, \phi_0^* \rangle}{\langle F \psi, \phi_0^* \rangle},\tag{7}
$$

$$
\overline{\beta} = \sum_{l} \overline{\beta}_{l},\tag{8}
$$

$$
c_l = \frac{\langle \chi_l C_l, \phi_0^* \rangle}{\langle I_{1/V} \psi, \phi_0^* \rangle}.
$$
\n(9)

The system of coupled equations satisfied by ψ and C_l is then obtained by replacing the factorization (1) in (3) :

$$
\begin{cases}\nI_{1/V} \left(\frac{\partial \psi}{\partial t} + \psi \frac{1}{N} \frac{\partial N}{\partial t} \right) = \mathscr{L} \psi + F_p \psi + \frac{1}{N} \sum_l \chi_l \lambda_l C_l \\
\frac{\partial C_l}{\partial t} = -\lambda_l C_l + \frac{N}{\chi_l} F_l \psi\n\end{cases}
$$
\n(10)

2.2. The two quasi-static approaches

2.2.1. The improved quasi-static method (IQM)

The IQM, introduced first in ([Ott and Meneley, 1969\)](#page--1-0), consists in solving the two coupled systems of equations (4) and (10) . The algorithm can be summarized as follows $(c$ is a convergence criterion chosen by user):

- Step 1: The point-kinetic equations (4) are solved first, over a macro time interval Δt , using micro time-steps dt.
- Step 2: Knowing the amplitude and its time derivative at $t + \Delta t$, the shape equations (10) are then solved, directly with Δt as a time-step.
- Step 3: The error ϵ on the normalization [\(2\)](#page-0-0) is evaluated and the shape is renormalized. This is necessary because the pointkinetic equations assume that the shape satisfies the normalization condition [\(2\).](#page-0-0)
- Step 4: Point kinetic parameters are computed.
- Step 5: Step 1-4 are iterated until $\varepsilon < c$.

Several numerical schemes are possible on this basis.

2.2.2. The predictor-corrector quasi-static method (PCQM)

In the PCQM ([Dulla et al., 2008](#page--1-0)) the flux equations (3) are coupled with the point-kinetic ones (4) . The algorithm can be summarized as follows:

- Step 1: The flux equations (3) are solved first, over a macro timestep Δt .
- Step 2: An approximate shape function is derived at $t + \Delta t$ from the flux of step 1 and condition (2) .
- Step 3: Point-kinetic parameters are evaluated and the pointkinetic equations (4) are solved, over the macro time interval Δt using micro time-steps dt.
- Step 4: Flux is reconstructed as the product of the shape of step 2 and of the amplitude of step 3.
- Step 5: Precursor concentrations are recomputed with this new flux.

Here again, details of the implementation can vary.

2.2.3. Block diagrams

Block diagrams for the two algorithms are sketched in Fig. 1:

Fig. 1. Block diagrams for IQM and PCQM calculation of a time-step.

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