

DYNAMICS OF MSR

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ABSTRACT

The Dynamics of the Molten Salt Reactor (MSR) - one of the 'Generation IV' concepts - was studied by means of a code developed at FZR. The graphite-moderated channel type MSR was selected for the numerical simulation of the reactor with liquid fuel. The MSR dynamics is very specific because the fuel flow influences the delayed neutrons distribution. Presently, there are only a few accessible numerical codes appropriate for the MSR simulation, therefore the DYN3D-MSR code was developed based on the FZR in-house LWR code DYN3D. It allows the 3D transient simulating of neutronics in combination with parallel channel type thermal-hydraulics. By means of DYN3D-MSR, several transients typical for the liquid fuel system were analyzed. Those transients were initiated by reactivity insertion, by overcooling of fuel at the core inlet, by the fuel pump start-up or coast-down, or by the blockage of selected fuel channels.

Keywords: Molten Salt, Liquid Fuel, Reactor Dynamics, MSR.

1. INTRODUCTION

The first projects for reactors with liquid fuel go back to the pioneers' time of nuclear energy, when various designs with different fuel-type have been proposed. From those the Molten Salt Reactor (MSR) could be considered as the most developed project [1]. The research on MSR was performed mainly in the Oak Ridge National Laboratory (ORNL), where the Molten Salt Reactor Experiment (MSRE) was realized in the sixties. This experiment has shown that the molten salt technology beside the energy production offers three main advantages:

- 1) excellent neutron economy,
- 2) continuous or in-batch reprocessing,
- 3) inherent safety features.

These three advantages make the MSR attractive also for the present Generation IV (GIF) initiative [2]. The aim of the GIF initiative is the development of technologies that achieve safety performance, waste reduction, and proliferation resistance while providing a nuclear energy option that is economically competitive. The MSR is studied also for spent fuel partitioning and transmutation (P&T) purposes. It is the aim of P&T to separate the different components of spent fuel (partitioning) and to reduce the radiotoxicity, mainly by the transmutation of actinides. A critical MSR can serve as transmutor with acceptable burning rates. The time period the radiotoxic inventory needs to reach the level of uranium ore can be shortened by transmutation from millions of years to less than one thousand years. Within the GIF initiative the graphite-moderated channel type MSR based on the previous ORNL research is considered (Fig. 1). In this MSR, the mixture of molten fluoride salts circulates through the whole primary circuit acting as fuel and coolant simultaneously. The fuel, in the form of uranium tetra fluoride, is dissolved in the carrier salt. The core is formed from hexagonal graphite blocks each with a central fuel channel (see Fig.2). The liquid fuel causes two physical peculiarities: the fission energy is released predominantly directly into the liquid fuel and the delayed neutrons (DN) precursors are convected (drifted) by the fuel flow out of the core. Therefore, specific tools are needed for the reactor dynamics analysis of the MSR.

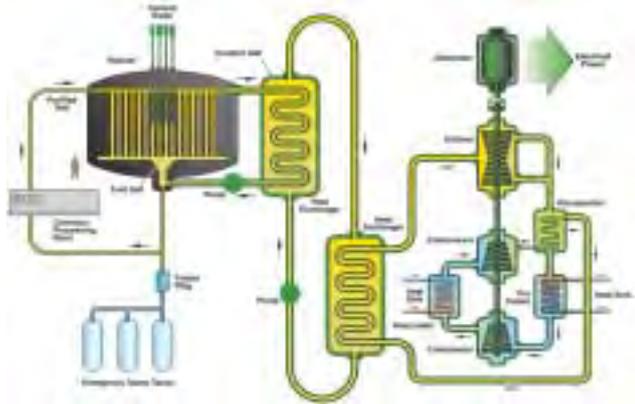


Fig. 1: Scheme of MSR [2].

Consideration of the MSR within the GIF initiative and in the EU-project MOST (Review of Molten Salt Technology) in which we participated was the motivation to develop DYN3D-MSR for MSR dynamics simulation at the Institute of Safety Research. The code DYN3D-MSR [3] is based on the FZR in-house code DYN3D [4] which was developed and validated initially for LWR applications. Worldwide, only one comparable 3D code is presently suitable for the spatial numerical simulation of MSR dynamics.

2. NEUTRONICS OF THE MSR

The DYN3D-MSR code consists of a neutron-kinetics and a thermal-hydraulics part [5]. The neutron-kinetics model is based on the neutron diffusion equation with two energy groups. The diffusion equation is solved in nodal approach. The prompt fission neutron source is calculated using the same nodal flux expansion method [6] as the basic DYN3D code. Part of the neutrons is released with time delay when the precursor nuclides decay. Due to the drift of these precursors with the liquid fuel along axial direction an extended delayed neutrons model had to be implemented. The drift is described by an additional convective term in the equation for precursors concentration. The drift must also be considered in the primary circuit loop outside the core.

The following set of equations is solved in each node n :

$$\frac{\partial \Phi_1^n(z,t)}{v_1^n \partial t} + \nabla J_1^n(z,t) + \Sigma_1^n(t) \Phi_1^n(z,t) = \frac{1}{k_{eff}} \sum_{g=1}^2 (1 - \beta_g^n) \Sigma_{f,g}^n(t) \Phi_g^n(z,t) + \sum_{j=1}^M \lambda_j^n C_j^n(z,t) + S_{ext}^n(z); \quad (1)$$

$$\frac{\partial \Phi_2^n(z,t)}{v_2^n \partial t} + \nabla J_2^n(z,t) + \Sigma_a^n(t) \Phi_2^n(z,t) = \Sigma_s^n(t) \Phi_1^n(z,t); \quad (2)$$

$$\frac{\partial C_j^n(z,t)}{\partial t} + \frac{\partial}{\partial z} (v^n C_j^n(z,t)) = \frac{1}{k_{eff}} \sum_{g=1}^2 \beta_{g,j}^n v \Sigma_{f,g}^n \Phi_g^n(z,t) - \lambda_j^n C_j^n(z,t); \quad (3)$$

standard notation is used. Assuming constant drift velocity within one node and during one time step the equation (3) can be integrated in each node. These nodal solutions are combined with the method of characteristics to find the precursors distribution in the whole primary circuit. This method involves a tracking of fictitious fluid particle during one time step Δt on its way through one or more nodes. The particle which trajectory ends on the outgoing edge of given node is selected. Once the starting point of the trajectory is found the initial precursors' concentration is found by the linear interpolation between two edge values, which are known from the pervious time step. In next the known nodal solutions relevant for the trajectory position are sequentially used to find the precursors concentration on the end of the trajectory. To obtain the overall precursors distribution this procedure is used for every node.

As a consequence of the drift, the DN precursors propagate through the entire primary circuit (see Fig. 2). Nevertheless, the final effect on the neutron-kinetics depends mainly on the flow velocity and on the volumetric fuel proportion in the core and in the primary circuit, respectively.

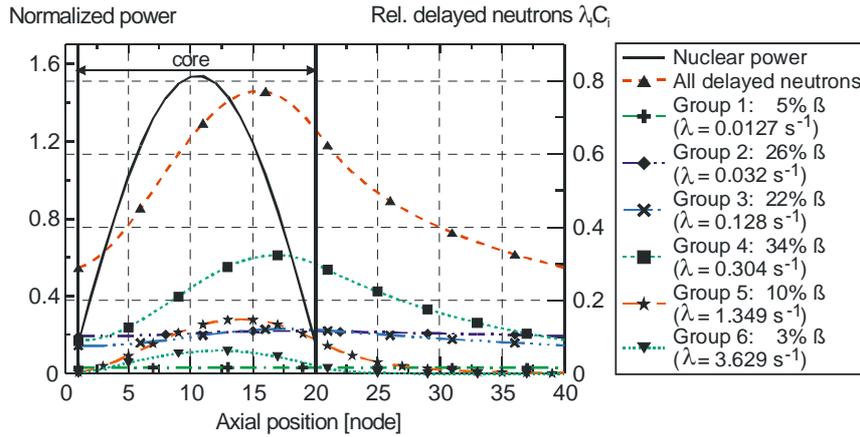


Fig. 2: Distribution of relative reactor power and 6 delayed neutrons groups for a salt velocity of $v=1m/s$. The core and the external loop consist each from 20 axial nodes.

3. THERMAL-HYDRAULICS OF THE MSR

The thermal-hydraulics of DYN3D-MSR is based on the DYN3D light water reactor model [7]. However, the energy balance equation must be adapted for the liquid fuel. As mentioned above, it is a unique feature of the MSR reactors that the liquid molten salt is flowing through the core and the whole primary circuit acting as fuel and coolant simultaneously. The energy from fission is predominantly released directly in the fuel and is immediately convected by the flow out of the core, while the graphite is heated-up by the gamma and neutron radiation [8]. Thus the energy balance equation will include three energy sources:

$$\frac{\partial}{\partial t} (\rho h) + \frac{\partial}{\partial z} (\rho v h) = Q_{fission} + Q_{delay} + Q_{graphite}, \quad (4)$$

standard notation is used. The three sources represent the direct heat release $Q_{fission}$, the decay heat Q_{delay} , and the heat exchange between graphite and salt $Q_{graphite}$. The heat source $Q_{graphite}$ is defined by heat release and heat conduction in the graphite and by heat exchange between graphite and liquid fuel. The heat conduction and heat transfer models had to be modified. Heat exchange correlations were implemented for the liquid fuel. The heat conduction model was adopted for the solid graphite structures. The method of effective heat transfer coefficients was used to calculate the temperature distribution in the graphite. To determine these transfer coefficients an

analytical solution of the heat conduction equation must be known in the graphite. Thus, the hexagonal graphite blocks were approximated by cylinders with an inner fuel channel and divided into segments.

When all the three sources are known, the equation (4) can be integrated in each node using the same assumption of constant drift velocity as mentioned in the previous chapter. Once the solution in each node is known, the method of characteristics can be used to get the overall temperature distributions. As an example, typical axial temperature distributions in one fuel element divided into 20 axial nodes and 4 radial graphite segments are shown in Fig. 3.

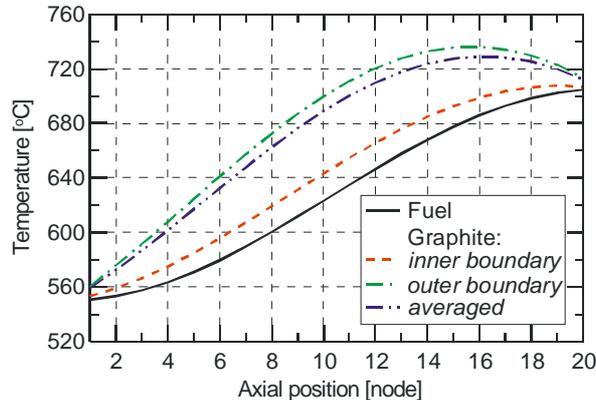


Fig. 3: Axial distributions of temperature in the MSR fuel channel. Inlet temperature of the fuel is 550°C. Flowing through the core; it is heated up by 150°C, approximately.

4. MOLTEN SALT REACTOR EXPERIMENT

The Molten Salt Reactor Experiment – MSRE was an 8-MW, fluid-fuelled test reactor that operated from 1965 through 1969 in the Oak Ridge National Laboratory – ORNL [9]. The fuel consisted of a mixture of lithium, beryllium, and zirconium fluoride salts, to which uranium was added as UF_4 . The molten salt mixture with the temperature above 600°C circulated through the reactor vessel, fuel pump, heat exchanger, and piping. These components have been constructed from Hastelloy-N (70%Ni, 18%Mo, 7%Cr, 5%Fe). The MSRE core consisted from square graphite stringers, which form the fuel channels. The original geometry can be seen from Fig. 4.

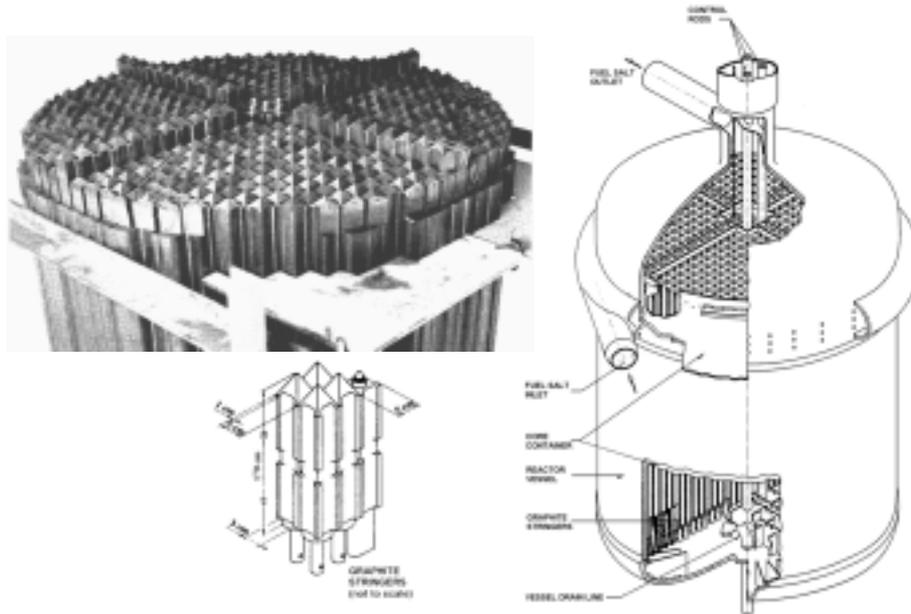


Fig. 4: 610 graphite stringers prepared for the installation into the reactor vessel (top-left), several graphite stringers which form the core in detailed view (lower-left), and the MSRE reactor vessel (right).

5. MSRE NODALIZATION

The MSRE core consisted from 610 square graphite stringers. The inner core diameter was 135 cm and the outer pressure vessel diameter was 142 cm. The width of each stringer was 5 cm and the inner core radius was filled by 577 full stringers. The remaining 33 elements were partly cut to fit into the cylindrical vessel (see Fig. 5). By this structure 1150 fuel channels were formed in the core. Through these channels the liquid fuel circulates by the rate 52 l/s. It corresponds to the drifting velocity of 19.7 cm/s. The nominal fuel pressure was 344.7 kPa and the nominal inlet and outlet temperature was 635°C and 663°C, respectively.

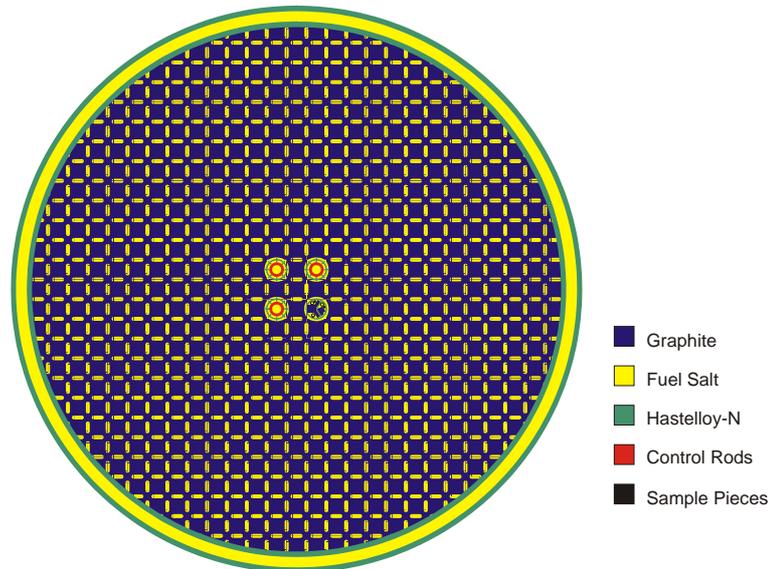


Fig. 5: Structure of the MSRE core with detailed view of the central part.

The code DYN3D-MSR is able to calculate with both Cartesian and hexagonal fuel assembly geometry. To reduce the calculating time the core was approximated with only about 200 channels while neglecting the central part construction and the control rods. The best approximation was obtained for hexagonal assembly geometry using only 199 channels (see Fig. 6). Along the axial direction, the core was divided into the 20 nodes. From those, 18 nodes represent the graphite-salt area. The graphite stringers were 166cm long, thus each from these 18 nodes was 9.2cm long. The first and the last node represent the upper and lower plena, respectively. The effective plena heights were estimated to be 17.2cm. The total core height was approximately 200cm.

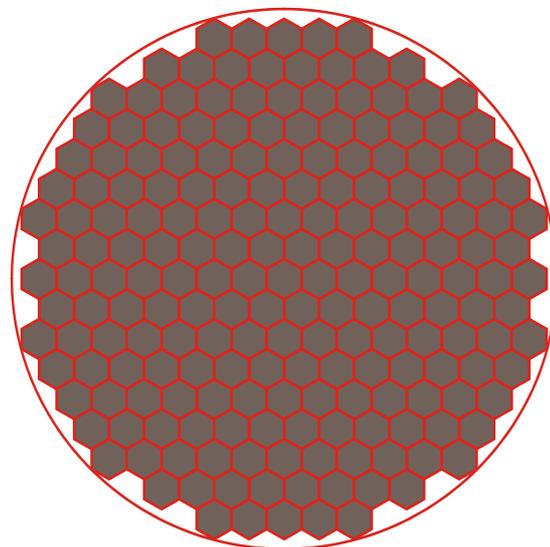


Fig. 6: Approximation of core by 199 hexagons.

The residual salt volume in the piping and in the fuel pump was represented by arbitrary number of uniform nodes. The size of these nodes was given by the drifting velocity and by the overall recirculation time (duration of one passage through the whole primary circuit). With the MSRE the recirculation time was 25s for the nominal fuel velocity.

6. DYN3D-MSR CALCULATIONS

In the frame of the MOST project, a validation benchmark was defined, based on the experimental data measured during the MSRE [10]. The cross-section libraries appropriate for the calculation were prepared by EdF using the Apollo code [11]. The benchmark was calculated by means of different codes in five institutions [12]. For the validation of DYN3D-MSR neutronics, especially the fuel pump start-up and coast-down transients are important. During these two transients the fuel flow velocity varies and the DN distribution is changing. However, constant power is maintained by means of control rods (Fig. 7).

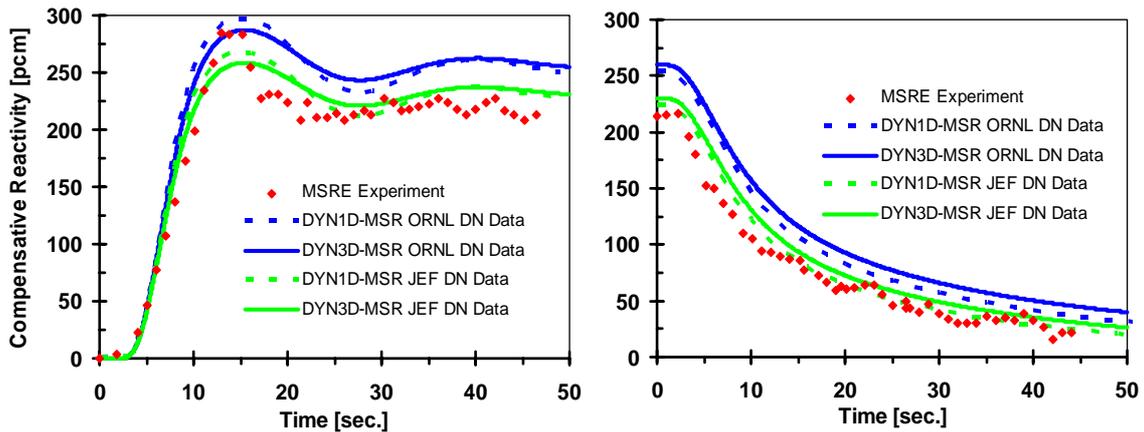


Fig. 7: Compensative reactivity inserted by control rods during the fuel pump start-up (left) and coast-down (right) transients for original ORNL and calculated JEF DN data.

After successful validation, the code was applied to several hypothetical transients in a typical MSRE core. The 300 pcm prompt reactivity jump is presented here as an example of different power transients. The power increases rapidly during the first few seconds of this transient. However, the response of fuel salt temperature is immediate and the power increase is slowed-down rapidly due to the negative reactivity introduced by the fuel temperature. The long-term behavior is determined by the slowly growing graphite temperature through the moderator temperature feedback coefficient (Fig. 8).

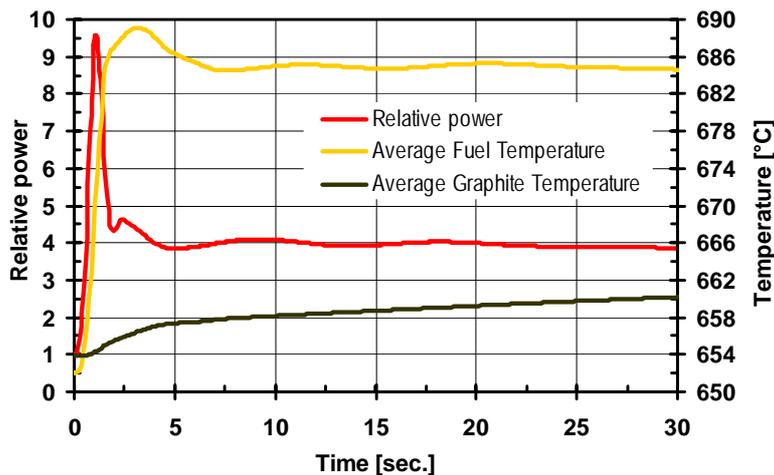


Fig. 8: Response of the MSRE power and average graphite and fuel temperature after the prompt reactivity jump (300 pcm).

7. SUMMARY AND OUTLOOK

The delayed neutrons drift and the direct release of fission energy into the liquid fuel are two attributes specific to MSR dynamics. The appropriate models have been integrated into the DYN3D-MSR code. The code was validated against the experimental results from the MSRE. Then it was applied to several hypothetical transients analyses, where space-dependent effects are relevant. In the considered transients, the response of the MSR is characterized by the immediate change of the fuel temperature with changing power and fast negative temperature feedback to the power.

The DYN3D-MSR code has been shown to be an effective tool for MSR dynamics studies. In the future the code is scheduled to be extended by modeling the heat exchanger between primary and secondary circuit and by considering the reprocessing unit, to cover a wider spectrum of transients.

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