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On the use of a dedicated ballast pellet for a prompt self-ejection mechanism after a temperature transient in lead-cooled fast reactors

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Abstract

The potential use of changes in buoyancy as a reactivity feedback mechanism during temperature transients in heavy liquid metal fast reactors (HLMFRs) is discussed. It is shown that with the use of ballast pellets ($\sim 15\%$ volume fraction) introduced in combination with fuel pellets, fuel rods will be endowed with a reliable self-ejection mechanism that is able to compensate temperature transients. Utilizing a simplified model, an estimate of the negative reactivity insertion expected from this mechanism is derived. The use of ballast pellets opens up the possibility of introducing greater amounts of actinides into the core, as well as providing a solution to the classical problem of positive coolant temperature reactivity coefficients in fast reactors.

Keywords: Heavy liquid metal fast reactors, buoyancy, temperature transient compensation, Generation IV reactors

1. Introduction

One of the unique features of heavy liquid metal fast reactors (HLMFRs) with lead or lead-bismuth eutectic coolant is the very high density of the coolant: the coolant density in HLMFRs is similar to that of the fuel. The potential use of this feature has either been overlooked by nuclear designers or seen as a “nuisance”, and, as a result, preventive measures such as the

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use of tungsten deadweight (ballast) to overcome buoyancy forces have been proposed [3, 7].

The objective of this study was to assess the potential for exploiting
10 changes in buoyancy forces as a control mechanism for fuel rod self-ejection
during HLMFR temperature transients, thereby providing a reliable solution
to the well-established problem of the positive coolant temperature reactiv-
ity coefficient exhibited by sodium fast reactors and also in lead-cooled fast
reactors depending on the size of the reactor core. This concept is expected
15 to represent a passive safety feature for the system but it does not represent
at all a control device to be used during reactor normal operation.

The effect of buoyancy forces in HLMFRs as a positive aspect in safety
analysis during a post-accident heat removal scenario was recently investi-
gated by Arias [4]. It was found that, because of the similar densities of the
20 fuel and the heavy liquid metal (HLM) coolant, an inherent passive safety
feedback self-removal mechanism governed by buoyancy is developed, pro-
pelling the packed bed away from the wall, and preventing temperatures
that could jeopardize the structural integrity of the vessel being reached,
as well as reducing the re-criticality potential by limiting the allowable bed
25 depth.

Thus, it is interesting to consider whether buoyancy forces, rather than
being regarded as a nuisance during nominal operating conditions, can be
harnessed as a mechanism for endowing fuel rods with unique safety prop-
erties only available in HLMFRs. In the sections that follow, this possibility
30 will be investigated and discussed. However, the reader should be aware that
the results reported in this preliminary analysis of the proposed concept are
based on idealizations, of the sort which are inevitable in preliminary the-
oretical assessments of concepts, and therefore should not be misconstrued
as definitive detailed analysis. The final verdict about the feasibility of the
35 proposed concept will only be reached following detailed analysis of the com-
plexities arising from the proposed solutions, the subject of future work.
Nonetheless, we feel that this preliminary assessment is appropriate at this
time, to encourage (or not) further careful investigation of the idea.

2. Buoyancy forces as a fuel rod ejection mechanism

40 Fig. 1 illustrates schematically the mechanism we seek to exploit. For the
envisaged mechanism to work as intended the density of the coolant needs

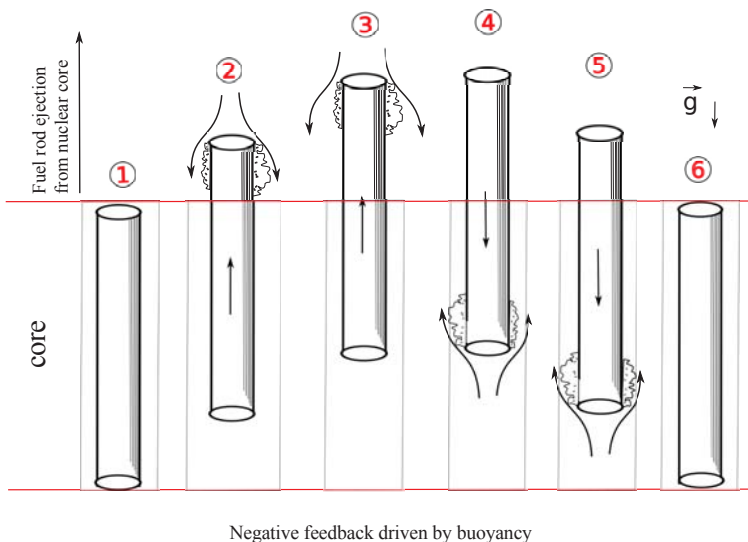


Fig. 1. Fuel rod ejection by buoyancy forces. Sequence: (1) Insertion of reactivity, leading to rising temperatures; (2) Due to relative changes in density with temperature, buoyancy effects act and the fuel rod is propelled upwards; (3) A subcriticality condition is reached, leading to falling temperatures; (4) Relative changes in density lead to loss of buoyancy and the fuel rod falls back down; (5) Fuel rod re-enters the core; (6) End of transient.

to become greater than the effective density of the fuel as the temperature increases.

Fig. 2 shows the variation of density as a function of temperature for mixed oxide (MOX) and UO_2 fuels and Pb-Bi eutectic and Pb coolants. This indicates that the relative changes of HLM coolant and fuel densities with temperature are not favorable. However, before deciding on the feasibility of the posited buoyancy mechanism, the fuel densities shown in Fig. 2 need to be corrected to account for the presence of stainless steel, mostly in the form of cladding. Thus, to take into account the effect of stainless steel on the total density of the fuel, a combined fuel-steel density may be defined as:

$$\bar{\rho}_f = F_f \rho_f + (1 - F_f) \rho_s \quad (1)$$

where F_f is the volume fraction of fuel and ρ_f and ρ_s are the densities of the fuel and stainless steel, respectively.

For practical purposes, the densities can be approximated as linear functions of temperature:

$$\rho_i = \rho_{i,0} - \alpha_i T_i \quad (2)$$

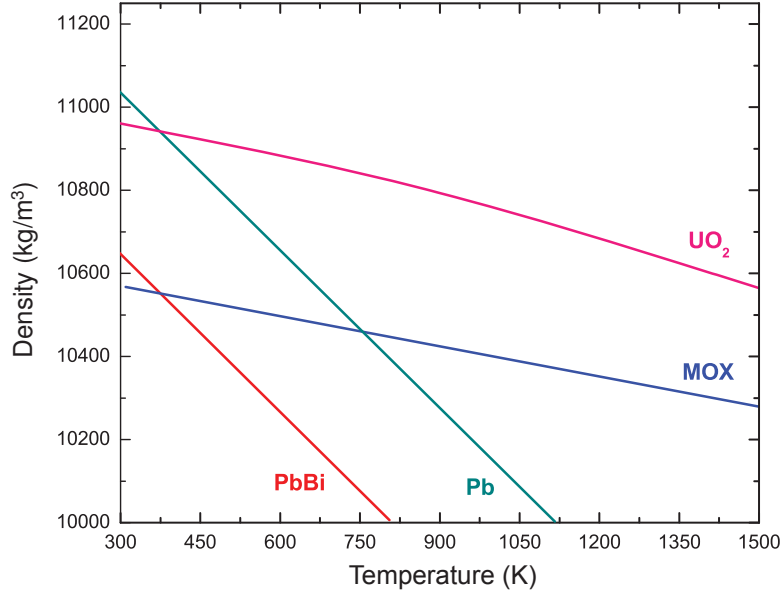


Fig. 2. Density variations of Pb-Bi eutectic and Pb coolants and MOX and UO₂ fuels as functions of temperature.

where the subscript i denotes the specific material, for example, $i = f$ for fuel, c for coolant, s for stainless steel, and $\rho_{i,0}$ is the density of material i at a temperature of 0 K, α_i is the rate of change of density of material i with temperature, and T_i is the temperature of material i in K. Then, the combined density given by Eq. (1) can be represented as a function of temperatures as:

$$\bar{\rho}_f = \bar{\rho}_{f,0} - \bar{\alpha}_f T_f \quad (3)$$

where

$$\bar{\rho}_{f,0} = [F_f \rho_{f,0} + (1 - F_f) \rho_{s,0}] \quad (4)$$

and

$$\bar{\alpha}_f = (1 - F_f) \frac{T_s}{T_f} \alpha_s \quad (5)$$

where T_s is the average temperature of the cladding and can be calculated as $T_s = T_f - \Delta T$, with ΔT being the temperature drop between fuel and cladding. A typical value of ΔT is ~ 200 K. This value has been assumed for the preliminary calculations in this paper.

From the available data in the literature, the linear relationships for fuels [17], coolants [16] and stainless steel [12] shown in Table 1 were formulated.

70 All densities are given in kg m^{-3} for temperatures in K. The corresponding relationships are depicted in Fig. 3, where a volume fraction of stainless steel of 45.6% (from Table 2) was assumed.

Table 1. Assumed density variations with temperature.

Material type	Material	Equation
Coolant	Pb	$\rho_{\text{Pb}} = 11478.69 - 1.32T_c$
Coolant	PbBi	$\rho_{\text{PbBi}} = 11093.71 - 1.33T_c$
Fuel	UO ₂	$\rho_{\text{UO}_2} = 11122.84 - 0.36T_f$
Fuel	MOX	$\rho_{\text{MOX}} = 10657.97 - 0.255T_f$
Cladding	Stainless steel SS-316	$\rho_s = 8077.729 - 0.42T_s$
Ballast	Tungsten	$\rho_w = 19300.0 - 0.22T_f$

Referring to Fig. 3, it can be seen the densities of the HLM coolants are consistently greater than those of the combined fuel-steel options. Thus, the desired buoyancy mechanism for self-ejection of a fuel rod will only be possible with the use of deadweight or ballast to increase the effective density of the fuel. The use of such ballast is discussed below.

2.1. The tungsten ballast pellet

Although the use of tungsten as ballast in lead-cooled reactors has been proposed previously, its application was for a totally different purpose: tungsten ballast is located outside the core and used to keep the fuel assemblies in their designated positions by providing a downward force exceeding the force due to buoyancy under refueling conditions [2]. In other words, buoyancy forces are not contemplated as the basis of a feedback mechanism but, rather, they are neutralized over all temperatures by the use of an excess of ballast.

The proposed use of tungsten ballast here is with a totally different purpose in mind. We want to neutralize buoyancy, but only in the nominal range of working temperatures of the reactor, and we want buoyancy forces to appear if the nominal operating temperature range is exceeded, for example during a temperature or power transient. So, by introducing a tungsten ballast pellet occupying just the right volume within the fuel rod (as depicted in Fig. 4) we will be able to endow the fuel rod with a reliable mechanism for self-ejection or self-disassembly, as depicted in Fig. 1.

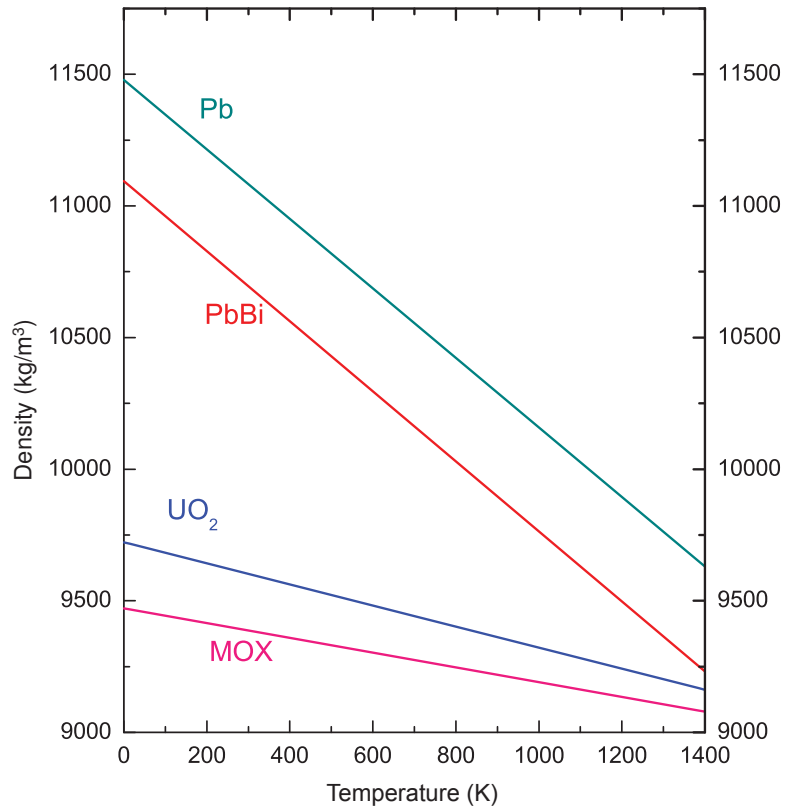


Fig. 3. Density variations as functions of temperature of Pb-Bi eutectic and Pb coolants and MOX- and UO₂-based fuels with a representative volume of stainless steel cladding.

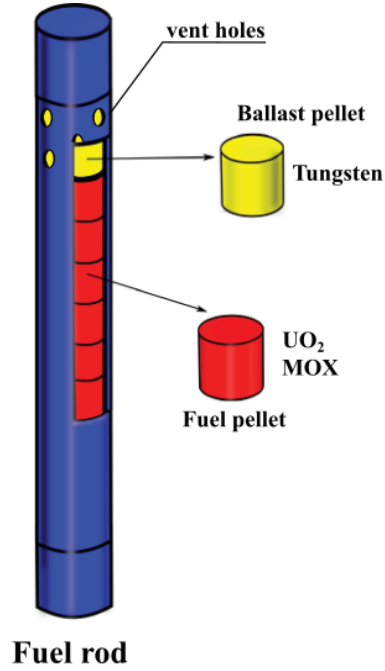


Fig. 4. The ballast pellet concept: by introducing a ballast pellet into the fuel rod it is possible to harness buoyancy forces to provide negative reactivity feedback.

95 Our first step, therefore, is to derive an expression that allows us to determine what tungsten ballast pellet fraction will be needed, and our second step is to establish an initial estimate of the negative reactivity insertion arising from the consequent fuel rod self-ejection.

100 First, we need to define an “effective density” taking into account the volume fraction occupied by tungsten ballast pellets. Proceeding as in our previous analysis, the effective fuel-steel-tungsten density is:

$$\rho_{f,\text{eff}} = (1 - F_b)\bar{\rho}_f + F_b\rho_w \quad (6)$$

where F_b is the volume fraction of tungsten ballast used and ρ_w its density. From our foregoing discussion, Eq. (6) yields the following relationship:

$$\rho_{\text{eff}} = \rho_{\text{eff},0} - \alpha_{\text{eff}}T_f \quad (7)$$

where

$$\rho_{\text{eff},0} = (1 - F_b)\bar{\rho}_{f,0} + F_b\rho_{w,0} \quad (8)$$

105 Design considerations dictate that the ballast pellet should be positioned at the top or bottom of the fuel element, thereby avoiding thermal stresses between fuel pellets, and also enabling the ballast pellet to act as a reflector (thanks to the high density of tungsten) and/or as a bottom- or top-cap, as schematically indicated in Fig. 4. The design of the ballast pellet will also

110 be influenced by the location of the gas plenum. If the gas plenum is at the same end of the fuel element as the ballast, then the ballast pellet should contain holes to allow the free flow of fission gas towards the plenum.

Thus, accounting for the contribution due to the expansion of the ballast, the effective rate of change of density is

$$\alpha_{\text{eff}} = (1 - F_b)\bar{\alpha}_f + F_b \frac{T_w}{T_f} \alpha_w \quad (9)$$

115 where T_w is the temperature of the ballast at the appropriate location. Because of the high thermal conductivity of tungsten ($\kappa_w \approx 173 \text{ WK}^{-1}\text{m}^{-1}$), T_w can be assumed to be approximately equal to the local temperature of the fuel. The fuel temperature falls by around 50% between its maximum axial value (close to the center of the fuel element) and the outermost axial
 120 levels where the ballast should be placed. Thus, to be on the safe side, a conservative preliminary value for the effective ballast temperature is taken as $T_w \approx \frac{1}{2}T_f$. In overestimating the temperature of tungsten, we are underestimating its density and thus overestimating the volume fraction needed. From the available literature [12], the density of tungsten fits the relationship
 125 given in Table 1.

Fuel rod ejection driven by buoyancy will only occur when the effective density of the fuel becomes lower than that of the surrounding coolant, or:

$$\rho_c > \rho_{\text{eff}} \quad (10)$$

To progress our analysis, we need an expression connecting the temperature of the fuel with the temperature of the coolant at the same instant
 130 in time. It should be noted, however, that even if the condition given by Eq. (10) is satisfied, this does not guarantee the feasibility of the proposed buoyancy mechanism: we must, additionally, be sure that this condition is accomplished at a power below the critical power that can jeopardise the structural integrity of the cladding. Thus, it is important to relate the fuel
 135 and coolant temperatures to the power being generated in the fuel. For transients in which reactivity ρ is much lower than the delayed neutron fraction $\rho \ll \beta$, the resulting reactor period would be considerably longer than the fuel thermal time constant τ given by [14]:

$$\tau \approx R_f M_f c_f \quad (11)$$

where M_f and c_f are the mass and specific heat capacity of the fuel, respectively, and R_f is the fuel thermal resistance given by:

$$R_f = \frac{1}{4\pi L \kappa_f} + \frac{1}{2\pi r_g L h_g} + \frac{1}{2\pi \kappa_s L} \ln \left(\frac{r_{s2}}{r_{s1}} \right) + \frac{1}{2\pi r_{s2} L h_c} \quad (12)$$

where κ_f is the thermal conductivity of the fuel, L the fuel rod length, r_g and h_g are the effective gap radius and heat transfer coefficient, respectively, r_{s2} and r_{s1} the outer and inner cladding radius, respectively, κ_s the thermal conductivity of the cladding, and h_c the coolant heat transfer coefficient.

It should be mentioned that Eq. (12) refers to the peak fuel temperature (centerline or hollow), not to the average fuel temperature. The latter is the temperature upon which density depends. A correction could be introduced by multiplying the first term in the right-hand side of Eq. (12) by $\frac{1}{2}$ [19]. However, the use of a peak fuel temperature, on one hand, results in an overestimation of the ballast pellet volume, but, on the other hand, in an underestimation of the power at which the buoyancy becomes effective. These effects will be somewhat compensatory, and, in view of the uncertainties in this preliminary assessment, let us use the peak fuel temperature in our calculations.

For the case where the reactor period is much longer than τ , the fuel and coolant temperatures can be expressed as functions of the power P as [14]:

$$T_f = \left[R_f + \frac{1}{2\dot{m}_c c_c} \right] P + T_i \quad (13)$$

and

$$T_c = \frac{1}{2\dot{m}_c c_c} P + T_i \quad (14)$$

where \dot{m}_c is the coolant mass flow rate and heat capacity, respectively; and T_i the coolant inlet temperature.

Thus, using the equations above, we find that the point at which the condition given by Eq. (10) is met occurs at a power given by:

$$P^* = \frac{\rho_{c,0} - \rho_{\text{eff},0} - T_i (\alpha_c - \alpha_{\text{eff}})}{\frac{\alpha_c - \alpha_{\text{eff}}}{2\dot{m}_c c_c} - \alpha_{\text{eff}} R_f} \quad (15)$$

To better understand the implications of these results, we assume some typical values for the relevant parameters. For the calculation of the thermal

Table 2. Design parameters of the HLMFR core concept considered, from [20].

Parameter	Value
Power	600 MW _e
Pellet outer radius	3.3 mm
Cladding inner radius	3.4 mm
Cladding outer radius	4.55 mm
Pitch-to-diameter ratio	1.5
Length of upper plenum	100 cm
Length of lower plenum	10 cm
Active pin length	100 cm
Pin-fuel volume fraction	54.4%
Pin-steel volume fraction	45.6%
Average linear pin power	11.5 kWm ⁻¹

resistance, we take: $\kappa_f = 2 \text{ Wm}^{-1}\text{K}^{-1}$, $\kappa_s = 15 \text{ Wm}^{-1}\text{K}^{-1}$; from [15], $h_g =$
165 $5678.26 \text{ Wm}^{-2}\text{K}^{-1}$, $h_c = 34069.58 \text{ Wm}^{-2}\text{K}^{-1}$; and, from Table 2, $r_{s2} = 4.55$
mm, $r_{s1} = 3.4$ mm, with a core length of $L = 100$ cm. These result in a
fuel thermal resistance of $R_f \approx 4.39 \times 10^{-2} \text{ KW}^{-1}$. For the coolant, we take
 $c_c = 160 \text{ JK}^{-1}\text{kg}^{-1}$. The maximum coolant velocity allowed for lead-based
coolants is in the range 2–3 ms^{-1} because of issues of erosion [21]. Then,
170 for the channel dimensions given in Table 2, the coolant mass flow rate is
 $\dot{m}_c = 2 \text{ kg s}^{-1}$.

For the average nominal linear pin power, we take a value of 115 W cm^{-1} ,
as in Table 2 [20]. Taking an inlet temperature of $T_i = 750 \text{ K}$, corresponding
with a nominal linear pin power of 100 W cm^{-1} , then results in fuel and
175 coolant temperatures that vary as functions of linear pin power as shown in
Fig. 5.

Fig. 6 shows how the densities of Pb-Bi eutectic and Pb coolants will
vary as functions of linear pin power according to these equations along with
the variation of the effective density of MOX (Fig. 6) fuels. In these figures,
180 the choice of the fraction of tungsten ballast pellets used was more or less
arbitrary, with the only purpose being to obtain an estimate of the amount of
ballast needed to stop the transient safely, i.e. to ensure that fuel rod ejection
occurs at a linear power significantly smaller than a certain design constraint,
for example, the 472 W cm^{-1} limit suggested by Hitachi [8]. However, as will
185 be apparent to the reader, the nuclear designer has a certain freedom of choice

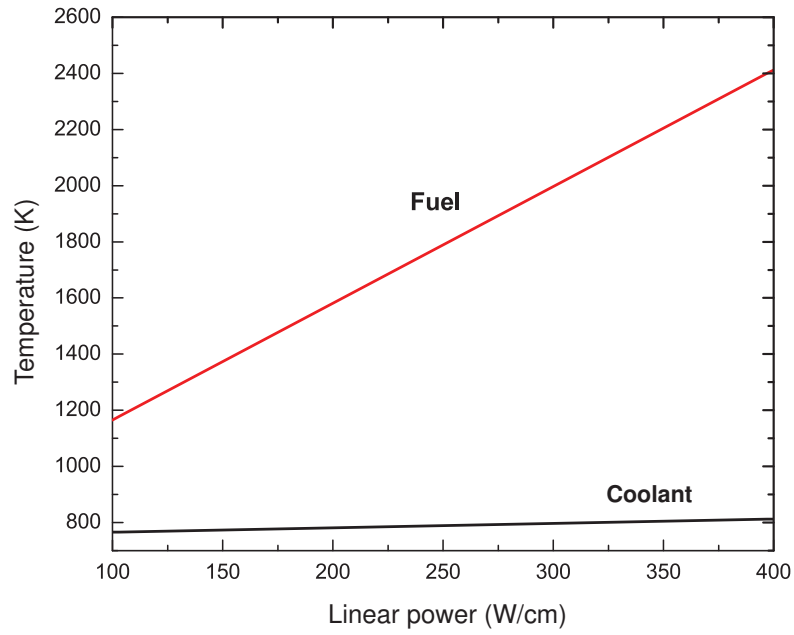


Fig. 5. Fuel and coolant temperatures as functions of linear pin power.

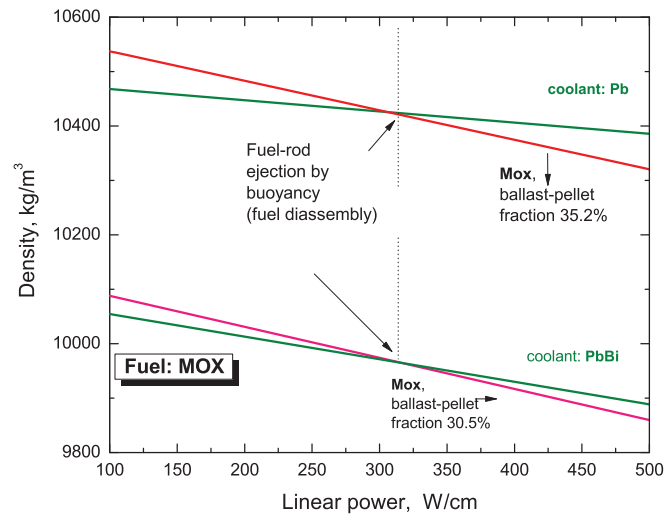


Fig. 6. Densities of coolants and MOX fuel as functions of linear pin power.

over the maximum power at which the fuel rod is ejected. If the fraction of ballast is reduced from the values used in Fig. 6, ejection will start at a lower linear power, but the system will also become more sensitive to small changes of temperature.

190 Next, we need to obtain a first estimate of the amount of negative reactivity insertion caused by the buoyancy-driven ejection of the fuel rod when the Eq. (10) condition is met. This will be our objective in the next section.

2.2. The negative reactivity insertion

195 In this section we will provide some first estimation on the negative reactivity insertion. This is a difficult calculation to accurately predict. Substantially uncertainties will be necessary introduced at every step of the analysis from the unavoidable idealizations required if analytical expressions are desired. If more precise calculation are desired more complex iterative methods will be required. For example, in view of several uncertainties, the simplest
 200 lumped model for heat transfer to arrive at the buoyancy-driven parameter and the terminal velocity seems for preliminary result preferable. But as mentioned before, this is very approximate and can cause significant error in the prediction of reactivity addition rate and the time scale involved for the buoyancy-driven mechanism to control the transient. Therefore, the negative reactivity insertion reported result from idealizations and is therefore
 205 not intended to typify negative reactivity estimates.

At the moment fuel rod ejection starts the maximum reactivity is given by

$$\rho_0 = \gamma_c \Delta T \quad (16)$$

210 where γ_c is the (positive) coolant temperature coefficient of reactivity and

$$\Delta T = T_c - T_c(0) \quad (17)$$

is the increase in coolant temperature from the initial value $T_c(0)$ to the temperature T_c when ejection occurs, i.e. at power $P = P^*$.

The negative reactivity insertion due to the sudden upward motion of the fuel rod over a small time-step Δt is:

$$\Delta \rho = - \left| \frac{\partial \rho}{\partial z} \right| \left| \frac{\partial z}{\partial t} \right| \Delta t = - \left| \frac{\partial \rho}{\partial z} \right| V_t \Delta t \quad (18)$$

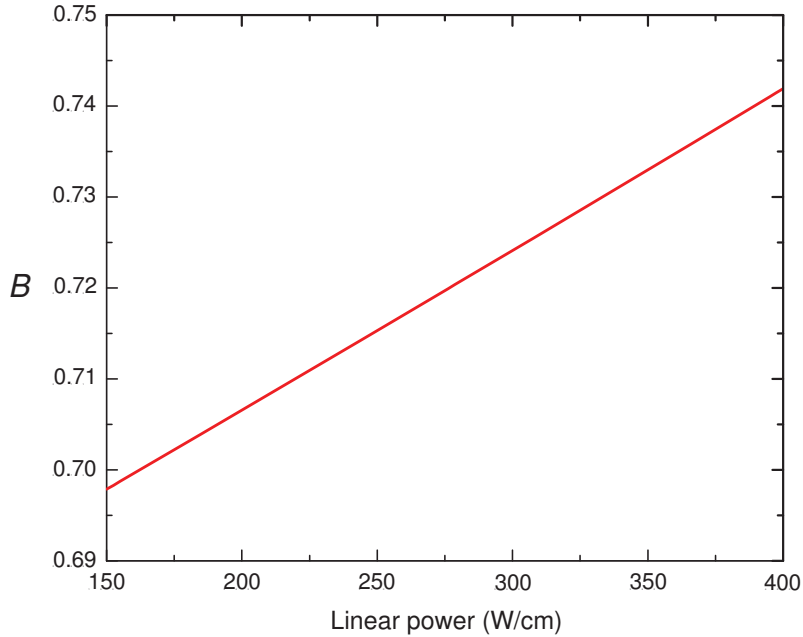


Fig. 7. The buoyancy-driving parameter B with lead coolant and UO_2 fuel as a function of linear pin power.

215 where V_t is approximately the terminal velocity of the cylindrical fuel rod, given by [13]:

$$V_t = \sqrt{\frac{2gL}{C_d} B} \quad (19)$$

where g is the acceleration due to gravity, L is the fuel rod length, C_d is the drag coefficient, and B is a buoyancy-driving parameter defined as:

$$B = \sqrt{\frac{\rho_c - \rho_{\text{eff}}}{\rho_{\text{eff}}}} \quad (20)$$

220 Using the representative values specified in the previous section, the relationship between B and linear pin power for lead coolant and UO_2 fuel is as shown in Fig. 7.

225 Now, in sufficiently slow transients, as soon as the condition given by Eq. (10) is satisfied, there will be a small prompt jump in power, but then the system will come to equilibrium as the rise in fuel and coolant temperatures and the increase in B compensate for ϱ_0 , sending $\varrho(t) \rightarrow 0$. Thus, neglecting the effect of the negative fuel temperature coefficient of reactivity and other

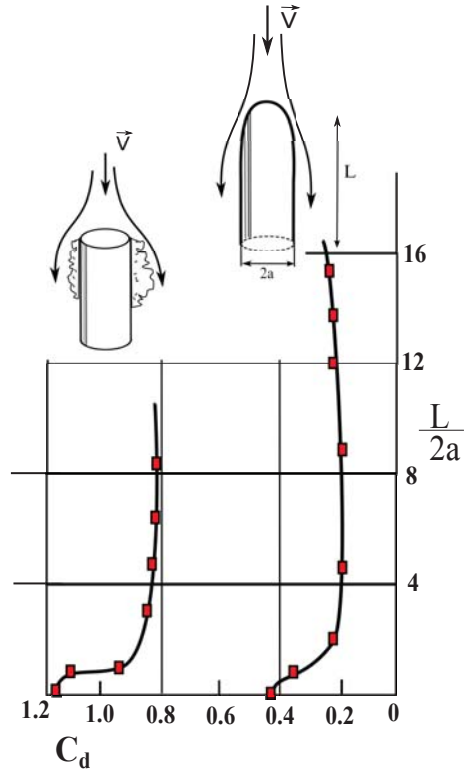


Fig. 8. Drag coefficients of blunt-nosed and rounded-nosed cylinders versus fineness ratio $L/2a$ [10].

negative feedback mechanisms, the new equilibrium power will be given by [14]:

$$P(\infty) = P^* + 2\dot{m}_c c_c \left[\frac{\rho_o}{\gamma_c} \right] \quad (21)$$

Using the power $P(\infty)$ we can calculate the coolant and fuel temperatures from Eqs. (13) and (14) and then their respective densities. This then allows us to find the value of the buoyancy parameter B given by Eq. (20). For example, taking $\Delta T = \rho_o/\gamma_c \approx 5$ K, we obtain an approximate value for $B \approx 0.1$. The drag coefficient C_d will be between 1.2 for a blunt-nosed cylinder or 0.2 for a rounded nose, as shown in Fig. 8 [10]. Thus, for an optimized fuel rod with a rounded end-cap, as depicted in Fig. 9, we can assume $C_d = 0.2$, and with a total fuel rod length (including plenums) of 210 cm (see Table 2), we obtain a terminal velocity of $V_t \approx 1.44$ m s⁻¹.

Finally, we need an estimate of the variation of reactivity with the dis-

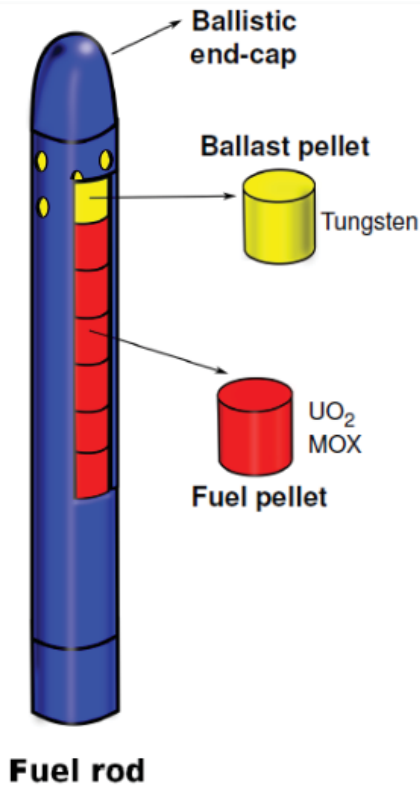


Fig. 9. A possible optimized fuel rod design for a lead or lead-bismuth cooled reactor. The end-cap is rounded to enhance the ejection velocity.

placement of the ejected fuel rod, i.e. $\partial\rho/\partial z$. Unfortunately, this is a highly
 240 uncertain parameter; its accurate computation requires knowledge of the
 specific location at which the fuel rod ejection occurs, as well as the spe-
 cific design of the rod. A calculation performed using the SCALE 6 software
 [6] for a typical fuel rod channel, using lead as the coolant and reflective
 245 boundary conditions is translate into assuming a homogeneous core in which all
 the elements of the core have ballast pellet and also all are moving which is
 within the linear fuel management model in which is assumed that the total
 core reactivity is the summation of the reactivity of each element. The use of
 this linear model, although admittedly is rather simplified, nevertheless will
 250 allow to provide with a conservative value of $\partial\rho/\partial z \approx -50 \text{ pcm cm}^{-1}$. Then
 using our previously calculated estimate of the fuel rod terminal velocity, we
 have a rate of negative insertion on -7200 pcm s^{-1} . Taking a typical posi-
 tive coolant temperature coefficient of reactivity to be 0.36 pcm K^{-1} [20] and
 $\Delta T = 5 \text{ K}$, the time needed for the buoyancy-driven mechanism to control
 255 this transient will be a tiny fraction of a second once the Eq.(10) is met.

Thus, the foregoing calculations indicate that by using a modest fraction
 ($\sim 15\%$) of tungsten ballast pellets the fuel rod will be endowed with a reliable
 self-ejection mechanism during temperature transients. It should be noted
 that, in these preliminary calculations, other components of the fuel rod

260 which can reduce its effective density even more were neglected, the most
important being the gas plenum chambers (if fission gases are not vented
directly into the coolant). However, the potential reduction in the effective
fuel rod density due to the gas plenums can be compensated by using tungsten
rather than stainless steel for the lower and upper plenums in the fuel rod.

265 3. Conclusions

In this paper we explored the possibility of taking advantage of buoy-
ancy forces in heavy liquid metal cooled fast reactors to endow the fuel rod
with a reliable and passive negative feedback mechanism through fuel rod
ejection (a fuel self-disassembly mechanism) during a temperature transient,
270 compensating the positive coolant temperature coefficient of reactivity that
some fast reactors feature. It was deduced that, through the use of tung-
sten ballast pellets introduced into the fuel rod design, such a mechanism is
feasible, with the volume occupied by the ballast pellets being less than 15%.

This preliminary assessment was based on unavoidable idealizations, some
275 conservative and others non-conservative. It should not be misconstrued
as a definitive, detailed analysis. Additional R&D is required to further
explore the possibilities of this concept, to seek optimal values for the design
variables, and to determine real practical applicability as details are refined.
Only then will the feasibility of the proposed concept be fully established.

280 4. Appendix

•The use of gas plenum.

In our previous calculations, the gas plenum was not taken into account. The
gas plenum length could be typically the same of the active length, and then
an extra fraction of dedicated ballast would be necessary to avoid buoyancy.
285 Because the density of the fission gas can be neglected in comparison with
the ballast, then it is easy to see that the fraction of volume of gas plenum
with ballast required should be $\approx \frac{\rho_c}{\rho_w}$ where ρ_c and ρ_w are the density of
the coolant and the ballast, respectively. Nevertheless, there is an actual
trend to eliminate fission gas plenums and to vent fission product gases to
the primary coolant system in a controlled manner,[1],[9]. Venting the fuel
290 pins enables deep burnups required to sustain the core, for over 40 years and
greatly reduces the probability of cladding failures. In addition, in lead cooled
and even more in lead bismuth cooled reactors the use of gas plenums and
the retention of fission gas plenums have been recently questioned,

295 **Nomenclature**

- B = buoyancy parameter defined by Eq. (20)
 C_d = drag coefficient
 c_i = heat capacity of material i
 F_f = volume fraction of fuel
300 g = acceleration due to gravity
 h = heat transfer coefficient
 L = length (of fuel pin or fuel rod)
 \dot{m}_c = coolant mass flow
 M_f = mass of fuel
305 P = pin power
 P^* = pin power at onset of rod ejection
 r = radius
 R_f = thermal resistance of fuel pin
 t = time
310 T = temperature
 T_i = inlet temperature of coolant
 V_t = terminal velocity
 z = vertical coordinate

315 **Greek symbols**

- α_i = rate of change of density of material i with temperature
 β = fraction of delayed neutrons
 γ_c = coolant temperature coefficient of reactivity
 κ_i = thermal conductivity of material i
320 ρ_i = density of material i
 ρ = reactivity

Subscripts

- c = coolant
325 f = fuel
 g = gap
 s = stainless steel
 w = tungsten

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