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LOCO - A LINEARISED MODEL FOR ANALYSING THE ONSET
OF COOLANT OSCILLATIONS AND FREQUENCY
RESPONSE OF BOILING CHANNELS

by

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(*CSIRO Mineral Physics Division)

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ABSTRACT

Industrial plant such as heat exchangers, thermosiphon reboilers, and nuclear and conventional boilers are prone to coolant flow oscillations (or density wave oscillations) which may not be detected

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(Continued)

until some catastrophic failure (e.g. boiler tube burnout, corrosion or rupture) terminates normal operation.

In this report, a hydrodynamic model is formulated in which the one-dimensional, non-linear, partial differential equations for the conservation of mass, energy and momentum are perturbed with respect to time, linearised (that is, terms higher than first order are neglected), and Laplace-transformed into the s-domain for frequency response analysis. A computer program has been developed to integrate numerically the resulting non-linear ordinary differential equations by finite difference methods. A sample problem demonstrates how the computer code is used to analyse the frequency response and flow stability characteristics of a heated channel.

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L CODES; FLOW MODELS; TWO-PHASE FLOW; HYDRODYNAMICS; DIFFERENTIAL EQUATIONS; FINITE DIFFERENCE METHOD; FREQUENCY ANALYSIS

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1. INTRODUCTION

The performance and reliability of nuclear and conventional steam generators, process heat exchangers, thermosiphon reboilers and other two-phase heat transfer equipment are dependent on the steady flow of coolant through their component heated channels. Under certain operating conditions (coolant pressures, temperatures, flow rates, heat fluxes etc.), a variety of fundamental and higher mode coolant flow instabilities can occur which may cause:

- a) failure of the channel structure or containment due to the high thermal stresses or corrosion induced by thermal cycling, dryout or burnout of the heat transfer surfaces;
- b) fatigue failure of the channel structure or containment due to the excitation of mechanical resonances;
- c) control problems in maintaining the desired operating conditions, particularly the steam conditions supplied to subsequent process equipment (e.g. steam turbine).

The following analysis formulates a linearised hydrodynamics model for calculating the frequency response of a 'boiling channel' and, in particular, for predicting the onset of static (flow excursions) and dynamic (flow or density wave oscillations) instabilities in the coolant flow. The analysis is prefaced by a review of the basic regions and two-phase flow regimes that are postulated to occur in the boiling channel, and the possible methods

for modelling the coolant flow behaviour.

The model is based on a linearised solution of the cross-section averaged equations for the conservation of mass, energy and momentum of the two-phase (vapour/liquid) mixture. The equations are perturbed with respect to time, linearised (that is, higher than first-order terms are neglected), and Laplace-transformed into the s-domain for frequency response analysis. The resulting non-linear ordinary differential equations are spatially integrated by finite difference methods. Finite difference equations are derived for the steady state and dynamic analysis of the thermal non-equilibrium two-phase flow through the subcooled boiling region of the channel. Specification of certain parameters and variables reduces these equations to simpler forms for analysis of the single phase liquid and saturated boiling regions.

The channel is divided into a number of volume elements, and the continuity in flow conditions between successive elements is provided by coupling equations at the boundary (or interface). In a similar manner, the flow conditions between the single phase and two-phase regions are coupled by mass, energy and momentum conservation equations at the boiling boundary. The steady state and dynamic relationships between variables at the channel inlet and exit are calculated by piecewise integration along the channel. The channel may be operated in parallel with a number of other channels, in parallel with a large bypass, or in a closed loop in series with a pump. Stability criteria are derived by specification of appropriate boundary conditions for the parallel channel, channel/bypass and

closed loop systems, and these enable the stability of the coolant flow through the boiling channel to be assessed.

A computer program (LOCO - a Linearised model for analysing the Onset of Coolant Oscillations in boiling channels) has been developed to perform the computations. The program data input specifies the channel geometry and operating conditions, and permits the calculation of the frequency response relationships between perturbations in channel inlet flow rate, temperature, pressure, average power (inputs) and the over pressure drop, exit void fraction (outputs). The inputs may be frequency independent ('white noise') for transfer function and stability calculations, or frequency dependent real or complex (experimental) data for frequency response calculations.

This analysis is part of a larger investigation which is described in more detail in the references by Romberg (1978) and Romberg and Rees (1980).

2. FLOW REGIONS AND REGIMES IN BOILING TWO-PHASE FLOW

During normal operation, the coolant enters the heated channel as subcooled liquid and discharges as a vapour/liquid mixture or superheated vapour from the channel exit. For modelling purposes the flow structure (vapour/liquid distribution) may be divided into five inter-connected hydrodynamic regions (Figure 1):

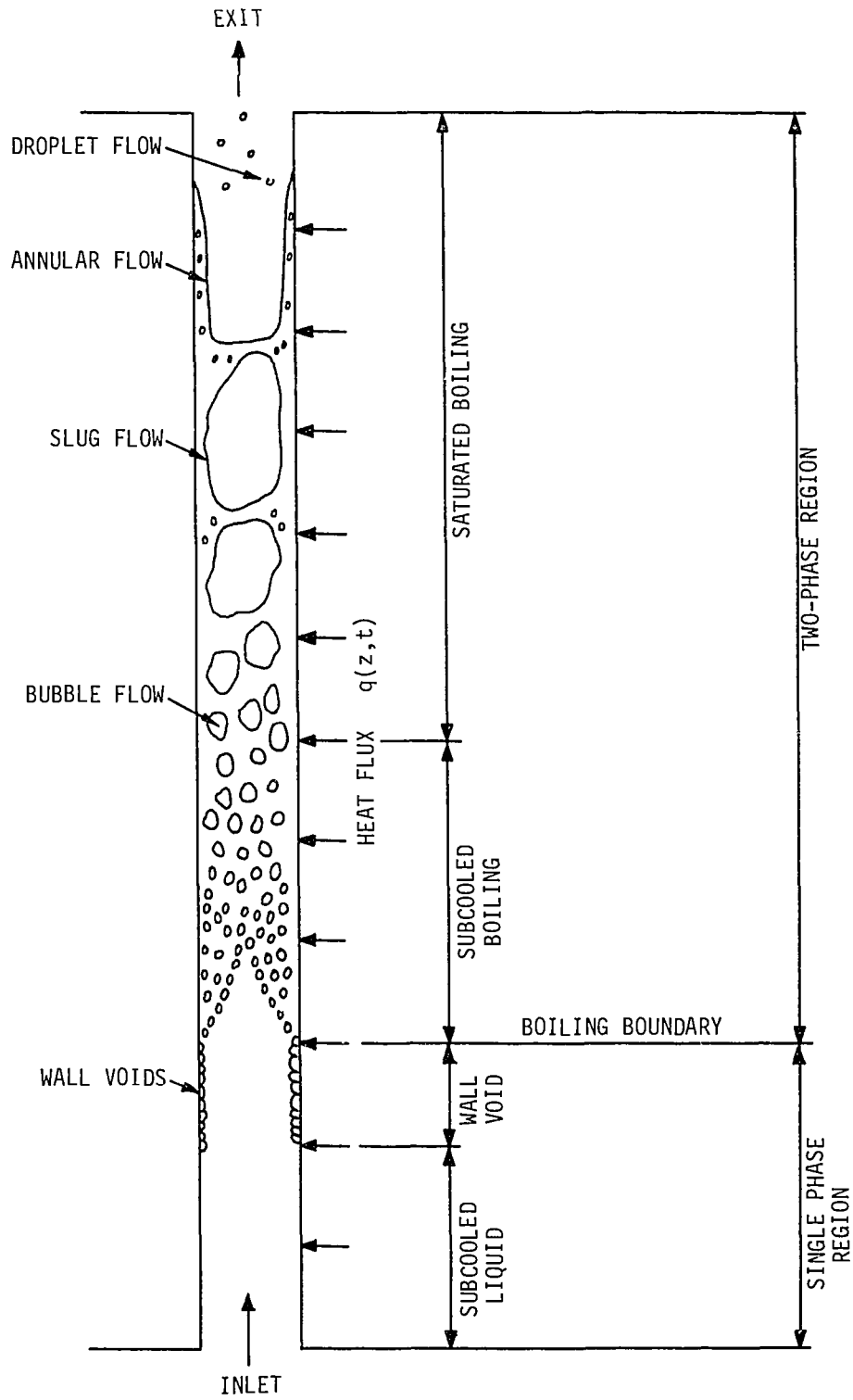


FIGURE 1 - SCHEMATIC DIAGRAM OF A BOILING CHANNEL SHOWING FLOW REGIONS AND TWO-PHASE REGIMES

1. single phase liquid region - no vapour bubbles, incompressible flow;
2. wall void region - the local wall temperature is greater than the local saturation temperature, vapour bubbles form and remain attached to the heated surface;
3. subcooled boiling region - the bulk (average) liquid temperature is less than the local saturation temperature, the vapour bubbles become detached from the wall due to thermal effects (low flow rates) or hydrodynamic drag effects (high flow rates);
4. saturated boiling region - the liquid temperature reaches saturation, violent boiling agitation at the heated surface;
5. superheated vapour region - no liquid phase, compressible vapour flow.

The liquid entering the channel has its temperature raised by heat transfer at the heat source/coolant interface. The rate of heat transfer depends on the cross-sectional velocity and temperature profiles. Since the bulk (cross-sectional average) liquid temperature is less than the saturation temperature, it is also called the subcooled liquid region. The local wall temperature is higher than the local liquid temperature, but not high enough to initiate nucleate boiling on the heated surface.

As a volume element of coolant moves further along the channel, the local conditions at the heated surface (heat fluxes, wall/coolant temperatures, surface asperities, etc.) are sufficient to initiate nucleate boiling, but the bubble diameters are small and are dominated by surface tension effects, so that they remain attached to the heated surface. These bubbles act as surface asperities to increase the frictional resistance to fluid motion, and improve the wall/coolant

heat transfer by their agitation and action as turbulence promoters. This region is termed the wall void region. Further downstream the bubbles grow in size and eventually become detached from the heated surface by thermal agitation or hydrodynamic action. The axial position at which this occurs denotes the subcooled boiling boundary, the beginning of the subcooled boiling region.

In the subcooled boiling region, the bubbles travel downstream near the wall or move, by turbulent interaction, into the centre of the channel, where they collapse due to thermal non-equilibrium with the subcooled liquid phase. The mean lifetime of the bubbles in this region is statistically random, and their growth and collapse can cause pressure transients (acoustic oscillations) which vibrate the channel structure. In general, the mean lifetime of a bubble ensemble is dependent on their mean free path, boundary layer profiles, number of nucleation sites etc. These are nonmeasurable quantities, and the axial void (vapour fraction) distribution in this region must be evaluated by empirical correlations based on experimental data. These correlations are a function of the channel geometry and operating conditions (pressures, temperatures, flow rates, power inputs etc.).

Further downstream, turbulent mixing of the vapour and liquid phases ultimately raises the liquid temperature to saturation, and this is the commencement of the saturated boiling region. In contrast to the subcooled boiling region, the vapour and liquid phases in the saturated boiling region are in thermal equilibrium (neglecting the slight superheat of the vapour phase), and the bubbles grow in size and no longer collapse. The heat input in this region is entirely dissipated in producing more voids, whereas the heat input to the subcooled boiling region is dissipated in raising the liquid temperature and producing more voids.

The flow structure in the two-phase region is divided into flow

regimes, principally bubble, slug, annular and droplet flow regimes (Tong 1965). Combinations of these basic regimes are possible. The vapour bubbles merge as they move downstream to form slugs of vapour. These slugs may also merge to form larger slugs, until the channel cross-section consists of a vapour core (with droplet entrainment), and liquid at the solid surface. This is termed annular flow. With further heating, the liquid phase at the wall disappears altogether, and is only present as droplets or mist in the vapour core. Ultimately the liquid phase disappears entirely, and only superheated vapour is left (superheat region). The superheat region is neglected in the present analysis, but may be approximately modelled with unit quality (vapour mass fraction) flow conditions.

3. ANALYSIS CONSIDERATIONS AND ALTERNATIVE METHODS

The previous description of the boiling channel flow structure gives some indication of the difficulties involved in the development of suitable models. Most analyses solve the partial differential equations for the conservation of mass, energy and momentum in the single phase and two-phase regions of the channel assuming one-dimensional turbulent flow. A detailed mathematical model of the two-phase flow structure requires six conservation equations (one for each phase), plus constitutive equations to define the fluid properties of the system, and mass, energy and momentum transfer relationships at the vapour/liquid boundaries (Bird et al. 1960, Govier and Aziz 1972, Boure et al. 1973, Boure 1975). However, since detailed transfer relationships (laws) are still not available, the use of six conservation equation models is not yet possible, and they are only useful in clarifying the assumptions of existing simplified models (Boure 1975). In recent years, there has been considerable emphasis on the development of mathematically consistent two-phase flow models, particularly for application to thermal non-equilibrium and non-linear

flow conditions (limit cycle and acoustic oscillations, rapid depressurisation transients), in which the use of empirical steady state correlations to replace the vapour/liquid transfer laws is suspect.

The traditional methods for analysing the excursive and oscillatory flow stability of boiling channels are schematically summarised in Figure 2. The mass, energy and momentum conservation equations are averaged over the channel flow area, and are spatially integrated by analytical or piecewise methods using constitutive equations to define local values for the fluid properties, heat transfer, subcooled boiling, two-phase friction and slip ratio. The analytical methods are dependent on simplifying assumptions (e.g. invariant fluid properties, homogeneous flow, no slip between vapour and liquid phases, etc.) to enable integration of the conservation equations in the two-phase region of the channel. These assumptions, however, restrict the models to a limited range of applications. The piecewise methods are less restrictive, and are based on the subdivision of the channel into small volume elements and the replacement of the spatial derivatives by their first order finite difference equivalents.

The alternative dynamic models depicted in Figure 2 result from the different methods for integrating the time dependent terms. The most important models are the non-linear time domain (NLTD) and linear frequency domain (LFD) distributed models, in which the time derivative is eliminated respectively by direct numerical integration of the non-linear differential equations, or by small perturbation methods and Laplace transformation of the linear equations for transfer function analysis in the frequency domain. The relative merits of NLTD and LFD models may be assessed from the following observations (see reviews by Neal and Zivi 1965 and Grumbach 1969):

a) Discrepancies between theoretical and experimental flow instability thresholds are primarily caused by inadequate mathematical

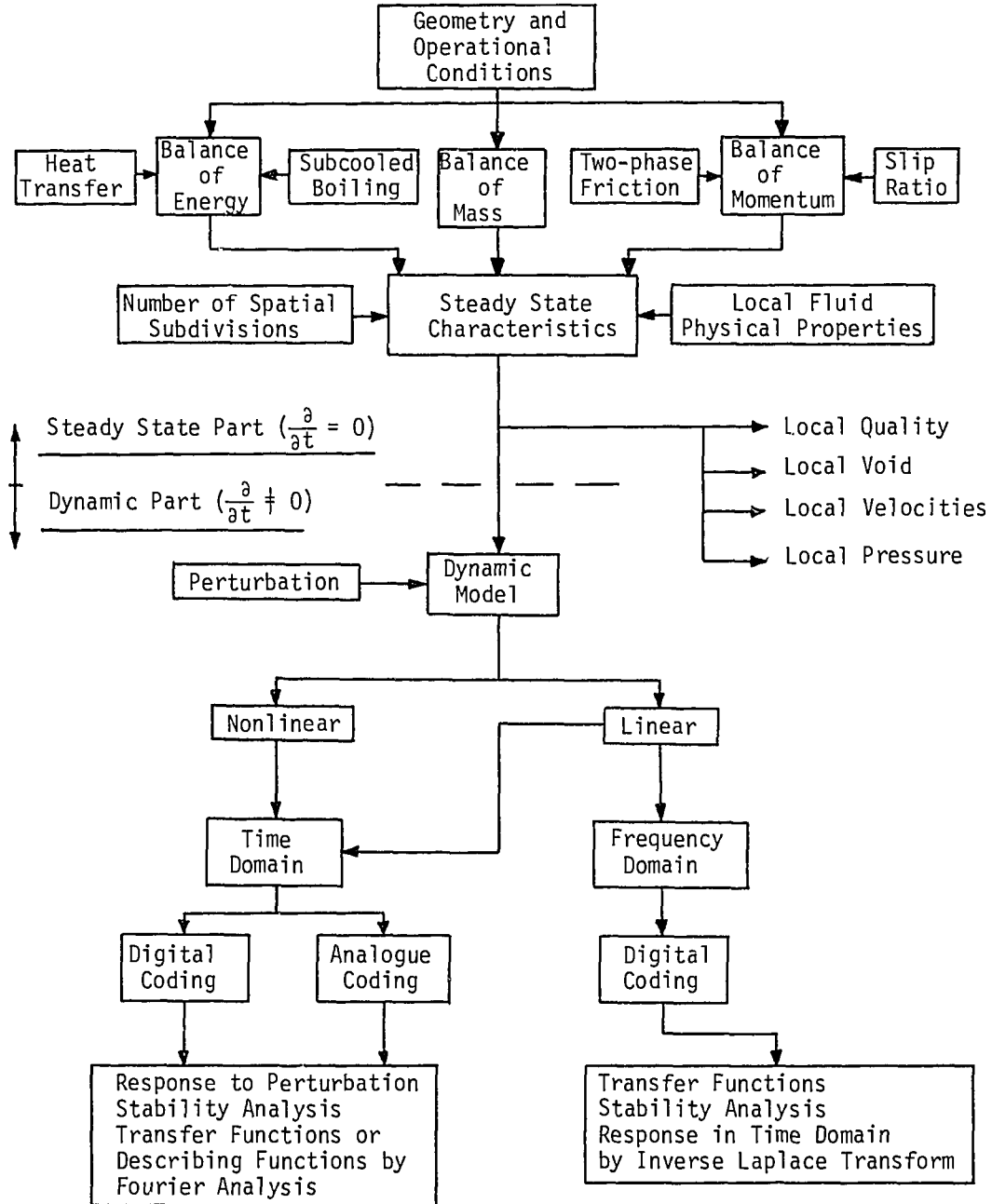


FIGURE 2 - POSSIBLE METHODS FOR BOILING CHANNEL CALCULATIONS

formulations, particularly numerical instabilities in the dynamic calculations. The NLTD models require small time steps to satisfy numerical stability constraints, and take between one and two orders of magnitude longer in computing time than LFD models (Spinks 1972).

b) Discrepancies between theoretical and experimental flow instability thresholds are reduced by matching the steady state characteristics (especially flow rate and/or pressure drop). In some models this can be achieved by adjustment of the coefficients in the empirical correlations.

c) Both models predict the local stability of the boiling channel, and NLTD models are also able to predict the amplitude and period of non-linear limit cycle oscillations beyond the instability threshold. However the empirical correlations are designed to model steady state conditions, and their validity for non-linear dynamic analysis is suspect. The LFD models assume the perturbed variables fluctuate about time invariant mean (steady state) values, and thus small perturbation of an arbitrary function or correlation $f(x,y)$ is linearly related to the steady state by the first order partial differential equation

$$f^*(x,y) = \left(\frac{\partial f^0}{\partial x^0} \right) x^* + \left(\frac{\partial f^0}{\partial y^0} \right) y^*$$

where the superscripts (o) and (*) denote respectively the steady state and perturbed variables.

Neal and Zivi (1965) found that the LFD model by Jones (1961) proved to be the most accurate in predicting the flow instability thresholds of experimental data, and for this type of analysis the greater computational efficiency of LFD models is an important consideration. Some NLTD models, however, are able to assess the channel flow stability during start-up and shut-down operation of the

plant, and are useful benchmark models for evaluation of non-linear lumped parameter control models.

4. CONSERVATION EQUATIONS AND MODEL ASSUMPTIONS

In the present analysis the mass, energy and momentum conservation equations for two-phase flow are solved assuming one-dimensional turbulent flow. The equations are perturbed, linearised and Laplace transformed for frequency domain analysis. The linearised differential equations are spatially integrated by finite difference methods. Constitutive equations define the local fluid properties, heat transfer, friction, subcooled boiling and slip correlations.

The two-phase conservation equations of mass, energy and momentum averaged over the channel flow area are respectively

$$\text{Mass: } \frac{\partial}{\partial z} \langle \rho_l (1-\alpha) U_l + \rho_g \alpha U_g \rangle + \frac{\partial}{\partial t} \langle \rho_l (1-\alpha) + \rho_g \alpha \rangle = 0 \quad (1)$$

$$\begin{aligned} \text{Energy: } & \frac{\partial}{\partial z} \langle \rho_l (1-\alpha) U_l h_l + \rho_g \alpha U_g h_g \rangle \\ & + \frac{\partial}{\partial t} \langle \rho_l (1-\alpha) h_l + \rho_g \alpha h_g - \alpha p \rangle = \langle Q \rangle \end{aligned} \quad (2)$$

$$\begin{aligned} \text{where } \langle Q \rangle &= - \frac{1}{A} \int_0^R \frac{1}{r} \frac{\partial}{\partial r} (r q_r) \cdot 2\pi r \, dr \\ &= \frac{2\pi R}{A} q_w = \frac{P_H}{A} q_w \quad (\text{single tube}) \\ &= \frac{1}{A} \sum_{j=1}^N P_{Hj} q_{wj} \quad (\text{fuel rod cluster}) \end{aligned}$$

and kinetic energy, potential energy, viscous heating, and axial conduction effects are neglected;

$$\begin{aligned} \text{Momentum: } & \frac{\partial}{\partial z} \langle \rho_l (1-\alpha) U_l^2 + \rho_g \alpha U_g^2 \rangle + \frac{\partial}{\partial t} \langle \rho_l (1-\alpha) U_l + \rho_g \alpha U_g \rangle \\ & + K_f \frac{\rho_l \langle U_l \rangle^2}{2} \phi^2 + \langle \rho_l (1-\alpha) + \rho_g \alpha \rangle g \cos \theta = - \frac{\partial}{\partial z} \langle p \rangle \end{aligned} \quad (3)$$

where the single phase friction factor (K_f) and the two-phase local friction multiplier (ϕ^2) are defined in Appendix A.

In the above equations the brackets <-----> denote the cross-section averaging operation

$$\langle F \rangle = F(z,t) = \frac{1}{A} \int_A F(r,z,t) dA \quad (4)$$

The product terms in the brackets <-----> cannot be readily determined from experimental data, and are usually assumed to be approximated by the product of the cross-section averaged individual variables, for example,

$$\begin{aligned} \langle \rho_l (1-\alpha) U_l^2 + \rho_g \alpha U_g^2 \rangle &= \langle G_l U_l + G_g U_g \rangle \\ &\approx \langle G_l \rangle \langle U_l \rangle + \langle G_g \rangle \langle U_g \rangle \end{aligned}$$

where the liquid, vapour and mixture mass fluxes are respectively

$$\left. \begin{aligned} \langle G_l \rangle &= \langle \rho_l (1-\alpha) U_l \rangle = \langle (1-x)G \rangle \\ \langle G_g \rangle &= \langle \rho_g \alpha U_g \rangle = \langle xG \rangle \\ \langle G \rangle &= \langle \rho_l (1-\alpha) U_l + \rho_g \alpha U_g \rangle = \langle G_l \rangle + \langle G_g \rangle \end{aligned} \right\} \quad (5)$$

This approximation is reasonable for fully developed turbulent flow in which the radial profiles of the variables are relatively uniform in the central region of the channel.

The conservation equations (1) to (3) can thus be simplified to

$$\text{Mass: } \frac{\partial}{\partial z} \{G\} + \frac{\partial}{\partial t} \{\rho_l (1-\alpha) + \rho_g \alpha\} = 0 \quad (6)$$

$$\text{Energy: } \frac{\partial}{\partial z} \{G_l h_l + G_g h_g\} + \frac{\partial}{\partial t} \{\rho_l (1-\alpha) h_l + \rho_g \alpha h_g - \alpha p\} = Q \quad (7)$$

$$\begin{aligned} \text{Momentum: } \frac{\partial}{\partial z} \{G_l U_l + G_g U_g\} + \frac{\partial}{\partial t} \{G\} + \frac{K_f}{2} \frac{G^2 \phi^2}{\rho_l} \\ + \{\rho_l (1-\alpha) + \rho_g \alpha\} g \cos \theta = - \frac{\partial p}{\partial z} \end{aligned} \quad (8)$$

in which the brackets <-----> have been omitted, and the symbols denote cross-section averaged variables, as in the notation in equation (4).

The mixture conservation equations (6) to (8) are solved as follows:

1. the channel variables (pressures, mass fluxes, void fraction, etc.) are perturbed with respect to time, and the product terms linearised (terms higher than first order are neglected);

2. the perturbed variables are transformed into the frequency domain by Laplace transformation,

$$F(z,s) = \tilde{F}(z) = \int_0^{\infty} F^*(z,t) e^{-st} dt \quad (9)$$

where $F(z,t) = F^0(z) + F^*(z,t)$, $s (= \sigma + j\omega)$ is the complex Laplace operator, and zero initial conditions are assumed;

3. the first order differential equations are integrated spatially over a small space node (axial length Δz x channel cross-sectional area) using the forward difference equation

$$\int_0^{\Delta z} \frac{\partial \tilde{F}(z)}{\partial z} dz \approx \tilde{F}_{i+1} - \tilde{F}_i \quad (10)$$

and the trapezoidal rule,

$$\int_0^{\Delta z} \tilde{F}(z) dz \approx \frac{\Delta z}{2} (\tilde{F}_{i+1} + \tilde{F}_i) \quad (11)$$

4. the fluid properties and heat flux at the wall/coolant interface are assumed spatially invariant within the space node, and all fluid properties except the vapour density (ρ_g) are assumed time invariant.

5. FINITE DIFFERENCE EQUATIONS

In the following analysis the finite difference equations for thermal non-equilibrium two-phase (vapour/liquid) flow are derived. The equations apply to a single space node in the subcooled boiling region of the channel. By specifying appropriate values for certain variables, the same equations are applicable to the single phase

liquid, wall void and saturated boiling regions of the channel. The analysis is based on a special form of the energy equation which is derived by manipulation of the mass and energy equations (6), (7) in conjunction with a subcooled boiling liquid enthalpy correlation. The mixture enthalpy is used as the fundamental variable rather than the vapour void fraction, which is usually obtained by solution of a void propagation equation (Zuber et al 1967, Zuber and Staub 1966, Zuber and Staub 1967, Hancox and Nicoll 1971). If the vapour void fraction is used as the fundamental variable then separate finite difference equations are required for the single phase and two-phase regions of the channel, rather than the single set of equations as derived in the present analysis.

5.1 Subcooled Boiling Region Energy Equation

The energy equation (7) may be rewritten as

$$\frac{\partial}{\partial z} \{Gh\} + \frac{\partial}{\partial t} \{\rho_l(1-\alpha) h_l + \rho_g \alpha h_g\} = Q + \alpha \frac{\partial p}{\partial t} \quad (12)$$

where the liquid and vapour mass fluxes are replaced by their respective equations from (5), and the mixture enthalpy is defined as

$$h = (1-x)h_l + x h_g \quad (13)$$

The mass equation (6) may be rewritten as

$$\frac{\partial}{\partial z} \{G\} - \frac{\partial}{\partial t} \{\Delta\rho\alpha\} = 0 \quad (14)$$

and substitution in the energy equation (12) yields

$$\begin{aligned} G \frac{\partial}{\partial z} \{h\} - \Delta\rho\alpha \frac{\partial}{\partial t} \{h\} + \frac{\partial}{\partial t} \{\rho_l h_l + (\rho_g h_g - \rho_l h_l + h\Delta\rho)\alpha\} \\ = Q + \alpha \frac{\partial p}{\partial t} \end{aligned} \quad (15)$$

The vapour void fraction is eliminated from the third term on the left side of (15) by noting that, from (13) and (C.16), Appendix C,

$$\rho_l h_l + (\rho_g h_g - \rho_l h_l + h\Delta\rho)\alpha = \rho_l \{(1-K_0)h_l + K_0 h\}$$

and thus (15) becomes

$$G \frac{\partial}{\partial z} \{h\} + (K_0 \rho_\ell - \Delta \rho \alpha) \frac{\partial}{\partial t} \{h\} + \rho_\ell (1-K_0) \frac{\partial}{\partial t} \{h_\ell\} = Q + \alpha \frac{\partial p}{\partial t} \quad (16)$$

where the Bankoff factor K_0 , (C.11) Appendix C, is assumed to be space and time invariant within the space node.

The liquid enthalpy in the subcooled boiling region is related to the mixture enthalpy by either an exponential or a hyperbolic tangent subcooled boiling correlation as follows (Zuber et al 1967, Kroeger and Zuber 1968):

Exponential correlation -

$$h_\ell^+ = \frac{h_\ell - h_0}{h_f - h_0} = 1 - \exp(-h^+) \quad (17)$$

Hyperbolic tangent correlation -

$$h_\ell^+ = \tanh(h^+) \quad (18)$$

where
$$h^+ = \frac{h - h_0}{h_f - h_0}$$

in which h_0 is the liquid enthalpy at the subcooled boiling boundary.

Equations (17) and (18) satisfy the boundary conditions

$h_\ell = h_0$ at the subcooled boiling boundary ($z = z_0$), and

$h_\ell = h_f$ at the saturated boiling boundary.

Hence, in equation (16), the liquid enthalpy is related to the mixture enthalpy by

$$\frac{\partial}{\partial t} \{h_\ell\} = \frac{\partial h_\ell^0}{\partial h^0} \frac{\partial}{\partial t} \{h\} = \frac{\partial h_\ell^+}{\partial h^+} \frac{\partial h}{\partial t} = F^+ \frac{\partial h}{\partial t} \quad (19)$$

$$\left. \begin{array}{l} \text{where } F^+ = \exp(-h^+) \quad (\text{Exponential correlation}) \\ \text{and } F^+ = 1 - \tanh^2(h^+) \quad (\text{Tanh correlation}) \end{array} \right\} \quad (20)$$

Hence equation (16) simplifies to

$$G \frac{\partial h}{\partial z} + \{\rho_\ell (K_0 + (1-K_0) F^+) - \Delta \rho \alpha\} \frac{\partial h}{\partial t} = Q + \alpha \frac{\partial p}{\partial t} \quad (21)$$

5.2 Subcooled Boiling Region Steady State Equations

Mass flux: Setting the time derivative term in (6) or (14) to zero yields

$$\frac{d}{dz} \{G^0\} = 0 \quad (22)$$

which, on integration, gives

$$G_2^0 = G_1^0 = G^0 \quad (\text{a constant}) \quad (23)$$

Enthalpy: The steady state form of (21) is

$$G^0 \frac{d}{dz} \{h^0\} = Q^0 \quad (24)$$

which, on integration, gives

$$h_2^0 = h_1^0 + \frac{Q^0 \Delta z}{G^0} \quad (25)$$

The liquid enthalpy h_l^0 is calculated from the correlations in equations (17) or (18), and the other variables of interest are calculated as follows:

$$(13) \rightarrow x_2^0 = \frac{h_2^0 - h_{l2}^0}{h_g^0 - h_{l2}^0} \quad (26)$$

(C.16), Appendix C \rightarrow

$$\alpha_2^0 = \frac{K_o \rho_l x_2^0}{\rho_l + \Delta \rho x_2^0} \quad (27)$$

$$(5) \rightarrow \left. \begin{aligned} G_{l2}^0 &= (1-x_2^0) G_2^0 \\ G_{g2}^0 &= x_2^0 G_2^0 \\ U_{l2}^0 &= \frac{G_{l2}^0}{\rho_l (1-\alpha_2^0)} \\ U_{g2}^0 &= \frac{G_{g2}^0}{\rho_g \alpha_2^0} \end{aligned} \right\} (28)$$

Pressure drop: From equation (8)

$$\begin{aligned} \frac{d}{dz} (G_l^0 U_l^0 + G_g^0 U_g^0) + \frac{K_f^0}{2} \frac{G^0 \phi^0}{\rho_l} + (\rho_l (1-\alpha^0) + \rho_g \alpha^0) g \cos \theta \\ = - \frac{dp^0}{dz} \end{aligned} \quad (29)$$

and integration gives

$$\Delta p_2^0 = p_1^0 - p_2^0 = \Delta p_{2m}^0 + \Delta p_{2f}^0 + \Delta p_{2g}^0 \quad (30)$$

where the momentum, friction and gravitational pressure drop components are respectively

$$\begin{aligned} \Delta p_{2m}^0 &= G_{l2}^0 U_{l2}^0 - G_{g2}^0 U_{g2}^0 \\ \Delta p_{2f}^0 &= \frac{K_f^0}{2} \frac{G_2^0}{\rho_l} \int_1^2 \phi^0{}^2 dz = \Delta p_{1f}^0 \frac{(\phi_1^0{}^2 + \phi_2^0{}^2)}{2} \\ \Delta p_{2g}^0 &= g \cos \theta \int_1^2 (\rho_l - \Delta p \alpha^0) dz \\ &= \Delta z \left\{ \rho_l - \Delta p \frac{(\alpha_1^0 + \alpha_2^0)}{2} \right\} g \cos \theta \end{aligned} \quad (31)$$

5.3 Subcooled Boiling Region Dynamic Equations

Perturbation, linearisation and Laplace transformation of equation (14) yields

$$\frac{d}{dz} \{ \tilde{G} \} = s \{ \Delta p \tilde{\alpha} - \alpha^0 \rho_g' \tilde{p} \} \quad (32)$$

The perturbed void fraction $\tilde{\alpha}$ is replaced by the perturbed mixture enthalpy \tilde{h} . From (C.18), Appendix C, the perturbed void fraction is

$$\tilde{\alpha} = \frac{1}{X} \{ \tilde{h} - x^0 (1-x^0) \frac{\rho_g'}{\rho_g} \tilde{p} \} \quad (33)$$

and, from equations (14) and (19), the perturbed flow quality in terms of the mixture enthalpy is

$$\tilde{x} = \frac{[1 - (1-x^0) F^+] \tilde{h}}{(h_g - h_l^0)} = \frac{\Gamma^0 \tilde{h}}{Q^0} \quad (34)$$

where Γ^0 is a vapour source term, and is given by

$$\Gamma^0 = \frac{Q^0}{h_g - h_l^0} [1 - (1-x^0) F^+] \quad (35)$$

Hence equation (32) becomes

$$\frac{d}{dz} [\tilde{G}] = sa\tilde{h} - sb\tilde{p} \quad (36)$$

where

$$a = \frac{\Delta\rho \Gamma^0}{X Q^0} = \frac{\Delta\rho}{X(h_g - h_l^0)} [1 - (1-x^0) F^+] \quad (37)$$

and

$$b = \rho_g' \left[\alpha^0 + \frac{\Delta\rho x^0 (1-x^0)}{X \rho_g} \right]$$

Integration of (36) gives

$$\tilde{G}_2 = \tilde{G}_1 + \frac{s\Delta z}{2} \{a_1 \tilde{h}_1 + a_2 \tilde{h}_2 - (b_1 + b_2) \tilde{p}_1\} \quad (38)$$

Similarly, perturbation, linearisation and Laplace transformation of the energy equation (21) yields

$$Q^0 \frac{d}{dz} \{\tilde{h}\} + c\tilde{h} = Q^0 \left\{ \frac{\tilde{Q}}{Q^0} - \frac{\tilde{G}}{G^0} \right\} + s\alpha^0 \tilde{p} \quad (39)$$

which has been simplified using (24), and where

$$c = \rho_l \{K_0 + (1-K_0) F^+\} - \Delta\rho\alpha^0 \quad (40)$$

From (B.19), Appendix B, and noting that $\tilde{h}_l = F^+ \tilde{h}$,

$$\frac{\tilde{Q}}{Q^0} = \frac{\tilde{q}_w}{q_w^0} = (1 + \tau_{HS})^{-1} \left[\frac{\tilde{Q}_s}{Q_s^0} + b_t \tau_{HS} \frac{\tilde{G}}{G^0} - F^+ \tau_{HS} \frac{\tilde{h}}{\Delta h_w^0} \right] \quad (41)$$

where

$$\Delta h_w^0 = C_{pl} \Delta T_w^0 = C_{pl} (T_w^0 - T_l^0)$$

Substitution of (41) in (39) and integrating gives, in conjunction with (38)

$$\tilde{h}_2 = \frac{\left(1 - sT_1(s)\right)\tilde{h}_1 + \left(\frac{\Delta h^0}{1 + \tau_H s}\right)\frac{\tilde{Q}_s}{Q_s^0} - P(s)\frac{\tilde{G}_1}{G^0} + s\bar{\tau}_p(s)\tilde{p}_1}{\{1 + sT_2(s)\}} \quad (42)$$

$$\begin{aligned} \text{where } P(s) &= \Delta h^0 \left[\frac{1 + (1 - b_t)\tau_H s}{1 + \tau_H s} \right] \\ T(s) &= \frac{\Delta z}{2G^0} \left[c + \frac{Q^0 F^+ \tau_H}{\Delta h_w^0 (1 + \tau_H s)} + \frac{a P(s)}{2} \right] \\ T_p(s) &= \frac{\Delta z}{2G^0} \left[(\alpha_1^0 + \alpha_2^0) + \frac{(b_1 + b_2)}{2} P(s) \right] \end{aligned} \quad (43)$$

The perturbed flow quality, void fraction, mass fluxes and velocities are given respectively by:

$$\begin{aligned} (13) \quad \tilde{x}_2 &= \frac{\{1 - (1 - x_2^0) F_2^+\} \tilde{h}_2}{(h_g - h_{l2}^0)} \\ (C.18), \text{ Appendix C} \quad \tilde{\alpha}_2 &= \frac{1}{x_2} \left[\tilde{x}_2 - x_2^0 (1 - x_2^0) \frac{\rho_g'}{\rho_g^0} \tilde{p}_1 \right] \\ (5) \quad \tilde{G}_{g2} &= x_2^0 \tilde{G}_2 + G^0 \tilde{x}_2 \\ \tilde{G}_{l2} &= \tilde{G}_2 - \tilde{G}_{g2} \\ \tilde{U}_{g2} &= U_{g2}^0 \left[\frac{\tilde{G}_{g2}}{G_{g2}^0} - \frac{\tilde{\alpha}_2}{\alpha_2^0} - \frac{\rho_g'}{\rho_g^0} \tilde{p}_1 \right] \\ \tilde{U}_{l2} &= U_{l2}^0 \left[\frac{\tilde{G}_{l2}}{G_{l2}^0} + \frac{\tilde{\alpha}_2}{(1 - \alpha_2^0)} \right] \end{aligned} \quad (44)$$

where \tilde{G}_2 is calculated from (38).

Perturbation, linearisation and Laplace transformation of the momentum equation (8) gives the nodal pressure drop as

$$\Delta \tilde{p}_2 = \Delta \tilde{p}_{2m} + \Delta \tilde{p}_{2f} + \Delta \tilde{p}_{2g} \quad (45)$$

where the perturbed momentum, friction and gravitational pressure drop components are respectively

$$\begin{aligned} \text{Momentum: } \Delta \tilde{p}_{2m} &= \int_1^2 d \left[U_l^0 \tilde{G}_l + G_l^0 \tilde{U}_l + U_g^0 \tilde{G}_g + G_g^0 \tilde{U}_g \right] \\ &\quad + s \int_0^{\Delta Z} \tilde{G} dz \\ \text{or } \Delta \tilde{p}_{2m} &= \Delta \left[U_l^0 \tilde{G}_l + G_l^0 \tilde{U}_l + U_g^0 \tilde{G}_g + G_g^0 \tilde{U}_g \right] + \frac{s \Delta Z}{2} (\tilde{G}_1 + \tilde{G}_2) \\ \text{Friction: } \Delta \tilde{p}_{2f} &= \frac{K_f^0 G^0{}^2}{2\rho_l} \int_0^{\Delta Z} \left[\left(\frac{\tilde{K}_f}{K_f^0} + \frac{2\tilde{G}}{G^0} \right) \phi^0{}^2 + \tilde{\phi}^2 \right] dz \\ \text{or } \Delta \tilde{p}_{2f} &= \frac{\Delta p_{1f}^0}{2} \left[(2-B_f) \frac{(\phi_1^0{}^2 \tilde{G}_1 + \phi_2^0{}^2 \tilde{G}_2)}{G^0} + (\tilde{\phi}_1^2 + \tilde{\phi}_2^2) \right] \end{aligned} \quad (46)$$

Gravitation:

$$\begin{aligned} \Delta \tilde{p}_{2g} &= \int_0^{\Delta Z} \left[\Delta \rho \tilde{\alpha} - \alpha^0 \rho_g' \tilde{p} \right] g \cos \theta. dz \\ \text{or } \Delta \tilde{p}_{2g} &= \frac{1}{2} \left[\Delta \rho (\tilde{\alpha}_1 + \tilde{\alpha}_2) - (\alpha_1^0 + \alpha_2^0) \rho_g' \tilde{p}_1 \right] \Delta Z g \cos \theta. \end{aligned}$$

The perturbed two-phase friction multiplier (ϕ^2) in equation (46) is calculated from equation (A.13), Appendix A.

5.4 Single Phase Region Equations

Since the single phase liquid region has no vapour, the variables relating to the vapour phase are zero, that is, for the steady state and dynamics

$$\alpha = 0, \quad x = 0$$

and

$$G_g = 0, \quad U_g = 0 \quad (47)$$

The coefficients F^+ and K_0 in the subcooled boiling region energy equation (21) assume the values

$$F^+ = 1$$

and

$$K_0 = 1 \quad (48)$$

With these modifications the subcooled boiling region equations automatically reduce to simpler single phase region equations.

The single phase region equations are also used to model the wall void region. The vapour bubbles attached to the wall act as surface asperities, and are hydrodynamically represented by increasing the relative roughness ratio (ϵ) in the single phase friction factor equation (A.3), Appendix A. This method, proposed by Staub (1968), results in an increased single phase friction pressure drop component. The equations to calculate the incipience of the wall void region and the relative roughness factors based on vapour bubble height (Romberg 1972) are not considered in the present model.

5.5 Saturated Boiling Region Equations

The temperature of the subcooled liquid ultimately reaches the local saturation temperature, and the fluid properties of the liquid phase are defined by their local saturation values. Of particular importance is the subcooled liquid enthalpy, which becomes

$$\begin{aligned} h_l^0 &= h_f \\ \tilde{h}_l &= \tilde{h}_f = 0 \end{aligned} \quad (49)$$

so that, from equations (17), (18) and (20),

$$F^+ = 0 \quad (50)$$

Equations (49) and (50) assume that the saturated liquid enthalpy h_f is time invariant (Jones 1963), that is, the subcooled liquid

enthalpy perturbations propagating through the single phase and subcooled boiling regions are completely attenuated in the saturated boiling region. Hence, perturbations in the mixture enthalpy, equation (13), are solely due to perturbations in the flow quality (x). The subcooled boiling region energy equation (21) and finite difference equations automatically apply to the saturated boiling region with these constraints.

6. SECTION BOUNDARY EQUATIONS

The spatial integration of the conservation equations (6) - (8) is performed by dividing a heated channel of arbitrary geometry into a number of sections of constant flow area. These sections are further subdivided into smaller nodes and subnodes to satisfy numerical stability requirements when the finite difference equations derived previously are applied. The boundary between any two sections may represent a localised discontinuity in area and/or obstruction (bend, orifice plate, valve etc.) in the coolant flow path. Boundary coupling equations are required to model the flow conditions due to the obstruction or restriction, and these are obtained by solution of the macroscopic conservation equations (Bird et al 1960, Govier and Aziz 1972). Of particular importance is the two-phase restriction pressure drop, for which a number of models have been developed (Tong 1965, Janssen 1966, Husain and Weisman 1975, Weisman et al 1976).

In the following analysis the macroscopic equations for the conservation of mass and energy are solved for the general case of a section boundary with an abrupt change in cross-sectional area.

6.1 Section Boundary Conservation Equations

The boundary interface may represent an abrupt expansion or contraction (Figure 3).

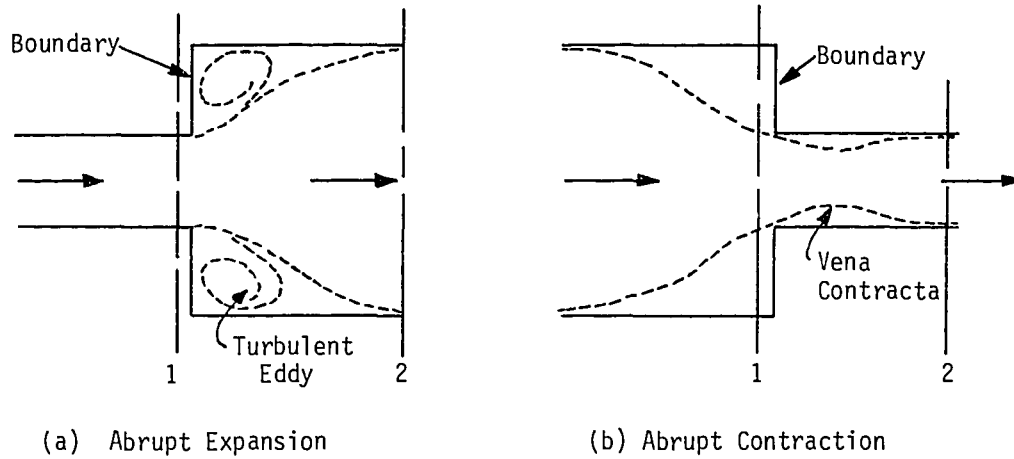


FIGURE 3 - SECTION BOUNDARY WITH AREA CHANGES

The macroscopic mass balance for the small control volume between stations (1) and (2) is, neglecting mass accumulation (Bird et al 1960),

$$\frac{dM}{dt} = W_1 - W_2 = \Delta W = 0 \quad (51)$$

$$\begin{aligned} \text{where } W &= W_\ell + W_g = AG \\ W_\ell &= A_\ell \rho_\ell U_\ell = A(1-\alpha)\rho_\ell U_\ell = AG_\ell = A(1-x)G \\ W_g &= A_g \rho_g U_g = A\alpha \rho_g U_g = AG_g = AxG \end{aligned} \quad (52)$$

The convected energy at stations (1) and (2) is, neglecting potential energy,

$$\begin{aligned} E &= W_\ell h_\ell + \frac{1}{2} W_\ell U_\ell^2 + W_g h_g + \frac{1}{2} W_g U_g^2 \\ &= W \{u + pv_m + e\} \end{aligned} \quad (53)$$

$$\begin{aligned}
 \text{where } h &= u + pv_m \\
 u &= (1-x)u_l + xu_g \\
 v_m &= \frac{1}{\rho_m} = v_l(1-x) + v_g x \\
 \rho_m &= \frac{\rho_l \rho_g}{\rho_g + \Delta\rho x} = \frac{1}{v_m} \\
 e &= \frac{1}{2} \{ (1-x) U_l^2 + x U_g^2 \}
 \end{aligned} \tag{54}$$

The macroscopic energy balance for the small control volume is, neglecting energy accumulation,

$$\therefore \frac{dE}{dt} = E_1 - E_2 = \Delta E = 0 \tag{55}$$

6.2 Steady State Equations

From (51) and (55) the steady state mass and energy equations are respectively

$$w_2^0 = w_1^0 \tag{56}$$

$$u_1^0 + \rho_1^0 v_{m1}^0 + e_1^0 = u_2^0 + \rho_2^0 v_{m2}^0 + e_2^0 \tag{57}$$

For no phase change ("flashing") across the boundary, the flow quality and mixture specific volume are respectively

$$\begin{aligned}
 x_2^0 &= x_1^0 \\
 v_{m2}^0 &= v_{m1}^0 = v_m^0
 \end{aligned} \tag{58}$$

From (52) and (56), the downstream mass flux is

$$G_2^0 = \frac{A_1}{A_2} G_1^0 \tag{59}$$

and from (27) the void fraction is

$$\alpha_2^0 = \frac{K_{02} \rho_l x_1^0}{\rho_l + \Delta\rho x_1^0} \quad (60)$$

The other variables (liquid and vapour mass fluxes and velocities) are calculated from equations (28).

From (57) the inlet restriction pressure drop is

$$\Delta p_R^0 = (p_1^0 - p_2^0) = \rho_m^0 \{ (e_2^0 - e_1^0) + (u_2^0 - u_1^0) \} \quad (61)$$

The change in internal energy is due to the irreversible friction loss, and is given by

$$\Delta p_L^0 = \rho_m^0 (u_2^0 - u_1^0) = K_R^0 \frac{G_2^0 \phi_R^0}{2\rho_l} \quad (62)$$

where the two-phase friction multiplier is (Beattie 1973)

$$\phi_R^2 = 1 + 1.284 \left\{ \frac{\rho_l}{\rho_g} - 1 \right\} x \quad (63)$$

and K_R is the single phase loss factor for the abrupt expansion or contraction (Bird et al 1960).

The momentum pressure drop component is

$$\begin{aligned} \Delta p_I^0 &= \rho_m^0 (e_2^0 - e_1^0) \\ &= \frac{\rho_m^0}{2} \{ (1-x_1^0)(U_{l2}^0 - U_{l1}^0)^2 + x_1^0 (U_{g2}^0 - U_{g1}^0)^2 \} \end{aligned} \quad (64)$$

and hence equation (61) becomes

$$\Delta p_R^0 = \Delta p_I^0 + \Delta p_L^0 \quad (65)$$

Note that for $A_2 > A_1$ (expansion) there is a deceleration of the coolant flow ($\Delta p_I^0 < 0$), and if $|\Delta p_I^0| > |\Delta p_L^0|$ there will be an overall pressure rise ($p_2^0 > p_1^0$, Δp_R^0 is negative). For no area change ($A_2 = A_1$), Δp_I^0 is zero so that $\Delta p_R^0 = \Delta p_L^0$ (e.g. bends, valves, etc.).

6.3 Dynamic Equations

Perturbation, linearisation and Laplace transformation of the mass and energy equations (51) and (53) gives

$$\tilde{w}_2 = \tilde{w}_1 \quad (66)$$

$$\text{and } \tilde{u}_1 + v_{m1}^0 \tilde{p}_1 + p_1^0 \tilde{v}_{m1} + \tilde{e}_1 = \tilde{u}_2 + v_{m2}^0 \tilde{p}_2 + p_2^0 \tilde{v}_{m2} + \tilde{e}_2 \quad (67)$$

For no flashing, the perturbed values of the flow quality and mixture specific volume are respectively

$$\tilde{x}_2 = \tilde{x}_1 \quad (68)$$

$$\text{and } \tilde{v}_{m2} = \tilde{v}_{m1} = \tilde{v}_m$$

where, from equation (54),

$$\begin{aligned} \tilde{v}_m &= \left(\frac{\partial v_m^0}{\partial x^0} \right) \tilde{x} + \left(\frac{\partial v_m^0}{\partial \rho_g^0} \right) \tilde{\rho}_g \\ &= \left(\frac{\Delta \rho}{\rho_l \rho_g^0} \right) \tilde{x}_1 - \left(\frac{x_1^0}{\rho_g} \frac{\rho_g'}{\rho_g} \right) \tilde{p}_1 \end{aligned} \quad (69)$$

The perturbed inlet restriction pressure drop is, from equation (67),

$$\Delta \tilde{p}_R = (\tilde{p}_1 - \tilde{p}_2) = \rho_m^0 \{ (\tilde{u}_2 - \tilde{u}_1) - \Delta p_R^0 \tilde{v}_m + (\tilde{e}_2 - \tilde{e}_1) \} \quad (70)$$

Now, from equation (62), the loss component of the restriction pressure drop is

$$\Delta p_L = \rho_m (u_2 - u_1) = K_R \frac{G_2^2 \phi_R^2}{2\rho_l} \quad (71)$$

so that

$$\begin{aligned}
\Delta \tilde{p}_L &= \rho_m^0 (\tilde{u}_2 - \tilde{u}_1) - \rho_m^0 \Delta p_L^0 \tilde{v}_m \\
&= \Delta p_L^0 \left[2 \frac{\tilde{G}_2}{G_2^0} + \frac{\tilde{\phi}_R^{-2}}{\phi_R^0{}^2} \right]
\end{aligned} \tag{72}$$

where, from equations (54) and (63) respectively,

$$\tilde{\rho}_m = -\rho_m^0{}^2 \tilde{v}_m \tag{73}$$

$$\text{and } \tilde{\phi}_R^{-2} = (\phi_R^0{}^2 - 1) \left[\frac{\tilde{x}_1}{x_1^0} - \left(\frac{\rho_g}{\Delta \rho} \quad \frac{\rho_g'}{\rho_g^0} \right) \tilde{p}_1 \right]$$

Similarly, from equation (64), the inertia component of the restriction pressure drop is

$$\Delta p_I = \rho_m (e_2 - e_1) \tag{74}$$

so that,

$$\Delta \tilde{p}_I = \rho_m^0 (\tilde{e}_2 - \tilde{e}_1) - \rho_m^0 \Delta p_I^0 \tilde{v}_m \tag{75}$$

where, from equation (54),

$$\tilde{e}_i = \frac{1}{2} \{U_{gi}^0{}^2 - U_{li}^0{}^2\} \tilde{x}_1 + (1-x_1^0)U_{li}^0 \tilde{u}_{li} + x_1^0 U_{gi}^0 \tilde{u}_{gi} \quad i=1,2$$

Substitution of equations (72) and (75) in (70) gives

$$\Delta \tilde{p}_R = \Delta \tilde{p}_I + \Delta \tilde{p}_L \tag{76}$$

The perturbed mass flux at station (2), Figure 3, is, from equations (52) and (66),

$$\tilde{G}_2 = \frac{A_1}{A_2} \tilde{G}_1 \tag{77}$$

The perturbed values of the other variables are calculated from equations (44).

These steady state and dynamic equations apply to both the sub-cooled and saturated boiling regions. The assumption of no flashing across the restriction is more inaccurate for

a) the subcooled boiling region, in which the temperature difference between the subcooled liquid and saturated (or superheated) vapour phases is greater, and

b) large restriction pressure drops (Δp_R), particularly those resulting in choked flow conditions. The equations automatically reduce to simpler forms for single phase flow ($\alpha=0$, $x=0$) and constant flow area ($A_2=A_1$).

7. SUBCOOLED AND SATURATED BOILING BOUNDARIES

The integration of the conservation equations (6) - (8) also requires coupling equations for the interface (boiling boundary) between the single phase and two-phase regions. The interface may denote either a subcooled or saturated boiling boundary depending on whether or not a subcooled boiling region is included in the model. The boundary between the subcooled and saturated boiling regions does not require special coupling equations, since the subcooled boiling region finite difference equations asymptote to the correct limits as defined by equations (49) and (50). The following analysis derives equations for the subcooled boiling boundary, and it is shown how these equations may be modified to model the saturated boiling boundary.

The subcooled boiling boundary (point of net vapour generation) is mathematically denoted by the position where the wall voids become detached from the heated surface. In practice, however, the statistical variations at the heated surface (heat fluxes, surface cavities, etc.) preclude a precise definition of the subcooled boiling position, and thus it is predicted from empirical correlations based on

experimental data. Kroeger and Zuber (1968) show that the ability to predict the void fraction in the subcooled boiling region depends on the accurate prediction of the boiling boundary position.

7.1 Subcooled Boiling Boundary Correlation

A number of methods have been developed to predict the subcooled boiling boundary position. These include

1. mathematical analyses which attempt to predict the local thermal-hydraulic conditions which cause bubble detachment (Bowring 1962, Levy 1967, Staub 1968), and
2. empirical correlations based on a postulated heat transfer coefficient (Ahmad 1970, Hancox and Nicoll 1971).

The mathematical models are too complicated and difficult for transient studies, and both methods, although agreeing with experimental data for high mass flow rates, give unsatisfactory results at low mass flow rates (Saha and Zuber 1974). Saha and Zuber (1974) postulate that at low mass flow rates the point of net vapour generation is dependent only on the local thermal conditions, and that condensation as the bubbles move away from the heated surface is governed by a diffusion process. They regard this as the 'thermally controlled region' in which the local Nusselt number is the dominant dimensionless parameter. At high mass flow rates the bubbles are detached by viscous shear effects, and they regard this as a 'hydrodynamically controlled' region in which, using Reynolds analogy (Knudsen and Katz 1958), the Stanton number becomes the dominant dimensionless parameter. The Nusselt and Stanton numbers are related by the Peclet number:

$$\text{Nu} = \text{Pe} \times \text{St} \quad (78)$$

$$\text{where } \text{Nu} = \frac{h_t D_e}{k_\ell}, \quad \text{Pe} = \frac{G D_e C_{p\ell}}{k_\ell}, \quad \text{St} = \frac{h_t}{G C_{p\ell}} \quad (79)$$

Saha and Zuber plotted the Peclet number versus the Stanton number for a range of experimental data, and obtained (Figure 4) a straight line of slope minus one for Peclet numbers less than 70,000 (thermally controlled region), which implies a constant value for the local Nusselt number. For Peclet numbers greater than 70,000 (hydrodynamically controlled region) the data has a constant Stanton number.

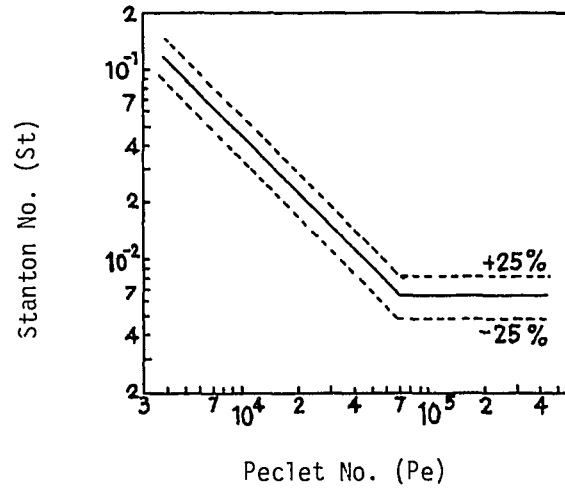


FIGURE 4 - PECLT NUMBER VERSUS STANTON NUMBER AT THE POINT OF NET VAPOUR GENERATION

Their correlation is given by

$$\begin{aligned} Nu &= \frac{h_{to} D_e}{k_\ell} = 455, & Pe < 70,000 \\ St &= \frac{h_{to}}{G C_{p\ell}} = 0.0065, & Pe > 70,000 \end{aligned} \quad (80)$$

Hence the heat transfer coefficient at the onset of subcooled boiling is

$$\left. \begin{aligned} h_{to} &= \frac{455 k_\ell}{D_e} & Pe \leq 70,000 \\ h_{to} &= 0.0065 G C_{p\ell} & Pe > 70,000 \end{aligned} \right\} \quad (81)$$

Their correlation is modified in the present analysis by replacing the equivalent hydraulic diameter (D_e) by the equivalent heated diameter (D_h), as suggested by Forgan and Whittle (1967) and Kreyger and Essler (1970). This improves the correlation of data by Evangelisti (1969) for annuli, which previously fell outside the $\pm 25\%$ tolerance band in Figure 4.

The liquid enthalpy for the onset of subcooled boiling is given by

$$h_o = h_f - \frac{C_{pl} q_{wo}}{h_{to}} \quad (82)$$

and the boiling boundary occurs when

$$h_l^o(z_o) = h_o \quad (83)$$

7.2 Subcooled Boiling Boundary Conservation Equations

The coupling of the steady state and perturbed variables between the single phase and two-phase regions is provided by the equations for the conservation of mass, energy and momentum at the boiling boundary interface (Jones 1963). Some analytical models (Davies and Potter 1967, Spinks 1972) express the perturbed variables in the two-phase region in terms of the movement in the boiling boundary position (perturbed boiling boundary). This method is impractical for finite difference models, since movement of the perturbed boiling boundary across a fixed space node boundary introduces numerical instabilities in the calculation of the perturbed variables in the two-phase region.

The mass, energy and momentum conservation equations at the boiling boundary interface are respectively:

$$\text{Mass:} \quad [G_l]' = [G_l + G_g]'' = [G]'' \quad (84)$$

$$\text{Energy:} \quad [G_l h_l]' = [G_l h_l + G_g h_g]'' = [Gh]'' \quad (85)$$

$$\text{Momentum:} \quad [G_l U_l]' = [G_l U_l + G_g U_g]'' \quad (86)$$

The superscripts (') and (") refer respectively to conditions upstream

and downstream of the boiling boundary interface.

Neglecting the voids attached to the wall, the steady state free-stream void fraction and flow quality are zero at the boiling boundary, that is

$$[\alpha^0]'' = 0, \quad [x^0]'' = 0 \quad (87)$$

$$\text{so that } [h_\ell^0]'' = [h^0]'' = [h_\ell^0]' = h_0 \quad (88)$$

$$\text{and, } [F^+]'' = 1$$

Hence it follows that

$$\begin{aligned} (5) \rightarrow [G_g^0]'' &= 0, \\ (84) \rightarrow [G_\ell^0]'' &= [G_\ell^0]' = [G^0]', \\ (86) \rightarrow [U_\ell^0]'' &= [U_\ell^0]' \end{aligned} \quad (89)$$

$$\text{and, implicitly, } [U_g^0]'' = 0$$

The perturbed, linearised and transformed conservation equations are respectively, from equations (84) - (86):

$$\underline{\text{Mass:}} \quad [\tilde{G}_\ell]' = [\tilde{G}_\ell + \tilde{G}_g]'' = [\tilde{G}]'' \quad (90)$$

$$\begin{aligned} \underline{\text{Energy:}} \quad [G_\ell^0 \tilde{h}_\ell + h_\ell^0 \tilde{G}_\ell]' &= [G_\ell^0 \tilde{h}_\ell + h_\ell^0 \tilde{G}_\ell + h_g^0 \tilde{G}_g]'' \\ &= [G^0 \tilde{h} + h^0 \tilde{G}]'' \end{aligned} \quad (91)$$

$$\underline{\text{Momentum:}} \quad [G_\ell^0 \tilde{U}_\ell + U_\ell^0 \tilde{G}_\ell]' = [G_\ell^0 \tilde{U}_\ell + U_\ell^0 \tilde{G}_\ell + G_g^0 \tilde{U}_g + U_g^0 \tilde{G}_g]'' \quad (92)$$

From (44) the perturbed flow quality at the boiling boundary is

$$[\tilde{x}]'' = \left[\frac{\{1 - (1-x^0) F^+\} \tilde{h}}{h_g - h_\ell^0} \right]'' = 0 \quad (93)$$

using equations (87) and (88), and hence it follows that

$$\begin{aligned}
 (44) \rightarrow \quad [\tilde{\alpha}]'' &= 0 \\
 (5) \rightarrow \quad [\tilde{G}_g]'' &= 0 \\
 (90) \rightarrow \quad [\tilde{G}_\ell]'' &= [\tilde{G}_\ell]' = [\tilde{G}]' \\
 (91) \rightarrow \quad [\tilde{h}]'' &= [\tilde{h}_\ell]'' = [\tilde{h}_\ell]' \\
 (92) \rightarrow \quad [\tilde{U}_\ell]'' &= [\tilde{U}_\ell]' \\
 \text{and, implicitly,} \quad [\tilde{U}_g]'' &= 0
 \end{aligned} \tag{94}$$

7.3 Saturated Boiling Boundary

In certain applications the subcooled boiling region may be neglected and the boiling channel hydrodynamics represented by a two region (single phase/saturated two-phase) model. The saturated boiling boundary denotes the position where the bulk liquid temperature reaches saturation. Hence (88) becomes

$$[\tilde{h}_\ell^0]'' = [\tilde{h}_\ell^0]' = h_f \tag{95}$$

The steady state equations (87) and (89) apply as previously for the subcooled boiling boundary, except that the subcooled liquid subscript (ℓ) is replaced by the saturated liquid subscript (f). Since the saturated liquid enthalpy is time invariant, equations (49), (3.50),

$$[\tilde{h}_f]'' = 0 \quad \text{and} \quad [F^+]'' = 0 \tag{96}$$

at the saturated boiling boundary, and the perturbed flow quality, (93), becomes

$$[\tilde{x}]'' = \frac{[\tilde{h}]''}{h_g - h_f} = \frac{[\tilde{h}_\ell]'}{h_{fg}} \tag{97}$$

Hence,

$$\begin{aligned}
 (44) \rightarrow [\tilde{\alpha}]'' &= \left[\frac{\tilde{x}}{X} \right]'' \\
 (5) \rightarrow [\tilde{G}_g]'' &= [G^0 \tilde{x}]'' \\
 (90) \rightarrow [\tilde{G}_f]'' &= [\tilde{G} - \tilde{G}_g]'' \\
 (92) \rightarrow [\tilde{U}_f]'' &= \frac{[G_l^0 \tilde{U}_l]'' + [U_f^0 \tilde{G}_g]''}{[G_l^0]''}
 \end{aligned}
 \quad \left. \vphantom{\begin{aligned} (44) \rightarrow [\tilde{\alpha}]'' \\ (5) \rightarrow [\tilde{G}_g]'' \\ (90) \rightarrow [\tilde{G}_f]'' \\ (92) \rightarrow [\tilde{U}_f]'' \end{aligned}} \right\} (98)$$

since, implicitly, $[\tilde{U}_g]'' = 0$.

A comparison of the above models shows that the essential difference between the subcooled and saturated boiling boundaries is the transfer of the perturbed energy flux across the boundary. The perturbed energy flux of the upstream liquid phase is transferred entirely to the downstream liquid phase at a subcooled boiling boundary, whereas it is transferred entirely to the downstream vapour phase at a saturated boiling boundary.

8. CHANNEL, BYPASS AND PUMP MODELS

A heated channel is usually one component of a larger system, and may be operated

- a) in parallel with a large number of channels of similar geometry,
- b) in parallel with a large bypass, or
- c) in closed circuit with a pump.

The following analysis shows how the finite difference equations derived previously are used to model the total channel. Equations are derived for the bypass and pump to permit analysis of the complete system.

8.1 Single Channel Model

The stepwise integration of the conservation equations using the finite difference, spatial boundary and boiling boundary equations

yields the local steady state and dynamic conditions along the channel. The steady state calculations are performed first to evaluate the coefficients in the dynamic equations and to obtain, in particular, the overall channel pressure drop, which may be written as

$$\Delta p^0 = \Delta p_1^0 + \Delta p_2^0 \quad (99)$$

where the single phase and two-phase steady state pressure drops are respectively

$$\Delta p_1^0 = \sum_{j=1}^{NSAT-1} \Delta p_{1j}^0 + \sum_{k=1}^{N1} \{ \Delta p_{1R}^0 \}_k$$

and

$$\Delta p_2^0 = \sum_{j=NSAT}^N \Delta p_{2j}^0 + \sum_{k=N1+1}^{NCH} \{ \Delta p_{2R}^0 \}_k$$

(100)

The nodal pressure drops (Δp_{1j}^0 , Δp_{2j}^0) have momentum, friction and gravitational components, equations (31). The channel is divided into NCH sections and N sub-nodes (see description of computer program, Section 10), and boiling is assumed to occur at the beginning of sub-node NSAT in section N1. Restriction pressure drops (Δp_{1R}^0 , Δp_{2R}^0) are evaluated at the beginning of each section in the single phase and two-phase regions of the channel.

The subdivision of the sections into nodes and sub-nodes is particularly important near the boiling boundary, where small node sizes are required to accurately resolve the boundary position and to ensure that the steady state and dynamic calculations in the two-phase region of the channel are numerically stable. The steady state piecewise integration is similar to that used for the dynamic analysis, Figure 5, in which the perturbed variables are replaced by their equivalent steady state values. The top blocks represent the nodal mass and energy conservation equations, and the lower blocks represent the nodal

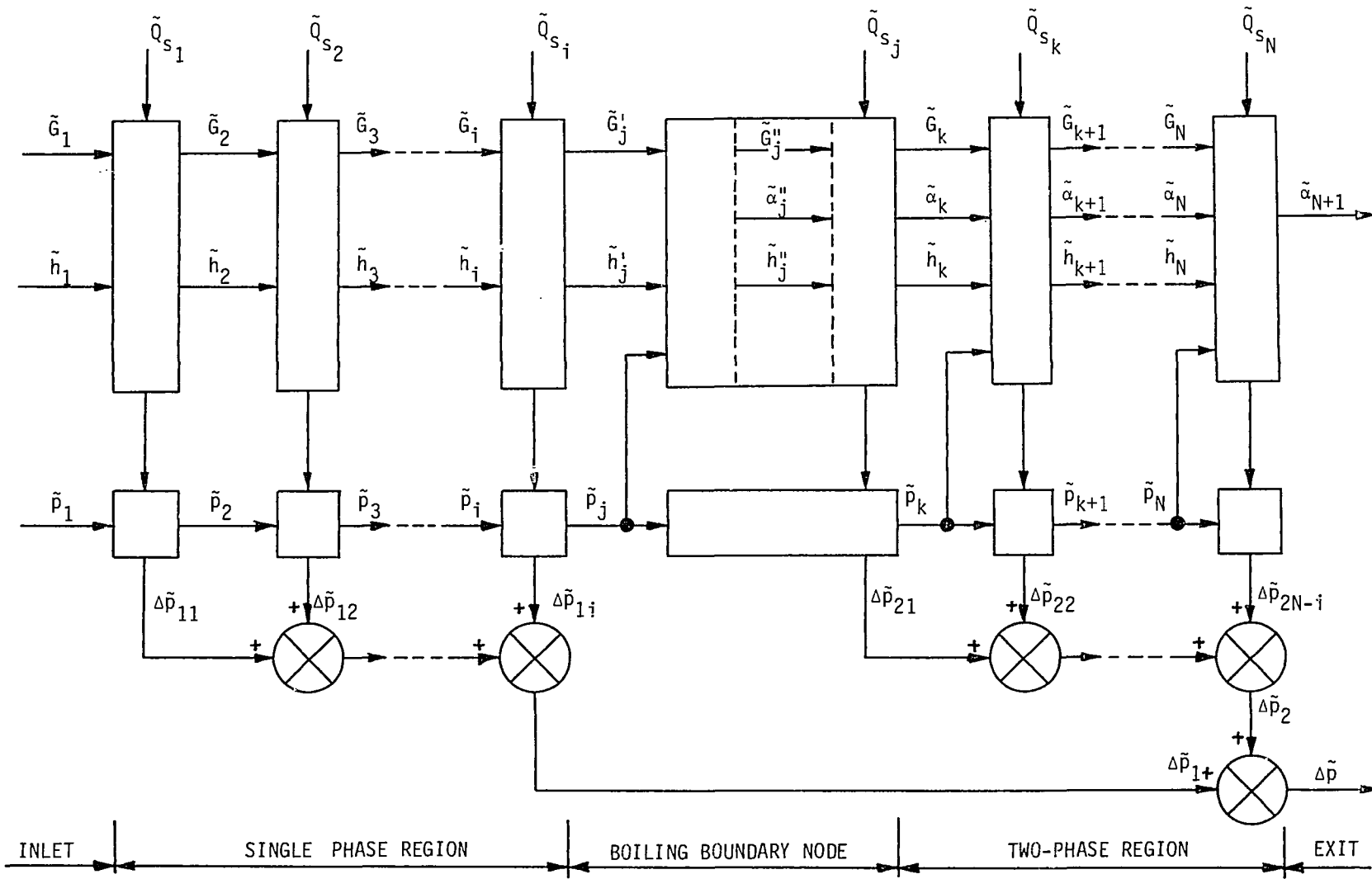


FIGURE 5 - SCHEMATIC DIAGRAM OF BOILING CHANNEL MODEL

momentum conservation equations. Only the transfer of principal variables between the blocks is shown.

In the dynamic analysis, the piecewise calculation of the local perturbed variables is performed for specified values of the complex Laplace operator. The mass flux (\tilde{G}_1), enthalpy (\tilde{h}_1) and pressure (\tilde{p}_1) perturbations at the channel inlet, and the nodal power density (\tilde{Q}_S) perturbations of the heat source are the system inputs, and the overall pressure drop perturbation ($\Delta\tilde{p}$) and the perturbed void fraction at the channel exit ($\tilde{\alpha}_{N+1}$) are the outputs. The perturbations at the channel inlet travel through the single phase region to the boiling boundary, where they are transmitted to the two-phase region. The perturbations of the vapour and liquid phases travel to the channel exit, where they interact with the external system.

The overall perturbed pressure drop is obtained by summation of the single phase and two-phase region pressure drops (Figure 5):

$$\Delta\tilde{p} = \Delta\tilde{p}_1 + \Delta\tilde{p}_2 \quad (101)$$

where $\Delta\tilde{p}_1$ and $\Delta\tilde{p}_2$ are given by equation (100) with the steady state values replaced by their respective perturbed values. From Figure 5, $\Delta\tilde{p}$ (via $\Delta\tilde{p}_1$, $\Delta\tilde{p}_2$) is also linearly related to the inlet mass flux, pressure, enthalpy and average power density by the equation

$$\begin{aligned} \Delta\tilde{p} = & \tilde{H}(\Delta p, G_1)\tilde{G}_1 + \tilde{H}(\Delta p, p_1)\tilde{p}_1 + \tilde{H}(\Delta p, h_1)\tilde{h}_1 \\ & + \tilde{H}(\Delta p, \langle Q_S \rangle)\langle\tilde{Q}_S\rangle \end{aligned} \quad (102)$$

where $\tilde{H}(\quad)$ is the frequency dependent transfer function relating the variables in the brackets.

The first term $\tilde{H}(\Delta p, G_1)$ in (102) is the channel 'hydraulic impedance' (Anderson 1970), and is the dominant term in the frequency range of interest (0.1 - 2 Hz). The second term $\tilde{H}(\Delta p, p_1)$ is due to the compressibility of the two-phase region, and is more influential at

higher frequencies (above 2 Hz). The third term, $\tilde{H}(\Delta p, h_1)$, results from the transfer of the inlet enthalpy perturbation to the boiling boundary, where it produces a marginal change in the two-phase region perturbed pressure drop. The fourth term, $\tilde{H}(\Delta p, \langle Q_s \rangle)$, is dependent on the heat source/coolant time constant τ_H , and is usually insignificant at frequencies above 0.1 Hz. Although the latter three terms have a marginal effect on the overall pressure drop $\Delta \tilde{p}$ in most heat transfer systems, they are retained in the present model for general system analysis.

8.2 Bypass Model

A bypass is sometimes placed in parallel with a test channel to control its overall pressure drop (e.g. Collins and Gacesa 1969). The steady state pressure drop of the bypass is given by

$$\Delta p_B^0 = \sum_{k=1}^{NB} \{ \Delta p_{Bf}^0 + \Delta p_{Bg}^0 + \Delta p_{BR}^0 \}_k \quad (103)$$

where NB is the number of bypass sections, and the friction, gravitation and restriction pressure drop components are calculated from equations (31) and (65) for single phase flow.

The perturbed pressure drop of the bypass is given by

$$\Delta \tilde{p}_B = \sum_{k=1}^{NB} \{ \Delta \tilde{p}_{Bm} + \Delta \tilde{p}_{Bf} + \Delta \tilde{p}_{BR} \}_k \quad (104)$$

where the momentum, friction and restriction pressure drop components are calculated from equations (46) and (76) for single phase flow. The perturbed pressure drop is linearly related to the perturbed mass flux at the bypass inlet by

$$\Delta \tilde{p}_B = \tilde{H}(\Delta p_B, G_B) \tilde{G}_{B1} \quad (105)$$

8.3 Pump Model

The pump operating in closed circuit with a test channel may have variable speed, throttle and/or bypass flow control. The net pump head as seen by the channel is assumed to be approximated by the quadratic equation

$$H_p = A_p + B_p Q_1 + C_p Q_1^2 \quad (106)$$

where Q_1 is the volumetric flow rate at the inlet to section 1 (pump discharge) of the channel, and is given by

$$Q_1 = \frac{A_1 G_1}{\rho_{\ell 1}} \quad (107)$$

The pressure rise across the pump is given by

$$\begin{aligned} -\Delta p_p &= H_p \rho_{\ell 1} g \\ &= \{A_p + B_p Q_1 + C_p Q_1^2\} \rho_{\ell 1} g \end{aligned} \quad (108)$$

Equation (108) gives the steady state pressure differential across the pump by insertion of the appropriate steady state values.

Perturbation linearisation and Laplace transformation of (108) yields

$$\begin{aligned} -\Delta \tilde{p}_p &= \{B_p + 2C_p Q_1^0\} A_1 g \tilde{G}_1 \\ &= \tilde{H}(\Delta p_p, G_1) \tilde{G}_1 \end{aligned} \quad (109)$$

$$\text{where } \tilde{H}(\Delta p_p, G_1) = \{B_p + 2C_p Q_1^0\} A_1 g \quad (110)$$

9. BOUNDARY CONDITIONS AND STABILITY CRITERIA

The analysis of the flow stability in parallel channel, channel/bypass and closed loop systems is performed by consideration of the mass and momentum exchange between the sub-systems (channel, bypass, pump), together with appropriate boundary conditions based on assumptions regarding the overall system behaviour. In this section,

the stability criteria for each type of system are derived.

9.1 Parallel Channel Stability

For N heated channels connected in parallel between common headers, the total inlet mass flow rate is

$$W_T = \sum_{n=1}^N W_n \quad (111)$$

and the overall pressure drop across all channels is

$$\Delta p = \Delta p_n, \quad n = 1, N \quad (112)$$

The mass and momentum equations (111) and (112) apply to both the steady state and dynamic behaviour of the system. At the onset of self-sustained oscillations in the coolant flow through a particular channel (n), flow oscillations are induced in the adjacent stable channels as per (111), and this redistribution is assumed to result in no change in the system pressure differential (or net pump head Δp) and, consequently, the pressure drop across the nth channel (Δp_n). Thus, the (linear) 'instability threshold' boundary condition for the parallel channel system is

$$\Delta p(f_{00}) = \Delta p_n(f_{00}) = 0 \quad (113)$$

where f_{00} is the frequency of the self-sustained oscillations (subscript (oo) = fundamental mode, (01) = first harmonic, etc.).

Dividing equation (101) by the perturbed mass flux at the channel inlet and invoking (113) results in the parallel channel stability criterion

$$\left. \begin{aligned} \tilde{H}_{00}(\Delta p, G_1) &= \tilde{H}_{00}(\Delta p_1, G_1) + \tilde{H}_{00}(\Delta p_2, G_1) > 0 \\ \text{that is, instability threshold occurs when} \\ \tilde{R}_{00}(\Delta p, G_1) &= 0 \\ \text{and} \\ \tilde{I}_{00}(\Delta p, G_1) &= 0 \end{aligned} \right\} \quad (114)$$

where the single phase, two-phase and overall hydraulic impedance functions are respectively

$$\left. \begin{aligned} \Delta\tilde{p}_1 &= \tilde{H}(\Delta p_1, G_1) \tilde{G}_1 \\ \Delta\tilde{p}_2 &= \tilde{H}(\Delta p_2, G_1) \tilde{G}_1 \\ \Delta\tilde{p} &= \tilde{H}(\Delta p, G_1) G_1 \end{aligned} \right\} \quad (115)$$

$\tilde{R}(\Delta p, G_1)$ and $\tilde{I}(\Delta p, G_1)$ are the real and imaginary components of $\tilde{H}(\Delta p, G_1)$, and the subscript (oo) denotes the frequency f_{oo} in equation (113). Equation (114) neglects the influence of the other terms in equation (101), and is consistent with most hydrodynamic models (Jones 1961, Davies and Potter 1967, Takahashi and Shindo 1971, Spinks 1972, Yadigaroglu and Bergles 1972).

When $\tilde{H}(\Delta p, G_1)$ is plotted on an Argand diagram, the channel stability is defined by the locus in the clockwise direction as follows:

- a) the channel is stable if the locus intersects the positive real axis, $\tilde{R}(\Delta p, G_1) > 0$, resulting in no encirclement of the origin;
- b) the channel is at the instability threshold if the locus passes through the origin, $\tilde{R}(\Delta p, G_1) = 0$;
- c) the channel is unstable if the locus intersects the negative real axis, $\tilde{R}(\Delta p, G_1) < 0$, resulting in an encirclement of the origin.

The locus of the hydraulic impedance function $\tilde{H}(\Delta p, G_1)$ also has the following important characteristics:

- 1) $\tilde{H}(\Delta p, G_1)$ has no poles, only zeros (see equation (46), Spinks 1972) and thus $\tilde{H}(\Delta p, G_1) \rightarrow j\infty$ as $f \rightarrow \infty$;
- 2) the number of encirclements of the origin is indicative of the number of unstable modes (fundamental and higher mode instabilities, Yadigaroglu and Bergles 1972);
- 3) the 'margin to instability' is given by the intercept of the positive real axis ($= \tilde{R}_c(\Delta p, G_1)$) at the crossover frequency f_c ;

4) a negative zero frequency value of $\tilde{H}(\Delta p, G_1)$, that is $\tilde{R}_0(\Delta p, G_1) < 0$, indicates an unstable steady state (excursive or Ledinegg flow instability, see Section 9.4);

5) for a given frequency (f_n) the vector $\tilde{H}_k(\Delta p, G_1)$ is the resultant of the single phase and two-phase region vectors $\tilde{H}_k(\Delta p_1, G_1)$ and $\tilde{H}_k(\Delta p_2, G_1)$, which, in turn, are vector sums of their respective perturbed pressure drop components.

The latter point is illustrated by the simplified block diagram of the boiling channel model given in Figure 6, and is important in diagnosing the inherent mechanisms tending to drive the channel unstable. The single phase region impedance function $\tilde{H}(\Delta p_1, G_1)$ has orthogonal resistive (friction and restriction) and inductive (inertia) components in the positive real and imaginary directions respectively, and is mostly stabilising. The two-phase region impedance function $\tilde{H}(\Delta p_2, G_1)$ has vector components whose phase relationships are dependent on the vapour transit time through the two-phase region, and which have a destabilising influence on the channel hydrodynamics. The magnitudes and phase angles of dominant destabilising vector components (usually two-phase friction and restriction) can be evaluated by alteration of the channel dimensions, operating conditions, correlation coefficients, etc., and gives important information for design assessment studies. It can be seen from Figure 6 that the boundary condition (113) is satisfied when the perturbed two-phase pressure drop ($\Delta \tilde{p}_2$) lags the perturbed single phase pressure drop ($\Delta \tilde{p}_1$) by 180° .

The block diagram in Figure 6 can be readily manipulated into a closed loop feedback diagram to permit application of conventional control theory. For example, Davies and Potter (1967) manipulate their analytical equations to formulate a feedback model with the single phase region dynamics in the forward loop, and the two-phase region dynamics in the feedback loop. Channel stability is determined by

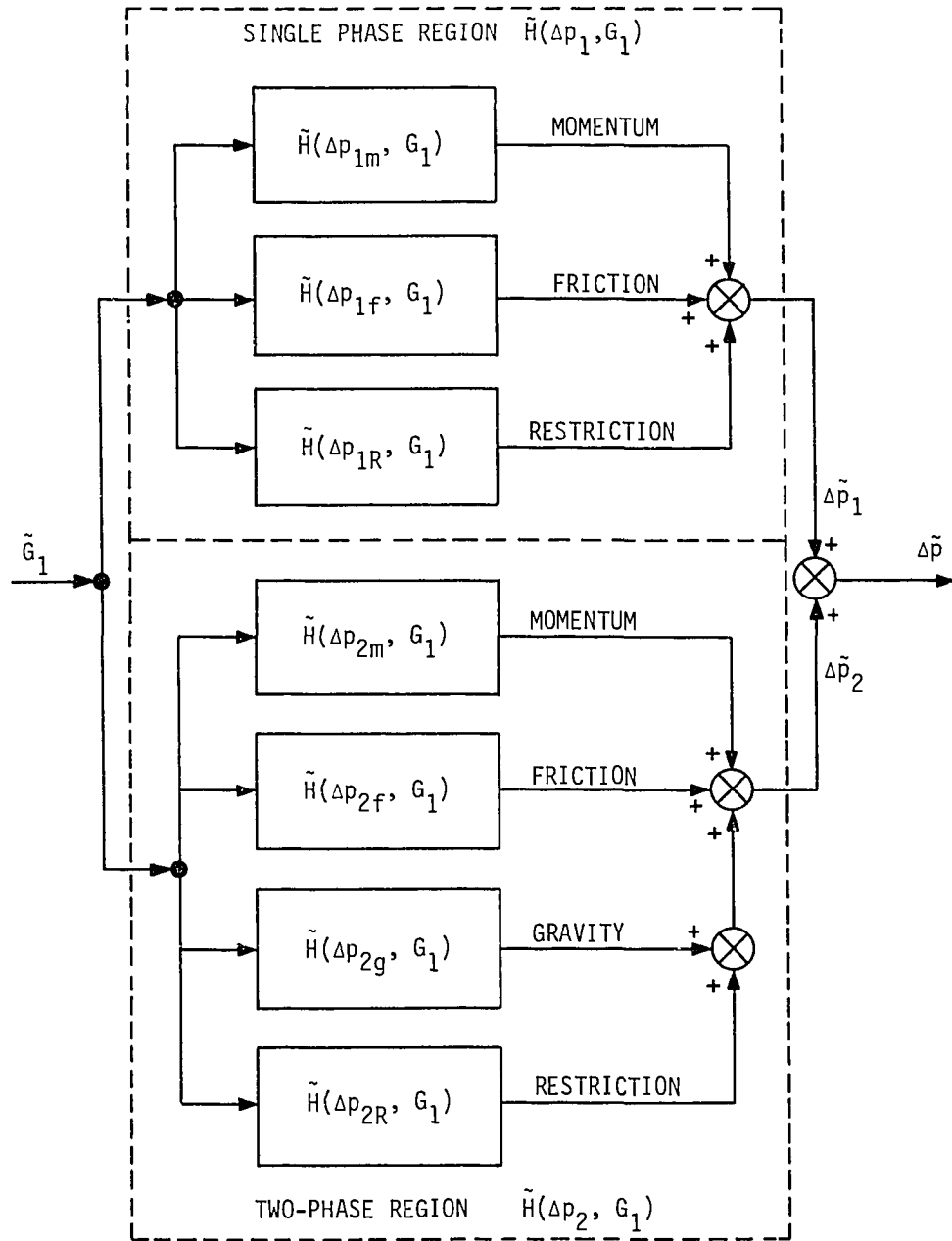


FIGURE 6 - SIMPLIFIED BLOCK DIAGRAM OF BOILING CHANNEL MODEL

applying the Nyquist stability criterion to the open loop transfer function

$$\frac{\Delta\tilde{p}_2}{\Delta\tilde{p}_1} = \tilde{H}(G_1, \Delta p_1) \tilde{H}(\Delta p_2, G_1) \geq (-1, 0) \quad (116)$$

using the same boundary condition (113) as the present model. Jones (1961) manipulated his hydrodynamic equations to obtain a feedback model with inlet pressure and mass flux perturbations as the input and output respectively, and the pressure perturbation at the boiling boundary (\tilde{p}_0) as the output of the feedback loop. Channel stability is determined by applying the Nyquist stability criterion to the open loop transfer function

$$\frac{\tilde{p}_0}{\Delta\tilde{p}_1} = \tilde{H}(G_1, \Delta p_1) \tilde{H}(p_0, G_1) \geq (-1, 0) \quad (117)$$

and results from the boundary condition that both the inlet and exit pressure perturbations are assumed to be zero, which is a particular solution of the boundary condition (113).

A major advantage of the present model is that the hydraulic impedance function $\tilde{H}(\Delta p, G_1)$ is defined in terms of measurable variables ($\tilde{G}_1, \Delta\tilde{p}$), and can be applied to system identification studies of experimental data from a test channel (e.g. Romberg and Rees 1975). By comparison, the feedback models (116) and (117) have the disadvantage that the perturbed variables $\Delta\tilde{p}_1$, $\Delta\tilde{p}_2$ and \tilde{p}_0 are non-measurable, since they are a function of an experimentally indeterminate boiling boundary position.

9.2 Channel/Bypass Stability

The bypass connected in parallel with a heated channel is designed to simulate the other (N-1) heated channels. For this configuration the mass equation (111) becomes

$$W_1 + W_B = W_T \quad (118)$$

and the momentum equation (112) becomes

$$\Delta p = \Delta p_B \quad (119)$$

If it is assumed that hydrodynamic interaction is between the channel and bypass only, that is, perturbations in the external system are negligible, then the dynamic boundary conditions are

$$\tilde{W}_1 + \tilde{W}_B = 0 \quad (120)$$

or $A_1 \tilde{G}_1 + A_{B1} \tilde{G}_{B1} = 0$

and $\Delta p (f_{00}) = \Delta p_B (f_{00}) \quad (121)$

From (105) and (120) the bypass pressure drop perturbation is

$$\Delta \tilde{p}_B = \tilde{H}(\Delta p_B, G_{B1}) \tilde{G}_{B1} = -\tilde{H}(\Delta p_B, G_1) \tilde{G}_1 \quad (122)$$

Substitution of (115) and (122) in (121) results in the channel/bypass stability criterion

$$\tilde{H}_{00}(\Delta p, G_1) + \tilde{H}_{00}(\Delta p_B, G_1) \geq 0 \quad (123)$$

that is, for the channel/bypass system to be stable the locus of equation (123) on an Argand diagram must not encircle the origin. A block diagram representation of the channel/bypass system is given in Figure 7.

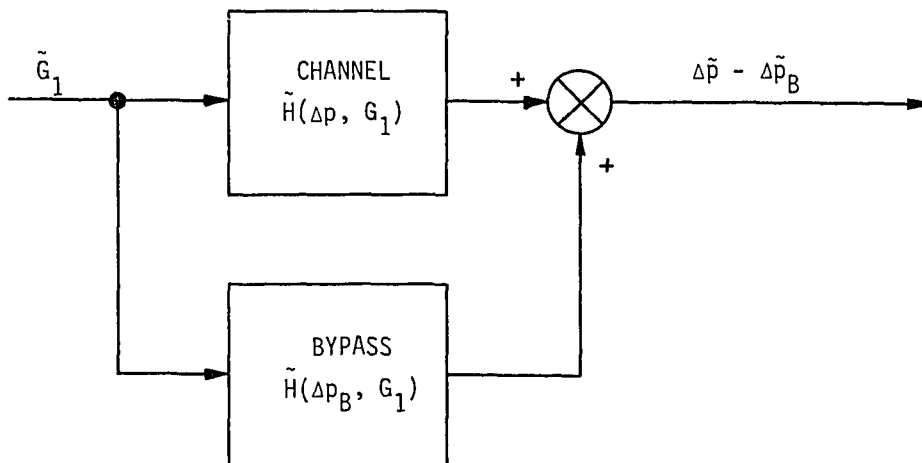


FIGURE 7 - BLOCK DIAGRAM OF CHANNEL/BYPASS SYSTEM

The bypass hydraulic impedance, $\tilde{H}(\Delta p_B, G_1)$, has single phase resistance and inertia components only, and has a stabilising effect on the channel flow. This was investigated experimentally by Collins and Gacesa (1969) and theoretically by Carver (1969), who showed that the threshold power (TP) for the onset of flow oscillations decreased asymptotically to the parallel channel ('infinite' bypass) threshold power when plotted as a function of the bypass ratio (W_B/W). Mathematically this means that, for large bypass ratio, in the frequency range of interest

$$\tilde{H}(\Delta p_B, G_1) \rightarrow 0 \quad (124)$$

and the channel/bypass stability criterion (123) converges to the parallel channel stability criterion (114).

9.3 Closed Loop Stability

For a closed loop system in which a channel is in series with a pump, summation of the pressure drops around the circuit results in the momentum boundary condition

$$\Delta p - \Delta p_p = 0 \quad (125)$$

$$\text{or} \quad \Delta p(f_{00}) - \Delta p_p(f_{00}) = 0$$

which is the same as the channel/bypass conditions (119) and (121) respectively. However, implicit in the boundary conditions (125) is the assumption that the channel exit and pump suction inlet conditions are separated by a large dispersive volume (e.g. steam drum), which decouples the input and output pressure perturbations in the frequency range of interest.

Substitution of equations (109) and (115) in (125) yields the "loop" stability criterion

$$\tilde{H}_{00}(\Delta p, G_1) + \tilde{H}_{00}(\Delta p_p, G_1) \geq 0 \quad (126)$$

From equation (110), the pump hydraulic impedance, $\tilde{H}(\Delta p_p, G_1)$, adds a single phase resistance to the channel hydraulic impedance, $\tilde{H}(\Delta p, G_1)$, and is stabilising. If the pump characteristic is fairly 'flat' ($B_p, C_p = 0$), or the system has 'natural circulation' flow due to a net gravitational head between inlet and outlet, the boundary condition (125) reduces to the parallel channel boundary condition (113).

9.4 Steady State Stability Criteria

The parallel channel, channel/bypass and closed loop boundary conditions (112), (119) and (125) must also be satisfied in the steady state,

$$\text{i.e.} \quad \Delta p^0 = \Delta p_s^0 \quad (127)$$

where Δp_s^0 is the net (supply) differential pressure imposed across the channel. Flow excursions (Ledinegg instabilities) occur when the channel pressure drop-flow characteristic has a negative slope region in which

$$\frac{\partial \Delta p^0}{\partial G_1^0} \ll \frac{\partial \Delta p_s^0}{\partial G_1^0} \quad (128)$$

Equation (128) can be replaced by the zero frequency transfer function equation

$$\tilde{H}_0(\Delta p, G_1) \ll \tilde{H}_0(\Delta p_s, G_1) \quad (129)$$

for the channel to be statically unstable, where

$$\left. \begin{aligned} \tilde{H}_0(\Delta p_s, G_1) &= 0 && \text{(parallel channel)} \\ &= -\tilde{H}_0(\Delta p_B, G_1) && \text{(bypass)} \\ &= -\tilde{H}_0(\Delta p_p, G_1) && \text{(pump)} \end{aligned} \right\} \quad (130)$$

As noted in Section 9.1, flow excursions occur when the resultant real vector of the zero frequency transfer functions in equations (129) and (130) is negative or zero.

10. DESCRIPTION OF COMPUTER PROGRAM (LOCO)

The finite difference equations derived in the previous sections have been incorporated in a digital computer program (LOCO - a Linearised model for analysing the Onset of Coolant Oscillations and frequency response of boiling channels), which performs the numerical calculations. This section describes the overall logic of the LOCO code. A complete listing of the code is given in Appendix D.

10.1 Program Data Input and Options

The hydraulic circuit between inlet and outlet headers is divided into an arbitrary number of sections ($NCH \leq 40$), and may optionally consist of an inlet region ($ND > 0$), a channel (or fuelled) region ($NC > 0$) and a riser region ($NR > 0$). The total number of sections is

$$NCH = ND + NC + NR \leq 40 \quad (131)$$

The geometry and operating data required for each section are specified by comment cards at the beginning of the program, and are listed in Appendix D. Each section has uniform flow area and power input or output. Total condensation of the two-phase flow is not permitted since no condensing boundary has been included in the present model. The two-phase slip correlation and friction multiplier coefficients are specified for each section, and may be varied between sections to account for changes in flow regime.

The program control parameters specify the various options available.

a) Bypass (NB) - the bypass is divided into NB sections; for no bypass $NB = 0$.

b) Flow rate (NDROP) - $NDROP = 0$: the channel pressure drop is calculated for a specified flow rate (W kg/s); $NDROP = 1$: the flow rate is calculated for a specified pressure drop ($PDROP$ kPa).

c) Subcooled boiling (NOSUB) - the liquid enthalpy in the subcooled boiling region is modelled using an exponential correlation (NOSUB = 1) or a hyperbolic tangent correlation (NOSUB = 2), equations (17), (18) respectively; subcooled boiling is neglected if NOSUB = 0.

d) Hydrodynamics (NDYN) - the program may be used for steady state calculations only (NDYN = 0), or steady state and transient calculations (NDYN > 0); the perturbed inputs (inlet mass flux, pressure, enthalpy and average power density) may be frequency independent (white noise, NDYN = 1) for stability calculations, frequency dependent and real (NDYN = 2) for frequency response calculations, or frequency dependent and complex (NDYN = 3) for multivariate spectral calculations.

e) Perturbed pressure drop components - these give information on the mechanisms tending to drive the channel unstable, and are calculated for

i) specified frequencies (CFREQ(N), N=1, NCFREQ) when $0 < \text{NCFREQ} < 10$;

ii) the initial frequency (FREQ) when the number of frequencies NF = 1;

iii) the 'crossover' frequency calculated by the program at which the polar plot of the perturbed pressure drop ($\Delta\bar{p}$) crosses the real axis in moving from the negative to positive imaginary half plane.

The print options (NPRINT), plot options (NPLOT), and fluid property options (NCARDS) are input device options as specified by the comment cards listed in Appendix D.

10.2 Program Operation

The LOCO program calculates the steady state and transient variables (mass fluxes, void fraction, component pressure drops etc.) for each space node in the channel as shown in the overall flow chart given in Figure 8. The MAIN subprogram reads and writes the input

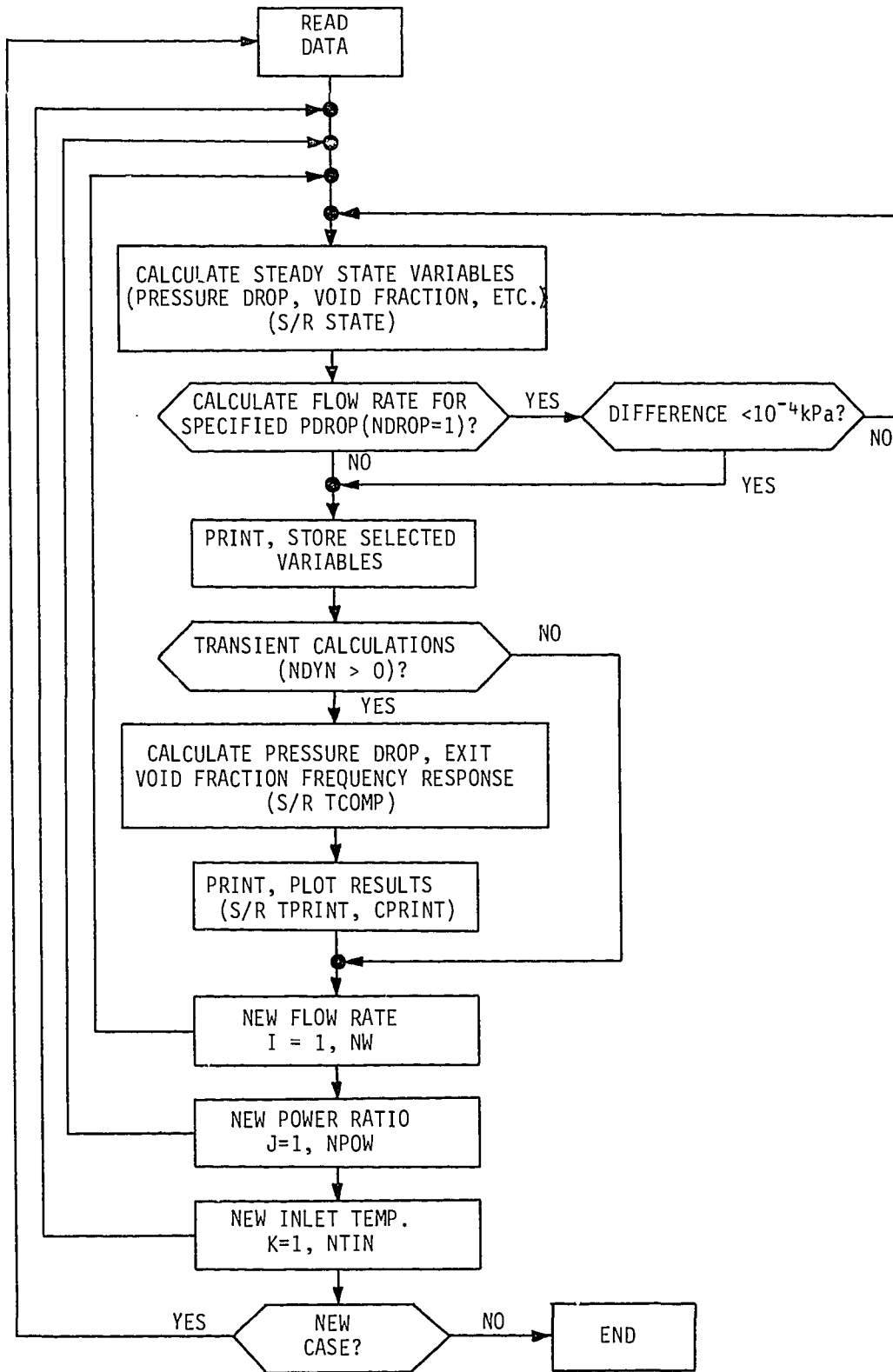


FIGURE 8 - 'LOCO' PROGRAM OVERALL FLOWCHART

data, converts the fluid properties to consistent units (subroutine TABLES), and transfers control to an executive subprogram (subroutine EXEC). This subroutine initialises the channel operating conditions (coolant inlet temperature, power input mass flow rate) and calls subroutine STATE for calculation of the channel steady state variables. These may be determined for a specified flow rate ($NDROP = 0$, W kg/s) or a specified pressure drop ($NDROP = 1$, $PDROP$ kPa). In the latter case $PDROP$ is specified as data input, or is calculated by the program if there is a bypass ($NB > 0$, subroutine *BYPASS*). The difference ($PDIFF$) between the calculated pressure drop ($PTOT$) and $PDROP$ is tested for a change in sign, and if this occurs, a regula falsi iteration procedure is used to calculate the mass flow rate (W) for the next repetitive calculation. The iteration is terminated when $|PDIFF| < 10^{-4}$ kPa. If no sign change in $PDIFF$ has occurred, the mass flow rate is incremented or decremented (DW kg/s) and a new $PTOT$ (and $PDROP$ if $NB > 0$) is calculated. The program prints W , $PTOT$, $PDROP$ and $PDIFF$ for each calculation, and terminates after NW flow rate steps if no sign change in $PDIFF$ has occurred ($NDYN$ is set to zero). Selected steady state variables are printed (subroutine *STATE*) for each node ($NPRINT = 0$) or section ($NPRINT > 0$) for the specified or calculated mass flow rate.

The program then proceeds to the transient calculations (if $NDYN > 0$) and calls subroutine *TFUNCN*, which edits the perturbed inputs (subroutine *DYNIN*) read in as data and calculates the value of the Laplace operator (s) for each frequency ($FREQ(J)$, $J = 1, NF$). Subroutine *TCOMP* is called at each frequency step to compute the channel pressure drop and exit void fraction frequency response. The channel pressure drop is modified by the perturbed pressure drop of the bypass ($NB > 0$) or pump. *TFUNCN* stores this information and tests whether the resultant perturbed pressure drop value has crossed the real axis of the polar

plot. If so, the subprogram calculates the crossover frequency by a regula falsi iteration procedure, and the value is stored, together with the single phase and two-phase perturbed pressure drop components. After completing the frequency calculations ($J = NF$), the channel pressure drop and exit void fraction frequency response values are printed (subroutine TPRINT, $0 \leq NPRINT \leq 3$) and plotted (subroutine NORGRA, $0 \leq NPRINT \leq 2$). This is followed by a printout of the single phase and two-phase perturbed pressure drop components of the calculated crossover frequencies, and the component frequencies (CFREQ(N), $N = 1, NCFREQ$) read in as data.

A list of the LOCO code subprograms and their respective functions is given in Appendix E. Spatial integration and default procedures are described in more detail in the following sections.

10.3 Spatial Integration Procedures

The NCH channel sections (131) are further subdivided by the program to ensure that the finite difference integration is numerically stable. In this respect, the spatial resolution of the subcooled or saturated boiling boundaries is particularly important. The shift in the boiling boundary position for changes in an operating variable (flow rate, power input etc.) must be continuous (or as smooth as possible), otherwise the steady state and, more particularly, the transient variables exhibit numerical instabilities.

The subdivision of the channel sections is performed in subroutine STATE, and the number of nodes (NODE) and subnodes (NBOIL) in each section is stored (COMMON/STATIC/..) and transferred to subroutine TCOMP, where it is used in the transient calculations. The steps in the spatial integration procedure are as follows:

- a) the sections in the single phase region are initially divided into 5 nodes (NODE = 5) and 1 subnode (NBOIL = 1);
- b) for each section in the single phase region, the program per-

forms a heat balance based on the fluid conditions at the section inlet:

c) if the section power input is 0.80 times the power required to boil the coolant, the number of section nodes (NODE) is calculated by

$$\text{NODE} = \text{DZ}(K)/0.50 + 1 < 50 \quad (132)$$

where DZ(K) is the Kth section length (cm);

d) for each node, another heat balance is performed, and if the nodal power input is 0.90 times the power required to boil the coolant, the node is subdivided into NBOIL subnodes, where

$$\text{NBOIL} = \text{DZ}(K)/\text{NODE}/0.10 + 1 \quad (133)$$

e) the boiling boundary is shifted to the beginning of the Ith subnode (labelled NSAT) where boiling occurs;

f) after the boiling boundary position has been calculated, the number of nodes and subnodes are respectively

$$\text{NODE} = 5 \quad (134)$$

and $\text{NBOIL} = (\text{NODES}-I)/\text{NODE}/(\text{NCH}-K+1) + 1$

where the maximum number of subnodes has been set at 600. (NODES <=600).

This procedure is designed to yield numerically stable results, but care must be exercised in deciding how the channel is sectioned, particularly in regions of high power input (channel region) or large pressure drop (valves, restrictions, etc.). The accuracy of the numerical calculation may be tested by increasing the number of sections in these regions and noting the change in numerical results for key variables (e.g. boiling boundary position, channel pressure drop, exit void fraction) in the steady state and transient calculations.

10.4 Abnormal Termination Procedures

Statements have been incorporated in subroutine STATE which will

terminate program execution when

a) the static pressure at the exit of a subnode, $P(I+1)$, is less than the minimum saturation pressure, $PRESS(1)$, in the fluid property tables;

b) sufficient heat is removed from a section to cause condensation of the two-phase mixture.

The first case only applies when $|NCARDS| > 0$. The subroutine prints the variables in the subnode where (a) applies, and returns control to subroutine EXEC, which proceeds to the next case or terminates execution.

In the second case the steady state calculation continues, but no transient calculations are performed (NDYN is set to zero). The present version of the LOCO code does not include a model for the condensation boundary.

11. SAMPLE TEST CASE

The transfer function of a test channel is calculated to demonstrate how the LOCO code may be used to assess its flow stability.

11.1 Test Channel

The test channel is connected between headers at the inlet and outlet, and consists of an inlet pipe with a venturi flowmeter, a heated section made of Pyrex glass tube with a concentric electric heater rod, and a glass riser section. The principal dimensions of the test channel are given in Appendix F. Details of the test channel instrumentation and hydrodynamic experimental results are given in the references by Romberg (1978) and Romberg and Rees (1980).

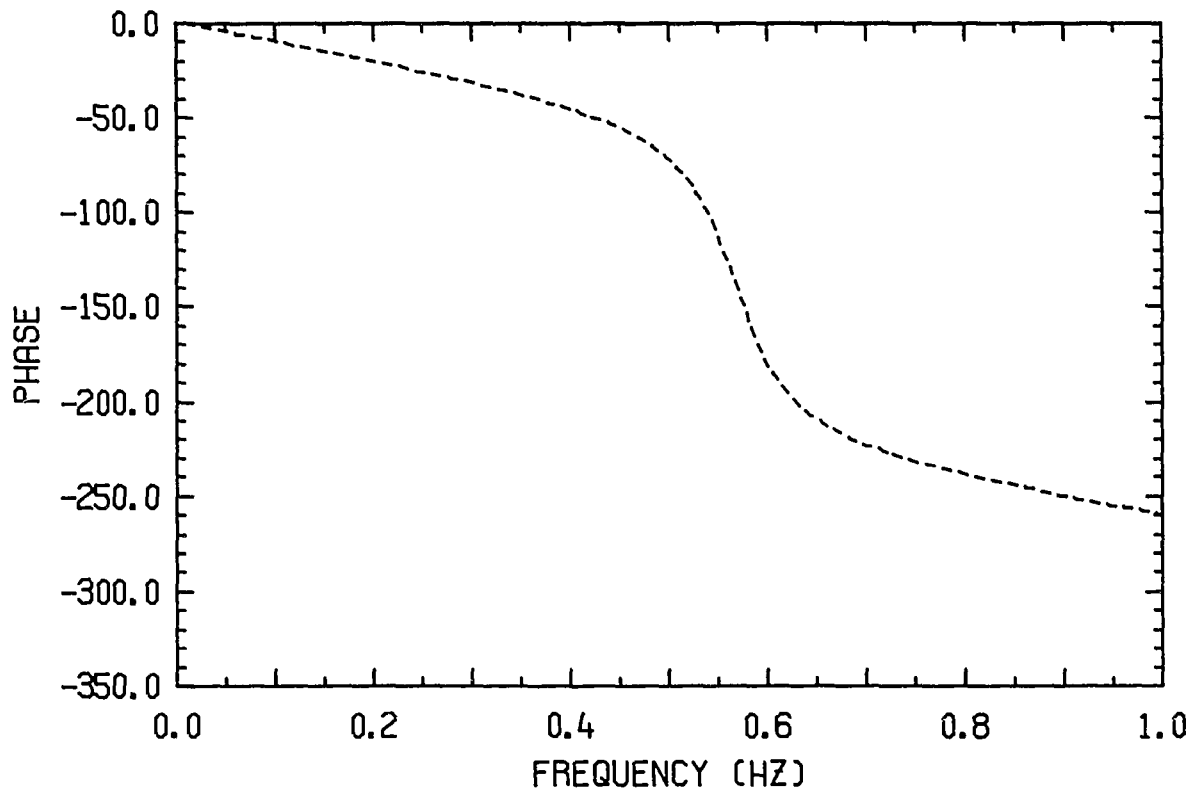
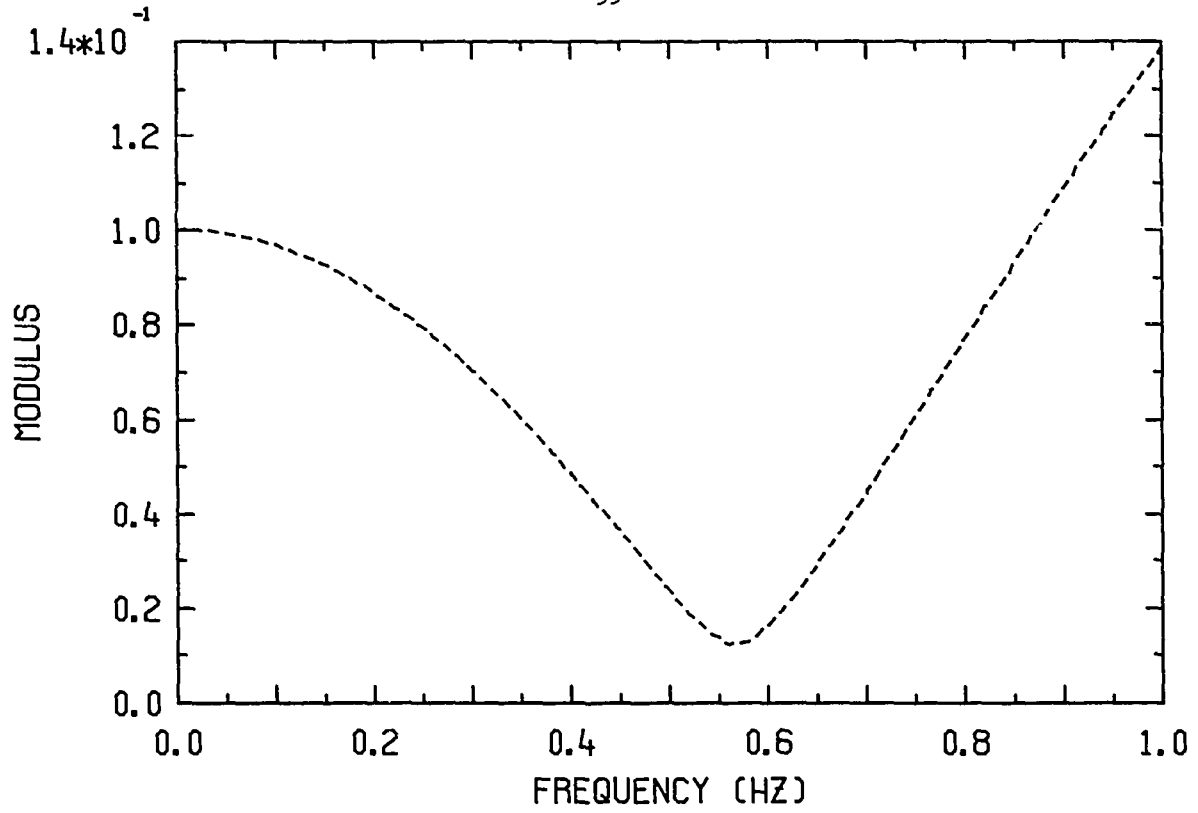


FIGURE 9 - FLOW - PRESSURE DROP TRANSFER FUNCTION

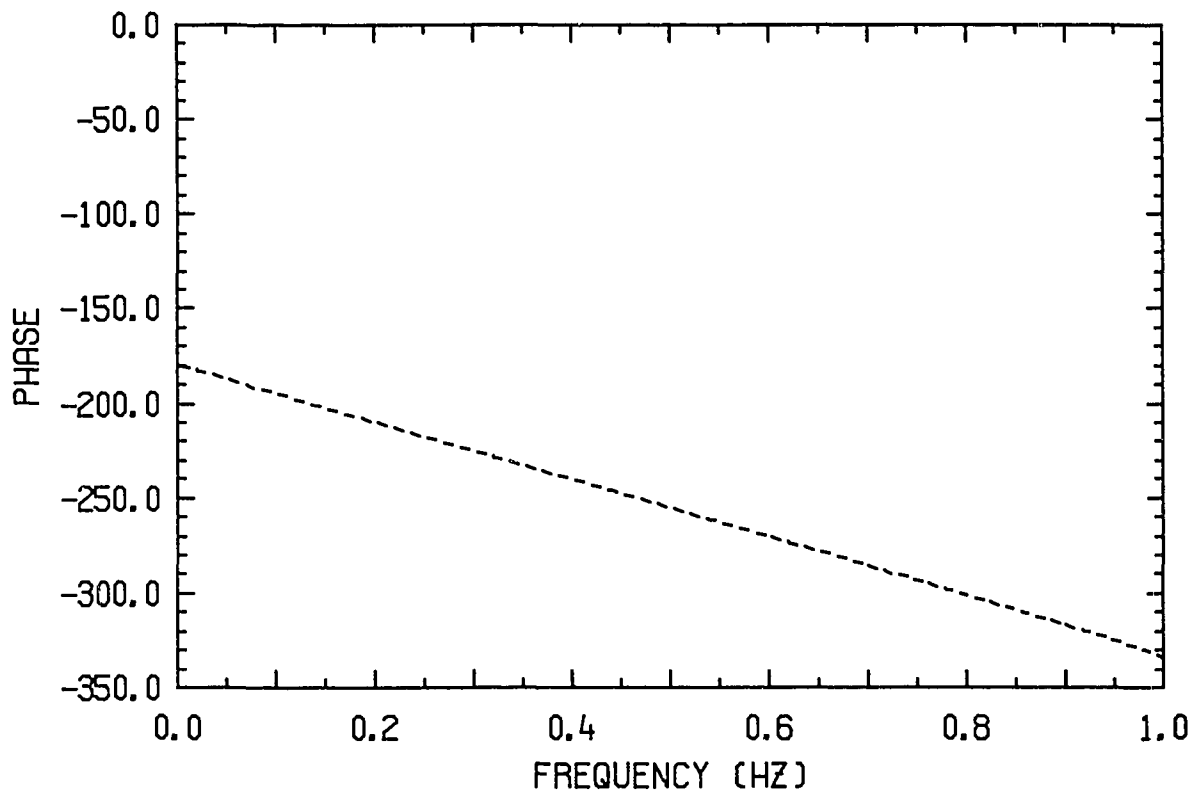
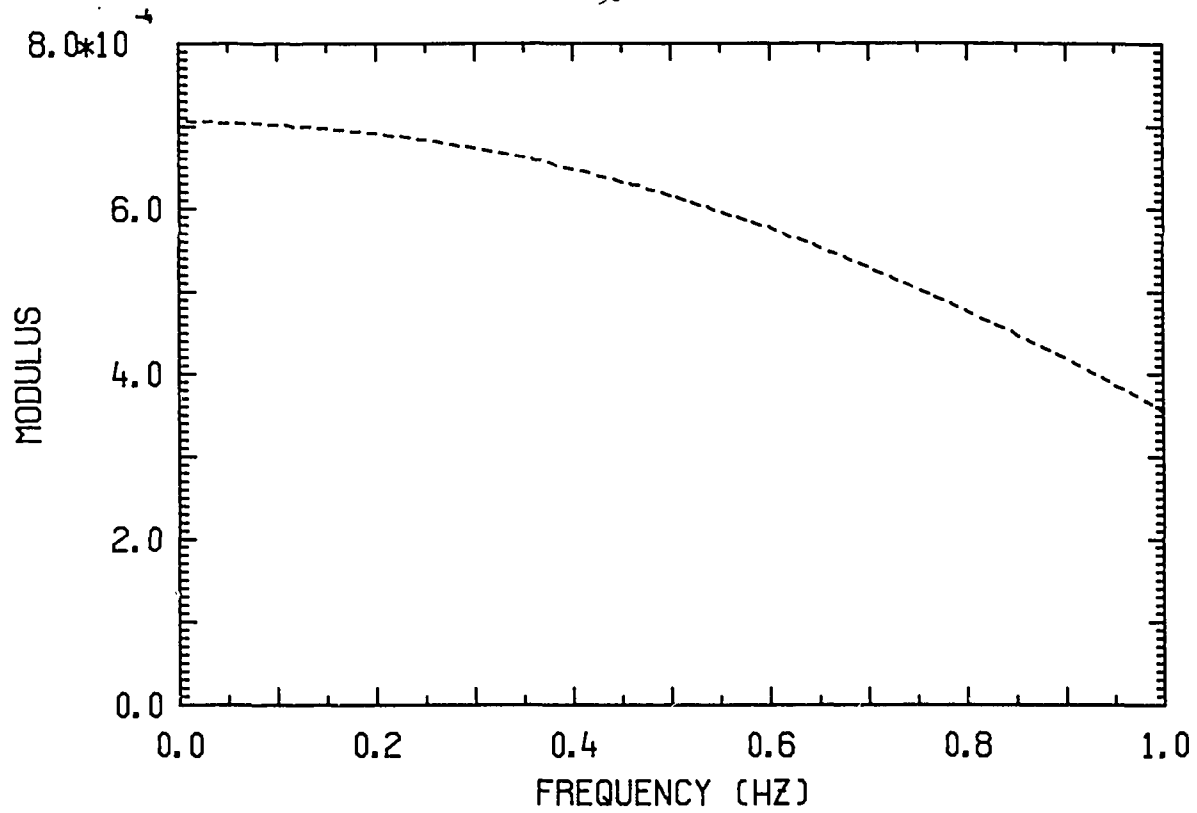


FIGURE 10 - FLOW - EXIT VOID FRACTION TRANSFER FUNCTION

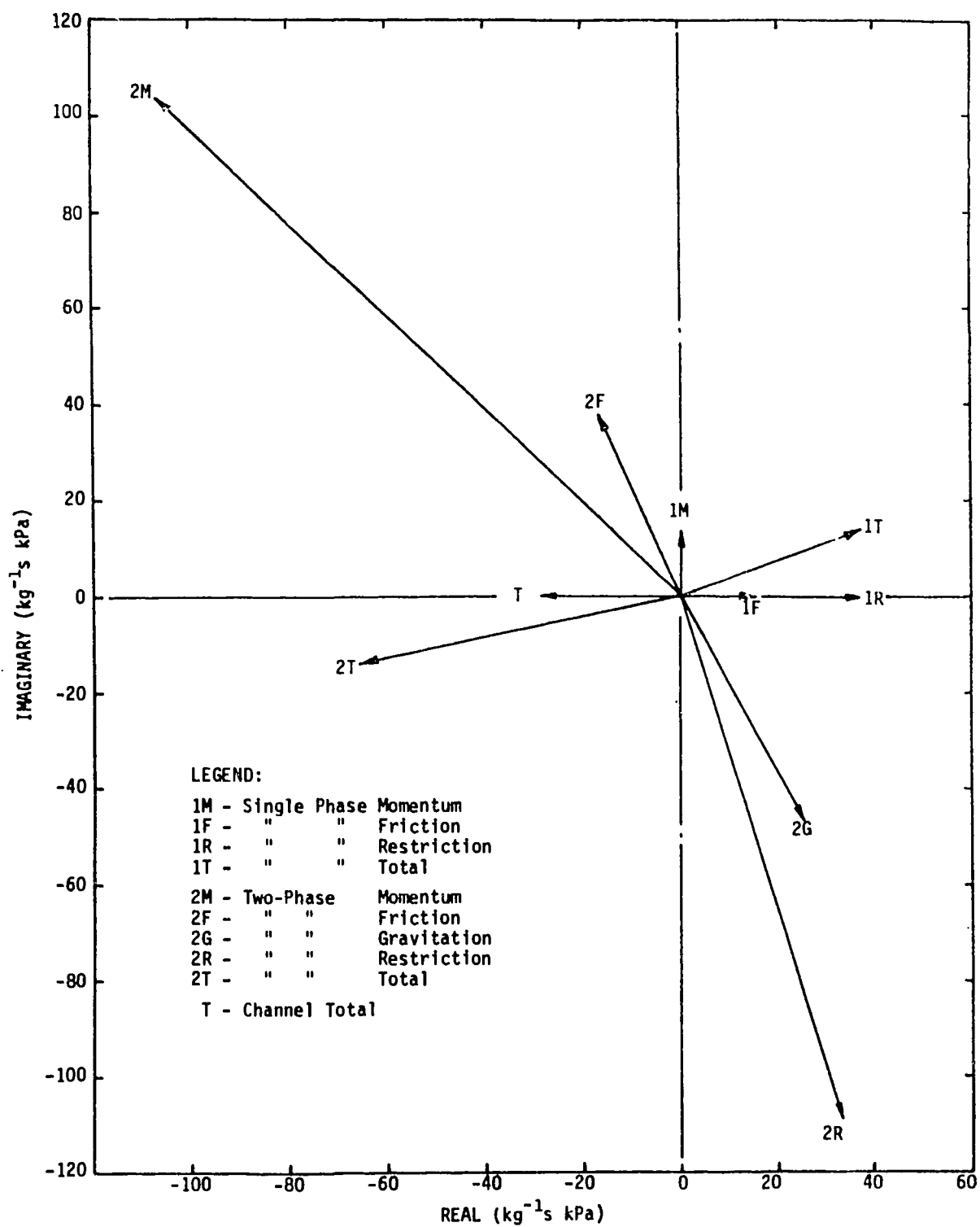


FIGURE 11 - VECTOR DIAGRAM OF TRANSFER FUNCTION COMPONENTS (1.84 kW)
(Crossover frequency = 0.595 Hz)

11.2 LOCO Results

A printout of the computer program input data is given in Appendix G. The inlet flow-pressure drop and inlet flow-exit void fraction transfer functions are plotted in Figures 9 and 10 respectively. The phase values of the inlet flow-pressure drop transfer function (hydraulic impedance) show that there is an encirclement of the origin, and thus, as discussed in Section 9.1, the channel is unstable. The code automatically prints the transfer function components at which it crosses the real axis, and these are plotted in Figure 11. The resultant vector (T) is to the left of the origin, which confirms that the channel is operating beyond the flow stability threshold.

Frequency response results, and their comparison with spectral estimates computed from the hydrodynamic measurements, are presented in the source references.

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NOMENCLATURE

<u>Symbol</u>	<u>Meaning</u>	<u>Units</u>
A	flow area or cross-sectional area	m^2
A_p	equation (152)	m
a	equation (A.11)	-
a	equation (37)	$kg^2 J^{-1} m^{-3}$
a_t	equation (B.8)	$W m^{-2} o_C^{-1}$
B_f	equation (A.9a)	-
B_p	equation (152)	$m^{-2} s$
b	equation (A.11)	-
b	equation (37)	$m^{-2} s^{-2}$
b_t	equation (B.8)	-
C_o	distribution parameter (C.6)	-
C_p	equation (106)	$m^{-5} s^2$
C	specific heat	$J kg^{-1} o_C^{-1}$
c	equation (A.11)	-
c	equation (40)	$kg m^{-3}$
D_e	equivalent hydraulic diameter (A.6)	m
D_h	equivalent heated diameter = $4A/P_h$	m
E	total energy per unit time (53)	W
E_{HC}	heat source/cladding thermal resistance (B.9)	-
e	mixture specific kinetic energy (54)	$m^2 s^{-2}$
$\tilde{F}, \tilde{F}(z)$	arbitrary perturbed, transformed variable (13)	-
F^+	subcooled boiling region correlation (20)	-
f	single phase friction factor (A.2)	-
G	mass flux	$kg m^{-2} s^{-1}$
g	gravitational acceleration	$m s^{-2}$
$\tilde{H}(y,x)$	transfer function between x (input), y (output)	-
H_p	pump head (106)	m

<u>Symbol</u>	<u>Meaning</u>	<u>Units</u>
h	specific enthalpy	J kg^{-1}
h_t	heat transfer coefficient (B.8)	$\text{W m}^{-2} \text{ } ^\circ\text{C}^{-1}$
h^+	dimensionless specific enthalpy (18)	-
J	mixture volumetric flux (C.4)	m s^{-1}
K_f	single phase friction loss factor (A.5)	m^{-1}
K_o	Bankoff factor (C.11)	-
K_R	single phase restriction loss factor	-
k	thermal conductivity	$\text{W m}^{-1} \text{ } ^\circ\text{C}^{-1}$
Nu	Nusselt number (79)	-
P	surface perimeter	m
\tilde{P}	transfer function (43)	m^{-1}
Pe	Peclet number (79)	-
p	static pressure	Pa
Q	power density	W m^{-3}
q	heat flux	W m^{-2}
Re	Reynolds number (A.7)	-
r	radius vector	m
S	Slip ratio	-
St	Stanton number (79)	-
s	Laplace transform operator	s^{-1}
T	temperature	$^\circ\text{C}$
\tilde{T}	transfer function (43)	s
\tilde{T}_p	transfer function (43)	$\text{kg}^{-1} \text{ m}^3 \text{ s}$
t	time	s
U	velocity	m s^{-1}
ΔU_g	vapour drift velocity (C.5)	m s^{-1}
u	specific internal energy	J kg^{-1}
V	weighted mean vapour drift velocity (C.12)	m s^{-1}
v	specific volume	$\text{m}^3 \text{ kg}^{-1}$

<u>Symbol</u>	<u>Meaning</u>	<u>Units</u>
W	mass flow	kg s ⁻¹
X	equation (C.19)	-
x	flow quality	-
z	axial coordinate	m
<u>Greek Symbols</u>		
α	void fraction	-
Γ	vapour source term (35)	kg m ⁻³ s ⁻¹
Δ	difference operator	-
ϵ	relative roughness ratio	-
θ	angle of inclination to vertical	-
μ	dynamic viscosity	Pa s
ρ	density	kg m ⁻³
σ	real component of Laplace operator	s ⁻¹
τ	shear stress	Pa
τ	time constant	s
ω	angular frequency	s ⁻¹
<u>Subscripts</u>		
B	bypass	
C	cladding	
H	heat source	
I	restriction inertia	
L	restriction loss	
P	pump	
R	restriction total	
T	total	
f	saturated liquid; friction component	
g	saturated vapour; gravitational component	
h	heated dimension or variable	

<u>Symbol</u>	<u>Meaning</u>	<u>Units</u>
l	subcooled liquid	
m	mixture; momentum component	
p	pressure	
r	radial component	
s	source	
w	wall; wetted surface	
0	subcooled boiling boundary	
1	space node inlet; channel inlet; single phase	
2	space node outlet; two-phase	

Superscripts

\sim	Laplace transformed variable; tensor
0	steady state variable
$'$	upstream of boiling boundary
$''$	downstream of boiling boundary

APPENDIX A.

SINGLE PHASE FRICTION FACTOR AND TWO-PHASE MULTIPLIER CORRELATIONSA1. Single Phase Friction Factor

Integration of the momentum conservation equation (3) over the flow area requires evaluation of the shear stress integral

$$\begin{aligned} \frac{1}{A} \int_A \frac{1}{r} \frac{\partial}{\partial r} (r \bar{\tau}_{rz}) dA &= \frac{1}{A} \int_0^R \frac{1}{r} \frac{\partial}{\partial r} (r \bar{\tau}_{rz}) 2\pi r dr \\ &= \frac{2\pi R}{A} \tau_w = \frac{P_w}{A} \tau_w \quad (\text{round tubes}) \\ &= \frac{1}{A} \sum_{j=1}^N P_{wj} \tau_{wj} \quad (\text{annuli, rod clusters}) \end{aligned} \quad (A1)$$

The wall shear stress (τ_w) is related to the single phase friction factor (f) by

$$\tau_w = f \frac{\rho_l \langle U_l \rangle^2}{2} \quad (A2)$$

where f is a function of the single phase Reynolds number (Re) and the relative roughness ratio (ϵ) of the channel walls, and is given by the Colebrook equation (Knudsen and Katz 1958)

$$\frac{1}{\sqrt{f}} = 4 \log \left[\frac{Re \sqrt{f}}{\epsilon Re \sqrt{f} + 4.67} \right] + 2.28 \quad (A3)$$

From (A1) and (A2) the single phase pressure gradient is

$$\frac{\partial p_{1f}}{\partial z} = K_f \frac{\rho_l \langle U_l \rangle^2}{2} = K_f \frac{G^2}{2\rho_l} \quad (A4)$$

in which the single phase friction loss factor is

$$K_f = \frac{P_w f}{A} = \frac{4f}{D_e} \quad (A5)$$

where D_e is the equivalent hydraulic diameter:

$$D_e = \frac{4 \times \text{flow area}}{\text{wetted perimeter}} = \frac{4A}{P_w} \quad (\text{A6})$$

A2. Transient Single Phase Friction Factor

From (A5) the perturbed and transformed friction loss factor is

$$\begin{aligned} \tilde{K}_f &= \frac{4}{D_e} \tilde{f} = \frac{4}{D_e} \left(\frac{\partial f^0}{\partial \text{Re}^0} \right) \tilde{\text{Re}} \\ &= \frac{4}{D_e} \text{Re}^0 \left(\frac{\partial f^0}{\partial \text{Re}^0} \right) \frac{\tilde{G}}{G^0} \end{aligned} \quad (\text{A7})$$

where $\text{Re} = \frac{G D_e}{\mu_\ell}$ and $G = \rho_\ell \langle U_\ell \rangle$

$$\text{From (A3)} \quad \frac{\partial f^0}{\partial \text{Re}^0} = - \left[\frac{2m}{m + B \text{Re}^0} \right] \frac{f^0}{\text{Re}^0} \quad (\text{A8})$$

$$\text{where} \quad m = \frac{4 \times 4.67}{2.3 \ln 10} = 8.11262 \quad (\text{A8a})$$

$$\text{and} \quad B = \epsilon + \frac{4.67}{\text{Re}^0 \sqrt{f^0}} \quad (\text{A8b})$$

Substitution of (A8) in (A7) gives

$$\frac{\tilde{K}_f}{K_f^0} = - B_f \frac{\tilde{G}}{G^0} \quad (\text{A9})$$

$$\text{where} \quad B_f = \frac{2m}{m + B \text{Re}^0} \quad (\text{A9a})$$

A3. Local Two-phase Friction Multiplier

The two-phase friction pressure gradient is expressed as a function of the single phase pressure gradient by

$$\frac{\partial p_{2f}}{\partial z} = \frac{\partial p_{1f}}{\partial z} \phi^2 \quad (\text{A10})$$

where ϕ^2 is the local two-phase friction multiplier, and is given by

the correlation (Beattie 1973)

$$\phi^2 = 1 + a \left[\left(\frac{\rho_l}{\rho_g} - 1 \right) x \right]^b G^c \quad (A11)$$

In the single phase region $\phi^2=1$, and for homogeneous two-phase flow $a=1$, $b=1$ and $c=0$.

A4. Transient Two-phase Friction Multiplier

From (A3.11) the perturbed two-phase friction multiplier is

$$\tilde{\phi}^2 = \left(\frac{\partial \phi^{0^2}}{\partial x^0} \right) \tilde{x} + \left(\frac{\partial \phi^{0^2}}{\partial G^0} \right) \tilde{G} + \left(\frac{\partial \phi^{0^2}}{\partial \rho_g^0} \right) \tilde{\rho}_g \quad (A12)$$

that is,
$$\tilde{\phi}^2 = (\phi^{0^2}-1) \left[b \left(\frac{\tilde{x}}{x^0} - \frac{\rho_l}{\Delta \rho} \frac{\rho_g^0}{\rho_g} \tilde{p} \right) + c \frac{\tilde{G}}{G^0} \right] \quad (A13)$$

APPENDIX B

HEAT SOURCE/COOLANT HEAT TRANSFERB1. Heat Source Cladding Equations

Neglecting axial conduction and the temperature dependence of the material properties, the heat source energy equation for an arbitrary space node is

$$\rho_H C_H \frac{\partial}{\partial t} [T_H(r,t)] + \frac{1}{r} \frac{\partial}{\partial r} [r q_r(r,t)] = Q_s(r,t) \quad (B1)$$

The energy equation for the adjacent cladding is

$$\rho_c C_c \frac{\partial}{\partial t} [T_c(r,t)] + \frac{1}{r} \frac{\partial}{\partial r} [r q_r(r,t)] = 0 \quad (B2)$$

Integration of (B1) and (B2) over their respective cross-sectional areas gives

$$\rho_H C_H \frac{d}{dt} T_H + \frac{(P_H q_H)_o - (P_H q_H)_i}{A_H} = Q_s \quad (B3)$$

$$\text{and } \rho_c C_c \frac{d}{dt} T_c + \frac{(P_c q_c)_o - (P_c q_c)_i}{A_c} = 0 \quad (B4)$$

where, in general,

$$\left. \begin{aligned} T &= T(t) = \frac{1}{A} \int_A T(r,t) dA \\ Q_s &= Q_s(t) = \frac{1}{A} \int_A Q_s(r,t) dA \\ \text{and } \frac{1}{A} \int_A \frac{1}{r} \frac{\partial}{\partial r} (r q_r) dA &= \frac{(P q)_o - (P q)_i}{A} \end{aligned} \right\} \quad (B5)$$

in which the subscripts i,o refer to the inner and outer perimeters respectively. In the last equation of (B5) the radius and flux vectors are assumed to act in the same direction. If they act in opposite directions then the right hand side has a negative sign.

Equations (B3) and (B4) are solved for an annular channel in which the inner or outer wall is heated. The boundary conditions at the heat source/cladding interface and the cladding/coolant interface are respectively

$$a) \quad (q_H)_i = 0; \quad (q_H)_o = (q_C)_i = q_{HC} \quad (B6)$$

$$\text{and } b) \quad (q_C)_o = q_w; \quad q_w = h_t (T_w - T) \quad (B7)$$

where h_t , the local heat transfer coefficient, is a function of the local Reynolds number (Re) and Prandtl number (Pr), and is given by the correlation

$$h_t = a_t (Re)^{b_t} (Pr)^{c_t} \quad (B8)$$

If it is assumed that

$$T_c = E_{HC} T_H = T_w \quad (B9)$$

where E_{HC} is a thermal resistance factor to account for temperature gradients across the heat source/cladding interface, then (B3) and (B4) become respectively

$$\frac{A_H \rho_H C_H}{E_{HC}} \frac{d}{dt} [T_w] + P_{HC} q_{HC} = A_H Q_S \quad (B10)$$

$$\text{and } A_C \rho_C C_C \frac{d}{dt} [T_w] + P_C q_w - P_{HC} q_{HC} = 0 \quad (B11)$$

Adding (B10) and (B11) yields:

$$\left[\frac{\rho_H C_H}{E_{HC}} + \frac{A_C \rho_C C_C}{A_H} \right] \frac{d}{dt} [T_w] + \frac{P_C q_w}{A_H} = Q_S \quad (B12)$$

Perturbation, linearisation and Laplace transformation of (B12) gives the steady state and transient equations respectively as

$$\frac{P_C q_w^0}{A_H} = Q_S^0 \quad (B13)$$

$$\left[\frac{\rho_H C_H}{E_{HC}} + \frac{A_C \rho_C C_C}{A_H} \right] s \tilde{T}_W + \frac{P_C}{A_H} \tilde{q}_W = \tilde{Q}_S \quad (B14)$$

The cladding/coolant boundary condition (B7) is used to eliminate the perturbed wall temperature \tilde{T}_W from (B14). Perturbation and transformation of (B7) gives:

$$\tilde{q}_W = h_t^0 (\tilde{T}_W - \tilde{T}) + \Delta T_W^0 \tilde{h}_t \quad (B15)$$

or

$$\tilde{T}_W = \frac{\tilde{q}_W}{h_t^0} - \Delta T_W^0 \frac{\tilde{h}_t}{h_t^0} + \tilde{T}$$

Substitute (B15) into (B14):

$$\frac{\tilde{q}_W}{q_W^0} = \frac{1}{(1+\tau_H s)} \frac{\tilde{Q}_S}{Q_S^0} + \frac{\tau_H s}{(1+\tau_H s)} \frac{\tilde{h}_t}{h_t^0} - \frac{\tau_H s}{(1+\tau_H s)} \frac{\tilde{T}}{\Delta T_W^0} \quad (B16)$$

where τ_H is the fundamental mode time constant, and is given by

$$\tau_H = \frac{A_H}{P_C h_t^0} \left[\frac{\rho_H C_H}{E_{HC}} + \frac{A_C \rho_C C_C}{A_H} \right] \quad (B17)$$

From (B8) the perturbed heat transfer coefficient is

$$\tilde{h}_t = \left\{ \frac{\partial h_t^0}{\partial Re} \quad \frac{\partial Re}{\partial G^0} \right\} \tilde{G} = b_t h_t^0 \frac{\tilde{G}}{G^0} \quad (B18)$$

and the perturbed liquid enthalpy is

$$\tilde{h}_\ell = \left(\frac{\partial h_\ell^0}{\partial T^0} \right) \tilde{T} = C_{P\ell} \tilde{T}$$

Hence (B16) becomes

$$\frac{\tilde{q}_W}{q_W^0} = \frac{1}{1+\tau_H s} \left[\frac{\tilde{Q}_S}{Q_S^0} + b_t \tau_H s \frac{\tilde{G}}{G^0} - \frac{\tau_H s}{C_{P\ell} \Delta T_W^0} \tilde{h}_\ell \right] \quad (B19)$$

This equation applies for the single phase and subcooled boiling regions of the channel. In the saturated boiling region

$$\tilde{h}_l = \tilde{h}_f = 0 \quad (\text{B20})$$

Equation (B19) is applicable to other channel geometries by alteration of equation (B17) for the overall time constant τ_H . An average 'thermal capacitance factor' (stored heat/unit surface area/unit temperature) is defined as

$$\langle \theta_T \rangle = \frac{\sum_n^N \left[\frac{A_H \rho_H C_H}{E_{HC}} + A_C \rho_C C_C \right]_n}{\sum_n^N [P_C]_n} \quad (\text{B21})$$

where the cladding values apply to those between the heat source (or sources) and the coolant only. The heat transfer coefficient implicitly becomes

$$\langle h_t^0 \rangle = h_t^0 = \frac{1}{N} \sum_n^N [h_t^0]_n \quad (\text{B22})$$

Hence the average fundamental mode time constant is

$$\langle \tau_H \rangle = \tau_H = \frac{\langle \theta_T \rangle}{\langle h_t^0 \rangle} \quad (\text{B23})$$

Note that heat source/coolant heat transfer is neglected if

$$\tau_H = 0 \quad (\text{B24})$$

APPENDIX C

TWO-PHASE SLIP CORRELATION AND VOID/QUALITY RELATIONSC1. Two-phase Slip Correlation

The slip ratio, or holdup ratio (Govier and Aziz 1972), is defined as the ratio of the weighted cross-section averaged vapour and liquid velocities:

$$S = \frac{\langle \bar{U}_g \rangle}{\langle \bar{U}_l \rangle} \quad (C1)$$

The weighted cross-section averaged vapour and liquid velocities are defined as

$$\left. \begin{aligned} \langle \bar{U}_g \rangle &= \frac{\int U_g dA_g}{\int dA_g} = \frac{\int \alpha U_g dA}{\int \alpha dA} = \frac{\langle \alpha U_g \rangle}{\langle \alpha \rangle} \\ \langle \bar{U}_l \rangle &= \frac{\int U_l dA_l}{\int dA_l} = \frac{\int (1-\alpha) U_l dA}{\int (1-\alpha) dA} = \frac{\langle (1-\alpha) U_l \rangle}{\langle 1-\alpha \rangle} \end{aligned} \right\} \quad (C2)$$

Hence (C1) becomes

$$S = \frac{\langle \alpha U_g \rangle}{\langle \alpha \rangle} \frac{\langle 1-\alpha \rangle}{\langle (1-\alpha) U_l \rangle} \quad (C3)$$

Zuber and Findlay (1965) define the volumetric flux (mixture velocity) by

$$J = (1-\alpha) U_l + \alpha U_g \quad (C4)$$

a vapour drift velocity by

$$\Delta U_g = U_g - J \quad (C5)$$

and a distribution parameter by

$$C_0 = \frac{\langle \alpha J \rangle}{\langle \alpha \rangle \langle J \rangle} \quad (C6)$$

From (C5) and (C6) it follows that

$$\langle \alpha U_g \rangle = C_0 \langle \alpha \rangle \langle J \rangle + \langle \alpha \Delta U_g \rangle \quad (C7)$$

and substitution in (C4) gives

$$\langle (1-\alpha)U_l \rangle = \langle J \rangle (1 - C_0 \langle \alpha \rangle) - \langle \alpha \Delta U_g \rangle \quad (C8)$$

Hence the Zuber and Findlay slip correlation is

$$S = \frac{\langle 1-\alpha \rangle}{\frac{1}{C_0 + \frac{\langle \alpha \Delta U_g \rangle}{\langle \alpha \rangle \langle J \rangle}} - \langle \alpha \rangle} \quad (C9)$$

Alternatively,

$$S = \frac{\langle 1 - \alpha \rangle}{K_0 - \langle \alpha \rangle} \quad (C10)$$

$$\text{where } K_0 = \frac{\langle \alpha \rangle \langle J \rangle}{\langle \alpha U_g \rangle} = \frac{1}{C_0 + \frac{V}{\langle J \rangle}} \quad (C11)$$

$$\text{and } V = \frac{\langle \alpha \Delta U_g \rangle}{\langle \alpha \rangle} \quad (C12)$$

The factor (K_0) is commonly termed the Bankoff factor (Bankoff 1960), and (V) is termed the weighted mean (cross-section averaged) vapour drift velocity (Zuber and Findlay 1965). The distribution parameter (C_0) and the weighted mean drift velocity (V) are a function of the two-phase flow regime (bubble, slug, annular, droplet regime - Zuber et al. 1967). In the present analysis the factor (K_0) and the weighted mean drift velocity are assumed to vary in space, but invariant in time. This assumption is not unreasonable in view of the indeterminate nature of the void and velocity profiles.

C2. Void-Quality Relations

From equation (C5), the weighted mean vapour and liquid velocities are also given by

$$\langle \bar{U}_g \rangle = \frac{\langle x \rangle \langle G \rangle}{\langle \rho_g \rangle \langle \alpha \rangle} \quad (C13)$$

$$\langle \bar{U}_l \rangle = \frac{\langle 1-x \rangle \langle G \rangle}{\langle \rho_l \rangle \langle 1-\alpha \rangle}$$

Substitution in (C1) gives

$$S = \frac{\langle 1-\alpha \rangle \langle x \rangle \langle \rho_l \rangle}{\langle \alpha \rangle \langle 1-x \rangle \langle \rho_g \rangle} \quad (C14)$$

Equating (C10) and (C14) gives:

$$x = \frac{\alpha \rho_g}{K_0 \rho_l - \Delta \rho \alpha}, \quad \Delta \rho = \rho_l - \rho_g \quad (C15)$$

or alternatively,

$$\alpha = \frac{K_0 \rho_l x}{(\rho_g + \Delta \rho x)} \quad (C16)$$

where cross-section averaged variables are assumed.

From (C16) the perturbed and Laplace transformed void fraction is

$$\tilde{\alpha} = \left(\frac{\partial \alpha^0}{\partial x^0} \right) \tilde{x} + \left(\frac{\partial \alpha^0}{\partial \rho_g^0} \right) \tilde{\rho}_g \quad (C17)$$

$$\text{or} \quad \tilde{\alpha} = \frac{1}{X} \left[\tilde{x} - x^0 (1-x^0) \frac{\rho_g^0}{\rho_g^0} \tilde{p} \right] \quad (C18)$$

$$\text{where} \quad X = \frac{(\rho_g^0 + \Delta \rho x^0)^2}{K_0 \rho_l \rho_g^0} = K_0 \left(\frac{\rho_l}{\rho_g^0} \right) \left(\frac{x^0}{\alpha^0} \right)^2 \quad (C19)$$

$$\text{and} \quad \tilde{\rho}_g = \left(\frac{\partial \rho_g^0}{\partial p^0} \right) \tilde{p} = \rho_g^0 \tilde{p} \quad (C20)$$

APPENDIX D
LISTING OF 'LOCO' COMPUTER PROGRAM

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C *****
C * LOCO - A LINEARISED MODEL FOR CALCULATING THE FREQUENCY RESPONSE *
C *      AND ONSET OF COOLANT OSCILLATIONS IN BOILING CHANNELS.      *
C * AUTHOR : T. M. ROMBERG, AUSTRALIAN ATOMIC ENERGY COMMISSION. *
C *****
C
C ***GENERAL NOTE : THE DATA CARDS ARE READ IN AS INDICATED BY THE FOLL-
C OWING COMMENT CARDS. SUBSEQUENT CASES ARE READ IN AFTER THE FLUID
C PROPERTIES CARDS, WHICH ARE READ IN AFTER THE FIRST CASE ONLY.
C
C -----
C ***TITLE CARD : FORMAT (20A4)
C -----
C ***CONTROL CARDS : FORMAT (16I5)
C ND      = NUMBER OF INLET PIPING SECTIONS )
C NC      = NUMBER OF CHANNEL SECTIONS      ) TOTAL = NCH = ND+NC+NR
C NR      = NUMBER OF RISER SECTIONS        )
C NB      = NUMBER OF BYPASS PIPING SECTIONS.
C NTIN    = NUMBER OF INLET TEMPERATURE INCREMENTS OR DECREMENTS.
C NPOW    = NUMBER OF CHANNEL POWER RATIO INCREMENTS OR DECREMENTS.
C NW      = NUMBER OF FLOW RATE INCREMENTS OR DECREMENTS.
C NF      = NUMBER OF FREQUENCY VALUES.
C NCFREQ  = NUMBER OF FREQUENCIES FOR COMPONENT PRESS. DROP CALCULATIONS
C -----
C NDROP   = 0 : NO ITERATION TO FIND FLOW RATE FOR GIVEN PRESSURE DROP
C NODES   = MAXIMUM NUMBER OF SPACE NODES (<=600)
C         = 1 : ITERATION TO FIND FLOW RATE FOR GIVEN PRESSURE DROP
C NOSUR   = 0 : NO SUBCOOLED BOILING REGION
C         = 1 : SUBCOOLED BOILING REGION (EXPONENTIAL CORRELATION)
C         = 2 : SUBCOOLED BOILING REGION (HYPERBOLIC TANGENT CORRELATION)
C NDYN    = CONTROL PARAMETER FOR DYNAMIC CALCULATIONS
C         = 0 : NO DYNAMIC CALCULATIONS
C         = 1 : REAL INPUTS (FREQUENCY INDEPENDENT)
C         = 2 : REAL INPUTS (FREQUENCY DEPENDENT)
C         = 3 : COMPLEX INPUTS (FREQUENCY DEPENDENT)
C NPRINT  = PRINTOUT CONTROL PARAMETER
C         = 0 : AXIAL NODAL VALUES+PRESSURE DROP AND EXIT VOID DYNAMICS
C         = 1 : AXIAL SECTION VALUES+PRESSURE DROP AND EXIT VOID DYNAMICS
C NPLOTT  = PLOT OPTION
C         = 0 : NO PLOTS
C         = 1 : BODE PLOTS OF PRESSURE DROP AND EXIT VOID DYNAMICS
C NCARDS  = NUMBER OF FLUID PROPERTY DATA CARDS.
C         = +VE : FLUID PROPERTIES ARE PRINTED (NOTE : NCARDS>3)
C         = -VE : FLUID PROPERTIES ARE NOT PRINTED (NOTE : NCARDS<-3)
C         = 0  : CONSTANT FLUID PROPERTIES.
C -----
C NGDYN   = INPUT DEVICE OPTION FOR MASS FLUX FREQUENCY DATA
C NPDYN   = INPUT DEVICE OPTION FOR INLET PRESSURE FREQUENCY DATA
C NHGYN   = INPUT DEVICE OPTION FOR INLET ENTHALPY FREQUENCY DATA
C NQDYN   = INPUT DEVICE OPTION FOR HEAT SOURCE FREQUENCY DATA
C NSPEC   = INPUT DEVICE FOR PRESS DROP+EXIT VOID FREQUENCY DATA
C NPROP   = INPUT DEVICE OPTION FOR FLUID PROPERTIES DATA
C ***NOTE : NXXXX - NO INPUT=0, CARDS=1, DATASET=N : NPROP=1 IF NCARDS=0
C -----

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C***THE FOLLOWING CARDS HAVE FORMAT (8F10.3)
C TIN    = INLET TEMPERATURE TO SECTION 1.....C
C DTIN   = +VE : INCREMENT IN INLET TEMPERATURE.....C
C        = -VE : DECREMENT IN INLET TEMPERATURE.....C
C -----
C POWR   = CHANNEL REGION POWER RATIO (POW(N)=PKW(N)*POWR)
C DPOWR  = +VE : INCREMENT IN CHANNEL REGION POWER RATIO
C        = -VE : DECREMENT IN CHANNEL REGION POWER RATIO
C -----
C W      = INITIAL MASS FLOW RATE.....KG/S
C DW     = +VE : INCREMENT IN MASS FLOW RATE.....KG/S
C        = -VE : DECREMENT IN MASS FLOW RATE.....KG/S
C -----
C PIN    = INLET PRESSURE TO SECTION 1.....KPA
C PDROP  = PRESSURE DROP BETWEEN INLET AND OUTLET.....KPA
C        = 0.0 FOR NDROP = 0 OR NB > 0
C WTOT   = TOTAL MASS FLOW RATE THROUGH CHANNEL AND BYPASS.....KG/S
C AP,BP, = COEFFICIENTS FOR PUMP HEAD-FLOW CHARACTERISTIC,
C CP     = (HEAD=AP+BP*Q+CP*Q*Q, HEAD=M, Q=LITRE/S)
C -----
C ASB    = LOCAL NUSSELT NUMBER FOR ONSET OF SUBCOOLED BOILING
C        (PECLET NO<=70000 : ASB=455.0 NORMALLY)
C BSB    = LOCAL STANTON NUMBER FOR ONSET OF SUBCOOLED BOILING
C        (PECLET NO >70000 : BSB=0.0065 NORMALLY)
C        NOTE : ASB AND BSB ARE RELATED BY THE EQUATION ASB=70000*BSB
C AHT,   = COEFFICIENTS FOR THE SUBCOOLED AND SATURATED BOILING REGION
C BHT,   = HEAT TRANSFER CORRELATION : HT = AHT*RE**BHT*PRL**CHT,
C CHT    = WHERE RE = REYNOLDS NUMBER, PRL = LIQUID PRANDTL NUMBER
C -----
C***FOLLOWING ELEVEN (11) CARDS ARE NOT REQUIRED IF NDYN=0.
C SIGMA  = REAL COMPONENT OF THE LAPLACE OPERATOR.....HZ
C FREQ   = INITIAL FREQUENCY FOR DYNAMIC CALCULATIONS.....HZ
C -----
C DFREQ  = FREQUENCY INCREMENT FOR DYNAMIC CALCULATIONS.....HZ
C -----
C CARD NOT REQUIRED IF NDYN = 0 OR NCFREQ = 0 : N=1,NCFREQ
C CFREQ(N)=FREQUENCIES FOR COMPONENT PRESSURE DROP CALCULATIONS.....HZ
C -----
C GDFAC  = NORMALISING FACTOR FOR INLET MASS FLUX PERTURBATION.
C PDFAC  = NORMALISING FACTOR FOR INLET PRESSURE PERTURBATION.
C HDFAC  = NORMALISING FACTOR FOR INLET ENTHALPY PERTURBATION.
C QDFAC  = NORMALISING FACTOR FOR CHANNEL POWER DENSITY PERTURBATION.
C -----
C***MODULUS (NDYN<3) OR REAL COMPONENT (NDYN=3).
C GDYN(N) = SPECTRUM OF INLET MASS FLUX PERTURBATION.....KG/(S*M**2)
C -----
C***COMPLEX COMPONENT (REQUIRED ONLY IF NDYN=3).
C GCDYN(N) = SPECTRUM OF INLET MASS FLUX PERTURBATION.....KG/(S*M**2)
C -----
C***MODULUS (NDYN<3) OR REAL COMPONENT (NDYN=3).
C PDYN(N) = SPECTRUM OF INLET PRESSURE PERTURBATION.....KPA
C -----
C***COMPLEX COMPONENT (REQUIRED ONLY IF NDYN=3).
C PCDYN(N) = SPECTRUM OF INLET PRESSURE PERTURBATION.....KPA
C -----
C***MODULUS (NDYN<3) OR REAL COMPONENT (NDYN=3).
C HDYN(N) = SPECTRUM OF INLET ENTHALPY PERTURBATION.....KJ/KG
C -----
C***COMPLEX COMPONENT (REQUIRED ONLY IF NDYN=3)
C HCDYN(N) = FREQUENCY SPECTRUM OF INLET ENTHALPY PERTURBATION.....KJ/KG
C -----

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C***MODULUS (NDYN<3) OR REAL COMPONENT (NDYN=3)
C QDYN(N) = FREQUENCY SPECTRUM OF NORMALISED CHANNEL POWER PERT.....W/W
C -----
C***COMPLEX COMPONENT (REQUIRED ONLY IF NDYN=3)
C QCDYN(N) = FREQUENCY SPECTRUM OF NORMALISED CHANNEL POWER PERT.....W/W
C -----
C***SECTION SPECIFICATIONS (N=1,NCH) :
C DZ(N) = LENGTH.....CM
C A(N) = FLOW AREA.....CM**2
C DE(N) = EQUIVALENT HYDRAULIC DIAMETER.....CM
C PH(N) = HEATED PERIMETER.....CM
C ANG(N) = ANGLE TO VERTICAL (UPFLOW=0.0 : DOWNFLOW=180.0).....DEG

C RK(N) = INLET RESTRICTION LOSS FACTOR
C E(N) = SURFACE RELATIVE ROUGHNESS RATIO
C -----
C PKW(N) = POWER INPUT (+VE, ZERO OR -VE).....KW
C TCAP(N) = THERMAL CAPACITANCE OF HEAT SOURCE AND CLADDING.....KJ/(M*S)
C CO(N) = DISTRIBUTION PARAMETER
C V(N) = WEIGHTED MEAN VAPOUR DRIFT VELOCITY.....M/S
C AA(N), = LOCAL TWO-PHASE FRICTION MULTIPLIER COEFFICIENTS :
C BB(N), =  $PHI = 1.0 + AA(N) * ((RHOL/RHOG - 1.0) * X) ** BB(N) * G ** CC(N)$ , WHERE
C CC(N) = RHOL,RHOG=LIQUID,VAPOUR DENSITIES:X=FLOW QUALITY:G=MASS FLUX
C -----
C***BYPASS SECTION SPECIFICATIONS (NB>0 ONLY : N=1,NB)
C ZB(N) = LENGTH.....CM
C AB(N) = FLOW AREA.....CM**2
C DB(N) = EQUIVALENT HYDRAULIC DIAMETER.....CM
C ANGB(N)= ANGLE OF INCLINATION TO VERTICAL.....DEG
C RKB(N) = INLET RESTRICTION FACTOR
C EB(N) = SURFACE RELATIVE ROUGHNESS RATIO
C -----
C TPL0T1 = PLOT TITLE CARD FOR PRESSURE DROP DYNAMICS (NPL0T,NE,0)
C TPL0T2 = PLOT TITLE CARD FOR EXIT VOID FRACTION DYNAMICS (NPL0T,NE,0)
C -----
C***SATURATION FLUID PROPERTY DATA CARDS (NCARDS,NE,0 : N=1,NCARDS)
C***FORMAT (F5.1,F9.2,F8.2,F8.1,F10.6,F10.5,F6.1,F8.4,F8.3)
C TEMP(N) = TEMPERATURE.....C
C PRESS(N) = PRESSURE.....KPA
C ENTHF(N) = LIQUID ENTHALPY.....KJ/KG
C ENTHG(N) = VAPOUR ENTHALPY.....KJ/KG
C SPVF(N) = LIQUID SPECIFIC VOLUME.....M**3/KG
C SPVG(N) = VAPOUR SPECIFIC VOLUME.....M**3/KG
C VISCOF(N)= LIQUID DYNAMIC VISCOSITY.....MICRO PA*S
C CONDF(N) = LIQUID THERMAL CONDUCTIVITY.....W/(M*K)
C SPHF(N) = LIQUID SPECIFIC HEAT.....KJ/(KG*K)
C -----
C***SATURATION FLUID PROPERTIES DATA CARD (NCARDS=0)
C***FORMAT (F5.1,F8.2,F8.1,F10.6,F10.5,F6.1,F8.4,F8.3)
C TSAT = TEMPERATURE.....C
C HF = LIQUID ENTHALPY.....KJ/KG
C HG = VAPOUR ENTHALPY.....KJ/KG
C VF = LIQUID SPECIFIC VOLUME.....M**3/KG
C VG = VAPOUR SPECIFIC VOLUME.....M**3/KG
C VISCF = LIQUID DYNAMIC VISCOSITY.....MICRO PA*S
C CONF = LIQUID THERMAL CONDUCTIVITY.....W/(M*K)
C CPF = LIQUID SPECIFIC HEAT.....KJ/(KG*K)
C -----

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C

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COMMON / CONT / ND,NC,NR,NTIN,NPOW,NW,NF,NDROP,NOSUB,NPRINT,NCH
COMMON /OPTION/ NGDYN,NPDYN,NHDYN,NQDYN,NODES
COMMON /DATA1 / TIN,DTIN,POWR,DPOWR,W,DW,
    PIN,PDROP,AP,BP,CP,ASB,BSB,AHT,BHT,CHT,
    TSAT,HF,HG,RHOF,RHOG,VISCF,CONF,CPF
COMMON /DATA2 / DZ(40),A(40),DE(40),PH(40),ANG(40),RK(40),F(40),
    PKW(40),TCAP(40),CO(40),V(40),AA(40),BB(40),CC(40)
COMMON /DATA3 / SIGMA,FREQ,DFREQ,FHZ(120),GDYN(120),HDYN(120),
    QDYN(120),PDYN(120),CFREQ(10),NDYN,NCFREQ
COMMON /DATA4 / GDFAC,PDFAC,HDFAC,QDFAC,GCDYN(120),PCDYN(120),
    HCDYN(120),QCDYN(120)
COMMON /BYPAS / ZB(20),AR(20),DB(20),ANGB(20),RKB(20),EB(20),
    CONL,CPL,DRHO,HFG,PRL,RHOL,VISCL,WTOT,NB
COMMON /TABLE / TEMP(100),PRESS(100),ENTHF(100),ENTHFG(100),
    ENTHG(100),SPVF(100),SPVG(100),DENF(100),
    DENG(100),VISCDF(100),CONDF(100),SPHF(100),
    PRF(100),FAC1,FAC2,ADRG,BDRG,NCARDS
COMMON /ERROR / CSPEC(80),CPHASE(80),ASPEC(80),APHASE(80),NERR
COMMON /CPLT / TPLT1(20),TPLT2(20),NPLT
DIMENSION STC(40),SCO(40),SV(40),SAA(40),SBB(40),SCC(40)
LOGICAL *4 TITLE(20)

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C

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NCASE =0
READ (1,210) (TITLE(I),I=1,20)
1100 READ (1,100,END=1000) ND,NC,NR,NB,NTIN,NPOW,NW,NF,NCFREQ
READ (1,100) NDROP,NODES,NOSUB,NDYN,NPRINT,NPLT,NCARDS
READ (1,100) NGDYN,NPDYN,NHDYN,NQDYN,NSPEC,NPROP
READ (1,110) TIN,DTIN
READ (1,110) POWR,DPOWR
READ (1,110) W,DW
READ (1,110) TCFAC,COFAC,VFAC,AAFAC,BBFAC,CCFAC
READ (1,110) PIN,PDROP,WTOT,AP,BP,CP
READ (1,110) ASB,BSB,AHT,BHT,CHT
BSB=ASB/70000.0
NERR=NDYN
IF (NDYN.EQ.0) NSPEC=0
NDYN=IABS(NDYN)
IF (NCASE.GT.0) GO TO 30
IF (NDYN) 15,15,10
10 READ (1,110) SIGMA,FREQ,DFREQ
IF (NCFREQ.GT.0) READ (1,110) (CFREQ(I),I=1,NCFREQ)
READ (1,110) GDFAC,PDFAC,HDFAC,QDFAC
NT=NF
IF (NDYN.EQ.1) NT=1
IF (NGDYN.EQ.0) GO TO 11
IF (NDYN.EQ.1.AND.NGDYN.NE.1) NGDYN=1
READ (NGDYN,110) (GDYN(I),I=1,NT)
IF (NDYN.EQ.3) READ (NGDYN,110) (GCDYN(I),I=1,NT)
11 IF (NPDYN.EQ.0) GO TO 12
IF (NDYN.EQ.1.AND.NPDYN.NE.1) NPDYN=1
READ (NPDYN,110) (PDYN(I),I=1,NT)
IF (NDYN.EQ.3) READ (NPDYN,110) (PCDYN(I),I=1,NT)
12 IF (NHDYN.EQ.0) GO TO 13
IF (NDYN.EQ.1.AND.NHDYN.NE.1) NHDYN=1
READ (NHDYN,110) (HDYN(I),I=1,NT)
IF (NDYN.EQ.3) READ (NHDYN,110) (HCDYN(I),I=1,NT)
13 IF (NQDYN.EQ.0) GO TO 14
IF (NDYN.EQ.1.AND.NQDYN.NE.1) NQDYN=1

```

```

      ;
      READ (NQDYN,110) (QDYN(I),I=1,NT)
      IF (NDYN.EQ.3) READ (NQDYN,110) (QCDYN(I),I=1,NT)
14  IF (NSPEC.EQ.0) GO TO 15
      READ (NSPEC,110) (CSPEC(I),I=1,NF)
      READ (NSPEC,110) (CPHASE(I),I=1,NF)
      READ (NSPEC,110) (ASPEC(I),I=1,NF)
      READ (NSPEC,110) (APHASE(I),I=1,NF)
15  NCH=ND+NC+NR
      DO 16 I=1,NCH
16  READ (1,110) DZ(I),A(I),DE(I),PH(I),ANG(I),RK(I),E(I)
      DO 20 I=1,NCH
20  READ (1,110) PKW(I),TCAP(I),CO(I),V(I),AA(I),BB(I),CC(I)
      IF (NB) 30,30,22
22  DO 24 I=1,NB
24  READ (1,110) ZB(I),AB(I),DB(I),ANGB(I),RKB(I),EB(I)
30  WRITE (3,140) (TITLE(I),I=1,20)
      WRITE (3,142)
      WRITE (3,150) ND,NC,NR,NB,NTIN,NPOW,NW,NF,NCFREQ
      WRITE (3,144)
      WRITE (3,150) NDROP,NODES,NOSUB,NERR,NPRINT,NPLOT,NCARDS
      WRITE (3,146)

      WRITE (3,150) NGDYN,NPDYN,NHDYN,NQDYN,NSPEC,NPROP
      WRITE (3,152)
      WRITE (3,160) TIN,DTIN
      WRITE (3,154)
      WRITE (3,160) POWR,DPOWR
      WRITE (3,156)
      WRITE (3,160) W,DW
      WRITE (3,157)
      WRITE (3,160) TCFAC,CQFAC,VFAC,AAFAC,BRFAC,CCFAC
      WRITE (3,158)
      WRITE (3,160) PIN,PDROP,WTOT,AP,BP,CP
      WRITE (3,159)
      WRITE (3,160) ASB,BSB,AHT,RHT,CHT
      IF (NCASE.GT.0) GO TO 42
      IF (NDYN) 40,40,31
31  WRITE (3,162)
      WRITE (3,160) SIGMA,FREQ,DFREQ
      IF (NCFREQ) 33,33,32
32  WRITE (3,163)
      WRITE (3,160) (CFREQ(I),I=1,NCFREQ)
33  WRITE (3,190)
      WRITE (3,160) GDFAC,PDFAC,HDFAC,QDFAC
      IF (NGDYN.EQ.0) GO TO 34
      WRITE (3,192)
      IF (NDYN.EQ.3) WRITE (3,200)
      WRITE (3,160) (GDYN(I),I=1,NT)
      IF (NDYN.NE.3) GO TO 34
      WRITE (3,202)
      WRITE (3,160) (GCDYN(I),I=1,NT)
34  IF (NPDYN.EQ.0) GO TO 35
      WRITE (3,194)
      IF (NDYN.EQ.3) WRITE (3,200)
      WRITE (3,160) (PDYN(I),I=1,NT)
      IF (NDYN.NE.3) GO TO 35
      WRITE (3,202)
      WRITE (3,160) (PCDYN(I),I=1,NT)

```

```

35 IF (NHDYN.EQ.0) GO TO 36
   WRITE (3,196)
   IF (NDYN.EQ.3) WRITE (3,200)
   WRITE (3,160) (HDYN(I),I=1,NT)
   IF (NDYN.NE.3) GO TO 36
   WRITE (3,202)
   WRITE (3,160) (HCDYN(I),I=1,NT)
36 IF (NQDYN.EQ.0) GO TO 37

   WRITE (3,198)
   IF (NDYN.EQ.3) WRITE (3,200)
   WRITE (3,160) (QDYN(I),I=1,NT)
   IF (NDYN.NE.3) GO TO 37
   WRITE (3,202)
   WRITE (3,160) (QCDYN(I),I=1,NT)
37 IF (NSPEC.EQ.0) GO TO 40
   WRITE (3,220)

   WRITE (3,160) (CSPEC(I),I=1,NF)
   WRITE (3,160) (CPHASE(I),I=1,NF)
   WRITE (3,222)
   WRITE (3,160) (ASPEC(I),I=1,NF)
   WRITE (3,226)
   WRITE (3,160) (APHASE(I),I=1,NF)
40 WRITE (3,164)
   DO 42 I=1,NCH
     IF (ND.GT.0.AND.I.EQ.1) WRITE (3,170)
     IF (NC.GT.0.AND.I.EQ.(ND+1)) WRITE (3,172)
     IF (NR.GT.0.AND.I.EQ.(ND+NC+1)) WRITE (3,174)
42 WRITE (3,168) DZ(I),A(I),DE(I),PH(I),ANG(I),RK(I),E(I),I
   ND1=ND+1
   DO 43 I=ND1,NCH
     IF (NCASE.EQ.0) STC(I)=TCAP(I)
     TCAP(I)=STC(I)*TCFAC
     IF (NCASE.EQ.0) SCO(I)=CO(I)
     CO(I)=SCO(I)*COFAC
     IF (NCASE.EQ.0) SV(I)=V(I)
     V(I)=SV(I)*VFAC
     IF (NCASE.EQ.0) SAA(I)=AA(I)
     AA(I)=SAA(I)*AAFAC
     IF (NCASE.EQ.0) SBB(I)=BB(I)
     BB(I)=SBB(I)*BBFAC
     IF (NCASE.EQ.0) SCC(I)=CC(I)
43 CC(I)=SCC(I)*CCFAC
   WRITE (3,166)
   DO 44 I=1,NCH
     IF (ND.GT.0.AND.I.EQ.1) WRITE (3,170)
     IF (NC.GT.0.AND.I.EQ.(ND+1)) WRITE (3,172)
     IF (NR.GT.0.AND.I.EQ.(ND+NC+1)) WRITE (3,174)
44 WRITE (3,168) PKW(I),TCAP(I),CO(I),V(I),AA(I),BB(I),CC(I),I
   IF (NB) 50,50,46

46 WRITE (3,176)
   WRITE (3,178)
   NDROP=1
   DO 48 I=1,NB
48 WRITE (3,160) ZB(I),AB(I),DB(I),ANGB(I),RKB(I),EB(I),I
50 FAC1=1.000000E+03
   FAC2=1.000000E-06
   NPRINT=IABS(NPRINT)

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      IF (NDYN.EQ.0) GO TO 53
      IF (NPLOT) 51,53,51
51  READ(1,210) (TPLT1(I),I=1,20)
      READ(1,210) (TPLT2(I),I=1,20)
      WRITE (3,215) (TPLT1(I),I=1,20)
      WRITE (3,215) (TPLT2(I),I=1,20)
53  IF (NCASE.GT.0) NCARDS=IABS(NCARDS)
      IF (NCASE) 54,54,60
54  NCARD=IABS(NCARDS)
      IF (NPROP.EQ.0) NPROP=1
      READ (NPROP,120) (TEMP(N),PRESS(N),ENTHF(N),ENTHG(N),SPVF(N),
      .           SPVG(N),VISCDF(N),CONDF(N),SPHF(N),N=1,NCARD)
      IF (NCARDS.LT.0) GO TO 56
      WRITE (3,180)
      WRITE (3,182)
      WRITE(3,184) (TEMP(N),PRESS(N),ENTHF(N),ENTHG(N),SPVF(N),
      .           SPVG(N),VISCDF(N),CONDF(N),SPHF(N),N=1,NCARDS)

56  CALL TABLES
      GO TO 60
58  READ (1,130) TSAT,HF,HG,VF,VG,VISCF,CONF,CPF
      RHOF=1.0/VF
      RHOG=1.0/VG
      WRITE (3,185)
      WRITE (3,186)
      WRITE (3,188) TSAT,HF,HG,VF,VG,VISCF,CONF,CPF
60  CALL EXEC (NCASE)
      NCASE=1
      GO TO 1100
100  FORMAT (16I5)
110  FORMAT (8F10.3)
120  FORMAT (F5.1,F9.2,F8.2,F8.1,F10.6,F10.5,F6.1,F8.4,F8.3)
130  FORMAT (F5.1,F8.2,F8.1,F10.6,F10.5,F6.1,F8.4,F8.3)
140  FORMAT ('1',' LOCO INPUT DATA : ',20A4,/2X,15('*'))
142  FORMAT ('0',6X,'ND',6X,'NC',6X,'NR',6X,'NB',5X,'NTIN',4X,
      .       'NPOW',5X,'NW',6X,'NF',4X,'NCFREQ')
144  FORMAT ('0',5X,'NDROP',2X,'NODES',4X,'NOSUB',3X,'NDYN',3X,
      .       'NPRINT',3X,'NPLOT',2X,'NCARDS')
146  FORMAT ('0',5X,'NGDYN',3X,'NPDYN',3X,'NHDYN',3X,'NQDYN',3X,
      .       'NSPEC',3X,'NPROP')
150  FORMAT (1X,16I8)
152  FORMAT ('0',1X,'TIN',7X,'DTIN')
154  FORMAT ('0',1X,'POWR',6X,'DPOWR')
156  FORMAT ('0',1X,'W',9X,'DW')
157  FORMAT ('0 TCFAC',5X,'COFAC',5X,'VFAC',6X,'AAFAC',5X,
      .       'BBFAC',5X,'CCFAC')
158  FORMAT ('0',1X,'PIN',7X,'PDROP',5X,'WTOT',6X,'AP',8X,'BP',8X,'CP')
159  FORMAT ('0 ASB',7X,'BSR',7X,'AHT',7X,'BHT',7X,'CHT')
160  FORMAT (1X,1P8E10.3)
162  FORMAT ('0',1X,'SIGMA',5X,'FREQ',6X,'DFREQ')
163  FORMAT ('0',1X,'CFREQ')
164  FORMAT ('0',1X,'DZ(N)',5X,'A(N)',6X,'DE(N)',5X,'PH(N)',5X,
      .       'ANG(N)',4X,'RK(N)',5X,'E(N)')
166  FORMAT ('0',1X,'PKW(N)',4X,'TCAP(N)',3X,'CO(N)',5X,'V(N)',6X,
      .       'AA(N)',5X,'BB(N)',5X,'CC(N)')
168  FORMAT (1X,1P7E10.3,16)
170  FORMAT (' INLET PIPING REGION')
172  FORMAT (' CHANNEL REGION')
174  FORMAT (' RISER REGION')

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176 FORMAT ('Ø BYPASS PIPING')
178 FORMAT (' ZB(N)',5X,'AB(N)',5X,'DB(N)',5X,'ANGB(N)',3X,'RKB(N)',
.
4X,'EB(N)')
180 FORMAT ('1',28X,'FLUID PROPERTIES')
182 FORMAT ('Ø',1X,'TEMP',4X,'PRESS',3X,'ENTHF',3X,'ENTHG',4X,'SPVF',
.
7X,'SPVG',2X,'VISCF',3X,'CONDF',4X,'SPHF')
184 FORMAT (1X,F5.1,F9.2,F8.2,F8.1,F10.6,F10.5,F6.1,F8.4,F8.3)
185 FORMAT ('Ø',24X,'FLUID PROPERTIES')
186 FORMAT ('Ø',1X,'TEMP',3X,'ENTHF',3X,'ENTHG',4X,'SPVF',7X,'SPVG',
.
2X,'VISCF',3X,'CONDF',4X,'SPHF')
188 FORMAT (1X,F5.1,F8.2,F8.1,F10.6,F10.5,F6.1,F8.4,F8.3)
190 FORMAT ('Ø GDFAC',5X,'PDFAC',5X,'HDFAC',5X,'QDFAC')
192 FORMAT ('Ø GDYN(N)')
194 FORMAT ('Ø PDYN(N)')
196 FORMAT ('Ø HDYN(N)')
198 FORMAT ('Ø QDYN(N)')
200 FORMAT (' REAL COMPONENT')
202 FORMAT (' IMAGINARY COMPONENT')

210 FORMAT (2ØA4)
215 FORMAT (1X,2ØA4)
220 FORMAT ('Ø CSPEC(N)')
222 FORMAT ('Ø ASPEC(N)')
224 FORMAT ('Ø CPHASE(N)')
226 FORMAT ('Ø APHASE(N)')
1000 STOP
END
SUBROUTINE EXEC (NCASE)
COMMON / CONT / ND,NC,NR,NTIN,NPOW,NW,NF,NDROP,NOSUB,NPRINT,NCH
COMMON /DATA1 / TIN,DTIN,POWR,DPOWR,W,DW,
.
PIN,PDROP,AP,ØP,CP,ASB,BSB,AHT,BHT,CHT,
.
TSAT,HF,HG,RHOF,RHOG,VISCF,CONF,CPF
COMMON /DATA2 / DZ(4Ø),A(4Ø),DE(4Ø),PH(4Ø),ANG(4Ø),RK(4Ø),E(4Ø),
.
PKW(4Ø),TCAP(4Ø),CO(4Ø),V(4Ø),AA(4Ø),BB(4Ø),CC(4Ø)
COMMON /DATA3 / SIGMA,FRFQ,DFREQ,FHZ(12Ø),GDYN(12Ø),HDYN(12Ø),
.
QDYN(12Ø),PDYN(12Ø),CFREQ(1Ø),NDYN,NCFREQ
COMMON /EXSTAT/ DPT,FAC,FAC3,GC,PI,PUMP,PY,NIT,NODE
COMMON /BYPAS / ZB(2Ø),AB(2Ø),DB(2Ø),ANGB(2Ø),RKB(2Ø),EB(2Ø),
.
CONL,CPL,DRHO,HFG,PRL,RHOL,VISCL,WTOT,NB
COMMON /TABLE / TEMP(1ØØ),PRESS(1ØØ),ENTHF(1ØØ),ENTHFG(1ØØ),
.
ENTHG(1ØØ),SPVF(1ØØ),SPVG(1ØØ),DENF(1ØØ),
.
DENG(1ØØ),VISCDF(1ØØ),CONDF(1ØØ),SPHF(1ØØ),
.
PRF(1ØØ),FAC1,FAC2,ADRG,ØDRG,NCARDS
FAC=1.ØØØØE-Ø2
FAC3=FAC1
NDYNS=NDYN
IF (NCASE.GT.Ø) GO TO 15
IF (NCARDS) 1Ø,1Ø,15
1Ø RHOL=RHOF
DRHO=RHOF-RHOG
HF=HF*FAC1
HG=HG*FAC1
HFG=HG-HF
CPF=CPF*FAC1
CPL=CPF
VISCF=VISCF*FAC2
VISCL=VISCF
CONL=CONF
PRL=CPF*VISCF/CONF
HSUB=CPF*(TSAT-TIN)/FAC1

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15 GC=9.807
   WI=W
   FCWRI=POWER
   DO 90 K=1,NTIN
   POWER=PCWRI
   DO 80 J=1,NPOW
   IT1=0
   IT2=0
   NIT=0
   W=WI
   IF (NDROP.EQ.1) WRITE (3,120)
   DO 70 I=1,NW
20 IF (NCARDS) 25,25,22
22 RHOL=TDF(TIN)
   VISCL=TVISCD(TIN)
   TSAT=SAT(TEMP,PRESS,PIN,NCARDS)
   HSUB=(THF(TSAT)-THF(TIN))/FAC1
25 PY=3.141592/180.0
30 IF (NB.GT.0) GO TO 35
   QT=W/RHOL*FAC1
   DPUMP=-RHOL*(AP+BP*QT+CP*QT*QT)*GC/FAC1
   PUMP=- (BP+2.0*CP*QT)*A(1)/FAC1
   GO TO 40
35 DPUMP=0.0
   PUMP=0.0
   PDROP=0.0
   CALL BYPASS(0)
40 IF (NDROP.EQ.0) WRITE (3,100) TIN,HSUB,POWER,W
   IF (NIT.EQ.1) WRITE (3,110) TIN,HSUB,POWER,W
   NDYN=NDYNS
   CALL STATE
   IF (NIT) 45,45,75
45 PTOT=DPT+DPUMP
   PDIFF=PTOT-PDROP
   IF (NDROP) 65,65,50
C
C   CALCULATION OF FLOW RATE FOR SPECIFIED PRESSURE DROP.
50 WRITE (3,130) W,PTOT,PDROP,PDIFF
   IF (PDIFF) 52,62,54
52 W1=W
   PDIFF1=PDIFF
   IT1=1
   GO TO 56
54 W2=W
   PDIFF2=PDIFF
   IT2=1
56 IF (IT1) 70,70,58
58 IF (IT2) 70,70,60
60 W=(PDIFF2*W1-PDIFF1*W2)/(PDIFF2-PDIFF1)
   CHECK=ABS(PDIFF)
   IF (PDROP.NE.0.0) CHECK=ABS(PDIFF/PDROP)
   IF (CHECK.LE.0.0050) GO TO 62
   GO TO 20
62 NIT=1
   GO TO 20
65 IF (NDYN.GT.0) CALL TFUNCN
70 W=W+DW
   GO TO 80

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6 HL1=H1
  HIN=H1
  ALFA(1)=0.0
  CK0(1)=1.0
  X(1)=0.0
  PHI(1)=1.0
  UL1=W/AC/RHOL
  UG1=0.0
  GM1=W*W/AC/AC
  TAU1=0.0
  TAU2=0.0

  Z=0.0
  ZBR=0.0
  IF (NDYN.EQ.0) DFREQ=0
  IF (NC.GT.1) GO TO 8
  NOSUR=0
  NPRINT=0
8  NBR=0
  NCD=0
  NSB=0
  NSUB=0
  NSAT=0
  NCDN=0
  I=0
  IF (NPRINT) 10,10,15
10 IF (NDROP.EQ.0.OR.NIT.EQ.1) WRITE (3,200)
  GO TO 20
15 IF (NDROP.EQ.0.OR.NIT.EQ.1) WRITE (3,205)
20 DO 190 K=1,NCH
  IF (NDROP.GT.0.AND.NIT.EQ.0) GO TO 22
  IF (ND.GT.0.AND.K.EQ.1) WRITE (3,210)
  IF (NC.GT.0.AND.K.EQ.(ND+1)) WRITE (3,220)
  IF (NR.GT.0.AND.K.EQ.(ND+NC+1)) WRITE (3,230)

  IF (NC.EQ.1) GO TO 22
  IF (NPRINT.EQ.0) WRITE (3,235) K
22 AR=A(1)/A(K)
  A1=AC
  PKWK=PKW(K)
  IF (K.GT.ND.AND.K.LE.(ND+NC)) PKWK=POWR*PKW(K)
  ZK=DZ(K)*FAC
  NWARN=0
  IF (NDROP.EQ.1.AND.NIT.EQ.0) NWARN=1
  AC=A(K)*FAC*FAC
  DC=DE(K)*FAC
  PC=PH(K)*FAC
  THETA=COS((ANG(K)*PY))
  DHE=4.0*AC/PC
  FLK=PKWK*FAC1/DZ(K)/PH(K)
  FLX=FLK/FAC/FAC
  G=W/AC
  G2=G*G
  DRGDK(K)=0.0
  IF (NC.GT.1) GO TO 23
  NODE=1
  NB0IL=1
  IF (K.GT.ND.AND.K.LE.(ND+NC)) NODE=2
  CHECK=PKWK*FAC1/W/(HBB-HIN)
  IF (CHECK.LE.1) NODE=1

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GO TO 26
23 IF (NSAT.GT.0.AND.NSAT.LT.1) GO TO 25
HBR=HF
IF (NOSUB.EQ.0) GO TO 24
PEC=G*DC*CPL/CONL
HTT=ASR*CONL/DC
IF (PEC.GT.7.00E+04) HTT=BSB*G*CPL
HSB=HF-CPL*FLX/HTT
HBR=HSB
24 NODE=ABS(10.0*PKWK*FAC1/W/(HBB-H1))+2
25 NCHECK=(NODES-I)/(NCH-K+1)+1
IF (NSAT.GT.0) NODE=2.0*DFREQ*ZK/UG1+1
IF (NODE.GT.NCHECK) NODE=NCHECK
26 MODE(K)=NODE
DO 180 J=1,NODE
IF (NC.EQ.1) GO TO 30
IF (NSAT.GT.0.AND.NSAT.LT.1) GO TO 28
HBR=HF
IF (NOSUB.GT.0) HBB=HSB
NBOIL=ABS(10.0*PKWK*FAC1/W/(HBB-H1))/NODE+2
LBOIL=(NODES-I)/NODE/(NCH-K+1)+1
IF (NBOIL.GT.LBOIL) NBOIL=LBOIL
IF ((NODE*NBOIL).GT.100) NBOIL=100/NODE
IF (K.LE.ND) NBOIL=5
JBOIL=2.0*DFREQ*ZK/NODE/UL1+1
GO TO 30
28 NBOIL=(NODES-I)/NODE/(NCH-K+1)+1
JBOIL=2.0*DFREQ*ZK/NODE/UG1+1
30 ZC=ZK/NODE/NBOIL
IF (NWARN.EQ.1) GO TO 31
IF (JBOIL.LT.NBOIL) GO TO 31
NWARN=1
NDIV=JBOIL/NBOIL+1
WRITE (3,295) K,NDIV
31 AS=PC*ZC
DH=PKWK*FAC1/W/NODE/NBOIL
IF (NC.GT.1) GO TO 38
IF (K.LE.ND.OR.K.GT.(ND+NC)) GO TO 38
DH=PKWK*FAC1/W
DHB=HBB-HIN
IF (DH-DHB) 32,32,34
32 ZC=ZK
GO TO 36
34 DH=DHB
IF (J.EQ.2) DH=PKWK*FAC1/W-DHB
ZC=DH*W*ZK/PKWK/FAC1
36 ZCL(J)=ZC
AS=PC*ZC
38 MBOIL(K,J)=NBOIL
DO 180 L=1,NBOIL
I=I+1
H2=H1+DH
HA=(H1+DH/2.0)/FAC1
IF (NCARDS) 45,45,40
40 TSAT=SAT(TEMP,PRESS,P1,NCARDS)
HF=THF(TSAT)
HG=THG(TSAT)
HFG=THFG(TSAT)

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CPI=TSH(TL1)
RHOL=TDF(TL1)
RHOG=TDG(TSAT)
DRHO=RHOL-RHOG
VISCL=TVISCD(TL1)
CONL=TCOND(TL1)
PRI=TPRN(TL1)
45 RE=G*DC/VISCL
   DRGDP(I)=0.0
   IF (NC.GT.1) GO TO 55
   IF (I.LE.(ND+1)) GO TO 70
   IF (J.EQ.2) GO TO 125
   IF (NBB.EQ.1) GO TO 130
   GO TO 70
50 NCD=1
   IF (NCON.EQ.0) NCON=I
   X(I+1)=0.0
   WRITE (3,270) ZBB
   NDYN=0
55 IF (NCD) 60,60,70
60 IF (NBB) 65,65,130
65 IF (NSB) 70,70,82
C
C SINGLE PHASE REGION
70 DHI=DH
   HL2=H2
   TL2=TL1+DH/CPL
   TLA=(TL1+TL2)/2.0
   HT(I)=0.023*CONL/DHE*RE**0.80*PRL**0.40
   HTB(I)=0.80
   PEC=RE*PRL
   HTT=ASB*CONL/DC
   IF (PEC.GT.7.00E+04) HTT=BSB*G*CPL
   HSB=HF-CPL*FLX/HTT
   ALFA(I+1)=0.0
   ALFM=0.0
   X(I+1)=0.0
   CKO(I+1)=1.0
   XM=0.0
   DX=0.0
   UL2=G/RHOL
   UG2=0.0
   GL2=G
   GG2=0.0
   GM2=G2/RHOL
   PHI(I+1)=1.0
   PHIA=1.0
   RHOA=RHOL
   IF (J.GT.1.OR.L.GT.1) GO TO 72
   ALFK(K)=0.0
   CKK(K)=1.0
   PHIK(K)=1.0
   E1=UL1*UL1/2.0
   UL1=G/RHOL
   GM1=G2/RHOL
72 IF (NC.EQ.1) GO TO 78
   IF (NOSUB) 74,74,76
74 IF (H2-HF) 78,78,125

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76 IF (H2-HSB) 78,79,80
78 TAU1=TAU1+ZC/UL1
   GO TO 150
C
C   SUBCOOLED BOILING REGION
80 NSB=1
   FHP(I)=1.0
   NSAT=I
   NSUB=I
82 IF (HF-HSB) 128,128,84
84 IF (H2-HSB) 50,50,88
88 IF (NOSUB-1) 90,90,94
90 FHP(I+1)=1.0/EXP((H2-HSB)/(HF-HSB))
   HL2=HF-(HF-HSB)*FHP(I+1)
   GO TO 96
94 FHP(I+1)=TANH((H2-HSB)/(HF-HSB))
   HL2=HSB+(HF-HSB)*FHP(I+1)
   FHP(I+1)=1.0-FHP(I+1)*FHP(I+1)
96 IF (HL2-HF) 100,128,128
100 X(I+1)=(H2-HL2)/(HG-HL2)
   IF (X(I+1)) 102,102,110
102 IF (I-NSAT) 104,104,106
104 NSB=0
   NSAT=0
   GO TO 150
106 X(I+1)=X(I)
110 DX=X(I+1)-X(I)
   DHL=HL2-HL1
   HLA=(HL1+HL2)/2.0
   TL2=TL1+DHL/CPL
   TLA=(TL1+TL2)/2.0
   IF (I-NSAT) 120,120,145
120 IF (NDROP.EQ.0.OR.NIT.EQ.1) WRITE (3,250) ZBB,PEC
   GO TO 145
125 FHP(I)=0.0
128 NBB=1
   NSAT=I
   IF (NDROP.EQ.0.OR.NIT.EQ.1) WRITE (3,260) ZBB
C
C   BULK BOILING REGION
130 HL2=HF
   TL2=TSAT-0.010
   TLA=TL2
   FHP(I+1)=0.0
   X(I+1)=(H2-HF)/HFG
   IF (X(I+1).GE.1.0) X(I+1)=0.999999
   IF (X(I+1)) 50,50,140
140 DX=X(I+1)-X(I)
145 GL2=(1.0-X(I+1))*G
   GG2=X(I+1)*G
   UJ=GL2/RHOL+GG2/RHOG
   CK0(I+1)=1.0/(CO(K)+V(K)/UJ)
   DUM=RHOG+DRHO*X(I+1)
   ALFA(I+1)=CK0(I+1)*RHOL*X(I+1)/DUM
   IF (ALFA(I+1).GE.1.0) ALFA(I+1)=0.999999
   ALFM=(ALFA(I)+ALFA(I+1))/2.0
   UL2=GL2/(1.0-ALFA(I+1))/RHOL
   UG2=GG2/ALFA(I+1)/RHOG

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GM2=GL2*UL2+GG2*UG2
PHI(I+1)=1.0+AA(K)*((RHOL/RHOG-1.0)*X(I+1))*BB(K)*G**CC(K)
PHIA=(PHI(I)+PHI(I+1))/2.0
IF (J.GT.1.OR.L.GT.1) GO TO 148
E1=((1.0-X(I))*UL1*UL1+X(I)*UG1*UG1)/2.0
GL1=(1.0-X(I))*G
GG1=X(I)*G
UJ=GL1/RHOL+GG1/RHOG
CKK(K)=1.0/(CO(K)+V(K)/UJ)
DUM=RHOG+DRHO*X(I)
ALFK(K)=CKK(K)*RHOL*X(I)/DUM
IF (ALFK(K).GE.1.0) ALFK(K)=0.999999
UL1=GL1/(1.0-ALFK(K))/RHOL
IF (I.EQ.NSUB.OR.I.EQ.NSAT) GO TO 146
UG1=GG1/ALFK(K)/RHOG
PHIK(K)=1.0+AA(K)*((RHOL/RHOG-1.0)*X(I))*BB(K)*G**CC(K)
GM1=GL1*UL1+GG1*UG1
GO TO 147
146 UG1=0.0
PHIK(K)=1.0
GM1=GL1*UL1
147 ALFM=(ALFK(K)+ALFA(I+1))/2.0
PHIA=(PHIK(K)+PHI(I+1))/2.0
148 XM=(X(I)+X(I+1))/2.0
RHOA=RHOL-DRHO*ALFM
TAU2=TAU2+2.0*ZC/(UG1+UG2)
HTL=0.023*((1.0-XM)*RE)**0.80*PRL**0.40*CONL/DHE
HTG=AHT*(XM*RE)**RHT*PRL**CHT*CONL/DHE
HT(I)=HTL+HTG
HTR(I)=(0.80*HTL+BHT*HTG)/HT(I)
C
C SINGLE PHASE AND TWO PHASE PRESSURE DROPS
150 DPG=RHOA*ZC*GC*THETA/FAC1
DPM=(GM2-GM1)/FAC1
FK=FRIC(RE,E(K))
BF(I)=16.22542/(8.11262+RE*E(K)+4.67/SQRT(FK))
DP1F(I)=4.0*FK*ZC/DC*G2/RHOL/2.0
DPF=DP1F(I)*PHIA/FAC1
DPK=0.0
IF (J.GT.1.OR.L.GT.1) GO TO 151
E2=((1.0-X(I))*UL1*UL1+X(I)*UG1*UG1)/2.0
IF (A1.EQ.AC) E2=E1
PHIR(K)=1.0+1.285*X(I)*(RHOL/RHOG-1.0)
DPR(K)=RK(K)*G2*PHIR(K)/RHOL/2.0
DPI(K)=RHOL*RHOG*(E2-E1)/(RHOG+DRHO*X(I))
DPK=(DPI(K)+DPR(K))/FAC1
DPRT=DPRT+DPK
151 DP=DPM+DPF+DPG+DPK

DPGT=DPGT+DPG
DPMT=DPMT+DPM
DPFT=DPFT+DPF
DPT=DPT+DP
P2=P1-DP
PA=(P1+P2)/2.0
IF (FLX) 154,152,154
152 HTW(I)=0.0
GO TO 156
154 HTW(I)=HT(I)/FLX/CPL

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270 FORMAT ('0 CONDENSING BOUNDARY POSITION =',F8.3,' CM FROM ',
.         'CHANNEL REGION INLET : NO CONDENSATION IS ALLOWED IN ',
.         'DYNAMIC CALCULATIONS'/)
280 FORMAT ('0 SINGLE PHASE REGION TRANSIT TIME =',1PE12.4,' SEC'/
.         '1X,' TWO PHASE REGION TRANSIT TIME =',E12.4,' SEC')
290 FORMAT ('0 TOTAL NUMBER OF CHANNEL NODES =',I3,
.         ' : SUBCOOLED NODES =',I3,' : SATURATED NODES =',I3)
295 FORMAT ('0 *** WARNING - SECTION',I3,' NUMERICALLY UNSTABLE : ',
.         'SUBDIVIDE LENGTH BY',I3/)
300 RETURN
END
SURROUTINE BYPASS (NBDYN)
COMMON /DATA1 / TIN,DTIN,POWR,DPOWR,W,OW,
.         PIN,PDROP,AP,BP,CP,ASB,BSB,AHT,BHT,CHT,
.         TSAT,HF,HG,RHOF,RHOG,VISCF,CONF,CPF
COMMON /DATA2 / D7(40),A(40),DE(40),PH(40),ANG(40),RK(40),E(40),
.         PKW(40),TCAP(40),C0(40),V(40),AA(40),BB(40),CC(40)
COMMON /EXSTAT/ DPT,FAC,FAC3,GC,PI,PUMP,PY,NIT,NODE
COMMON /BYPAS / ZB(20),AB(20),DB(20),ANGB(20),RKB(20),EB(20),
.
.         CONL,CPL,DRHO,HFG,PRL,RHOL,VISCL,WTOT,NB
COMMON /FUNCNC/ ALFA2,DELP1M,DELP1F,DELP1R,DELP1T,DELP2M,DELP2F,
.         DELP2G,DELP2R,DELP2T,DELPT,H,S,GD,HD,QSJ,PDIN
COMMON /BYPASD/ DELPTB
DIMENSION   RFR(20),DPFB(20),DPRB(20),GB(20)
COMPLEX*16   ALFA2,DELP1M,DELP1F,DELP1R,DELP1T,DELP2M,DELP2F,
.         DELP2G,DELP2R,DELP2T,DELPT,H,S,GD,HD,QSJ,PDIN
COMPLEX*16   DELPTB,GBD
C
IF (NBDYN) 5,5,15
5  WB=WTOT-W
   RA1=AB(1)*FAC*FAC
   DO 10 M=1,NB
     BZ=ZB(M)*FAC
     BA=AB(M)*FAC*FAC
     BDH=DB(M)*FAC
     THETB=COS((ANGB(M)*PY))
     BRK=RKB(M)
     BE=EB(M)
     GB(M)=WB/BA
     DPGB=RHOL*BZ*GC*THETB/FAC3
     REL=GB(M)*BDH/VISCL
     FKB=4.0*FRIC(REL,BE)*BZ/BDH
     BFR(M)=16.22542/(8.11262+REL*BE+4.67/SQRT(FKB))
     DPFB(M)=FKB*GB(M)*GB(M)/RHOL/2.0/FAC3
     RB=BA/RA1
     DPRB(M)=GB(M)*GB(M)*(1.0-RB*RB+BRK)/RHOL/2.0/FAC3
     BA1=BA
10  PDROP=PDROP+DPGB+DPFB(M)+DPRB(M)
   RETURN
15  A1=AB(1)
   GBD=CMPLX(1.0,0.0)*A(1)/A1
   DELPTB=0.0
   DO 40 I=1,NB
     GBD=GBD*A1/AB(I)
     DELPTB=DELPTB+(DPFB(I)*(2.0-BFR(I))+DPRB(I)+S*GB(I)*ZB(I)*
       FAC)*GBD/GB(I)/FAC3

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```

40 A1=AB(I)
RETURN
END

SUBROUTINE TFUNCN
COMMON / CONT / ND,NC,NR,NTIN,NPOW,NW,NF,NDROP,NOSUB,NPRINT,NCH
COMMON /DATA1 / TIN,DTIN,POWR,DPOWR,W,DW,
.
.
.
COMMON /DATA3 / SIGMA,FREQ,DFREQ,FHZ(120),GDYN(120),HDYN(120),
.
.
COMMON /DATA4 / GDFAC,PDFAC,HDFAC,QDFAC,GCDYN(120),PCDYN(120),
.
.
COMMON /EXSTAT/ DPT,FAC,FAC3,GC,P1,PUMP,PY,NIT,NODE
COMMON /BYPAS / ZB(20),AB(20),DB(20),ANGB(20),RKB(20),EB(20),
.
.
COMMON /BYPASD/ DELPTB
COMMON /FUNCNC/ ALFA2,DELPM,DELP1F,DELP1R,DELP1T,DELP2M,DELP2F,
.
.
COMMON /FUNIN / DGD,DHD,DPD,DOD,GOLD,HOLD,POLD,QOLD
COMMON /PRINT / AA(60),AT(60),GA(60),GADB(60),GT(60),GTDB(60),
.
.
.
DIMENSION CDP1M(5),CDP1F(5),CDP1R(5),CDP1T(5),CDP2M(5),
.
.
.
COMPLEX*16 ALFA2,DELPM,DELP1F,DELP1R,DELP1T,DELP2M,DELP2F,
.
.
.
COMPLEX*16 DGD,DHD,DPD,DOD,GOLD,HOLD,POLD,QOLD
COMPLEX*16 DELPTB,PUMPD
COMPLEX*16 CDP1M,CDP1F,CDP1R,CDP1T,CDP2M,CDP2F,CDP2G,CDP2R,
.
.
.
CDP2T,CDPT
PY=6.28319
DOMEG=0.1000*PY*DFREQ
NA=0
NM=0
NT=0
PUMPD=PUMP
IF (NB.EQ.0) DELPTB=0.0
SREAL=SIGMA*PY
DO 100 J=1,NF
FHZ(J)=FREQ+(J-1)*DFREQ
OMEGA=PY*FHZ(J)
S=CMPLX(SREAL,OMEGA)
CALL DYNIN (J)
CALL TCOMP
DELPT=DELPT/FAC3
IF (NB.GT.0) CALL BYPASS (1)
DELPT=DELPT+(DELPTB+PUMPD)*GD
GT(J)=CDABS(DELPT)
GTDB(J)=20.0*ALOG10(GT(J))
RT(J)=DREAL(DELPT)
QT(J)=DAIMAG(DELPT)
IF (J-1) 30,30,35
30 IF (RT(J).LT.0.0) NT=1
GO TO 60
35 IF (QT(J-1).LT.0.0.AND.(QT(J-1)*QT(J)).LT.0.0) GO TO 40
GO TO 60
40 OMEGA1=PY*FHZ(J-1)
QT1=QT(J-1)
OMEGA2=OMEGA

```

```

QT2=QT(J)
IT=1
42 OMEGM=(QT2*OMEGA1-QT1*OMEGA2)/(QT2-QT1)
S=CMPLX(SREAL,OMEGM)
IF (NDYN.EQ.1) GO TO 44
DOM=(OMEGM/PY-FHZ(J-1))/DFREQ
GD=GOLD+DGD*DOM
PDIN=POLD+OPD*DOM
HD=HOLD+DHD*DOM
QSJ=QOLD+DQD*DOM
44 CALL TCOMP
DELPT=DELPT/FAC3
IF (NB.GT.0) CALL BYPASS(1)
DELPT=DELPT+(DELPTB+PUMPD)*GD
ORD=DAIMAG(DELPT)
IF (ABS(ORD)-0.0010) 55,46,46
46 IT=IT+1
IF (IT.GT.2) GO TO 55
IF (ORD) 48,55,50
48 OMEGA1=OMEGM
QT1=ORD
GO TO 42
50 OMEGA2=OMEGM
QT2=ORD
GO TO 42
55 NM=NM+1
DPMIN=DREAL(DELPT)
IF (DPMIN.LT.0.0) NT=NT+1
OM(NM)=OMEGM
CDP1M(NM)=DELP1M
CDP1F(NM)=DELP1F
CDP1R(NM)=DELP1R
CDP1T(NM)=DELP1T
CDP2M(NM)=DELP2M
CDP2F(NM)=DELP2F
CDP2G(NM)=DELP2G
CDP2R(NM)=DELP2R
CDP2T(NM)=DELP2T
CDPT(NM)=DELPT
IF (IT.GT.20) WRITE (3,210) NM
60 AT(J)=ATAN2(QT(J),RT(J))*57.2958-NT*360.0
GA(J)=CDABS(ALFA2)
GADB(J)=20.0*ALOG10(GA(J))
RA(J)=DREAL(ALFA2)
QA(J)=DAIMAG(ALFA2)
IF (J-1) 90,90,65
65 IF (RA(J).LT.0.0.AND.(QA(J-1)*QA(J)).LT.0.0) GO TO 70
GO TO 90
70 OMEGA1=PY*FHZ(J-1)
QA1=QA(J-1)
OMEGA2=OMEGA
QA2=QA(J)
IT=1
72 OMEGM=(QA2*OMEGA1-QA1*OMEGA2)/(QA2-QA1)
S=CMPLX(SREAL,OMEGM)
IF (NDYN.EQ.1) GO TO 74
DOM=(OMEGM/PY-FHZ(J-1))/DFREQ
GD=GOLD+DGD*DOM

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PDIN=POLD+DPD*DOM
HD=HOLD+DHD*DOM
QSJ=QOLD+DQD*DOM
74 CALL TCOMP
ORD=DAIMAG(ALFA2)
IF (ABS(ORD)-1.0E-8) 82,82,76
76 IT=IT+1
IF (IT.GT.2) GO TO 82
IF (ORD) 78,82,80
78 OMEGA1=OMEGM
QA1=ORD
GO TO 72
80 OMEGA2=OMEGM
QA2=ORD
GO TO 72
82 AMIN=DREAL(ALFA2)
IF (AMIN.LT.0.0) NA=NA+1
90 IF (J.EQ.1.AND.RA(J).LT.0.0) NA=1
AA(J)=ATAN2(QA(J),RA(J))*57.2958-NA*360.0
100 CONTINUE
CALL TPRINT
IF (NF.EQ.1) CALL CPRINT
IF (NM) 140,140,110
110 DO 130 J=1,NM
WRITE (3,200)
S=CMPLX(SREAL,OM(J))
DELP1M=CDP1M(J)
DELP1F=CDP1F(J)
DELP1R=CDP1R(J)
DELP1T=CDP1T(J)
DELP2M=CDP2M(J)
DELP2F=CDP2F(J)
DELP2G=CDP2G(J)
DELP2R=CDP2R(J)
DELP2T=CDP2T(J)
DELP2=CDPT(J)
CALL CPRINT
130 CONTINUE
140 IF (NCFREQ) 170,170,150
150 DO 160 J=1,NCFREQ
WRITE (3,220)
OMC=CFREQ(J)*PY
S=CMPLX(SREAL,OMC)
JC=CFREQ(J)/DFREQ+1
CALL DYNIN (JC)
CALL TCOMP
DELPT=DELPT/FAC3
IF (NB.GT.0) CALL BYPASS (1)
DELPT=DELPT+(DELPTB+PUMPD)*GO
CALL CPRINT
160 CONTINUE
170 CONTINUE
200 FORMAT (1H1,////41X,'MINIMUM PURELY REAL PRESSURE DROP',
.      ' COMPONENTS'/41X,44('-'))
210 FORMAT ('0TOO MANY ITERATIONS FOR CONVERGENCE : NM =',I2)
220 FORMAT (1H1,////38X,'PRESSURE DROP COMPONENTS FOR SPECIFIED',
.      ' FREQUENCY'/38X,48('-'))

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300 RETURN
END
SURROUTINE DYNIN (J)
COMMON /OPTION/  NGDYN,NPDYN,NHDYN,NQDYN,NODES
COMMON /DATA3 /  SIGMA,FREQ,DFREQ,FHZ(120),GDYN(120),HDYN(120),
. QDYN(120),PDYN(120),CFREQ(10),NDYN,NCFREQ
COMMON /DATA4 /  GDFAC,PDFAC,HDFAC,QDFAC,GCDYN(120),PCDYN(120),
. HCDYN(120),QCDYN(120)
COMMON /EXSTAT/  DPT,FAC,FAC3,GC,PI,PUMP,PY,NIT,NODE
COMMON /FUNCNC/  ALFA2,DEL P1M,DELP1F,DELP1R,DELP1T,DELP2M,DELP2F,
. DELP2G,DELP2R,DELP2T,DELPT,H,S,GD,HD,QSJ,PDIN
COMMON /FUNIN /  DGD,DHD,DPD,DQD,GOLD,HOLD,POLD,QOLD
COMPLEX*16      ALFA2,DELP1M,DELP1F,DELP1R,DELP1T,DELP2M,DELP2F,
. DELP2G,DELP2R,DELP2T,DELPT,H,S,GD,HD,QSJ,PDIN
COMPLEX*16      DGD,DHD,DPD,DQD,GOLD,HOLD,POLD,QOLD
GD=0.0
PDIN=0.0
HD=0.0
QSJ=0.0
IF (NDYN-2) 10,20,30
10 IF (NGDYN.EQ.1) GD=CMPLX(GDYN(1),0.0)*GDFAC
IF (NPDYN.EQ.1) PDIN=CMPLX(PDYN(1),0.0)*PDFAC*FAC3
IF (NHDYN.EQ.1) HD=CMPLX(HDYN(1),0.0)*HDFAC*FAC3
IF (NQDYN.EQ.1) QSJ=CMPLX(QDYN(1),0.0)*QDFAC
RETURN
20 GCDYN(J)=0.0
PCDYN(J)=0.0
HCDYN(J)=0.0
QCDYN(J)=0.0
30 GOLD=GD
POLD=PDIN
HOLD=HD
QOLD=QSJ
IF (NGDYN.GT.0) GD=CMPLX(GDYN(J),GCDYN(J))*GDFAC
IF (NPDYN.GT.0) PDIN=CMPLX(PDYN(J),PCDYN(J))*PDFAC*FAC3
IF (NHDYN.GT.0) HD=CMPLX(HDYN(J),HCDYN(J))*HDFAC*FAC3
IF (NQDYN.GT.0) QSJ=CMPLX(QDYN(J),QCDYN(J))*QDFAC
DGD=-GOLD
DPD=-POLD
DHD=-HOLD
DQD=-QOLD
IF (CDABS(GD).GT.1.0E-30) DGD=GD-GOLD
IF (CDABS(PDIN).GT.1.0E-30) DPD=PDIN-POLD
IF (CDABS(HD).GT.1.0E-30) DHD=HD-HOLD
IF (CDABS(QSJ).GT.1.0E-30) DQD=QSJ-QOLD
RETURN
END
SUBROUTINE TCOMP
COMMON /CONT /  ND,NC,NR,NTIN,NPOW,NW,NF,NDROP,NOSUB,NPRINT,NCH
COMMON /DATA1 /  TIN,DTIN,POWR,POWR,W,DW,
. PIN,PDROP,AP,BP,CP,ASB,BSB,AHT,BHT,CHT,
. TSAT,HF,HG,RHOF,RHOG,VISCF,CONF,CPF
COMMON /DATA2 /  DZ(40),A(40),DE(40),PH(40),ANG(40),RK(40),E(40),
. PKW(40),TCAP(40),CO(40),V(40),AA(40),BB(40),CC(40)
COMMON /DATA3 /  SIGMA,FREQ,DFREQ,FHZ(120),GDYN(120),HDYN(120),
. QDYN(120),PDYN(120),CFREQ(10),NDYN,NCFREQ
COMMON /EXSTAT/  DPT,FAC,FAC3,GC,PI,PUMP,PY,NIT,NODE
COMMON /STATIC/  ALFA(601),BF(600),CKO(601),DP1F(600),DRGDP(600),

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.           HS(600),HT(600),HTW(600),HW(600),PHI(601),
.           FHP(601),RHOS(600),RHOW(600),HTB(600),ZCL(2),
.           X(601),ALFK(40),CKK(40),DP1(40),DPR(40),DRGDK(40),
.           PHIK(40),PHIR(40),NSUB,NSAT,NCON,NTOT,MODE(40),
.           MBOIL(40,50)
COMMON /FUNCNC/ ALFA2,DELP1M,DELP1F,DELP1R,DELP1T,DELP2M,DELP2F,
.           DELP2G,DELP2R,DELP2T,DELPT,H,S,GD,HD,QSJ,PDIN
COMPLEX*16 ALFA2,DELP1M,DELP1F,DELP1R,DELP1T,DELP2M,DELP2F,
.           DELP2G,DELP2R,DELP2T,DELPT,H,S,GD,HD,QSJ,PDIN
COMPLEX*16 ALFA1,E1,E2,GD1,GD2,GF1,GF2,GG1,GG2,GM1,GM2,H1,H2,
.           PD1,PHI1,PHI2,PHIR1,PS,PT,PTEX,UF1,UF2,UG1,UG2,
.           X1,X2,QSD,TAU1,TAU2,TAUP,TAUS,TSUB
COMPLEX*16 ALFA4

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C
C
C

CALCULATION OF TRANSFER FUNCTION COMPONENTS FOR A SPECIFIED FREQUE

```

NPD=1
PI=3.14159/180.0
EHC=1.0
A1=A(1)
ULI=W/RHOW(1)/A1/FAC/FAC
GD1=GD
H=HD
PD1=PDIN
ALFA1=0.0
ALFA2=0.0
X1=0.0
GG1=0.0
UG1=0.0
PHI1=0.0
PHIR1=0.0
E1=ULI*GD1
DELP1M=0.0
DELP1F=0.0
DELP1R=0.0
DELP2M=0.0
DELP2F=0.0
DELP2G=0.0
DELP2R=0.0
I=0
DO 200 K=1,NCH
NODE=MODE(K)
GD1=GD1*A1/A(K)
THETA=COS(PI*ANG(K))
TAUF=TCAP(K)*FAC3/PH(K)/FAC
PKWK=PKW(K)*FAC3
IF (K.GT.ND,AND.K.LE.(ND+NC)) PKWK=POWR*PKWK
AC=A(K)*FAC*FAC
G=W/AC
Q=PKWK/AC/DZ(K)/FAC
DO 160 J=1,NODE
NBOIL=MBOIL(K,J)
ZC=DZ(K)*FAC/NODE/NBOIL
IF (NC.GT.1) GO TO 10
IF (K.IE.ND,OR.K.GT.(ND+NC)) GO TO 10
ZC=ZCL(J)
10 DO 150 L=1,NBOIL

```

```

I=I+1
TAUH=TAUF/HT(I)
PS=1.0/(1.0+TAUH*S)
QSD=0.0
IF (NPD.EQ.0) PD1=0.0
IF (K.GT.ND.AND.K.LE.(ND+NC)) QSD=POWR*QSJ
IF (NSAT.GT.0.AND.I.GE.NSAT) GO TO 100
C
C SINGLE PHASE REGION
IF (CDABS(S)) 20,20,15
15 PT=S/G*(RHOW(I)+Q*HTW(I)*TAUH*PS)
PTEX=CDEXP((-PT*ZC))
H=H*PTEX+Q/G*(PS*QSD-(0.20*TAUH*S+1.0)*PS*GD1/G)*(1.0-PTEX)/PT
GO TO 25
20 H=H+Q*ZC*(QSD-GD1/G)/G
25 DELP1M=S*ZC*GD1+DELP1M
DELP1F=DP1F(I)*(2.0-BF(I))*GD1/G+DELP1F
ULI=G/RHOW(I)
IF (J.GT.1.OR.L.GT.1) GO TO 30
E2=ULI*GD1
IF (A(K).EQ.A1) E2=E1
DELP1R=E2-E1+2.0*DPR(K)*GD1/G+DELP1R
30 E1=ULI*GD1
DELP1T=DELP1M+DELP1F+DELP1R
DELP1T=DELP1T
GO TO 150
C
C***SUBCOOLED AND SATURATED BOILING REGIONS.
100 RHOD=RHOW(I)-RHOS(I)
HWS=HS(I)-HW(I)
FHP1=FHP(I)
FHPO=FHP(I+1)
XI=X(I)
X0=X(I+1)
DX=X0-XI
ALFAI=ALFA(I)
GGI=XI*G
GLI=(1.0-XI)*G
PHII=PHI(I)
CKI=CKO(I)
DRGDI=DRGDP(I)
IF (J.GT.1.OR.L.GT.1) GO TO 105
ALFAI=ALFK(K)
PHII=PHIK(K)
CKI=CKK(K)
DRGDI=DRGDK(K)
XE1=XI*(1.0-XI)*DRGDI/RHOS(I)
105 UGI=0.0
ULI=GLI/(1.0-ALFAI)/RHOW(I)
ALFAO=ALFA(I+1)

ALFM=(ALFAI+ALFAO)/2.0
GGO=X0*G
GLO=(1.0-X0)*G
UGO=GGO/ALFAO/RHOS(I)
ULO=GLO/(1.0-ALFAO)/RHOW(I)
PHIO=PHI(I+1)

```

```

XXD=RHOW(I)*DRGDPI/RHOD/RHOS(I)
XXI=(RHOS(I)+RHOD*XI)**2.0/CKI/RHOW(I)/RHOS(I)
XXO=(RHOS(I)+RHOD*XO)**2.0/CKO(I+1)/RHOW(I)/RHOS(I)
XXP=2.0*ALFM*DRGDPI
XXR=XO*(1.0-XO)*DRGDPI/RHOS(I)
XNI=RHOW(I)*(CKI+(1.0-CKI)*FHPI)-RHOD*ALFAI
XNO=RHOW(I)*(CKO(I+1)+(1.0-CKO(I+1))*FHP0)-RHOD*ALFAO
XAI=RHOD*(1.0-(1.0-XI)*FHPI)/XXI/HWS
XAO=RHOD*(1.0-(1.0-XO)*FHP0)/XXO/HWS
XBI=DRGDPI*(ALFAI-XI*(1.0-XI)*RHOD/XXI/RHOS(I))
XBO=DRGDPI*(ALFAO-XO*(1.0-XO)*RHOD/XXO/RHOS(I))

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C
C***ROILING BOUNDARY EQUATIONS.
IF (I.GT.NSAT) GO TO 110
H1=H
X1=(1.0-FHPI)*H1/HWS
ALFA1=X1/XXI
GG1=G*X1
GF1=GD1-GG1
UF1=GF1/RHOW(I)+ULI*ALFA1
UG1=0.0
GM1=2.0*ULI*GD1
GO TO 120
110 UGI=GG1/ALFA1/RHOS(I)
IF (J.GT.1.OR.L.GT.1) GO TO 120
GG1=XI*GD1+G*X1
GF1=GD1-GG1
ALFA1=(X1-XE1*PD1)/XXI
UG1=UGI*(GG1/GG1-ALFA1/ALFAI-DRGDPI*PD1/RHOS(I))
UF1=ULI*(GF1/GLI+ALFA1/(1.0-ALFAI))
GM1=ULI*GF1+GLI*UF1+UGI*GG1+GG1*UG1
PHI1=(PHI1-1.0)*(BB(K)*(X1/XI-XXD*PD1)+CC(K)*GD1/G)
PHIR1=(PHIR(K)-1.0)*(X1/XI-XXD*PD1)
120 TAUS=ZC*Q*(1.0+(1.0-HTR(I))*TAUH*S)*PS/G
TSUB=Q*TAUH*PS*HTW(I)
TAU1=ZC*(XNI+FHPI*TSUB+XAI*TAUS/2.0)/2.0/G
TAU2=ZC*(XNO+FHP0*TSUB+XAO*TAUS/2.0)/2.0/G
TAUP=ZC*(ALFM+(XBI+XBO)*TAUS/4.0)/G
H2=((1.0-S*TAU1)*H1+0*ZC*PS/G*QSD-TAUS*GD1/G+S*TAUP*PD1)/
(1.0+S*TAU2)
X2=(1.0-(1.0-XO)*FHP0)*H2/HWS
ALFA2=(X2-XXR*PD1)/XXO
IF (K.EQ.(NCH-1))ALFA4=ALFA2
GD2=GD1+S*ZC*(RHOD*(ALFA1+ALFA2)-XXP*PD1)/2.0
GG2=G*X2+XO*GD2
GF2=GD2-GG2
UG2=UGO*(GG2/GGO-ALFA2/ALFAO-DRGDPI*PD1/RHOS(I))
UF2=ULO*(GF2/GLO+ALFA2/(1.0-ALFAO))
GM2=ULO*GF2+GLO*UF2+UGO*GG2+GGO*UG2
PHI2=(PHI0-1.0)*(BB(K)*(X2/XO-XXD*PD1)+CC(K)*GD2/G)
DELP2M=GM2-GM1+S*ZC*(GD1+GD2)/2.0+DELP2M
DELP2F=DP1F(I)*((2.0-BF(I))*(PHI0*GD2+PHI1*GD1)/G
+PHI1+PHI2)/2.0+DELP2F
DELP2G=-(RHOD*(ALFA1+ALFA2)-XXP*PD1)*ZC*GC*THETA/2.0+DELP2G
IF (J.GT.1.OR.L.GT.1) GO TO 125
E2=(UGI*UGI-ULI*ULI)*X1/2.0+(1.0-XI)*ULI*UF1+XI*UGI*UG1
IF (A(K).EQ.A1) E2=E1
RHOM=RHOD/(RHOS(I)+RHOD*XI)

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RH0=RH0M*RHOW(I)*RHOS(I)/RHOD
XEK=XE1/(1.0-XI)*RHOW(I)/RHOD
DELP2R=RHO*(E2-E1)-RHOM*DPI(K)*(X1-XEK*PD1)+
      DPR(K)*(2.0*GD1/G+PHIR1/PHIR(K))+DELP2R
125 DELP2T=DELP2M+DELP2F+DELP2G+DELP2R
    DELPT=DELPT1+DELP2T
    E1=(UG0*UG0-ULO*ULO)*X2/2.0+(1.0-X0)*ULO*UF2+X0*UG0*UG2
    ALFA1=ALFA2
    GD1=GD2
    GG1=GG2
    GF1=GF2
    GM1=GM2
    H1=H2
    UF1=UF2
    UG1=UG2
    PHI1=PHI2
    X1=X2
150 PD1=PUIN-DELPT
160 CONTINUE
200 A1=A(K)
    ALFA2=ALFA4
    RETURN
    END
SUBROUTINE TPRINT
COMMON / CONT / ND,NC,NR,NTIN,NPOW,NW,NF,NDROP,NOSUB,NPRINT,NCH
COMMON /DATA1 / TIN,DTIN,POWR,DPOWR,W,DW,
      PIN,PDROP,AP,BP,CP,ASB,BSB,AHT,BHT,CHT,
      TSAT,HF,HG,RHOF,RHOG,VISCF,CONF,CPF
COMMON /DATA3 / SIGMA,FREQ,DFREQ,FHZ(120),GDYN(120),HDYN(120),
      QDYN(120),PDYN(120),CFREQ(10),NDYN,NCFREQ
COMMON /PRINT / AA(60),AT(60),GA(60),GADB(60),GT(60),GTDB(60),
      QA(60),QT(60),RA(60),RT(60)
COMMON /CPLOT / TPLT1(20),TPLT2(20),NPLOT
COMMON /ERROR / CSPEC(80),CPHASE(80),ASPEC(80),APHASE(80),NERR
REAL XX(2),YY(2)
15 WRITE (3,200)
   DO 20 J=1,NF
20 WRITE (3,210) FHZ(J),RT(J),QT(J),GT(J),GTDB(J),AT(J),RA(J),QA(J),
      GA(J),GADB(J),AA(J)
   IF (NERR,GE,0) GO TO 30
   SUM=0.0
   SUMA=0.0
   WRITE (3,300)
   DO 25 J=1,NF
   ERR=CSPEC(J)-GT(J)
   ERR=ERR*ERR
   SUM=SUM+ERR
   ERRA=ASPEC(J)-GA(J)
   ERRA=ERRA*ERRA
   SUMA=SUMA+ERRA
25 WRITE (3,290) FHZ(J),CSPEC(J),GT(J),ERR,ASPEC(J),GA(J),ERRA
   SUM=SUM/NF
   SUMA=SUMA/NF
   WRITE (3,310) SUM,SUMA
30 IF (NF.EQ.1) GO TO 100
   NLOG=1
   IF (NPLOT) 35,100,35

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35 XX(1)=0.0
   XX(2)=CSPEC(1)
   YY(1)=CPHASE(1)
   YY(2)=YY(1)
   CALL PLIM (NF,CSPEC,CPHASE,XX,YY)
   GO TO 37
36 XX(1)=0.0
   XX(2)=GT(1)
   YY(1)=AA(1)
   YY(2)=YY(1)
37 CALL PLIM (NF,GT,AA,XX,YY)
   CALL TPLLOT (FHZ,GT,CSPEC,AT,CPHASE,XX,YY,NF,NPLOT)
   CALL XYPEXT (0.,-1.0,TPLT1,80,2,0)
   XX(1)=0.0
   IF (NPLOT,GT,0) GO TO 44
   XX(2)=ASPEC(1)
   YY(1)=APHASE(1)
   YY(2)=YY(1)
   CALL PLIM (NF,ASPEC,APHASE,XX,YY)
   GO TO 46
44 XX(2)=GA(1)
   YY(1)=AA(1)
   YY(2)=YY(1)
46 CALL PLIM (NF,GA,AA,XX,YY)
   CALL TPLLOT (FHZ,GA,ASPEC,AA,APHASE,XX,YY,NF,NPLOT)
   CALL XYPEXT (0.,-1.0,TPLT2,80,2,0)
   CALL XYEND
200 FORMAT ('1',23X,'TOTAL PRESSURE DROP PERTURBATION',31X,'EXIT ',
.         'VOID FRACTION PERTURBATION'/24X,32('-',),31X,31('-',)/3X,
.         'FREQ',5X,'REAL',8X,'IMAG',8X,'GAIN',8X,'GAIN DB',5X,
.         'PHASE',10X,'REAL',8X,'IMAG',8X,'GAIN',8X,'GAIN DB',5X,
.         'PHASE')
210 FORMAT (1X,F6.4,1P5E12.3,2X,'*',5E12.3)
290 FORMAT (1X,F6.4,1P3E12.3,2X,'*',3E12.3)
300 FORMAT ('1 ERROR ANALYSIS IN CROSS-SPEC DENSITY MODULUS'/
.         3X,'FREQ',5X,'EXPT',8X,'THEORY',6X,'SQ ERR',
.         9X,'EXPT',8X,'THEORY',6X,'SQ ERR'/)
310 FORMAT ('0',10X,'MEAN SQUARE ERROR =',1PE12.4,8X,'MEAN SQUARE ',
.         'ERROR =',E12.4)
100 RETURN
   END
   SUBROUTINE PLIM (NF,YA,YB,X,Y)
   DIMENSION YA(NF),YB(NF),X(2),Y(2)
   DO 10 J=2,NF
   IF (YA(J).GT.X(2)) X(2)=YA(J)
   IF (YB(J).LT.Y(1)) Y(1)=YB(J)
   IF (YB(J).GT.Y(2)) Y(2)=YB(J)
10 CONTINUE
   RETURN
   END
   SUBROUTINE TPLLOT (X,YAA,YAB,YBA,YBR,XX,YY,NF,NPLOT)
   DIMENSION X(NF),YAA(NF),YAB(NF),YBA(NF),YBR(NF),XX(2),YY(2)
   COMMON /XYCOM$ / IDUM(3),HOR(2),VER(2)
   REAL*4 ZERO(2)/2*0.0/
   DO 20 I=1,2
   IF (I.EQ.2) GO TO 10
   CALL XYSKEW (1,1.0,6.0)

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CALL XYPAPE (-6.,-4.,X(1),X(NF),XX(1),XX(2))
CALL XYNAMY ('MODULUS',7)
CALL XYNAMX ('FREQUENCY (HZ)',14)
CALL XYSETL (2)
CALL XYLINE (X,YAA,NF)
IF (NPLOT.GT.0) GO TO 20
CALL XYSETL (1)
CALL XYLINE (X,YAB,NF)
GO TO 20
10 CALL XYSKEW (-1,0.,-5.0)
CALL XYPAPE (-6.,-4.,X(1),X(NF),YY(1),YY(2))
CALL XYNAMY ('PHASE',5)
CALL XYNAMX ('FREQUENCY (HZ)',14)
IF (VER(1)*VER(2).GT.0.0) GO TO 15
CALL XYSETL (1)
CALL XYLINE (HOR,ZERO,2)
15 CALL XYSETL (2)
CALL XYLINE (X,YRA,NF)
IF (NPLOT.GT.0) GO TO 20
CALL XYSETL (1)
CALL XYLINE (X,YBB,NF)
20 CONTINUE
RETURN
END
SURROUTINE CPRINT
COMMON /EXSTAT/ DPT,FAC,FAC3,GC,PI,PUMP,PY,NIT,NODE
COMMON /FUNCNC/ ALFA2,DELP1M,DELP1F,DELP1R,DELP1T,DELP2M,DELP2F,
.      DELP2G,DELP2R,DELP2T,DELPT,H,S,GD,HD,QSJ,PDIN
.
COMPLEX*16 ALFA2,DELP1M,DELP1F,DELP1R,DELP1T,DELP2M,DELP2F,
.      DELP2G,DELP2R,DELP2T,DELPT,H,S,GD,HD,QSJ,PDIN,COMP
WRITE (3,100)
PY=6.28319
PI=57.2958
SREAL=DREAL(S)/PY
SIMAG=DAIMAG(S)/PY
WRITE (3,120) SREAL,SIMAG
WRITE (3,110)
DO 20 J=1,10
IF (J.EQ.1) COMP=DELP1M/FAC3
IF (J.EQ.2) COMP=DELP1F/FAC3
IF (J.EQ.3) COMP=DELP1R/FAC3
IF (J.EQ.4) COMP=DELP1T/FAC3
IF (J.EQ.5) COMP=DELP2M/FAC3
IF (J.EQ.6) COMP=DELP2F/FAC3
IF (J.EQ.7) COMP=DELP2G/FAC3
IF (J.EQ.8) COMP=DELP2R/FAC3
IF (J.EQ.9) COMP=DELP2T/FAC3
IF (J.EQ.10) COMP=DELPT
CREAL=DREAL(COMP)
CIMAG=DAIMAG(COMP)
CGAIN=CDABS(COMP)
IF (J.GT.1) GO TO 5
IF (CDABS(S)) 10,10,5
5 CDB=0.0
IF (CGAIN.GT.0.0) CDB=20.0*ALOG10(CGAIN)
CPHASE=0.0
IF (CREAL.EQ.0.0.AND.CIMAG.GT.0.0) CPHASE=90.0
IF (CREAL.EQ.0.0.AND.CIMAG.LT.0.0) CPHASE=-90.0

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IF (CREAL.NE.0.0) CPHASE=ATAN2(CIMAG,CREAL)*PI
GO TO 15
10 CDB=0.0
CPHASE=0.0
IF (CREAL.EQ.0.0.AND.CIMAG.GT.0.0) CPHASE=90.0
IF (CREAL.EQ.0.0.AND.CIMAG.LT.0.0) CPHASE=-90.0
15 IF (J.EQ.1) WRITE (3,130) CREAL,CIMAG,CGAIN,CDB,CPHASE
IF (J.EQ.2) WRITE (3,140) CREAL,CIMAG,CGAIN,CDB,CPHASE
IF (J.EQ.3) WRITE (3,150) CREAL,CIMAG,CGAIN,CDB,CPHASE
IF (J.EQ.4) WRITE (3,160) CREAL,CIMAG,CGAIN,CDB,CPHASE
IF (J.EQ.5) WRITE (3,170) CREAL,CIMAG,CGAIN,CDB,CPHASE
IF (J.EQ.6) WRITE (3,180) CREAL,CIMAG,CGAIN,CDB,CPHASE
IF (J.EQ.7) WRITE (3,190) CREAL,CIMAG,CGAIN,CDB,CPHASE
IF (J.EQ.8) WRITE (3,200) CREAL,CIMAG,CGAIN,CDB,CPHASE

IF (J.EQ.9) WRITE (3,210) CREAL,CIMAG,CGAIN,CDB,CPHASE
IF (J.EQ.10) WRITE (3,220) CREAL,CIMAG,CGAIN,CDB,CPHASE
20 WRITE (3,110)
100 FORMAT (/41X,'REAL',13X,'IMAGINARY',13X,'GAIN',14X,'GAIN DB',
.          15X,'PHASE'/41X,4(' '),13X,9(' '),13X,4(' '),14X,7(' '),
.          15X,5(' '))
110 FORMAT(1X,125(' '))
120 FORMAT (/ ' LAPLACE OPERATOR (HZ)',14X,1PE12.4,8X,E12.4)
130 FORMAT (' 1-PHASE MOMENTUM',19X,4(1PE12.4,8X),0PF9.2)
140 FORMAT (' 1-PHASE FRICTION',19X,4(1PE12.4,8X),0PF9.2)
150 FORMAT (' 1-PHASE RESTRICTION',16X,4(1PE12.4,8X),0PF9.2)
160 FORMAT (' 1-PHASE TOTAL',22X,4(1PE12.4,8X),0PF9.2)
170 FORMAT (' 2-PHASE MOMENTUM',19X,4(1PE12.4,8X),0PF9.2)
180 FORMAT (' 2-PHASE FRICTION',19X,4(1PE12.4,8X),0PF9.2)
190 FORMAT (' 2-PHASE GRAVITY',20X,4(1PE12.4,8X),0PF9.2)
200 FORMAT (' 2-PHASE RESTRICTION',16X,4(1PE12.4,8X),0PF9.2)
210 FORMAT (' 2-PHASE TOTAL',22X,4(1PE12.4,8X),0PF9.2)
220 FORMAT (' CHANNEL TOTAL',22X,4(1PE12.4,8X),0PF9.2)
RETURN
END
FUNCTION FRIC (REW,E)
RE=ABS(REW)
FRICL=16.0/RE
FF=0.00080
DO 30 N=1,50
F=1.0/(4.0*ALOG10(RE*FF/(4.67+E*RE*FF))+2.28)
IF (ABS(1.0-FF*FF/(F*F)).LE.0.0010) GO TO 50
FF=ABS(F)
30 CONTINUE
FRIC=F*F
WRITE (3,40) RE,E,FRIC
40 FORMAT (1X,'FRICTION FACTOR ITERATION DID NOT CONVERGE : RE =',
.E12.4,' E =',E12.4,' F =',E12.4)
50 FRIC=F*F
IF (FRIC.LT.FRICL) FRIC=FRICL
RETURN
END
SUBROUTINE TABLES
COMMON /TABLE / TEMP(100),PRESS(100),ENTHF(100),ENTHFG(100),
.          ENTHG(100),SPVF(100),SPVG(100),DENF(100),
.          DENG(100),VISCDF(100),CONDF(100),SPHF(100),
.          PRF(100),FAC1,FAC2,ADRG,BDRG,NCARDS

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```

NCARD=NCARDS
NCARDS=IABS(NCARDS)
SP=0.0
SPP=0.0
SVG=0.0
SVP=0.0
SVV=0.0
DO 10 I=1,NCARDS
ENTHF(I)=ENTHF(I)*FAC1
ENTHG(I)=ENTHG(I)*FAC1
ENTHFG(I)=ENTHG(I)-ENTHF(I)
DENF(I)=1.0/SPVF(I)
DENG(I)=1.0/SPVG(I)
VISCDF(I)=VISCDF(I)*FAC2
SPHF(I)=SPHF(I)*FAC1
PRF(I)=SPHF(I)*VISCDF(I)/CONDF(I)
PLOG=ALOG(PRESS(I))
VGLLOG=ALOG(SPVG(I))
SP=SP+PLOG
SPP=SPP+PLOG*PLOG
SVG=SVG+VGLLOG
SVP=SVP+PLOG*VGLLOG
10 SVV=SVV+VGLLOG*VGLLOG
A=SVP-SP*SVG/NCARDS
B=SPP-SP*SP/NCARDS
C=SVV-SVG*SVG/NCARDS
BDRG=A/B
ADRG=EXP((SVG-BDRG*SP)/NCARDS)
ADRG=-BDRG/ADRG
BDRG=-(1.0+BDRG)
CORR=ABS(A)/SQRT(ABS(B*C))
IF (NCARD.GT.0) WRITE (3,100) ADRG,BDRG,CORR
100 FORMAT ('POWER LAW FIT FOR DRHOG/DP = ADRG*PRESS(I)**BDRG :'/
.          1X,'COEFFICIENTS ADRG =',1PE14.6,' : BDRG =',E14.6,
.          ' : CORRELATION COEFFICIENT =',E14.6)
RETURN
END
FUNCTION P(T)
COMMON /TABLE / TEMP(100),PRESS(100),ENTHF(100),ENTHFG(100),
.              ENTHG(100),SPVF(100),SPVG(100),DENF(100),
.              DENG(100),VISCDF(100),CONDF(100),SPHF(100),
.              PRF(100),FAC1,FAC2,ADRG,BDRG,NCARDS
CALL POLATE (TEMP,PRESS,T,FUNT,NCARDS)
P=FUNT
RETURN
ENTRY THF(T)
CALL POLATE (TEMP,ENTHF,T,FUNT,NCARDS)
THF=FUNT
RETURN
ENTRY THFG(T)
CALL POLATE (TEMP,ENTHFG,T,FUNT,NCARDS)
THFG=FUNT
RETURN
ENTRY THG(T)
CALL POLATE (TEMP,ENTHG,T,FUNT,NCARDS)
THG=FUNT
RETURN
ENTRY TVF(T)

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```

CALL POLATE (TEMP,SPVF,T,FUNT,NCARDS)
TVF=FUNT
RETURN
ENTRY TVG(T)

CALL POLATE (TEMP,SPVG,T,FUNT,NCARDS)
TVG=FUNT
RETURN
ENTRY TDF(T)
CALL POLATE (TEMP,DENF,T,FUNT,NCARDS)
TDF=FUNT
RETURN
ENTRY TDG(T)
CALL POLATE (TEMP,DENG,T,FUNT,NCARDS)
TDG=FUNT
RETURN
ENTRY TVISCD(T)
CALL POLATE (TEMP,VISCDF,T,FUNT,NCARDS)
TVISCD=FUNT
RETURN
ENTRY TCOND(T)
CALL POLATE (TEMP,CONDF,T,FUNT,NCARDS)
TCOND=FUNT
RETURN
ENTRY TSH(T)
CALL POLATE (TEMP,SPHF,T,FUNT,NCARDS)
TSH=FUNT
RETURN
ENTRY TPRN(T)
CALL POLATE (TEMP,PRF,T,FUNT,NCARDS)
TPRN=FUNT
RETURN
END
FUNCTION SAT(Y,X,P,NCARDS)
DIMENSION Y(NCARDS),X(NCARDS)
DO 10 J=1,NCARDS
IF (P.LE.X(1)) GO TO 15
IF (P.LT.X(J)) GO TO 20
10 CONTINUE
.15 WRITE (3,100) J,X(J),X(NCARDS),P
1000 FORMAT (1X,I5,5X,'FUNCTION SAT - PRESSURE OUTSIDE DATA RANGE :',
.      ' MINIMUM =',1PE10.3,' MAXIMUM =',E10.3,' PRESSURE =',
.      E10.3)
SAT=Y(1)
RETURN
20 GRAD=(Y(J)-Y(J-1))/(X(J)-X(J-1))
SAT =Y(J-1)+GRAD*(P-X(J-1))
RETURN
END
SUBROUTINE POLATE (X,Y,T,FUNT,NCARDS)
DIMENSION X(NCARDS),Y(NCARDS)
DO 10 J=1,NCARDS
IF (T.LT.X(1)) GO TO 15
IF (T.LT.X(J)) GO TO 20
10 CONTINUE
15 WRITE (3,100) J,X(J),X(NCARDS),T
1000 FORMAT(1X,I5,5X,'POLATE - TEMPERATURE OUTSIDE DATA RANGE: MINI',
.      ' MUM =',1PE10.3,' MAXIMUM =',E10.3,' TEMPERATURE =',E10.3)

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```
FUNT=Y(J)
RETURN
20 GRAD=(Y(J)-Y(J-1))/(X(J)-X(J-1))
FUNT=Y(J-1)+GRAD*(T-X(J-1))
RETURN
END
```

APPENDIX E
'LOCO' CODE SUBPROGRAM FUNCTIONS

SUBPROGRAM	FUNCTION
MAIN	Reads, writes input data; calls subroutines TABLES (NCARDS >0), EXEC.
EXEC	Control subroutine; calls subroutines BYPASS (NB>0), STATE, TFUNCN (NDYN>0); calculates flow rate for specified pressure drop (NDROP = 1).
STATE	Calculates steady state variables (mass fluxes, velocities, void fraction, pressure drops, etc.) in each space node between channel inlet and outlet.
BYPASS	Calculates the steady state and dynamic pressure drop in the bypass (NB sections).
TFUNCN	Control subroutine for dynamic calculations; calls subroutines DYNIN, TCOMP, BYPASS (NB>0), TPRINT, CPRINT; calculates crossover frequency when perturbed pressure drop crosses real axis of polar plot.
DYNIN	Edits the dynamic inputs read in as data.
TCOMP	Calculates the single phase and two-phase region perturbed pressure drop components and the perturbed void fraction (channel exit) for a specified value of the Laplace operator (s).
TPRINT	Prints the total pressure drop and channel exit void fraction frequency response functions; calls NORGRA (NPRINT ≤ 2).
CPRINT	Prints the single phase and two-phase region perturbed pressure drop components for a specified value of the Laplace operator (s).

<u>SUBPROGRAM</u>	<u>FUNCTION</u>
TPLOT	Plots the total pressure drop and channel exit void fraction frequency response functions (magnitude, phase and polar plots).
FRIC	Calculates single phase friction factor (STATE, BYPASS).
TABLES	Calculates the complete fluid properties data from the input data (NCARDS >0).
SAT	Calculates local saturation temperature from the local static pressure.
POLATE	Linear interpolation subroutine for fluid property calculations.

APPENDIX F
SAMPLE TEST CHANNEL DIMENSIONS

REGION	LENGTH (cm)	DIAMETER (cm)	FLOW AREA (cm ²)	HYDRAULIC DIAMETER (cm)	
Inlet*	52.7	2.58 ID	5.238	2.58	Horizontal pipe with venturi flowmeter 45 dia. from inlet. Vertical annulus with T-junction at inlet.
	17.8	0.953 ID 1.352 OD	0.722	0.399	
Channel*	91.5	0.953 ID	0.722	0.399	Vertical annulus with heated inner wall, glass outer wall.
		1.352 OD			
Riser	91.5	2.54 ID	5.067	2.54	Vertical glass tube.

* Test region for input-output measurements = inlet + channel regions

APPENDIX G
COMPUTER PROGRAM INPUT DATA FOR SAMPLE TEST CASE

ND	NC	NR	NB	NTIN	NPOW	NW	NF	NCFREQ
2	13	0	0	1	1	1	51	0
NDROP	NODES	NOSUB	NDYN	NPRINT	NPLOT	NCARDS		
0	600	2	1	1	-1	-21		
NGDYN	NPDYN	NHDYN	NDDYN	NSPEC	NPROP			
1	0	0	0	0	1			
TIN	DTIN							
4.138E+01	2.000E+00							
POWR	DPOWR							
1.835E+00	0.0							
W	DW							
8.830E-02	2.000E-02							
TCFAC	COFAC	VFAC	AAFAC	BBFAC	CCFAC			
0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0			
PIN	PDRPF	WTOT	AP	BP	CP			
1.628E+02	3.863E+01	2.100E+00	0.0	0.0	0.0			
ASB	BSB	AHT	BHT	CHT				
1.930E+02	2.757E-03	1.640E-01	0.0	1.000E+00				
SIGMA	FREQ	DFREQ						
0.0	0.0	2.000E-02						
GDFAC	PIFAC	HDFAC	QDFAC					
1.000E+00	1.000E+00	1.000E+00	1.000E+00					
GDYN(N)								
1.000E+00								
DZ(N)	A(N)	DE(N)	PH(N)	ANG(N)	RK(N)	E(N)		
INLET PIPING REGION								
5.270E+01	5.238E+00	2.581E+00	8.108E+00	9.000E+01	0.0	0.0		1
1.780E+01	7.222E-01	3.991E-01	2.992E+00	0.0	1.000E+00	0.0		2
CHANNEL REGION								
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		3
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		4
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		5
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		6
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		7
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		8
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		9
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		10
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		11
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		12
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		13
9.150E+00	7.222E-01	3.991E-01	2.992E+00	0.0	0.0	0.0		14
5.080E+00	4.355E+00	1.588E+00	2.992E+00	0.0	2.369E+01	0.0		15
PKW(N)	TCAP(N)	CO(N)	V(N)	AA(N)	BB(N)	CC(N)		
INLET PIPING REGION								
0.0	0.0	1.000E+00	0.0	1.000E+00	1.000E+00	0.0		1
0.0	0.0	1.000E+00	0.0	1.000E+00	1.000E+00	0.0		2
CHANNEL REGION								
1.000E-01	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		3
1.000E-01	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		4
1.000E-01	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		5
1.000E-01	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		6
1.000E-01	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		7
1.000E-01	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		8
1.000E-01	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		9
1.000E-01	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		10
1.000E-01	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		11
1.000E-01	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		12
0.0	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		13
0.0	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		14
0.0	0.0	9.835E-01	0.0	7.557E-01	1.000E+00	0.0		15
FLOW - PRESSURE DROP TRANSFER FUNCTION								
FLOW - EXIT VOID FRACTION TRANSFER FUNCTION								