

A Condensed Review of Nuclear Reactor
Thermal-Hydraulic Computer Codes
for Two-Phase Flow Analysis

by

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FOR TWO-PHASE FLOW ANALYSIS

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ABSTRACT

A review is made of the computer codes developed in the U.S. for thermal-hydraulic analysis of nuclear reactors. The intention of this review is to compare these codes on the basis of their numerical method and physical models with particular attention to the two-phase flow and heat transfer characteristics. A chronology of the most documented codes such as COBRA and RELAP is given. The features of the recent codes as RETRAN, TRAC and THERMIT are also reviewed. The range of application as well as limitations of the various codes are discussed.

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Abbreviations

The following abbreviations are referenced in this report:

ATWS: Anticipated Transients Without Scram
COBRA: Coolant Boiling in Rod Arrays
DBA: Drift Flux Model
DSM: Dynamic Slip Model
EPRI: Electric Power Research Institute
FLECHT: Full Length Emergency Cooling Heat Transfer
HEM: Homogeneous Equilibrium Model
LOFT: Loss of Flow Transient
LOCA: Loss of Coolant Accident
MWR: Method of Weighted Residuals
NSSS: Nuclear Steam Supply System
RIAs: Reactivity Insertion Accidents
RETRAN: RELAP4 - TRANsient
TRAC: Transient Reactor Analysis Code
UHI: Upper Heat Injection
WREM: Water Reactor Evaluation Model

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1. Introduction

Numerous computer codes have been written to calculate the thermal-hydraulic characteristics of the reactor core and the primary loop under steady-state and operational transient conditions as well as hypothetical accidents. New versions of some of these codes are still to come. The main purposes of the continuing effort in the development of such computer codes have been improved computational effectiveness and improved ability to predict the response of the core and the primary loop. Therefore, efforts have been continued to incorporate the recent models and methods of analysis in the areas of both hydrodynamics and heat transfer in two-phase flow to the extent that their prediction are reasonably reliable. For example, such a step by step development has been effected in the various versions of COBRA and RELAP Computer Programs.

The code users are therefore confronted with the need to develop criteria to choose the most appropriate version to handle a specified case. This is a two pronged decision since it requires not only an evaluation of the models and methods used in each code but also a comparison between the results and experimental data to observe how well these data are predicted.

An attempt is made here to address the first step, i.e., comparison of the models and methods. To accomplish this, a study was made on the physical models and numerical methods which have been employed in the WOSUB, RETRAN, TRAC and THERMIT as well as various versions of COBRA and RELAP as listed in Table 1-1.

Table 1-1

List of reviewed thermal hydraulic codes.

<u>Name of Code</u>	<u>Reference Number</u>
COBRA-I	1
COBRA-II	2
COBRA-III	3
COBRA-IIIC	4
COBRA-IIIP	5
COBRA-IV-I	6
COBRA-DF	7
COBRA-TF	8
RELAP2	9
RELAP3	10
RELAP3B-MOD101	11
RELAP4	12
RELAP4-MOD5	13
RELAP4-MOD6	14
RELAP4-MOD7	14
RELAP4-EM	15
RELAP5	16
WOSUB	17
RETRAN	18
TRAC	19
THERMIT	20

These codes, especially COBRA and RELAP series, are well known thermal hydraulic computer codes and have been extensively used in the nuclear industry. WOSUB and RETRAN introduce a new treatment for the hydrodynamics modeling. TRAC and THERMIT have gone further by applying the most advanced existing treatment of the two-phase flow, namely, three-dimensional, two-fluid, non-equilibrium model.

In the comparison that follows, both the advantages and drawbacks are noted in each code and ultimately it is attempted to assess the capability of each code for handling a specified case.

2. Classification of Nuclear Reactor Thermal-Hydraulic Codes

The existing thermal hydraulic codes may be classified under several categories as follows:

1) Capability of the system analysis

This contains two different classes of codes, namely, system component codes and loop codes. Basically, the hot channel or the fuel behavior codes are system component codes; however, some of these codes are extended to other situations far removed from subchannel (one channel) geometry. Integration of the down comer, jet pumps (in BWR's), bottom flooding, UHI and the like models into a component codes, makes it a vessel code. As distinct from the loop codes which are devised to analyze the whole primary side including reactor core and the secondary side, a variety of codes ranging from hot channel to vessel codes are called system component codes in this report.

2) Type of two-phase flow modeling

This part deals with the mathematical models used in thermal hydraulic codes to calculate the characteristics of the two-phase flow either in the reactor core or in the primary loop. The two pertinent methods in this respect, namely, the homogeneous equilibrium model and the two-fluid model fall in this category.

3) Range of application

Since the capability of each code to handle flow and fuel rod calculations depends upon the mathematical models used to represent the physical situations as well as the numerical methods employed, codes can be classified in these respects into steady-state, transient and accident analysis (such as LOCA) codes. Naturally, the more demanding codes in this respect are ATWS and LOCA codes.

4) Type of application

Codes may also be classified based upon their types, i.e., Best Estimate (BE) type and Evaluation Model (EM) type. The latter group are basically devised for the purpose of licensing.

The type of nuclear reactor for which thermal hydraulic codes are devised (such as PWR, BWR and LMFBR) may be another category. A detailed discussion concerning each mentioned category is presented in the following sections.

2.1 Classification According to System Analysis Capability

2.1.1 Component Codes

Core thermal hydraulic assessments necessitates analysis

of fluid passing axially along the parallel rod arrays. Such analysis is difficult to conduct due to the degree of freedom associated with parallel rod array and the two-phase flow and heat transfer involved in nuclear reactors. In addition, radial and axial variations of the fuel rod power generation exacerbates this situation.

Assumptions have been made to simplify the task of modeling the hydrodynamics and heat transfer characteristics of the rod arrays.

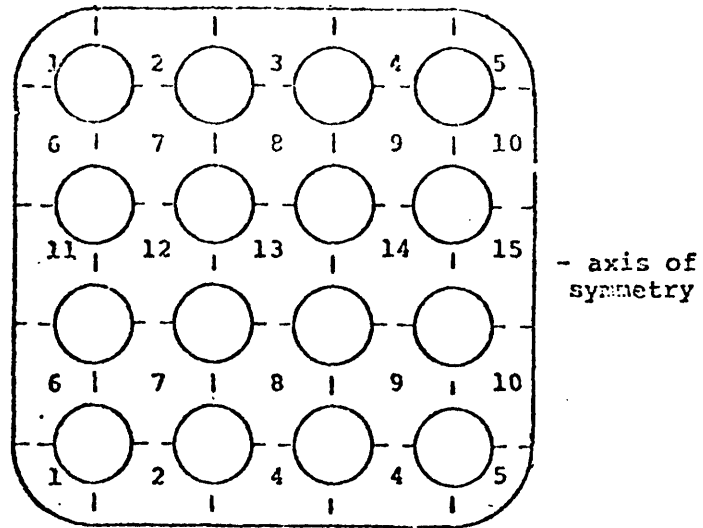
Generally, there are three pertinent methods⁽²¹⁾ used in rod bundle thermal hydraulic analysis of the nuclear reactor core as well as heat exchangers, namely, (a)-subchannel analysis, (b)-porosity and distributed resistance approach and finally (c)-benchmark rod-bundle analysis which uses a boundary fitted coordinate system.

The first approach is widely used in the subchannel codes such as COBRA, FLICA, HAMBO and THINC. Whereas the second approach is employed in THERMIT.

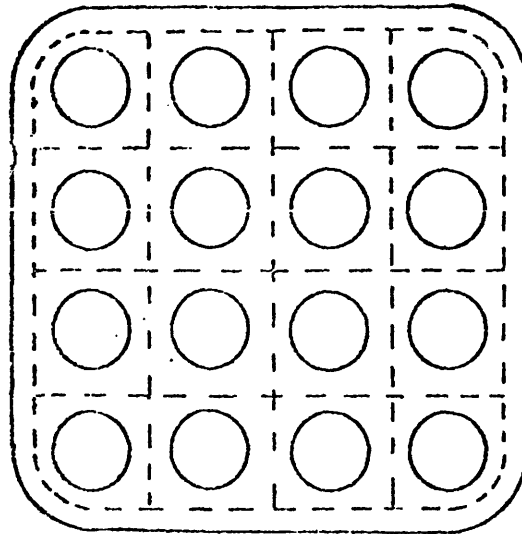
The subchannel approach will be more elaborated upon here, while a discussion in detail of these three concepts is presented in Ref 21.

In the subchannel approach, the rod array is considered to be subdivided into a number of parallel interacting flow subchannels between the rods. The fluid enthalpy and mass velocity is then found by solving the field or conservation equations for the control volume taken around the subchannel.

Although a rod-centered system with subchannel boundaries



(a)



(b)

Fig. 1^[22] (a) Coolant centered subchannel and conventional subchannel numbering.
(b) Rod center system with subchannel boundaries.

defined by lines of "zero-shear stress" between rods (Fig. 1-b) seems to be a well-defined control volumes,^{*} it has become customary to consider a coolant centered subchannel as a control volume (Fig. 1-a). The number of the above-mentioned control volumes axially is as many as the number of the channel length intervals.

Unlike the benchmark rod-bundles approach, the subchannel approach does not take into account the fine structure of both velocity and temperature within a subchannel.^{**} In other words, there are no radial gradients of flow and enthalpy in the subchannels but only across subchannel boundaries. Therefore, the flow parameters such as velocity, void fraction, and temperature are averaged over the subchannel area. Furthermore, the averaged values are assumed to be located at the subchannel centroid. The following example elaborates the latter assumption.⁽⁴⁸⁾

The transverse heat conduction in the fluid passing through the subchannels shown in Fig. 2-a becomes

$$q''_{ij} = k_{ij} \frac{\bar{T}_i - \bar{T}_j}{l_{ij}} \quad [1.a]$$

* This model was first introduced in the Italian subchannel, code (CISE (23)). It is especially preferred in modeling the strict annual two-phase flow condition, due to its resemblance to the annual geometry.

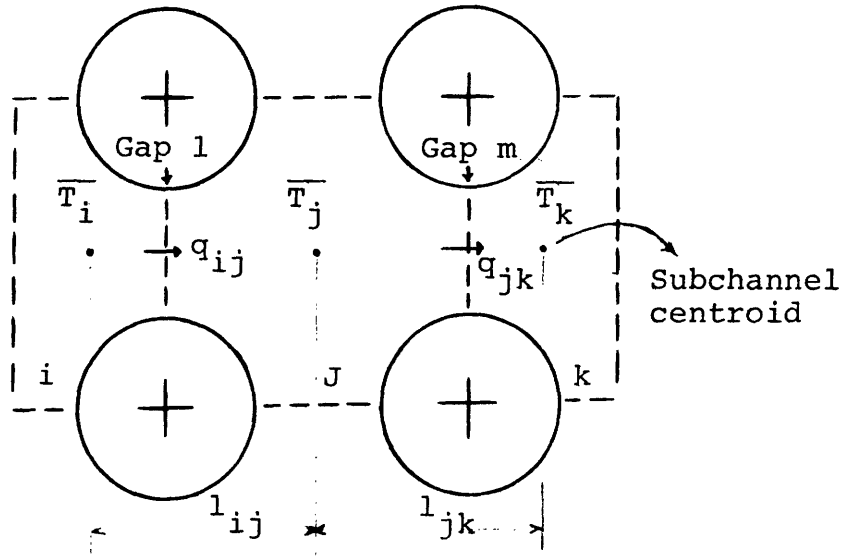
** An excellent discussion concerning the fine structure of the flow field within the coolant region is presented in Ref. (24).

and

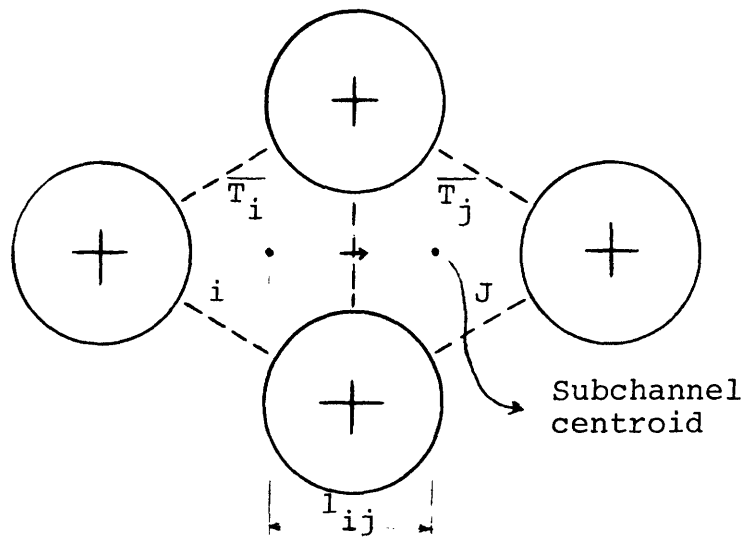
$$q''_{jk} = k_{jk} \frac{\bar{T}_j - \bar{T}_j}{l_{jk}} \quad [1.b]$$

where q'' , k , \bar{T} and l are heat flux, thermal conductivity, averaged temperature and finally centroid-to-centroid distance between the adjacent subchannels respectively. Assuming identical fuel rods, the centroid located averaged subchannel temperature seems to be a valid assumption for subchannel j . However, for subchannels i and k , it is expected that the averaged temperatures are located closer to the gap l and gap m respectively. This is also the case for the temperatures shown in Fig. 2-b. The centroid located averaged temperature is a valid assumption for low conductivity coolants and high $\frac{P}{D}$ ratios, whereas, it is a dramatic assumption for high conductivity coolants and tight rod bundles.*

* A discussion in detail and a suggested method to correct the centroid located averaged values are presented in reference 25.



(a)



(b)

Fig. 2 Lateral heat condition.

2.1.2 Loop Codes

Analysis of the whole primary system during transient conditions and hypothetical accidents such as loss of coolant, pump failure and nuclear excursions, necessitates modeling the whole loop components such as pipes, pressurizer (in PWR's), pumps, steam generator, jet pumps (in BWR's), valves and reactor vessel. Also, the effects of the secondary system need to be considered.

The thermal hydraulic behavior of the reactor core during the course of a transient is tied to the core nuclear characteristics through the reactor kinetics. Hence, the reactivity feedback should be considered in the process of the primary loop modeling.

The RELAP series of computer programs are the well-known transient loop codes which have been extensively used in the nuclear industry. These codes are basically devised to analyze transients and hypothetical accidents in the nuclear reactor loop of LWR's and mainly consist of four major parts as follows:

- (1) a thermal hydraulic loop part,
- (2) a thermal hydraulic core part,
- (3) a heat conduction part,
- (4) a nuclear part.

In these codes, the primary system is divided into volumes and junctions. The fluid volumes serve as control volumes, describe plenums, reactor core, pressurizer, pumps and heat exchangers. Each connection between volumes may be specified as a normal junction, a leak or a fill junction. A fill junction as its name implies, injects water into a well-specified volume. By definition, volumes specify a region of fluid within

a given set of fixed boundaries, whereas junctions are the common flow areas of connected volumes.⁽¹⁰⁾ Any fluid volumes may be associated with a heat source or a heat sink, such as fuel rods or the secondary side of a heat exchanger, respectively.

While RELAP2 is able to handle only three control volumes with a fixed set of pipes connecting these volumes, representing the whole primary loop, RELAP3B and RELAP4 are capable of handling as many as 75 volumes and 100 junctions or even more, at the expense of more computer core.*

2.2 Classification According to Two-Phase Model

2.2.1 Homogeneous Equilibrium Model

Flow characteristics in component and loop codes are calculated through solving the field or conservation equations written for the well specified control volumes. The basic assumption made in modeling the two-phase flow is representing the two-phase by a pseudo single phase. This method of modeling is also known as homogeneous equilibrium model (HEM). The HEM is extensively used in the thermal hydraulic codes. The homogeneous assumption implies that the phase velocities are equal and flow in the same direction, also the phase distribution is uniform throughout the control volumes. The equilibrium assumption requires the phase to be at the same pressure and temperature.

The one dimensional HEM codes use an approximate set of

*To do this, only the array sizes in the COMMON blocks should be increased.

field or conservation equations for the mixture in conjunction with the constitutive relations. The differential form of the conservation equations written for a mixture is as follows: (26)

Local mixture continuity equation:

$$\frac{\partial \rho_m}{\partial t} + \vec{\nabla} \cdot (\rho_m \vec{V}_m) = 0 \quad [1]$$

Local mixture momentum equations:

$$\frac{\partial}{\partial t} \rho_m \vec{V}_m + [\vec{\nabla} \cdot \rho_m (\vec{V}_m \vec{V}_m)] + \vec{\nabla} \cdot \bar{\bar{T}}_m - \rho_m \vec{g} = 0 \quad [2]$$

where the product $\vec{V} \vec{V}$ gives an array of nine components.

This product can be written as

$$\vec{V}_m \vec{V}_m = (V_i)_m (V_k)_m \quad (i, k = 1, 2, 3,)$$

The surface stress tensor, $\bar{\bar{T}}_m$, is made up of the pressure and the normal and the shear stresses

$$\bar{\bar{T}}_m = P_m \bar{\bar{I}} - \tau_m$$

where $\bar{\bar{T}}_m$ is the viscous stress tensor and $\bar{\bar{I}}$ is a unit tensor.

Local mixture energy equation:

$$\begin{aligned} \frac{\partial \rho_m}{\partial t} (U_m + 1/2 \vec{V}_m \cdot \vec{V}_m) + [\vec{\nabla} \cdot \rho_m (U_m + 1/2 \vec{V}_m \cdot \vec{V}_m) \vec{V}_m] = & [3] \\ -[\vec{\nabla} \cdot (\vec{q}_m - [\bar{\bar{T}} \cdot \vec{V}])] + \rho_m \vec{g} \cdot \vec{V}_m + \dot{Q}_m \end{aligned}$$

where \vec{q}_m is heat flux, \dot{Q}_m is the body heating term and U_m is the internal energy.

These balance equations need to be accompanied by the constitutive equations for τ_m , \vec{q}_m , and \dot{Q}_m , the equation of state, and the mixture properties.

2.2.1.1 Approximation to the field equations -
Component Codes

Approximations which are made in solving the conservation equations in the component codes using the homogenous equilibrium model for the two-phase flow will be discussed in this section.

Basically, none of the existing subchannel codes use such a generalized three dimensional set of field equations as are given by Equations 1, 2 and 3. Rather simplifying assumptions are made in these equations. For example, in most of the COBRA versions, flow is assumed to have a predominantly axial direction and all the "lateral" flow is lumped into one lateral momentum equation. The reason for such treatment may be justified by considering the non-orthogonal characteristics of subchannel arrangement (Fig. 3) which do not allow treatment of the lateral or transverse momentum equations as rigorously as the axial momentum equation. It is assumed that the interaction between two adjacent subchannels in the transverse direction is through two distinct processes,* namely, diversion cross-flow and turbulent mixing. Axial turbulent mixing between nodes is ignored.

The first process, diversion cross-flow is assumed to exist due to local transverse pressure difference in the adjacent subchannels. Such a process transfers mass, momentum

*A more general classification is given in Ref. (27) and is referenced in the model making process of WOSUB (17). Also see Ref. (47) for basic notation in subchannel analysis.

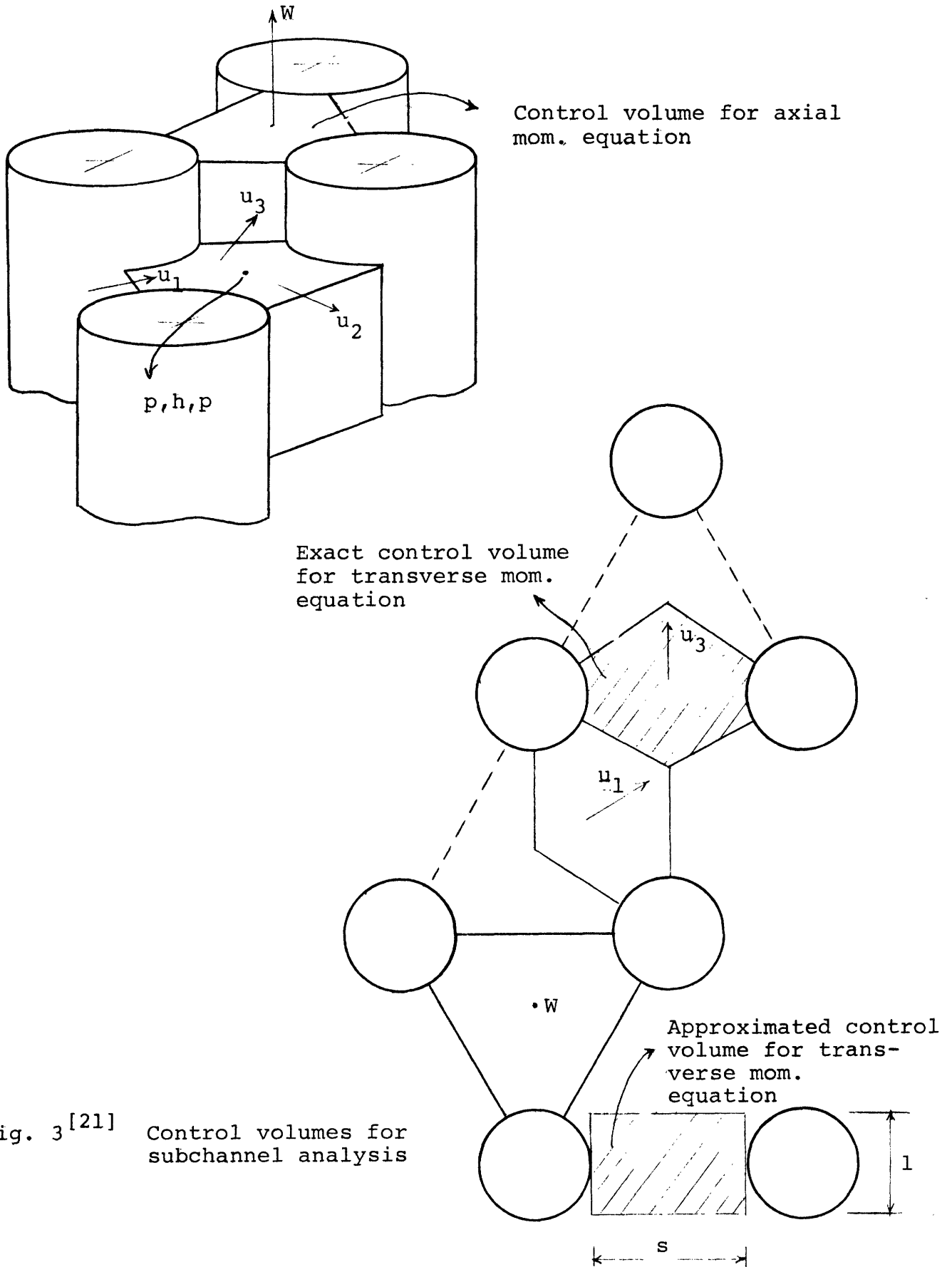


Fig. 3^[21] Control volumes for subchannel analysis

and energy with the assumption that the cross flow loses its sense of direction when it enters a subchannel⁽⁴⁾. Unlike the HEM versions of COBRA, WOSUB which is essentially devised for analysis of ATWS in BWRs does not account for diversion cross flow.

The second process, turbulent mixing is assumed to be caused by both pressure and flow fluctuation. In this process, no net mass transfers, only energy and momentum are involved. This is due to the assumption of the equi-mass model.* The magnitude of the turbulent mixing term is determined either by some correlations or by a physical model that includes empirical constant.

All the COBRA versions account for a single phase turbulent mixing while the two phase turbulent mixing term was added in the versions following COBRA-II, since COBRA-I does not account for this term.

It should be mentioned that forced flow mixing which is caused by some rod spacing methods such as a wire wrap or diverter vanes is taken into account, especially in those codes which are capable of analyzing fluid flow in LMFBRs such as COBRA-IIIC and COBRA-IV-I. Recently, a wire wrap model has been added⁽²⁸⁾ to COBRA-III-P which makes it capable of handling LMFBR flow analysis.

*The equi-volume model which is based on the change of same volume of flow is used in the MIXER code. For further detail see Ref. (22).

The steady-state versions of COBRA, namely, COBRA-I, COBRA-II and COBRA-III do not have any model for forced diversion cross flow.

A more complete form of transverse momentum equation is employed in COBRA-IIIC, COBRA-IIIP and COBRA-IV-I which includes the time and space acceleration of the diversion cross flow.

As a correction to the homogenous flow assumption, a one dimensional slip flow model which accounts for nonequal phase velocities, is considered in all the COBRA series up to and including COBRA-IV-I. A subcooled void calculation is also added to these codes. However, COBRA-I and the explicit scheme of COBRA-IV-I (to be described) do not have a subcooled void option.

In the COBRA codes, the energy equation has been further simplified by assuming the turbulent mixing and convection heat transfer as the unique mechanisms for internal energy exchange. In such treatment, it is assumed that⁽²⁹⁾

- no heat is generated within the fluid,
- changes in kinetic energy is small,
- no work against the gravity field.

Neglecting the time change of local pressure, $\frac{\partial P}{\partial t}$, limits these codes to transients with times that are longer than the sonic propagation time through the channel.⁽⁴⁾

Unlike the previous versions, the COBRA-IV-I momentum equations account for the momentum flux term.

Further simplifications to the axial momentum equation have been made by neglecting surface tension contribution. This requires equal phase pressures. This basic assumption in addition to the assumed equal phase temperatures are the result of the thermodynamic equilibrium assumption.

2.2.1.2 Approximations to the field equations -- Loop Codes

Assumptions made to solve the field equations in the loop codes using HEM are discussed here. Except for RELAP5, which is the latest publicly available version of RELAP series, the remaining versions use the HEM for their hydrodynamic modeling. Therefore, a set of conservation equations written for a mixture (Equations 1, 2, & 3) is applicable for theoretical considerations. For the practical purposes, approximations have been made to this generalized set. The RELAP codes, generally have a lumped parameter structure in which the spatial effects are integrated over the control volume for the conservations of mass and energy. For example, the mass balance in its differential form is

$$-\frac{\partial \rho}{\partial t} = \nabla \cdot (\rho \nabla) \quad [4]$$

Integrating over the control volume

$$-\iiint_V \frac{\partial \rho}{\partial t} d\tau = \iiint_V \nabla \cdot (\rho \vec{V}) d\tau \quad [5]$$

Now applying the divergence theorem to the right hand side of Equation [5], we get:

$$- \iiint_V \frac{\partial \rho}{\partial t} d\tau = \iiint_V \nabla \cdot (\rho \vec{V}) d\tau = \iint_S \rho \vec{V} \cdot \vec{n} ds$$

or

$$\iint_S \vec{n} \cdot (\rho \vec{V}) ds + \iiint_V \frac{\partial \rho}{\partial t} d\tau = 0 \quad [6]$$

Using M_j as the existing mass in the volume J at time t_i and considering the term $-\vec{n}_i \cdot (\rho S \vec{V})_i = W_{ij}$ which is equal to the inflow from side "i" into volume J (Fig. 4), the mass balance reduces to

$$\frac{d}{dt} M_j = \sum_{i=1}^n W_{ij} \quad [7]$$

Similarly, a simplified form of the energy equation which has been used in RELAP2 and various MOD's of RELAP3 is as follows:

$$\frac{dU_j}{dt} = \sum_{i=1}^n W_{ij} h_{ij} + Q_j \quad [8]$$

where U_j is the internal energy of volume J , h_{ij} is the enthalpy of fluid flowing from side i into volume j and finally Q_j is the heat input to volume j .

The effect of kinetic, potential and frictional energies are neglected in Equation [8]. However, the RELAP4 energy equation accounts for kinetic and potential energy changes.

Unlike the mass and energy equations, the momentum equation is written for a shifted control volume as shown in Fig. 4. This method minimizes the extrapolation of boundary conditions. The final form of the momentum equation used in various MOD's of RELAP3 is as follows:

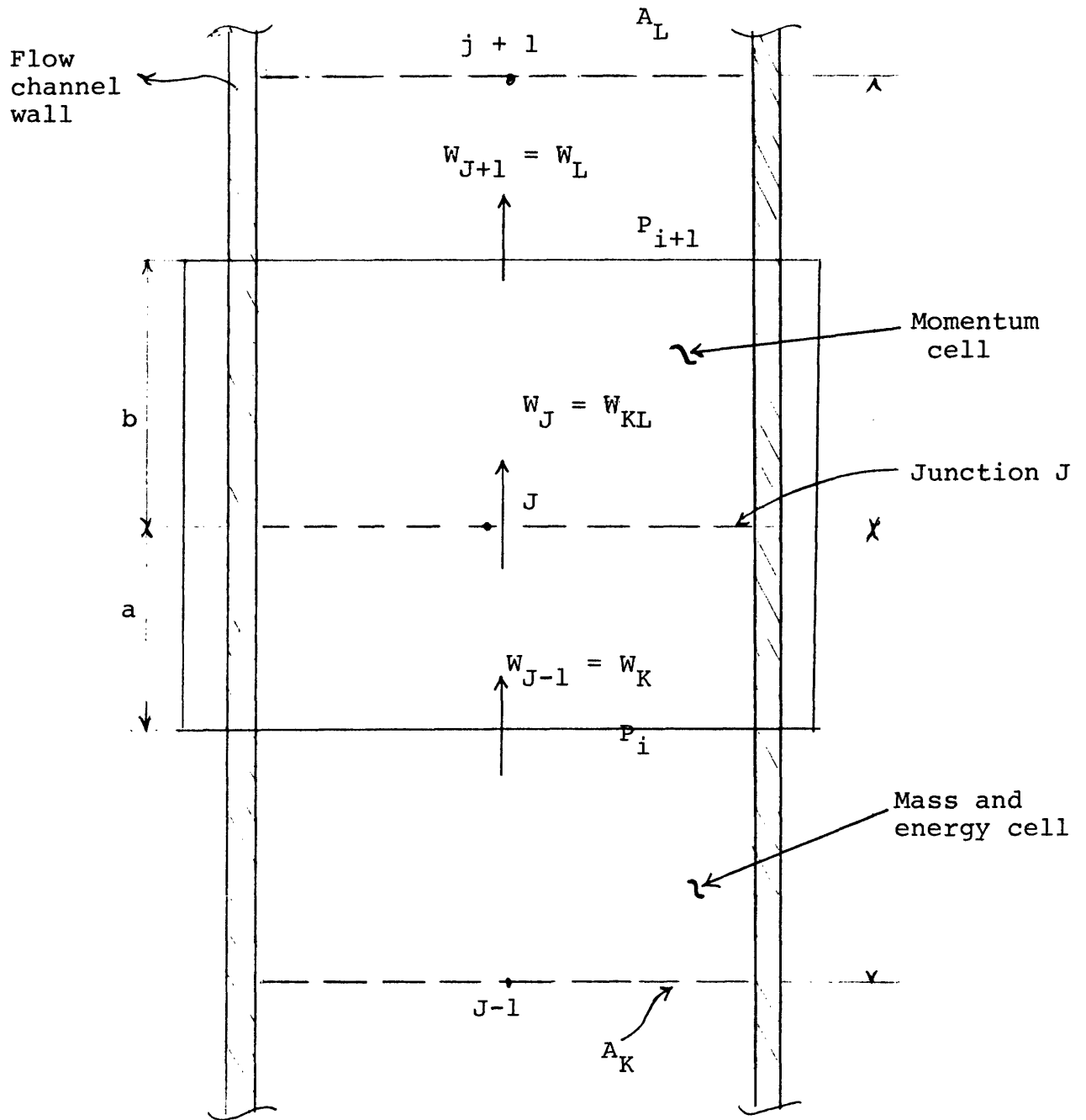


Fig. 4 Geometry for mass, momentum and energy equations.

$$\frac{1}{144g_c} \left(\frac{L}{A}\right) \frac{dW_j}{dt} = P_i - P_{i+1} + \Delta P_P + \int_{V_j} \frac{Pdz}{144} - \frac{k_j W_j |W_j|}{\rho_j} \quad [9]$$

However, like the energy equation, an improved form of the momentum equation is implemented in RELAP4 which takes the form: (30)

$$\begin{aligned} \underbrace{I_j \frac{dW_j}{dt}}_{\leftarrow 1 \rightarrow} &= \underbrace{(P_k + P_{kgj})}_{\leftarrow 2 \rightarrow} - \underbrace{(P_L + P_{Lgj})}_{\leftarrow 4 \rightarrow} - \underbrace{F_{fK} - F_{fL} - F_{fr}}_{\leftarrow 6 \rightarrow} \\ &- \underbrace{\int_{K_0}^{L_i} dF}_{\leftarrow 7 \rightarrow} - \underbrace{\int_{K_1}^{L_i} dP}_{\leftarrow 8 \rightarrow} - \underbrace{\int_{K_0}^{L_i} \frac{d(vW)}{A}}_{\leftarrow 9 \rightarrow} \end{aligned} \quad [10]$$

It is assumed that the junction inertia term, $I_j \frac{dW_j}{dt}$ in equation [10] or the corresponding term $\left(\frac{L}{A}\right) \frac{dW_i}{dt}$ in equation [9] represents the rate of change of momentum everywhere in the selected control volume J. (Fig. 5) In equation 10, I_j is the geometric "inertia" for the flow path and also,

W_j = flow rate in junction J,

P_k = Pressure in volume K,

P_{kgj} = gravity head contribution for volume K,

F = friction terms,

v = velocity,

A = flow area.

The significance of each term in equation [10] is as follows:

Term 1 represents the rate of change of momentum,

Term 2 and 4 represent the pressure drop between two volumes,

Terms 3 and 5 represent gravity,

Terms 7 and 8 represent the friction and pressure drop associated with expansion and contraction.

Term 6 represents the fanning friction terms.

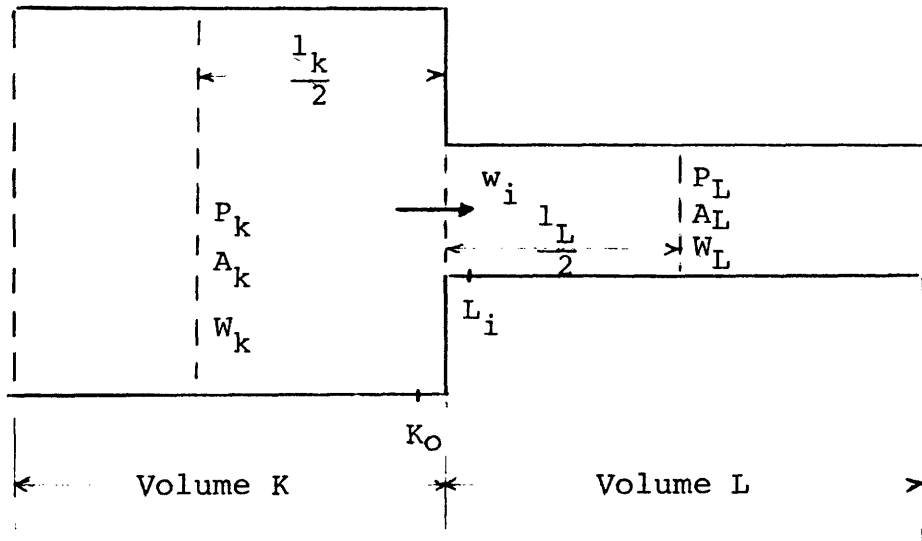


Fig. 5 Flow path control volume.

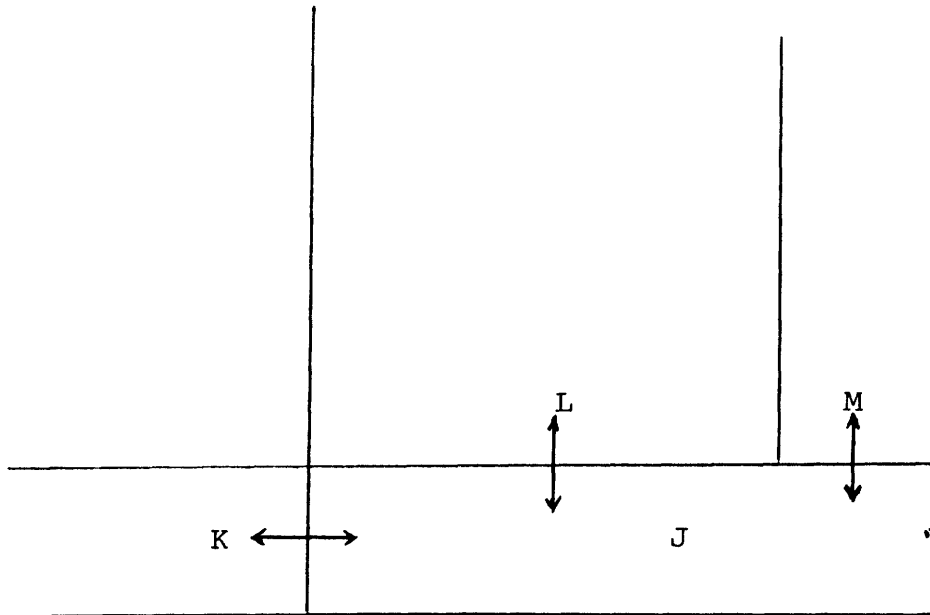


Fig. 6^[26] Lumped parameter simplification of multidimensional flow regimes.

Term 9 represents the momentum flux or spatial acceleration term. Comparing Equations 9, used in RELAP3, and 10, used in RELAP4, it is obvious that the major difference is inclusion of the momentum flux term in Equation 10. The momentum flux term for a homogeneous volume becomes

$$\int_k^L \frac{d(VW)}{A} = \frac{V_L W_L - W_k V_k}{A}$$

Although the inclusion of the momentum flux term has improved the momentum equation, there are some cases in the loop modeling that the code user has to ignore this term. (For example, when the given control volume, J, connects into more than two control volumes, as illustrated in Fig. 6; where the double-ended arrows indicate junction). A similar case might be encountered in the lower plenum modeling when the flow is highly three dimensional. In such cases, it is quite difficult to define the control volume boundaries for the momentum balance at each junction. These examples clarify the inability of a lumped parameter approach to model multidimensional regions. This is also the case in calculating the friction factor and heat transfer coefficient which by definition are⁽⁴⁸⁾

$$f_w = \int (-\mu \frac{du}{dr} |_w) dz \quad [11]$$

and

$$q'' = -k \frac{dT}{dr} |_w = h(T_w - T_b) \quad [12]$$

where μ and u are the viscosity and velocity of the flow and subscripts w and b represent values evaluated at wall and bulk

respectively,

It is clear from equations 11 and 12 that the derivatives are evaluated at the channel wall. However, since a lumped parameter approach doesn't account for velocity and temperature profile, therefore, the above mentioned derivatives do not make sense. It is this reason which necessitates an input specified friction factor for the codes using this approach.

The junction inertia term is another term in the momentum equation (Equations 9 and 10) which becomes rather ambiguous in modeling the complex geometries. The junction inertia arises from an approximation in the momentum equation to the temporal inertia term as follows: (30)

$$\int_{x_1}^{x_2} \frac{1}{A(x)} \frac{dw}{dt} dx \approx \frac{dw}{dt} \int_{x_1}^{x_2} \frac{dx}{A(x)} = I \frac{dw}{dt} \quad [13]$$

where x_1 = center of control volume 1 and x_2 = center of control volume 2, and I is the geometric inertia for the flow path defined as:

$$I = \int_{x_1}^{x_2} \frac{dx}{A(x)} \quad [14]$$

The geometric inertia for a homogeneous volume, Fig. 4, becomes $I = \frac{a}{2A_K} + \frac{b}{2A_L}$, however for complex geometrics, the inertia term may be determined by using a simplified assumption. The basic assumption which is introduced in this

respect⁽³⁰⁾ is that the inertia of a junction is composed of two independent contributions, one from each connecting volume. For example, Fig. 7 could represent a downcomer region. If we assume that Junction 1 communicates primarily with Junction 3, then with respect to the mentioned basic assumption, the geometric inertia will be:

$$I_j = I_{j1} + I_{j2} + I_{j3} \quad [15]$$

or

$$I_j = \frac{L_1}{2A_1} + \frac{L_2}{A_2} + \frac{L_1}{2A_3}$$

Where L_1 is the effective length of both Junctions 1 and 3 and the effective length of junction 2 is assumed to be $2L_2$.

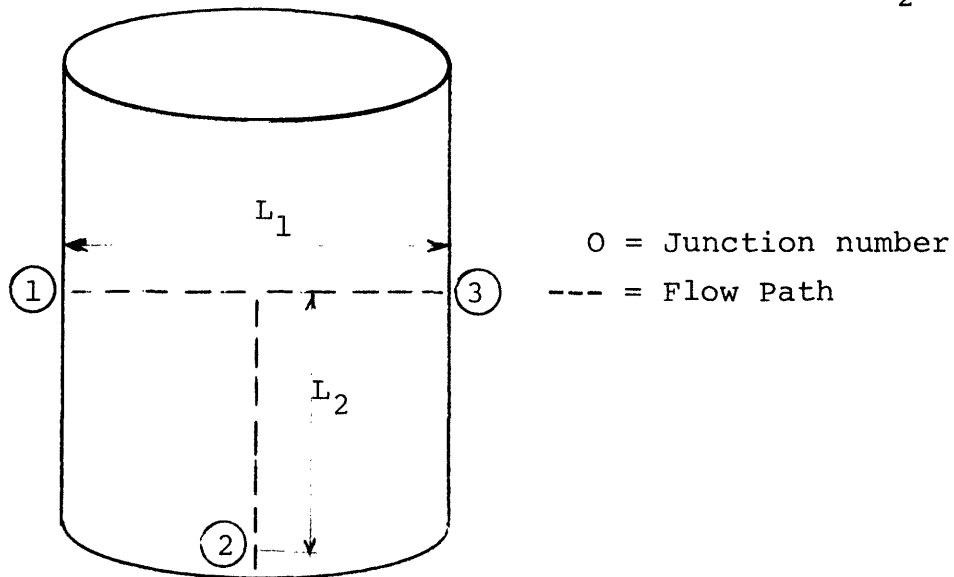


Fig. 7⁽³⁰⁾ - A Downcomer Representation in RELAP4

A schematic of RELAP4 model of a PWR is presented in Fig. 8. It illustrates the complexity of accounting for all volumes and junctions in a LWR plant.

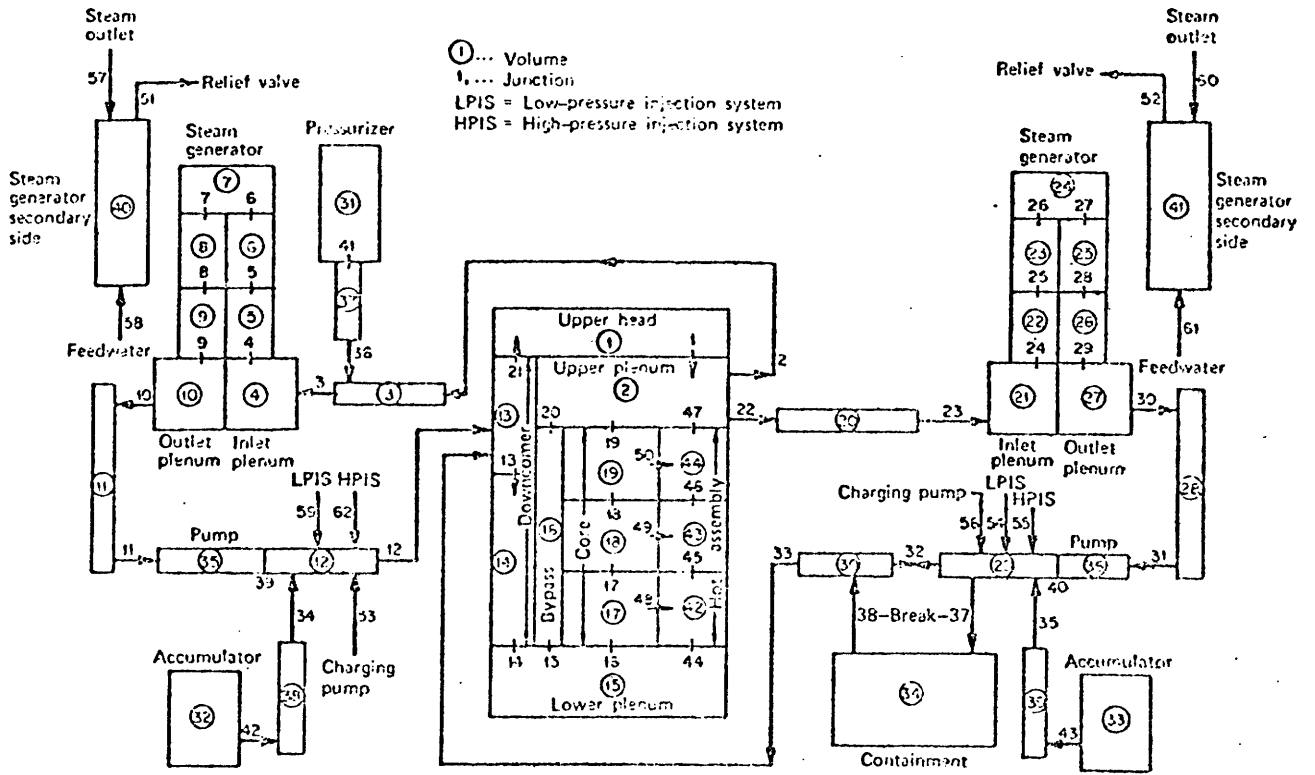


Fig. 8 [30] Schematic of RELAP4 model of a large PWR.

2.2.2 Improvements on the Homogeneous Equilibrium Model

By retaining more field equations, a more realistic approach to analysis of severe transients has become possible in recent codes. The increased number of field equations enable a code to analyse transients in which the situation is far beyond the capability of the rigid assumption of equal phase velocities and temperatures. In this respect, countercurrent two-phase flow and vapor-liquid phase separation during small break transients and emergency core coolant delivery are notable examples.

Since in a non-homogenous flow slip exists between the two phases, there is a relative motion of one phase with respect to the other. This relative motion arises due to density and/or viscosity differences between phases where usually the less dense phase will flow at a higher local velocity than the more dense phase, except for the gravity dominated flow⁽¹⁸⁾. The general effect of slip is to lower the void fraction below the homogeneous value.

The slip or hold up ratio, $s = \frac{v_g}{v_L}$, should not be confused with the slip velocity $v_{sL} = v_L - v_g$, or drift velocity, a concept which is used in the drift flux model.

Unlike the one-dimensional HEM, a non-homogeneous, one dimensional flow calculation for a two-phase flow in thermodynamic equilibrium, involves the solution of one

equation of state and five differential equations: a mixture energy equation and one continuity and one momentum equations for each phase as it is done in RELAP5.

2.2.2.1 Dynamic Slip Model

Simplified assumptions have been made to reduce the number of conservation equations while retaining the improvements over HEM codes. This is done in WOSUB and COBRA-DF by using the concept of diffusion or drift flux model, and in RETRAN by introducing the dynamic slip model.

RETRAN computer code is basically developed from the RELAP series of codes. It is a one-dimensional code which solves four field equations written for a fluid volume as follows: Mixture continuity, Momentum, Energy equation, and time dependent behaviour of the velocity difference, obtained by subtracting the momentum equations written for each phase. This additional momentum equation reads:

$$\frac{\partial V_{sl}}{\partial t} - V_g \frac{\partial V_g}{\partial X} + V_l \frac{\partial V_l}{\partial X} + \left(\frac{1}{\rho_l} + \frac{1}{\rho_g} \right) \frac{\alpha P}{\alpha X} + \left(\frac{1}{\alpha_l \rho_l} + \frac{1}{\alpha_g \rho_g} \right) \overline{A_{gL}} B_{gL} V_{sL} = 0 \quad [16]$$

where V, ρ, α, P represent velocity, density, void fraction and pressure respectively. Also $\overline{A_{gL}}$ represents the surface area between vapor and liquid phase per unit volume and B_{gL} represents the friction coefficient between vapor and liquid

phases.

In deriving equation (16), the following assumptions have been made:

- 1) The wall friction is nearly equal for the two-phase.
- 2) The momentum exchange between phases due to mass exchange is small.

In addition to inclusion equation (16) in RETRAN, some improvements have been made in the field equations used in RELAP4, as follows:

- 1) Additional term in the mixture momentum equation with respect to the momentum flux. Mixture momentum flux:

$$\frac{\partial}{\partial x} [A (\alpha_g \rho_g (V^2)_g) + \alpha_l \rho_l (V^2)_l] =$$

$$\frac{\partial}{\partial t} (vw) + \frac{\partial}{\partial x} \left(\frac{v_{sL}^2 \alpha_l \alpha_g \rho_l \rho_g A}{\rho} \right)$$

← ----- →
Additional Term

- 2) Additional term in energy equation which accounts for the time rate of change of kinetic energy,

$$\frac{\partial}{\partial t} \left[\rho A \left(\frac{U^2}{2} \right) \right].$$

3) Using a flow regime dependent two-phase flow friction multiplier.

2.2.2.2 Drift Flux Model

Unlike RETRAN, COBRA-DF which is a vessel code and uses the drift flux model, employs five field equations to determine phase enthalpy, density and velocity*. This code is used exclusively for examination of upper heat injection of water during a LOCA in a PWR.

Vapor diffusion or drift flux model is another step toward modeling a non-homogeneous non-equilibrium flow. The basic concept in this model is to consider the mixture of the two-phase as a whole, rather than treating each phase separately. The DFM is more appropriate for the mixture where dynamics of two components are closely coupled, however, it is still adequate where the relatively large axial dimension of the systems gives sufficient interaction time (26).

In this model in addition to the three field equations written for the mixture, there is a diffusion

* THOR which is developed at BNL (31) uses the DFM and accounts for thermal non-equilibrium of the dispersed phase only.

equation written for the dispersed phase which reads:

$$\frac{\partial \rho_g \alpha_g}{\partial t} + \vec{\nabla} \cdot (\alpha_g \rho_g \vec{V}_m) = \Gamma_g - \vec{\nabla} \cdot (\alpha_g \rho_g \vec{V}_{gj}) \quad [17]$$

where Γ_g is the phase change mass generation, \vec{V}_{gj} and \vec{V}_m are the drift velocity and mixture velocity respectively.

WOSUB which is a BWR rod bundle computer code uses the DFM and solves four field equations written for a subchannel control volume, as follows:

- 1) continuity equation for mixture
- 2) continuity equation for vapor. This equation reads:

$$A \frac{\partial}{\partial t} (\rho_g \alpha_i)_i + A \frac{\partial}{\partial z} (p_g J_g)_i = A \rho_{gi} \Psi_i + \rho_{gi} q_{gi} \quad [18]$$

where J_g = vapor flux

q_{gi} = vapor volume flow to subchannel i

Ψ_i = vapor volume generation in subchannel i
per unit volume

Equation [18] indicates the fact that the temporal and spatial increase in the mass of vapor in subchannel i is due to vapor generation in the subchannel and vapor addition from the adjacent subchannels. The vapor volume generation term, Ψ , appears in Equation [16] due to using the DFM. This term is part of the code constitutive package. It is modeled in WOSUB based on the Bowring's equation which relates

Ψ to the heat flux:

$$\Psi = \tau \frac{P_h}{A \rho_v h_{fg}} q'' \quad [19]$$

where τ is a coefficient depending upon coolant condition, P_h is the heated perimeter and q'' is heat flux in the fully developed nucleate boiling region, and A is the flow area. The effect of subcooled boiling non-equilibrium condition is considered in the final form of Ψ .

3) Mixture axial momentum equation: This is the only momentum equation considered in the code. Therefore it is clear that WOSUB is strictly one dimensional. This may be justified by considering the fact that the intention of creating WOSUB, has been analysing the flow characteristics in encapsuled PWR bundles as well as BWR bundle geometry⁽¹⁷⁾ in which, based on a channelwise node, the flow is predominately one dimensional. Nevertheless, the transverse effects are not totally forgotten. In fact a natural turbulence exchange mechanism is considered. Furthermore, vapor diffusion accounts for the tendency of diffusion vapor in the higher velocity regions.

These effects are considered in the momentum equation which is given by⁽¹⁷⁾:

$$\frac{\partial P}{\partial Z} = \left(\frac{\partial P}{\partial Z}\right)_{el} + \left(\frac{\partial P}{\partial Z}\right)_{ac} + \left(\frac{\partial P}{\partial Z}\right)_{fr} + \frac{\partial G}{\partial t} - \left[\xi + \left(\frac{\partial P}{\partial Z}\right)_{td}\right] \quad [20]$$

The last two terms stand for the axial momentum transferred into subchannel i and the turbulent shear stress, respectively. It is also evident that these two terms which connect the subchannel to its neighboring subchannels stem only from flow and pressure fluctuation and not transverse pressure difference as was discussed in Section 2.2.1.1.

4) Mixture energy equation which contains the inflow of enthalpy from adjacent subchannels.

Generally, the dynamic slip model, as it is used in RETRAN, has advantages over both slip ratio correlations, as used in most versions of COBRA, and DFM, as used in WOSUB, as follows⁽¹⁸⁾:

- 1) The slip correlations are based on steady-state data whereas the application is for transients.
- 2) They highly rely on empiricism which may eliminate many mechanistic effects.

3) The slip velocity $V_{sL} = V_g - V_L$ can only assume positive values, hence the possibility of rising liquid and falling vapor cannot be predicted.

2.2.3 The Two-Fluid Model

The inability of the simplified methods to treat the multidimensional, non-equilibrium separated and dispersed flows necessitates a better modeling of the two-phase flow. Anticipated reactor transients and postulated accidents like LOCA specially require a more realistic treatment.

Those cases in which a one-dimensional HEM is not acceptable are tabulated in Table 2-1.

TABLE 2-1⁽³²⁾
CASES WHERE 1-D HEM IS NOT
ACCEPTABLE

<u>Multidimensional Effects</u>	<u>Non-Equilibrium Effects</u>	<u>Phase Separation</u>
Downcomer region	ECC injection	Small breaks
Break flow entrance	Subcooled boiling	Steam generator
Plena	Post-CHF transfer	Horizontal pipe flow
Steam separators	ECC heat transfer	Counter current flow
Steam generators	Low-quality blowdown	PWR ECC bypass
Reactor core	Reflood quench front	BWR CCFL

The most flexible approach in modeling these cases is through using a two-fluid, full non-equilibrium concept, which is the most sophisticated model employed so far in treating the two-phase flow.

The derivation of the field equations in their general tensor form is quite involved. A detailed derivation is presented in Ref. 33. A short-hand representation for the two-fluid model is 2V2T or UVUT which stands for unequal phase velocities and temperatures - whereas 1V1T or EVET is used for HEM.

The unknowns and equations in this model are summarized in Table 2-2.

Table 2-2

Two Phase and Single Phase Comparison
with Respect to the Flow
Equations

Case	Unknowns	# of Unkn.	Type of Equations	# of Equ.
Single	\bar{V}	3	Conservation of Mass	1
Phase	P	1	" of Momentum	3
Flow	T	1	" of Energy	1
	ρ	1	Equation of State	1
		6		6
	α (void fra.)	1	Liquid Balance Equ.	1
	\vec{V}_g	3	Vapor " "	1
Two*	\vec{V}_l	3	Liquid Mom. "	3
Phase	ρ_l	1	Vapor " "	3
Flow	ρ_g	1	Liquid Energy	1
	P	1	Vapor " "	1
	T_l	1	Equation of State	2
	T_g	1	in Each Phase	
		12		12

* Table 2-3 gives a more detailed description of various approaches to modeling the two phase flow.

The two-fluid concept is employed in the advanced thermal hydraulic codes such as COBRA-TF, from BNWL, TRAC, KACHINA*, SOLA-FLX, SOLA-DF from LASL and finally THERMIT, which is developed at MIT under EPRI sponsorship.

TRAC is the state-of-the art primary loop analysis code. It employs a three-dimensional 2V2T model for the vessel and a one-dimensional drift flux model for the rest of the primary loop. The reactivity feedback is accounted for through coupling the point kinetic equations to the thermal hydraulic model. The same concept of volume and junction defined for RELAP series is used in TRAC as well. A cylindrical coordinate system is used in TRAC for modeling the three-dimensional reactor vessel. This doesn't satisfy the purpose of a common reactor core analysis with its square array pattern governed by the bundle design. THERMIT which is a vessel code, is basically the cartesian version of TRAC. Hence, the same field and constitutive equations used in TRAC is employed in THERMIT as well. A core or a fuel pin analysis in THERMIT essentially is based on treating a whole bundle cross-section as one node where the local details have been smeared throughout the cross-section. Therefore, neither TRAC nor THERMIT account for a turbulent mixing process. Devising a subchannelwise version for THERMIT using a coolant centered control volume

* K-FIX and K-TIF are two versions of KACHINA developed at LASL. The first stands for fully Implicit Exchange numerics and the second for Two Incompressible Fields.

has been initiated at MIT. This version will include a turbulent mixing model.

A summary of the aforementioned two-phase flow models is present in Table II-3. The notations and a description for specifications used in this Table are as follows:

- a) A partial non-equilibrium model, T_k T_{sat} , assumes one phase is at saturation, temperature of the other (k) phase computed.
- b) The notations τ, q, Γ, M and E stand for viscous stress, conduction heat transfer, interphase mass, momentum and energy exchange respectively.
- c) The notations: $V_r, V_G - V_m, V_G - J$ stand for relative velocity, diffusional velocity and the drift term respectively.

A glance at this table shows clearly that although the 2V2T model imposes no restriction on the flow condition such as velocity or enthalpy, however it contains the largest number of constitutive equations and it seems that the empiricism which enters in these equations is introduced at a more basic level than the less complicated models such as 1V1T approach.

Table 2-3 [23]
Two-Phase Flow Model Description

Model Designation	Characteristics	Restrictions		No. of Field Eqs.			No. of Inter-face Transf. Eq.	External Constitutive Eqs.		Final, Verifiable Results
		No.	Imposed on	Mass.	Mom. (vector)	Energ.		No.	Type	
1V1T	Homogeneous Equilib. (HEM)	3	$V_G=V_L, h_L$ $\frac{V_G}{V_L}$, and h_G V_G-V_L	1	1	1	0			
1VSLT	Slip, Equilibrium									
1DV1T	Drift, Equilibrium									
1VT _K ^T sat	Homogeneous, Partial Non-Equilib.	2	$V_G=V_L, h_L$ $\frac{V_G}{V_L}$ or h_G V_G-V_L	1	1	2	1			
1VST _K ^T sat	Slip, Partial Nonequilibrium									
1VDT _K ^T sat	Drift Flux (DF), Partial Non-Equilibrium			2						
2V1T	Two-Fluid, Equilibrium		h_L, h_G	1	2					

External Constitutive Eqs.		Final, Verifiable Results
No.	Type	
2		p \bar{v} α
3	$\frac{V_G}{V_L}$ V_G-V_L or V_G-J q_G, E	
4	τ \bar{q} $q_G, E, \frac{V_G}{V_L}$ Γ, V_r (or V_{Gm}, V_{Gj})	τ_L or τ_G
		P, V_L, V_G, α

Table 2-3 [23]
Two-Phase Flow Model Description
(continued)

Model Designation	Characteristics	Restrictions			No. of Field Eqs.		No. of Inter-face Transf. Eq.	External Constitutive Eqs.		Final, Verifiable Results
		No.	Imposed on	Mass.	Mom.	Energ.		No.	Type	
1V2T	Homogeneous, Full Non-Equilibrium		$V_G = V_L$		1	2		5	$\left[\begin{array}{l} \bar{\tau} \\ q_L \\ q_G \end{array} \right]$	$\left\{ p, \bar{V}, T_L, T_G, \alpha \right\}$
1VD2T	Drift, Full Non-Equilibrium	1	$V_G - V_L$	2			6	$\left[\begin{array}{l} \Gamma \\ E \\ V_r \end{array} \right]$		
2VT _k ^T sat	Two Fluid, Partial Non-Equilibrium		h_L or h_G			1	2	5	$\left[\begin{array}{l} \tau_L \\ \tau_G \\ M \end{array} \right]$	$\left\{ p, V_L, V_k, \alpha, T_L \text{ or } T_G \right\}$
2VT _k ^T sat	Two Fluid Partial Non-Equilibrium		h_L or h_G	1	2		6	$\left[\begin{array}{l} \bar{q}, \Gamma \\ \bar{q}, q_k, E \end{array} \right]$		
2V2T	Two Fluid (TF), Full Non-Equilibrium	0	None	2		2	3	7	$\left[\begin{array}{l} q_L, q_G, \Gamma, E \end{array} \right]$	$p, V_L, V_k, \alpha, T_L, T_G$

2.3 Classification According to Range of Application

2.3.1 LOCA CODES

The major task of thermal-hydraulic LOCA codes is analysis of the severe cases that are encountered by the reactor core or the primary loop during the period of a loss of coolant accident. The four phases of a LOCA, for a PWR double ended cold leg break, in order of occurrence are as follows:

- 1)*- A blowdown phase which generally lasts for 30 seconds, with 2200 psia initial pressure, and ends when ECCS starts to work.
- 2)- About sixty seconds after break initiation ECC fills the lower plenum and reaches the bottom of the core (Refill Phase).
- 3)- The refill phase is followed by a REFLOOD phase which lasts for about 150 seconds, during which the core is fully flooded and quenched by the coolant.
- 4)- Long-term cooling then follows.

It has become customary to call a code a LOCA code even if it is capable of describing only the first phase of the four aforementioned phases. At the same time two codes that are capable of handling the **blowdown**, may be entirely different with respect to their type. For example, one can be a component code whereas the other a loop code. To avoid any

* This step is divided in two periods according to Ref. 31, namely: a) Adiabatic Liquid Depressurization, b) The Blowdown Period.

confusion, a classification is necessary with respect to the code's application and type, in addition to the physical models and numerical methods. A usual way to classify the LOCA codes is by categorizing them into two groups as follows.

2.3.1.1 Evaluation Model Codes

The first group contains those codes which employ a conservative basis for their physical models. Such conservatism is mandated to satisfy the NRC acceptance criteria. These codes are called the EM-codes for Evaluation Model. They constitute the WREM package which has the capability to analyse the postulated LOCA with ECC injection in accordance with current commission acceptance criteria⁽³⁵⁾. The codes which constitute the WREM package⁽³⁶⁾ are the existing computer programs which have been modified to comply with the USNRC criteria. Most of the RELAP series of computer codes are a LOCA Licensing code such as RELAP4-MOD5 and RELAP5-MOD7. Whereas RELAP3B-MOD101 is essentially devised to analyse ATWS and RELAP4-MOD6 and RELAP5 are not based on conservative correlations. RELAP4-EM is the only version of RELAP which is specifically modified to comply with acceptance criteria.

The present EM codes comprise an assembly of codes run sequentially. Each member of the sequence is a stand-

alone code developed for some special application. With this respect, RELAP4-EM in conjunction with RELAP4-FLOOD⁽³⁶⁾ and TOODEE2⁽³⁷⁾ constitute the WREM package which perform the PWR LOCA analysis. Also a combination of RELAP4-EM and MOXY-EM⁽³⁸⁾ constitute the WREM package for a BWR LOCA analysis. The respected procedure for the above mentioned analysis are presented in Fig. 8 and 9.

2.3.1.2 Best Estimate Codes

Most of the LOCA codes lay in this category. The basic physical models used in these codes rely on the best estimate assessment rather than conservative correlations. Unlike the EM-codes, the BE-codes are mostly devised as one large system code consisting of various functions previously performed via the separated stand-alone codes. This guarantees the proper compatibility and continuity between the various calculational phases. As an example, the multi-purpose loop code TRAC can be used for the analysis of the whole phases of a LOCA namely, blowdown, refill and reflood. Unlike the EM-codes which mostly use a homogeneous equilibrium model in conjunction with a lumped parameter approach for their analysis, the best-estimate codes are much more demanding and the most recent BE-codes employ the state-of-the-art physical models. Therefore they

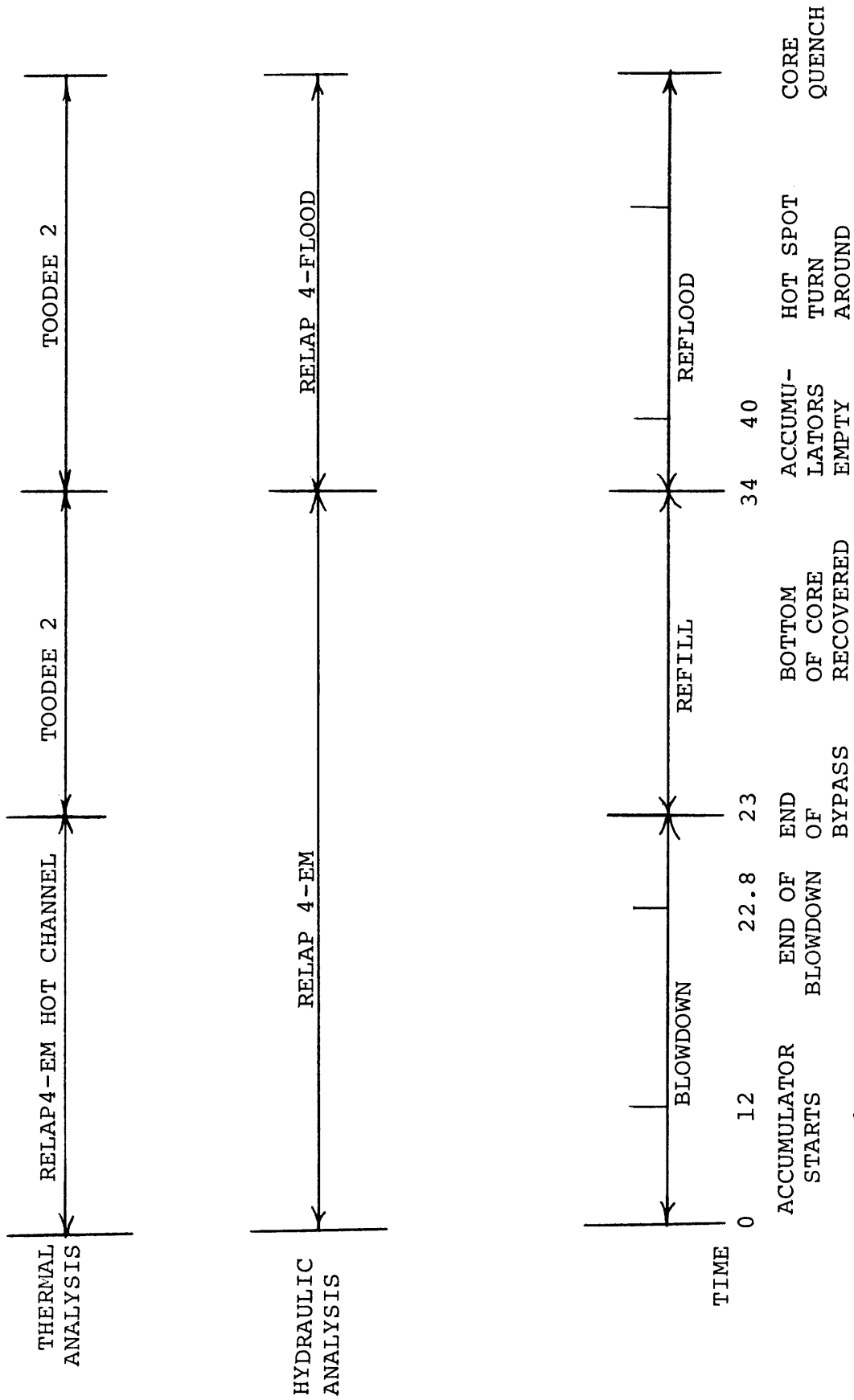


Fig.9.[15] PWR LOCA Analysis

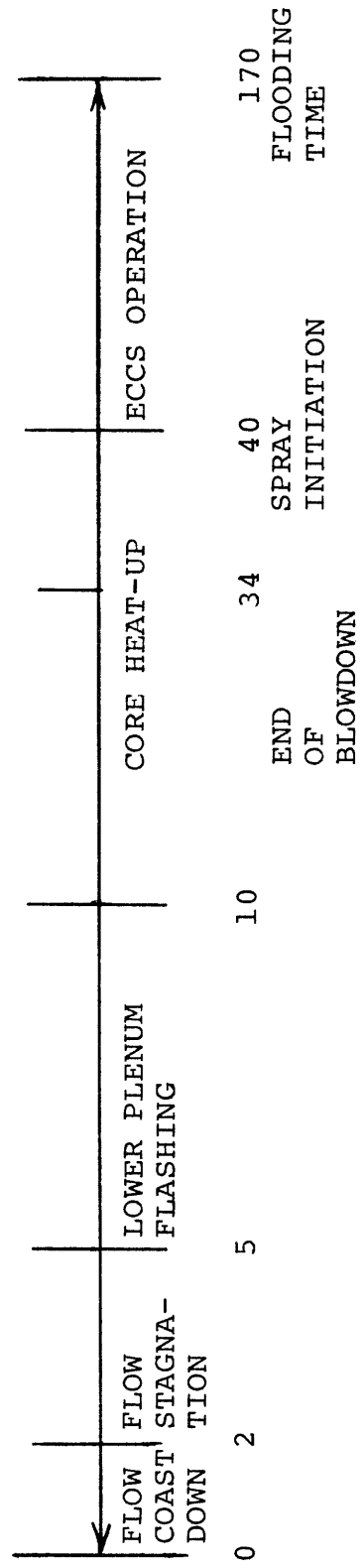
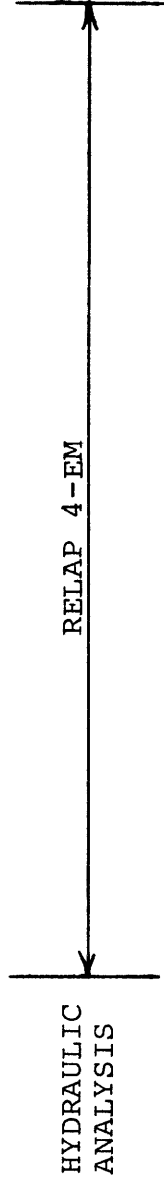


Fig. 10 [15] BWR LOCA Analysis

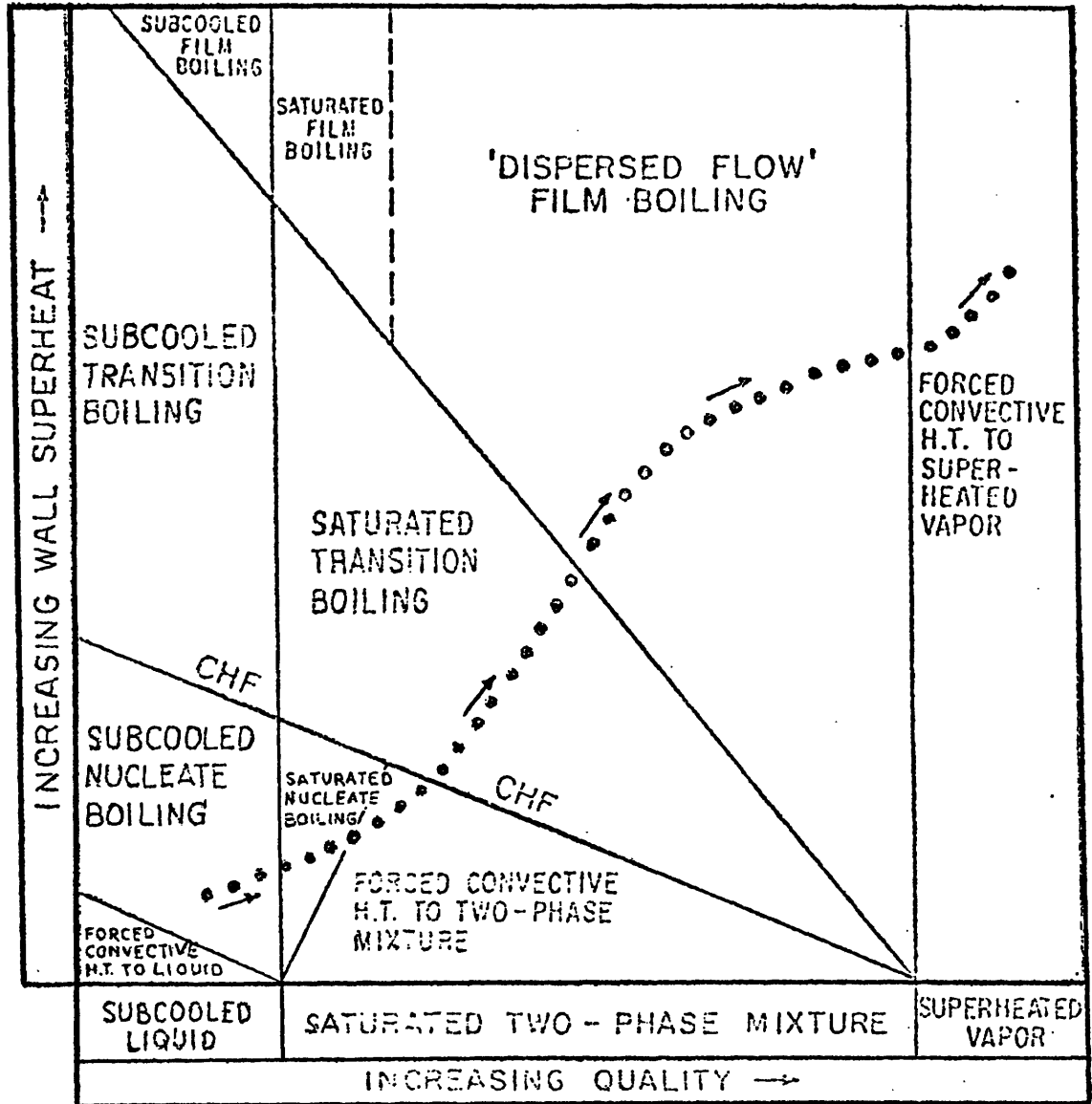
can be used to evaluate the degree of conservatism employed in the licensing (EM) calculations.

No codes have specifically been devised to handle the ATWS type of transients. In Table II-4 the causes and consequences of ATWS transients in both PWR's and BWR's are shown.

3. Two-Phase Heat Transfer Model

The energy balance of the field equations contains the contribution of the so-called wall heat transfer which accounts for the amount of heat transferred into or out of the control volume through a combination of convection and conduction heat transfer. This requires models for the wall heat transfer.

During hypothetical LOCA's, nearly all the two-phase heat transfer regimes are experienced by the coolant in the core of the NSSS, the steam generators, and the pipe of the hydraulic loop (see Fig. 11). This interdependence of the hydrodynamics and the wall heat transfer, as shown in Fig. 12, is accounted for in the thermal hydraulic LOCA codes through using a two-phase heat transfer package. These regimes are elaborated on in a pool boiling curve drawn for a fixed pressure, and shown in Fig. 13. According to Fig. 13, the path ABCDEF is obtained in a temperature



••••• APPROXIMATE ROD TRAJECTORY IN BLOWDOWN TRANSIENT

Fig. 11^[40] Heat transfer regimes traversed in blowdown.

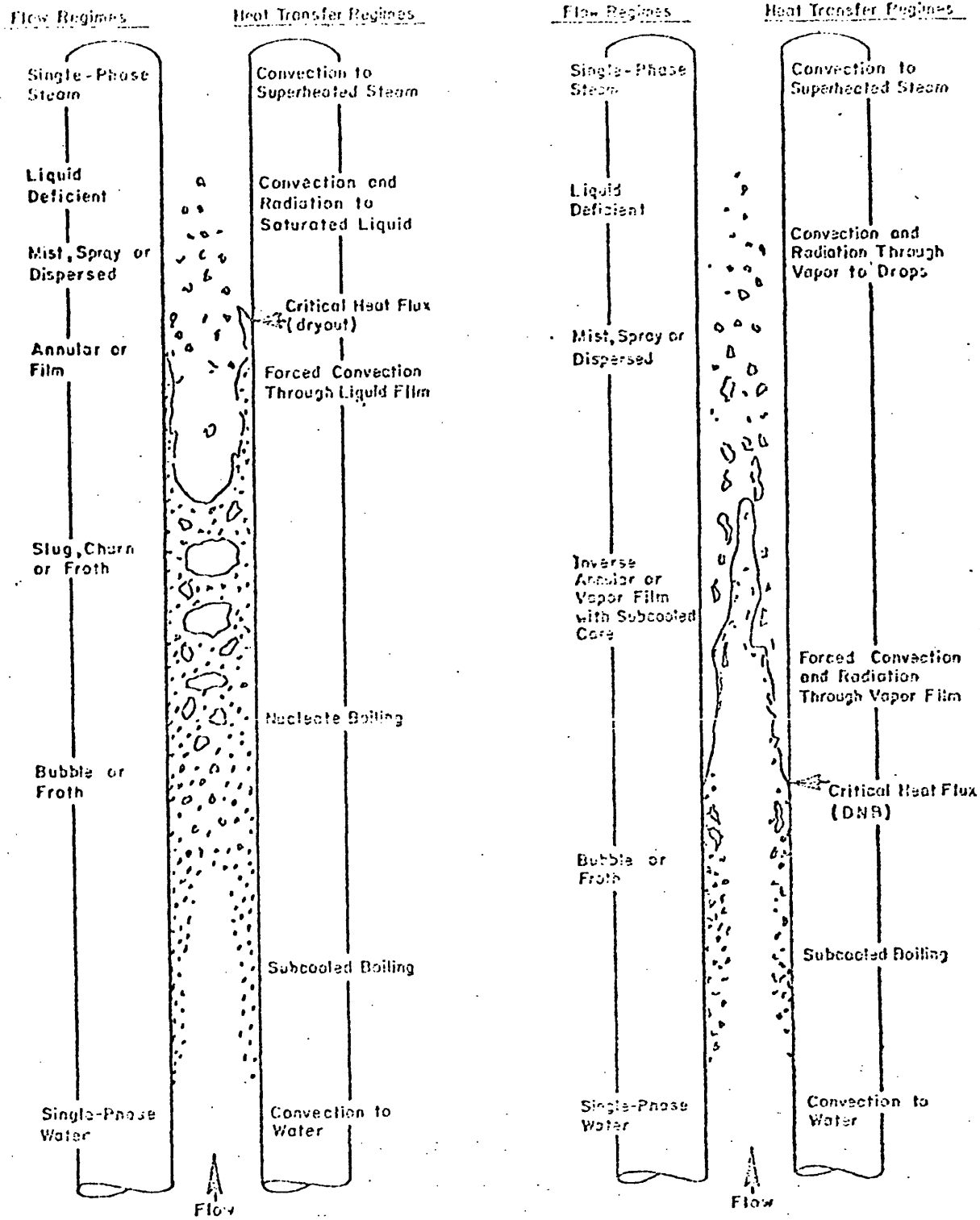


Fig. 12 [19] Flow and heat transfer regimes in rod array with vertical upflow.

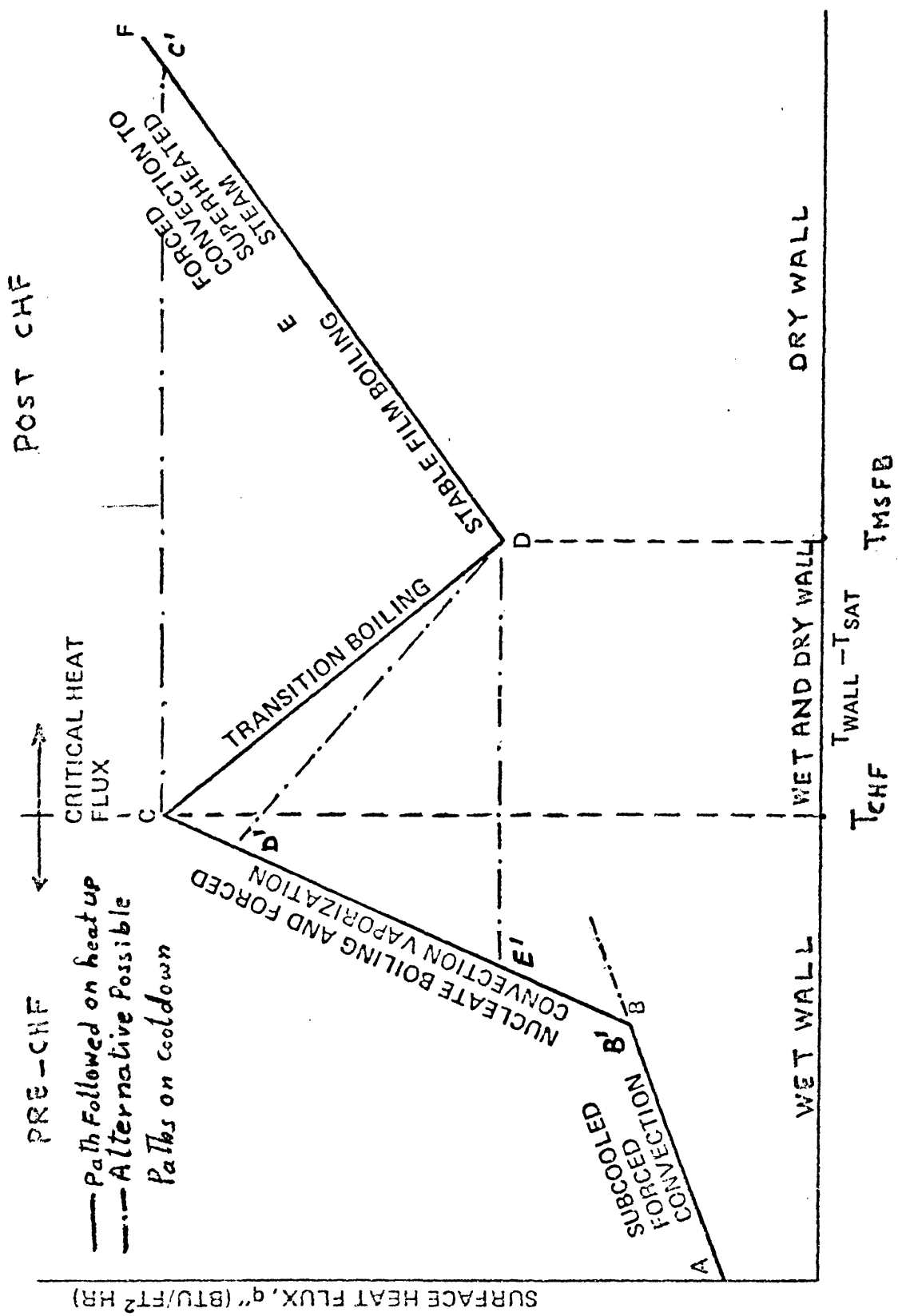


Figure 13 -Heat Transfer Regimes at Fixed Pressure

controlled surface as the temperature is increased. In general, the same path on cooldown process is not followed in the heat-up process by the boiling mechanics. For example in a heat flux controlled surface, the path ABCC'F will be followed in which point c' indicates the new equilibrium state of the surface at the heat flux value q_{CHF} .

At steady-state operation conditions, a NSSS fuel rod is a heat flux-controlled surface with a non-uniform axial heat flux distribution. In this case a reduction of the heat flux may be traced on the curve of Fig. 13 by the path EC'D E'BA⁽¹⁸⁾.

Unlike the steady-state conditions, during a hypothetical LOCA it is not clear which mechanism prevails, since the fuel rods of NSSS's may behave as heat flux-controlled surfaces for some parts of the transient and as temperature controlled surfaces for other parts of the transient.

3.1 Heat Transfer Regimes and Correlations

The recent thermal hydraulic LOCA codes have increased their capability of the two-phase heat transfer assessment by inclusion of more distinct heat transfer regimes and using more realistic correlations for calculations of the heat transfer coefficient in each regime.

As was discussed in section 3.1, a stand-alone LOCA code is capable of handling only one phase of hypothetical LOCA's, whereas integrated LOCA codes such as RELAP4-MOD7 and TRAC have the capability of calculating both blowdown and refill/reflood phases of a LOCA. This capability is made possible through inclusion of the unique features of bottom flooding (in PWR) and top spray (in BWR) of reflood heat transfer, in the blowdown heat transfer package. Such features are quench front, rewetting and liquid entrainment. Also thermal radiation and dispersed flow film boiling are specially pronounced in reflood heat transfer and are treated explicitly in the reflood heat transfer packages*.

The heat transfer package which was used in the early versions of the RELAP series such as RELAP2, is used extensively in the thermal hydraulic codes**. This package is used in various versions of RELAP3 as well as RELAP3B-MOD101. Later it was modified by replacing the quality by void fraction to determine the pre-CHF heat transfer regimes and by treating the transition boiling explicitly in which case the heat transfer coefficient is calculated using the MC DOUNOUCH, MILICH and KING correlation. Also

* See REFLUX⁽³⁹⁾ package which is developed at MIT to analyse the reflood phase of a LOCA.

**This package is essentially adopted from the THETA hot-channel code.

the Berenson and Groeneveld correlations were added to the formerly existing Dougall-Rohsenow correlation in the film boiling regime. RELAP4-MOD5 and RETRAN use this modified version and a simplified form of this new version was implemented in COBRA-IV-I. RELAP4-EM employs the new version with further modifications to satisfy the acceptance criteria. For example return to nucleate boiling is precluded once CHF happens. Also the GE correlation is added to the CHF correlations as an option to replace the Barnett correlation for BWR analysis.

There are however several disadvantages associated with this package⁽⁴⁰⁾.

- 1) There is no CHF scheme to consider CHF during flow reversal or stagnation, which are characteristic of blowdown in large, cold leg breaks in PWR's.
- 2) Use of Thom's correlation up to a void fraction equal to 0.8 which corresponds to a quality equal to 0.42 at 2250 Psia, which is above the quality range for which this correlation was verified.
- 3) Extensive use of correlations whose data base rely on tube or annular geometry, while their application is for rod bundle geometry.

- 4) Using the correlations which have a steady-state data base for transient conditions.

For this reason a heat transfer package called BEEST developed at MIT to overcome these drawbacks. BEEST⁽⁴⁰⁾ stands for BEst ESTimate heat transfer analysis. It is based on best estimate assessment rather than conservative correlations. Several tests of BEEST showed that it is able to construct the complete boiling curve where different heat transfer regimes are smoothly connected (Fig. 14). The heat transfer selection logic in this package is based on the comparison of the clad surface temperature with the two distinct temperatures on the boiling curve, namely the temperature at the minimum stable film boiling point, T_{MSFB} , and the temperature at the critical heat flux point, T_{CHF} (Fig. 13). This is certainly an unambiguous, efficient and valid criterion for selecting the appropriate heat transfer regime. Once the regime is identified, the second step is to apply a chosen correlation for the heat transfer regime selected. The upflow and downflow heat transfer are treated separately through using the void fraction. The transition boiling in this package is treated in a unique way. This treatment is based upon an interpolation between the Q''_{MSFB} and Q''_{CHF} (which are the heat flux corresponding to the T_{MSFB} and T_{CHF}) with

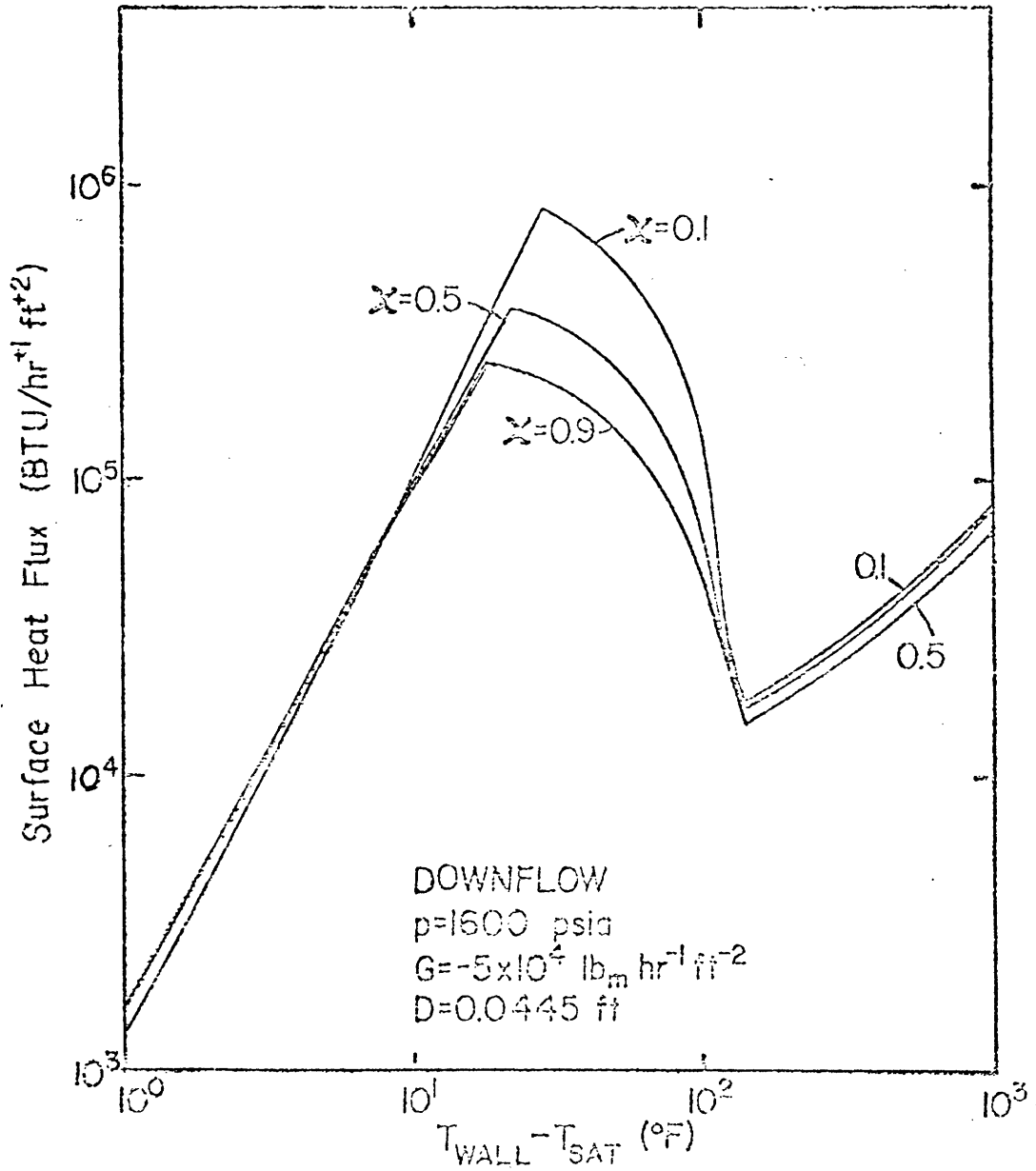


Fig. 14^[40] Effect of quality on calculated boiling curve for low flow downflow, predicted by BEEST package.

respect to the temperature ratio as follows*:

$$Q''_{TB} = E Q''_{CHF} + (1-E) Q''_{MSFB} \quad [21]$$

where

$$E = (T_{wall} - T_{MSFB}) / (T_{CHF} - T_{MSFB})^2 \quad [22]$$

In equation (22), E may be interpreted as the fraction of wall area that is wet. BEEST uses the Biasi correlation for the CHF calculations. The Biasi correlation is essentially a dry-out correlation. Therefore it is appropriate for high flows and qualities where the vapor is a dominant factor leading to dry-out. For low flows and qualities the void-CHF correlation developed at MIT is used. The RELAP heat transfer package which was discussed earlier uses the Barnett correlation as well as

* This concept was first introduced by W. Kirchner (see Ref. 41), in the form of a Log-Log interpolation:

$$h_{TB} = \frac{Q''_{CHF}}{T_{wall} - T_{sat}} \left(\frac{T_{wall}}{T_{CHF}} \right)^{\lambda} \text{ where}$$

$$\lambda = \frac{\text{Log } Q''_{CHF} - \text{Log } Q''_{MSFB}}{\text{Log } T_{CHF} - \text{Log } T_{MSFB}}$$

Kirchner then applied his model in the heat transfer package of TRAC.

the modified Barnett and the B&W-2 correlations for the CHF calculations. In the pre-CHF regimes, the Chen correlation is used in the subcooled nucleate boiling, saturated nucleate boiling as well as forced convection vaporization. This correlation has predicted the existing data with reasonable agreement⁽³⁸⁾ as compared to the other correlations such as: Dengler-Addams, Schrock-Grossman, Bennett et al, Sani and finally Guerrieri - Talty. The Chen correlation is applicable to flow regimes from slug flow through annular flow. While its data base is for low pressures⁽⁴²⁾, in most applications it is used at elevated pressures. Also, its dependence on the wall temperature which necessitates an alternative procedures, makes it less desirable.

The advantages of the BEEST heat transfer package namely, treating the upflow and downflow separately, using a once through heat transfer regime selection logic, using wall temperature as a heat transfer regime selection tool, using a best estimate assessment and incorporating the new improvements in heat transfer, has made it acceptable to the state-of-the-art LOCA codes. THERMIT uses BEEST with some modifications such as replacement of the void fraction calculated from DFM by that calculated in THERMIT. TRAC heat transfer packages is also very similar to BEEST. In fact it can be considered as an improved version of BEEST

with the following additions:

- 1) Adding two options to the CHF correlation namely, the Bowring and the Zuber Pool boiling correlations.
- 2) Inclusion of the thermal radiation contribution in the film boiling regime.
- 3) Using a horizontal film condensation to represent the low flow rates.
- 4) Inclusion of a vertical film condensation regime.
- 5) Considering laminar and turbulent flow correlations in steady-state calculation for forced convection to two-phase mixture.

A comparison of heat transfer selection logic and correlations used in different thermal-hydraulic codes is present in Tables 1.1, 1.2 and 1.3. The notations and specifications used in these tables are as follows:

- 1) Thermodynamic quality is represented by

$$x = \frac{h-h_f}{h_{fg}} \quad \text{where } h \text{ represents enthalpy, whereas}$$
$$X \text{ represents the true quality } X = \frac{W_g}{W_f + W_g}, \text{ where } W$$

is the mass flow rate.

2) The notations α , T_w , T_f , T_{sat} , P , G represent the void fraction, wall, fluid and saturated temperatures, pressure and finally mass flux, respectively. The dimension of the pressure and mass flux are in terms of "Psia" and "lbm/hr-ft²" respectively.

3) The terms "High" and "Low" flow used in these tables are in accordance with the flooding correlation which read

$$J_f^* \frac{1}{2} + m J_g^* \frac{1}{2} = K \quad [23]$$

where for turbulent flow m is equal to unity and J_f^* and J_g^* are dimensionless velocities:

$$J_f^* = J_f \rho_f^{\frac{1}{2}} [gD(\rho_f - \rho_g)]^{-\frac{1}{2}} \quad [24]$$
$$J_g^* = J_g \rho_g^{\frac{1}{2}} [gD(\rho_f - \rho_g)]^{-\frac{1}{2}}$$

where D is pipe diameter and K is the flow criteria. For example, in low flow region according to Ref. 40, this criterion is

$$J_f^* \frac{1}{2} + J_g^* \frac{1}{2} < 1.36 \quad \text{for upflow}$$
$$J_f^* \frac{1}{2} + J_g^* \frac{1}{2} < 3.5 \quad \text{for downflow} \quad [25]$$

4) The letter x and α in the parentheses in Table 3.2, in front of the Thom and Schrock-Grossman correlations imply that an interpolation is made with respect to the quality and void fraction respectively, i.e. quality or void fraction weighted heat transfer coefficient.

4. Fuel Rod Model

Temperature excursions of the fuel rod in case of any transient or accident are a major point of concern in the reactor safety analysis. A high temperature rise following severe transients is a threat to the cladding material whose integrity must be guaranteed in order to prevent any release of radioactive materials. There are four barriers preventing the release of radioactive fission gases to the environment under normal operating conditions namely, the UO_2 fuel, the fuel rod cladding, the reactor primary systems, and finally, the reactor containment building⁽⁴⁰⁾. Accordingly fuel melting, threatens the first barrier, and clad rupture violates the second barrier. The ECCS final acceptance criteria requires that failure of these barriers must be avoided under any circumstances. This necessitates a realistic fuel rod modeling, specially for the LOCA codes.

In general a fuel rod model consists of an approach to solution of the general three dimensional, time dependent

Table 3-1

Heat Transfer Regime Selection Logic

Heat Transfer Regime Code Name	Forced Convection Single Phase Flow	Fully Developed Subcooled Nucleate Boiling	Saturated Fully Developed Nucleate Boiling	Forced Convection Vaporization	Pool Boiling	Forced Convection Stable Film Boiling	Transition Stable Film Boiling	Forced Convection Single Phase Boiling	Low Pressure Film Boiling	High Pressure Film Boiling	Pool Film Boiling	Forced Convection Single Phase Boiling	Natural Convection to Vapor	
RELAP2	$x < 0.0$	$n < 0.0$	$0.0 < n < 0.1$	$0.1 < n < 0.6$	$q'' > q_{CHF}$								$x > 1.0$	-
RELAP3		$T_w > T_{NB}$	$0.0 < n < 0.1$	$0.6 < n < 1.0$	$q'' > q_{CHF}$									
RELAP3B	$T_w < T_{NB}$	$T_w > T_{NB}$			$q'' > q_{CHF}$									
RELAP4		$\alpha < 0.8$			$q'' > q_{CHF}$									
COBRA-IV-I	$T_w < T_{sat}$	$T_w > T_{sat}$	$0.8 < \alpha < 0.9$	$\alpha > 0.9$ steady-state	$\alpha > 0.9$	$P < 500$	$G > 2 \times 10^5$	$P < 500$	$G > 2 \times 10^5$	$P < 500$	$G > 2 \times 10^5$	$h > h_g$	-	
RETRAN		$T_w > T_{sat}$			$q'' > q_{CHF}$									
WOSUB	$x < 0.0$	$x > 0.0$ $h < h_{sat}$	$x > 0.0$ $h < h_{sat}$									$h > h_g$	-	
TRAC	$\alpha < 0.05$	$T_w < T_{CHF}$	$\alpha < 0.9995$ $T_w < T_{CHF}$	$T_w < T_{CHF}$	$T_w < T_{CHF}$	$T_w < T_{CHF}$	$T_w < T_{CHF}$	$T_w < T_{CHF}$	$T_w < T_{CHF}$	$T_w < T_{CHF}$	$T_w < T_{CHF}$	$\alpha > 0.9995$	$\alpha > 0.9995$	
	$T_w < T_{sat}$	steady-state	steady-state		$T_{CHF} < T_w < T_{MSFB}$	$T_w > T_{MSFB}$		High Flow	Low Flow			High Flow	Low Flow	
Thermit	$T_w < T_{sat}$	$T_F < T_{sat}$	$T_w < T_{CHF}$	$T_F < T_{sat}$	$T_{CHF} < T_w < T_{MSFB}$	$T_w > T_{MSFB}$		High Flow	Low Flow			$x > 0.99$	$x > 0.99$	

Table 3-2
Heat Transfer Correlations
Used in Each Regime

Pre-CHF Heat Transfer Regimes				CHF	
Code Name	Forced Convection Single Phase Flow	Fully Developed Subcooled Nucleate Boiling	Saturated Fully Developed Nucleate Boiling	Forced Convection Vaporization	
RELAP2	Sieder-Tate	Thom	Thom, Schrock-Grossman (x) Schrock-Grossman	M. Barnett	
RELAP3		Thom, Schrock-Grossman (α)	Schrock-Grossman, Dittus-Bolter	Barnett	
RELAP3B					
RELAP4	Dittus-Bolter	M. Chen	Chen	Schrock-Grossman	B&W-2
RETRAN					
COBRA-IV-5					
RELAP4-MOD6					
WOSUB					
TRAC	M. Chen	Dittus-Bolter, Chen			Low G M. Zuber High G M. W-3 Barnett Israel Jenssen-Levy Cise Zuber/Biasi Biasi Bowring Biase Void-CHF
THERMIT	Sieder-Tate				

Table 3-3

Heat Transfer Correlations Used in Each Regime

Post-CHF Heat Transfer Regimes

	Forced Convection Transition Boiling	Pool Transition Stable Film Boiling	Forced Convection Stable Film Boiling	Low Pressure Film Boiling	Pool Film Boiling	Forced Convection Single Phase Vapor	Natural Convection to Vapor	Normal Radiation	Condensation
AP2									
AP3			Dougall-Rohsenow						
AP3B						Dittus-Boelter			
AP4									
RA-V-I	McDonough-Millich-King	Berenson	Groenveld, Dougall-Rohsenow						
RAN									
AP4-MOD6	Low Flow: Hsu High Flow: M. Tong-Young M. Condie-Bengston	Low Flow: Brohley-Pomeranz, Hsu High Flow: Groenveld, Condie-Bengston				→ α -weighted → Max. of them		yes	
C		Bjornard	Bromely, Dougall-Rohsenow					yes	Horizontal: Chats, Collier Vertical: Carpenter-Colburn
RMIT	Griffith			High G: Low Pressure: M. Dittus-Boelter High Pressure: Groenweld 5.7	Low G: M. Bromely	Dittus-Boelter	McAdams		

Poisson's equation (heat conduction equation)

$$\rho c_p \frac{dT}{dt} = \vec{\nabla} \cdot (\underline{K} \vec{\nabla} T) + Q''' \quad [26]$$

where T, t, ρ and c_p represent the temperature, time, material density and specific heat respectively. In this equation \underline{K} represents the conductivity tensor, $\underline{K} = K_{ij}$, $i=1,2,3$, $j=1,2,3$ and Q''' is the heat source density which represents the amount of heat generated in the material per unit volume per unit time. Generally as the cylindrical shape of fuel rod dictates, a cylindrical coordinate system is chosen to expand the first term in the RHS of Equation 26.

4.1 Fuel Region

The expanded form of equation 24 in the fuel region is:

$$\rho c_p \frac{\partial T}{\partial t} = \frac{1}{r} \frac{\partial}{\partial r} \left(r k \frac{\partial T}{\partial r} \right) + \frac{1}{r^2} \frac{\partial}{\partial \theta} \left(k \frac{\partial T}{\partial \theta} \right) + \frac{\partial}{\partial z} \left(k \frac{\partial T}{\partial z} \right) + Q''' \quad [27]$$

where k in equation 25 is no longer a tensor but a time dependent scalar. This simplification is made possible through the valid assumption of homogeneous,

isotropic solid for UO_2 and fuel rod cladding material. The first term in the right hand side of Equation 27 is considered in the more general form in COBRA-IV-1 as follows:

$$r - \text{direction: } \frac{1}{R^2 r^{a-1}} \frac{\partial}{\partial r} (r^{a-1} k \frac{\partial T}{\partial r})$$

By assuming $a=2$, the cylindrical and $a=1$, the planar fuel can be treated.

The total derivative in the left hand side of Equation 26 is changed to a partial derivative in Equation 27. This simplification is possible as long as a stationary solid is treated. This in turn is a valid assumption since the fuel centerline melting is to be prohibited by design under any circumstances.

The azimuthal, or θ -direction, conduction is ignored in all the reviewed fuel rod models. This implies an assumption of infinite circumferential heat conduction.

The axial conduction, Z -direction, is only considered in COBRA-IV-1 and is ignored in the other codes.

Further simplification to Equation 27 is possible by assuming that all physical properties are temperature independent, in addition to the isotropic assumption. This is done for example in WOSUB. However, the temperature

dependence of thermal conductivity and heat capacity is considered in TRAC and THERMIT. The latter uses a chebyshev polynomial fitted to the MATPRO⁽⁴³⁾ expressions which represent fits to experimental data for fuel and clad material properties. For example a cubic and a quadratic polynomial is used to fit the temperature dependence of ρ , c_p and k of the fuel, respectively.

The Kirchoff's transformation is used in COBRA-IV-1 to reduce Equation 27 to a linear partial differential equation. By using this method the temperature dependence of k is taken into account.

As for the RELAP series, RELAP2, RELAP3, RELAP3B and RELAP4 use a simplified lumped model for their heat conduction calculation. In these codes heat generation is determined by reactor kinetics routines or by input specified values for power versus time. The fuel rod model used in these codes is patterned after the model used in the HEAT1 code. The final form of the heat conduction equation is presented in Table 4.1, equation 28. In this equation, the average temperature is

defined by (11).

$$\bar{T} = \frac{\int_V \rho c_p T dV}{\int_V \rho c_p dV} = \frac{\sum_{n=1}^n (\rho V)_n (c_p)_n T_n}{\sum_{n=1}^n (\rho V)_n (c_p)_n}$$

where V and N represent the fuel volume and fuel pin annulus number respectively. Also s in Equation 27 is the fuel pin surface area.

Equation 27 represents the U-tube as well as once-through steam generator modeled in RELAP3B-MOD101. Naturally Equation 27 does not include any heat source density term.

Equation 28 reflects the cartesian geometry used in the RETRAN fuel rod model. Thermal conductivity and heat source density temporal and spatial dependency are accounted for.

In Table 4.1, Equation 31, Equation 32 and Equation 33 represent the COBRA IV-1, COBRA-III P, WOSUB, TRAC and THERMIT, one dimensional heat conduction equation. Thermal source density Q''' in COBRA-IIIC & COBRA-III P is calculated as follows: The total power is

$$Q = \pi D \Delta Z Q'' \quad [35]$$

where D, ΔZ and Q'' are the fuel rod diameter, the axial interval and the heat flux respectively. Now dividing by the fuel volume gives

$$Q''' = Q'' \frac{\pi D \Delta Z}{\frac{\pi D_f^2}{4} \Delta Z} = Q'' \frac{4D}{D_f^2} \quad [36]$$

where D_f represents the fuel pellet diameter.

The thermal source density in WOSUB is assumed to be spatially uniform but time dependent, whereas in TRAC and THERMIT it depends both on position and time. Equation 34 demonstrates the COBRA-IV-I fuel rod model. This equation is found by using the Kirchoff's transformation.

$$\theta = \frac{1}{k_o} \int_{T_o}^T k(T) dt \quad [37]$$

where k_o is the conductivity at reference temperature T_o . Differentiating Equation 37 with respect to r and t and substituting in Equation 27 we will come up with Equation 34.

4.2 Fuel-Clad Gap

The fuel-clad gap heat transfer coefficient is implicitly treated in those models in which clad and the fuel-clad gap, are lumped together.

This is done for example in the COBRA-IIIC fuel rod model. However, upon the importance of the fuel-clad gap resistance to the heat flow, it is treated explicitly in RELAP3B-MOD101, WOSUB, TRAC and THERMIT.

The gap heat transfer coefficient depends upon the fission gas product in the gap, the radiation heat

Table 4-1

Approximations Made to the Equation of Heat Conduction

Code Name	Characteristics	Fuel Region	Clad Region	Eq. No.
RELAP2 RELAP3 RELAP3B RELAP4 (Fuel rod)	Radial lumped Parameter approach	$V \rho C_p \frac{dT}{dt} = Q''' + Sk \frac{dT}{dr}$	Included in the fuel region except for RELAP3B	(28)
RELAP3B (Heat ex. and Steam ge.)	One-dimensional cylindrical coo. system	$\rho C_v \frac{\partial T}{\partial t} = k \left[\frac{\partial^2 T}{\partial r^2} + \frac{1}{r} \frac{\partial T}{\partial r} \right]$	-	(29)
RETRAN	One-dimensional cartesian coo. system (x-direction)	$\rho C_p \frac{\partial T(x,t)}{\partial t} = \frac{\partial}{\partial x} \left[k(x,T) \frac{\partial T(x,t)}{\partial x} \right] + Q'''(x,t)$		(30)
COBRA-IIIC COBRA-IIIP	One-dim. cylindrical coo. system temp. Independent properties and thermal source density	$\rho C_p \frac{\partial T(r,t)}{\partial t} = k \left(\frac{\partial^2 T}{\partial r^2} + \frac{1}{r} \frac{\partial T}{\partial r} \right) + Q'''$	Fuel-clad gap and clad are lumped together	(31)
WOSUB	One-dim. cylindrical coo. system temp. independent properties	$\rho C_p \frac{\partial T(r,t)}{\partial t} = k \left(\frac{\partial^2 T}{\partial r^2} + \frac{1}{r} \frac{\partial T}{\partial r} \right) + Q'''(t)$	$Q'''(t) = 0$	(32)
TRAC THERMIT	One-dim. cylindrical coo. system	$\rho C_p \frac{\partial T(r,t)}{\partial t} = \frac{1}{r} \frac{\partial}{\partial r} \left[rk(r,t) \frac{\partial T(r,t)}{\partial r} \right] + Q'''(r,t)$	Q''' Metal-Water Reaction	(33)
COBRA-IV-I	Two-dim. planar and cylindrical geometry	$\rho C_p k(T) \frac{\partial \theta}{\partial t} = \frac{k_0}{R^2} \frac{\partial}{\partial r} \left[r a - 1 \frac{\partial \theta}{\partial r} \right] + \frac{\partial}{\partial z} \left[a - 1 \frac{\partial \theta}{\partial z} \right] + Q'''$	Fuel-clad gap and clad are lumped together	(34)

transfer across the gap as well as the fuel-clad contact and fuel-clad pressing⁽²⁰⁾. These are modeled in the GAPCON, MATPRO, FRAP-S and FRAP-T codes. It's MATPRO model which is implemented in THERMIT. In this model all the above mentioned factors are considered. The model used in TRAC ignores the effect of fuel pressing against the clad, whereas it is correlated in THERMIT in terms of the fuel contact pressure against the clad.

An effective gap heat transfer coefficient is used in WOSUB. Although this is not as realistic as the models used in TRAC and THERMIT, it still allows nodalization in the clad. COBRA-IIIC and COBRA-IV-1 assume the outer fuel surface and the inner clad surface are in a single node. In this case the conduction equation is written between the fuel and clad exterior surface. The heat transfer coefficient used for this purpose is defined as⁽⁴⁾:

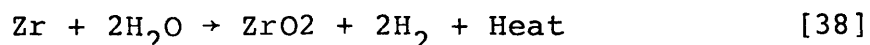
$$\frac{1}{H} = \frac{1}{H_g} + \frac{Y_c}{k_c}$$

where Y_c and k_c are the cladding thickness and conductivity and H_g is the fuel-clad gap conductance.

4.3 Clad Region

As mentioned in the previous section 4.2, only WOSUB, TRAC and THERMIT permit clad nodalization. While

the clad region is assumed to be heat source free in WOSUB, a metal-water reaction is considered as a heat source in that region in the TRAC fuel rod model as well as RELAP4. This reaction happens at elevated temperatures, below the cladding material melting point, between zirconium and steam and is expressed as⁺:



Both TRAC and RELAP4 use the parabolic rate law of Baker and Just to represent the rate of this reaction but in a different system of units⁺⁺. The mathematical statement of the parabolic rate law reads:

$$-\frac{dr}{dt} = \left(\frac{a}{R_0 - r}\right) \exp\left(-\frac{b}{T}\right) \quad [39]$$

where r , R_0 , t and T represent the radius at each moment, the initial clad exterior radius, time and temperature respectively. In this equation a and b are constant values. By integrating Equation 39 between the initial and final radii of a time step, the mass of zirconium reacted per unit length during the time step will be found. The amount of heat generated in the clad is then proportional to this reacted mass, and it will be considered as the internal heat source in the clad region.

+ This exothermic reaction, results in hydrogen gas which poses a threat to the fuel rods in case of accidents by excluding the upper part of the rods to be covered by the coolant.

++ TRAC and THERMIT are the only thermal hydraulic codes using the SI units.

5. Numerical Methods

The mathematical models which were discussed in the previous sections are solved numerically in the computer codes, because analytical solutions are impractical.

In the subchannel codes, the axial length of the reactor core is divided into several intervals which make each interval the computational control volume. The set of field equations in a finite difference form in conjunction with the constitutive equations are solved for the central volumes. The boundary conditions at the inlet of the core are, uniform or nonuniform pressure and coolant densities and enthalpies. The axial and radial heat flux profile must be specified. The solution is based on reaching a uniform pressure at the core outlet. For this purpose, a marching technique may be used. In a step by step, or marching technique, the calculation starts from the bottom of the core for all subchannels and moves upward. Inlet velocities are first assumed to be known and then solved alternatively through the external iteration loop. At each axial node, the cross-flow is guessed which allows solving the energy equation. A new value for cross-flow is calculated from pressure drop in each subchannel, which in turn is calculated from the momentum equation.

This internal iteration on cross-flow is continued until an acceptable pressure balance is reached. By knowing the heat addition into the axial cell and calculated cross-flow for the axial cell, the values of coolant density, velocity, enthalpy, and pressure can be determined at the exit of each computational cell. In turn, these values will be used as the information for the next axial cell. This procedure continues until the top of the core is reached where the criterion of uniform exit pressure is checked. If this criterion is not met, the external iteration loop which covers the whole channel length must be continued, using improved guesses of the flow division among the subchannels at the inlet. This procedure was employed in COBRA-II and HAMBO. The number of external iterations over the core length depends upon the coupling between subchannels. If this is weak (e.g. if the cross-flows are small), a single pass marching solution technique is adequate⁽²²⁾, otherwise a multipass marching solution is necessary. This concept is used in COBRA-IIIC, in which a pattern of subchannel boundary pressure differentials for all mesh points is guessed simultaneously and then the corresponding pattern of cross-flow is completed using a marching technique up the channel. By updating the pressure differentials during each external

iteration loop, the effects of downstream will be propagated upstream.

The procedure used in COBRA-IIIP is somewhat different. A new treatment is introduced for the transverse momentum equation which couples the adjacent computational cells. This includes the spatially semi-implicit treatment of the pressure field.⁺ Using this method guarantees the diagonal dominance of the matrices governing the pressure fields⁽⁴⁹⁾. The computed pressure field is then used in the transverse momentum equation to determine the cross-flow distribution. Applying the new concept in COBRA-IIIP has made it capable of increasing the number of computational subchannels markedly, i.e. from 15, in COBRA-IIIC, to 625 in COBRA-IIIP⁽⁴⁴⁾. Furthermore, it has increased the computational effectiveness resulting in a shorter running time.⁺⁺

⁺ By introducing a scalar, value θ , having an arbitrary value between 0 and 1 the pressure field is written:

$$[P] = \theta [P_j] + (1-\theta) [P_{j-1}]$$

By introducing this concept into transverse momentum equation, allows the cross-flow distribution to be driven by any combination of the pressure fields that exist at the top and the bottom of each plane of computational cells.

⁺⁺ Unlike COBRA-IIIC, a double precision is used in computation of pressure field and gradients which is specially pronounced in cross-flow distribution calculation in the vicinity of grids. This in turn will increase the computer running time.

While the marching technique determines the flow condition under steady-state situations, in transients the whole procedure will be repeated for each time increment, implicitly in both COBRA-IIIC and COBRA-IIIP. The marching technique is also employed in WOSUB, where for the sake of numerical stability, a backward finite difference form is used in space and time. The lack of transverse equation and cross-flow is compensated by the concept of recirculation loop which is based on the assumption that the net volumetric flow recirculation around closed loops connecting communicating subchannels is zero⁽¹⁷⁾.

Simultaneous solutions of the finite difference form of the field equations written for the previously defined computational cell, is another solution technique used in some subchannel codes such as SABRE and COBRA-IV-1. Since the calculated values will be advanced in each time step explicitly, this additional option in COBRA-IV-1 be called the "explicit solution scheme". The following possibilities are available in COBRA-IV-1 solution algorithm:

- 1) Steady-state and transient calculation using the COBRA-IIIC implicit solution scheme.

2) Implicit steady-state and explicit transient with either a ΔP or inlet flow boundary condition.

3) Explicit transient calculation based specified initial values with either a ΔP or inlet flow boundary condition and a zero flow initially.

The addition of the explicit numerical scheme with a ΔP boundary condition makes COBRA-IV-1 capable of handling flow reversal, recirculation and coolant expulsions as well as severe flow blockage.

These additional capabilities stem from solving a true boundary value problem rather than dealing with an initial value problem in the marching type solution technique.

The additional numerical scheme in conjunction with the boiling curve package and the improved fuel pin-model makes COBRA-IV-1 capable of assessing accidents such as a LOCA, where due to the heat transfer package inability of analysing reflood, the code capability limits to the blowdown phase of a LOCA. It should be realized that due to numerical instabilities and convergence difficulties which mostly result from discontinuities introduced by the physical models, the "explicit scheme" of COBRA-IV-1 uses the strict HEM, i.e., it does not contain the Levy subcooled boiling model or any slip correlation. Furthermore the lack of computational effectiveness, as

compared to COBRA-IIIC and IIIP, and using the reference pressure concept* which excludes any effect of compressibility and also large heat flux oscillations observed in prediction of depressurization transients⁽⁴⁶⁾ have made COBRA-IV-1 less desirable. Overcoming these deficiencies has been presumably the motivation of creating COBRA-DF and TF.

As for the loop codes, a fully implicit solution scheme, temporally, is employed in all the RELAP series as well as RETRAN. An automatic time step variation is built in RELAP3B-MOD101. Using this feature, the time-step size increases automatically during slowly varying portions of a transient case of a computer run and vice versa. Both implicit and explicit solution schemes are employed in RETRAN.

Unlike the implicit method which is unconditionally stable, the explicit method is conditionally stable in which the so-called "courant criterion" must be respected. This criterion reads:

$$\left| U \frac{\Delta T}{\Delta X} \right| < 1 \quad \text{or} \quad \Delta T < \left| \frac{\Delta X}{U} \right| \quad [40]$$

where U is fluid velocity, ΔT and ΔX are time step and

*The concept of reference pressure which ignores the sound wave propagation effects is employed in all the HEM versions of COBRA as well as WOSUB. This limitation is circumvented in COBRA-DF by using the ACE method, also see Ref. [45] in which this method is applied to the COBRA-IV-1 field equations.

mesh spacing respectively. It is clear that for fast transients which involve rapidly changing flow, very small time steps may be needed to resolve the flow evaluation. An explicit numerical scheme may be used for this purpose.

Longer time steps are desirable in calculating the mild transients, which necessitates using the implicit numerical scheme. This scheme has not been used in the three dimensional thermal hydraulic codes, because, the fully implicit difference equations are very difficult to solve in more than one space dimension. A marching solution method may be applied to circumvent this difficulty, but as it was discussed earlier, no true boundary value problem can be handled by this method, only initial value problems in which general boundary conditions and local flow reversal cannot be treated.

A compromise between the above mentioned techniques has been made in THERMIT by using a "semi-implicit" numerical scheme. As the name implies, both implicit and explicit schemes are employed in such a way that by differencing terms involving sonic propagation implicitly, limitations $\left| \frac{(U \pm C) \Delta T}{\Delta X} \right| < 1$ have been eliminated, whereas the liquid and vapor convection are treated explicitly.

Therefore the limitation imposed on the time increment to satisfy courant criteria (Equation 40) still exists. A default value is usually built in the codes which use a temporal explicit scheme to exclude the computational instability.

6. Summary and Conclusions

6.1 Summary

A summary of the aforementioned methods and models used in each reviewed code and the range of each code application are presented in Table 6-1 through 6-5.

Some of the terms which are used in these tables are further explained as follows:

Small breaks (Table 6-1): postulated breaks that are smaller than about 10% of the double-ended break in the discharge flow area.

Licensing codes (Table 6-1): notations are in accordance with the notations defined in Table 6-3.

Homologous model (Table 6-4): described the centrifugal pumps and specifies relations, connections head, torque, flow rate, and rotational speed.

Table 6-1
Component and Loop Codes Comparisons

Code Name	TYPE							APPLICATION				CHARACTERISTICS			
	PWR	BWR	LMFBR	Licensing Best Estimate System	Component	Blowdown	Reflood	Hot Channel	Small Break	Hydraulic Model	Geometry	Numerics	Case	Reference	
COBRA I	X	X	X	X	X			X		1V1T	Axial Lateral	Forward Marching	Steady State	1	
COBRA II	X	X	X	X	X			X		1V1T	Axial Lateral	Forward Marching	Steady State	2	
COBRA III	X	X	X	X	X			X		1V1T	Axial Lateral	Pseudo Boundary Condition	Steady State	3	
COBRA III-C	X	X	X	X	X			X		1V1T	Axial Lateral	Implicit	Steady State & Transient	4	
COBRA III-P	X	X	X	X	X			X		1V1T	Axial Lateral	Semi-Implicit Forcing Function for cross flow	Steady State Transient	5	
COBRA IV-I	X	X	X	X	X	X		X		1VS1T	Axial Lateral	Implicit Explicit	Steady State & Transient	6	
COBRA-DF	X	X		X	X	X	X			1VD1T			Steady State & Transient	7	
COBRA-TF	X	X		X	X	X	X			2V2T			Steady State & Transient	8	

Table 6-1
Component and Loop Codes Comparisons
(continued)

Code Name	TYPE							APPLICATION				CHARACTERISTICS				
	PWR	BWR	LMBFR	Licensing	Best Estimate	System	Component	Blowdown	Reflood	Hot Channel	Small Break	Hydraulic Model	Geometry	Numerics	Case	Reference
WOSUB		X			X	X				X		1VD1T	Axial Lateral	Implicit Forward Marching (1)	ATWS	17
THERMIT	X	X	X		X	X	X	X	X			2V2T	3-D x,y,z	Semi-Implicit		20
RELAP2	X	X		X	X	X						1V1T	Lumped Parameters	Implicit		9
RELAP3	X	X		X	X	X						1V1T	Lumped Parameters	Implicit		10
RELAP3B (MOD101)	X	X		X	X	X						1V1T	Lumped Parameters	Implicit		11
RELAP4	X	X		X	X	X			X			1V1T	Lumped Parameters	Implicit		12
RELAP4-EM ⁽²⁾	X	X		X	X	X	X	X	X			1V1T	Lumped Parameters	Implicit		15
RELAP4-FLOOD	X			X	X			X				1V1T	Lumped Parameters	Implicit		36
RELAP4 (MOD5)	X			X	X	X		X	X			1VD1T	Lumped Parameters	Implicit		13
RELAP4 (MOD6)	X			X	X	X	X	X	X			1VD1T	Lumped Parameters	Implicit		14

¹ No diversion cross flow

² In conjunction with TODDEE (for PWR) and MOXY (for BWR).

Table 6-1
Component and Loop Codes Comparisons
(continued)

Code Name	TYPE						APPLICATION						CHARACTERISTICS			
	PWR	BWR	Licensing	Best Estimate	System	Component	Blowdown	Reflood	Hot Channel	Small Break	Hydraulic Model	Geometry	Numerics	Reference		
RELAP4 (MOD7)	X	X	X	X	X		X	X	X	X	1VD1T	Lumped Parameters	Implicit	14		
RELAP 5	X	X	X		X	X	X		X	X	2VT _k T _{sat}	1-D	Semi-Implicit	16		
RETRAN	X	X		X	X		X	X	X	X	1VDS1T ⁽¹⁾	1-D	Implicit Explicit	18		
TODEE	X		X			X	X	X						37		
MOXY		X	X			X	X	X						38		
TRAC-P1	X			X	X		X	X	X	X	2V2T 1VD1T _k T _{sat}	3-D r, θ, z	Semi-Implicit	19		

¹ Dynamic slip model from the two-fluid theory

Table 6-2
Models and Methods Used in Some
Thermal Hydraulic Codes

CAPABILITY OR MODEL		CODE NAME						
		COBRA I BNWL 1967	COBRA II BNWL 1970	COBRA III BNWL 1971	COBRA IIIC BNWL 1973	COBRA IIIP MIT 1977	COBRA IV-I BNWL 1976	WOSUB MIT 1978
<u>Conservation Equation</u>								
1	Homogeneous equilibrium model, HEM(1V1T)	X	X	X	X	X	X	
2	One dimensional mass, momentum, energy equation	X	X	X	X	X	X	
3	Drift flux model						X	
4	Separate continuity equation for liquid and vapor phases						X	X
5	Separate momentum and energy equation for liquid and vapor phases							X
6	Turbulent liquid-liquid mixing in the subcooled region, in energy equation						X	X
7	Turbulent shear stress in mixture momentum equation						X	
<u>Numerical Scheme</u>								
Flow Solution-Steady State:								
8	Marching method (forward marching)	X	X				X	
9	Pseudo boundary value method			X	X	X	X	
10	True boundary value method						X	
11	New treatment of transverse momentum equation					X		

Table 6-2
Models and Methods Used in Some
Thermal Hydraulic Codes
(continued)

CAPABILITY OR MODEL		CODE NAME							
		COBRA I	COBRA II	COBRA III	COBRA IIIC	COBRA IIIP	COBRA IV-I	WOSUB	THERMIT
Flow Solution - Transient:									
12	Fully implicit				X	X	X	X	
13	Semi-implicit								X
14	Explicit, arbitrary flow field and boundary conditions (ACE method)						X		
Flow Energy Solution:									
15	Spatially explicit	X	X	X	X	X			
16	Spatially implicit						X	X	
<u>Equation of State</u>									
17	Reference pressure	X	X	X	X	X	X	X	
18	Local Pressure								X
19	Superheated steam properties						X		X
20	Steam table that contains the derivative of fluid properties								X
<u>Transverse Transport</u>									
Cross Flow Model:									
21	Pressure resistance only	X	X	X					
22	Transient momentum equation				X	X	X		
23	Forced diversion cross flow				X	X	X		

Table 6-2
Model and Methods Used in Some
Thermal Hydraulic Codes
(continued)

	CAPABILITY OR MODEL	CODE NAME							
		COBRA I	COBRA II	COBRA III	COBRA IIIC	COBRA IIIP	COBRA IV-I	WOSUB	THERMIT
24	Lateral momentum flux						X	X	
25	Two dimensional transverse flow								X
	<u>Turbulent Mixing:</u>								
26	Single phase turbulent mixing		X	X	X	X	X	X	X
27	Two phase turbulent mixing			X	X	X	X	X	X
28	Vapor drift on a volume to volume exchange basis							X	
	<u>Accident Analysis</u>								
29	Severe flow blockage, coolant expulsion, flow reversal						X		X
30	Recirculation loop						X	X	X
	<u>Single Phase Flow</u>								
31	Nonuniform channel friction		X	X	X	X	X	X	X
32	Laminar and turbulent friction correction						X	X	X
33	Hot wall friction correction				X	X	X		
	<u>Two Phase Flow</u>								
34	One-dimensional slip flow		X	X	X	X	X		

Table 6-2
Models and Methods Used in Some
Thermal Hydraulic Codes
(continued)

	CAPABILITY OR MODEL	CODE NAME							
		COBRA I	COBRA II	COBRA III	COBRA IIIC	COBRA IIIP	COBRA IV-I	WOSUB	THERMIT
47	One-dimensional heat con- duction equ. (r-direc.)				X	X		X	X
48	Two-dimensional heat con- duction equ. (r,z-direc.)						X		
49	Implicit finite difference solution scheme				X	X			X
50	Collocation method							X	
51	Orthogonal collocation technique (MWR)						X		
52	Temperature dependent thermal conductivity						X		X
53	Transient (time dependent) heat source density							X	X
54	Constant fuel-clad gap heat transfer coefficient				X	X	X	X	
55	Thermal radiation and interfacial contact in the gap heat transfer coefficient								X
56	Planar or cylindrical fuel						X		
57	Axial fuel zone						X		

Table 6-2
 Models and Methods Used in Some
 Thermal Hydraulic Codes
 (continued)

	CAPABILITY OR MODEL	CODE NAME					
		RELAP 2	RELAP 3	RELAP 3B	RELAP 4	RETRAN	TRAC-P1
		INEL 1968	INEL 1970	INEL 1976	INEL 1973	INEL 1977	LASAL 1978
	<u>Conservation Equations</u>						
1	Homogeneous equilibrium model, HEM (1V1T)	X	X	X	X	X	
2	One-dimensional mass, momentum, energy equation	X	X	X	X	X	
3	Inclusion of K.E. and P.E. in energy equation				X	X	X
4	Consideration of area and density change in momentum equation				X	X	X
5	Dynamic slip model from the two fluid theory to account for nonhomogeneous flow					X	
6	Three dimensional (r,θ,z) flow for vessel						X
7	One-dimensional flow with drift flux model for the rest of the primary loop						X
	<u>Numerical Scheme</u>						
8	Fully implicit solution scheme temporally	X	X	X	X	X	X
9	Factor to modify the fully implicit scheme				X	X	X
10	Automatic time step variation			X			
11	Explicit scheme					X	

Table 6-2
Models and Methods Used in Some
Thermal Hydraulic Codes
(continued)

CAPABILITY OR MODEL		CODE NAME					
		RELAP 2	RELAP 3	RELAP 3B	RELAP 4	RETRAN	TRAC-P1
<u>Equation of State</u>							
12	Local Pressure	X	X	X	X	X	X
13	Steam table that contains the derivative of fluid properties				X		X
14	Extension of the steam table above the critical pressure			X			
<u>Physical Model</u>							
Pump Characteristics:							
15	Homologous pump			X	X	X	X
16	Only one pump coastdown curve	X	X				
17	Independent tripping on the independent signals			X	X	X	
18	Pump is at a junction	X	X				
19	Pump is in a volume			X	X	X	X
20	Consideration of inertial effect				X	X	X
21	Consideration of frictional torque				X	X	X
22	Consideration of bearing and windage torque						X
23	Option for two phase pump				X	X	X
24	Motor torque option, pump stop option, & dimensionless head ratio difference data in two phase pump					X	

Table 6-2
 Models and Methods Used in Some
 Thermal Hydraulic Codes
 (continued)

CAPABILITY OR MODEL		CODE NAME					
		RELAP 2	RELAP 3	RELAP 3B	RELAP 4	RETRAN	TRAC-P1
Choked Flow:							
25	Moody's two-phase choked flow model	X	X	X	X	X	
26	Henry-Fauske and extended Henry-Fauske					X	
27	Sonic choking				X	X	
Heat Exchanger:							
28	Non-conduction model (input specified secondary temperature & a constant effective heat transfer coefficient)	X	X	X	X	X	
29	Time dependent heat exchanger (Input specified table of normalized power versus time)				X	X	
30	Time dependent secondary temperature			X			
31	U-tube steam generator (one dimension heat conduction equation)			X			
32	Once through steam generator			X			
Single Phase Friction Factor:							
33	Laminar friction factor				X	X	X
34	Turbulent friction factor				X	X	X
35	Input specified friction factor	X	X	X			

Table 6-2
 Model and Methods Used in Some
 Thermal Hydraulic Codes
 (continued)

	CAPABILITY OR MODEL	CODE NAME					
		RELAP 2	RELAP 3	RELAP 3B	RELAP 4	RETRAN	TRAC-P1
	Two Phase Frictional Multipliers:						
36	Modified Baroczy correlation			X	X	X	
37	New correlation based on modified Baroczy			X			
38	Beattie correlation using Bennet flow regime map					X	
39	CISE model						X
40	Annular flow model						X
41	Chisholm model						X
42	Homogeneous correlation					X	X
43	Armand model						X
	Bubble Rise Model:						
44	Linear approximation for the density of bubbles versus height	X	X	X	X	X	
	Heat Transfer Package:						
45	Correlations for pre and post CHF	X	X	X	X	X	X
46	Ability to construct boiling curve				X	X	X
47	Treatment of transition boiling explicitly				X	X	X
48	Condensation calculation						X
49	Reflood heat transfer package					X	X

Table 6-2
 Model and Methods Used in Some
 Thermal Hydraulic Codes
 (continued)

CAPABILITY OR MODEL	CODE NAME					
	RELAP 2	RELAP 3	RELAP 3B	RELAP 4	RETRAN	TRAC-P1
Heat Transfer Regime Selection Tool:						
50 Quality and CHF	X	X	X			
51 Void fraction and CHF				X	X	
52 Local clad surface temp.						X
Fuel Rod Model:						
53 Lumped approach	X	X	X	X		
54 One-dimensional heat conduction equation					X	X
55 Variable gap size during a transient			X			X
56 Thermal radiation and interfacial contact in the gap heat transfer coefficient				X		X
57 Fuel-clad gap heat transfer coefficient burn up dependent						X
58 Thermal conductivity temperature dependent					X	X
59 Exothermic metal-water reaction considered as a heat source in the cladding material				X	X	X
60 Explicit numerical scheme				X	X	
61 Implicit numerical scheme	X	X	X	X	X	X

Table 6-3

CODE NAME	Liquid Continuity Equation Vapor Cont. Eq. Mixture Mom. Eq. Mixture Energy Equation	Mixture Cont. Equation Mixture Mom. Eq. Differential Mom. Eq. Mixture Energy Eq.	Liquid Cont. Eq. Vapor Cont. Eq. Mixture Mom. Eq. Liquid Energy Eq. Vapor Energy Eq.	Liquid Cont. Eq. Vapor Cont. Eq. Liquid Mom. Eq. Vapor Mom. Eq. Mixture Energy Equation	Liquid Cont. Eq. Vapor Cont. Eq. Liquid Mom. Eq. Vapor Mom. Eq. Liquid Energy Eq. Vapor Energy Eq.	TOTAL NUMBER OF CONSERVATION EQUATIONS
WOSUB	X					4
RETRAN		X				4
COBRA-DF			X			5
RELAP5				X		5
COBRA-TF					X	6
TRAC					X	6
THERMIT					X	6

6.2 Conclusions

A study was made of a number of well-known thermal hydraulic codes. This study attempted to cover both models and methods used in these codes summarizing their basic elements in various tables. The following results are drawn:

6.2.1 Component Code

The COBRA series as well as WOSUB and THERMIT fall in this category. As for the COBRA codes, the steady-state versions namely COBRA-I, II, III are certainly obsolete by now and are not suggested for further considerations. COBRA-IIIC and its MIT version COBRA-IIIC/MIT*, which utilize the HEM model of two-phase flow, can be used for both normal operation and transient conditions. The great flexibility of COBRA-IIIC/MIT to simulate core regions, bundles and subchannels at the same time, makes it more desirable to use. An alternative marching solution using implicit numerical scheme, which converges on the cross-flow is used in this code. The fuel rod model in conjunction with Pre-CHF heat transfer correlations make it a fast running code for steady-state as well as mild transients. On the other hand COBRA-IIIP/MIT is capable of handling a larger number of computational subchannels with considerable computational effectiveness, since it deals with the diagonally

*The major difference between these two codes is that the latter uses a dynamic data management subroutine which allows the dimensions of the principal arrays as well as the total computer storage requirement to be a function of the problem size.

dominant pressure matrix. However it should be realized that these codes, COBRA-IIIC/MIT and IIIP/MIT, are not devised for the purpose of accident analysis such as a LOCA. In fact neither the physical models nor the numerical methods have such capability. For example, lack of a heat transfer package and use of the marching solution technique do not allow any extreme flow as well as reliable fuel rod temperature calculations. A step toward the analysis of severe transients and/or accidents is taken in COBRA-IV-I in which more realistic physical models, with respect to the 2-D fuel rod model and heat transfer package, are implemented. More importantly, a field equation solution technique, explicit solution, is employed in addition to the COBRA-IIIC implicit type solution scheme. It is the simultaneous set of differential equations using explicit solution technique which makes COBRA-IV-I capable of handling severe flow blockages, flow reversal, coolant expulsions and other extreme flow situations. Also, the field equations solution using ACE technique allows specifying a ΔP boundary condition which relaxes the impractical specification of inlet flow boundary condition in severe transients such as blowdown.

Despite these advantages, it should be recalled that several disadvantages are associated with this code which are as follows:

- 1) It relies extensively on the HEM,
- 2) Computational ineffectiveness,
- 3) The time change of local pressure is ignored,
- 4) Subcooled void is excluded in the explicit scheme.

A more realistic two-phase flow model that relaxes the assumption of the HEM is used in the newer versions of COBRA namely, COBRA-DF and TF, still under development. Except for the highlights of the models used in these codes that are presented in the summary tables, there is little additional information available on these codes for the time being.

Since analysis of the BWR normal operation and transient conditions is more demanding with respect to the two-phase flow modeling, the recent efforts in the BWR models have been focused on using more realistic assumptions with regard to vapor-liquid momentum exchange, or phase distribution as for example the WOSUB code.

Unfortunately, the marching type solution technique, simplifying assumptions such as the reference pressure concept and lack of a complete boiling curve package, in the current WOSUB version, put a limit on its application. For example, no reliable blowdown calculations can be performed with WOSUB in its present form. Furthermore, WOSUB is not supposed to be used for very fast transients.

It is now clear that extreme flow situations which are a point of concern in severe transients can only be evaluated by using more physically accurate field equations. The two-fluid concept provides the potential for increased accuracy in modeling the two-phase flow. By implementing this concept in the most recent codes, such as THERMIT, a number of limitations imposed on the flow characteristics are relaxed. Now the motion of two-phase in different directions, having different temperatures, velocity and pressure can be realistically analysed in three dimensions. The best estimate heat transfer package, BEEST, and an improved fuel rod model, specially with respect to the material temperature dependent and fuel-clad gap modeling, which are included in THERMIT, provide a reliable fuel rod temperature calculation and DNBR prediction. A semi-implicit numerical

method is used to circumvent the instabilities associated with the explicit numerical scheme.

In spite of the above mentioned advantages, it should be remembered (see section 1.3) that the need for mathematical models representing physical phenomena increases with the degree of sophistication of the two-phase flow modeling. Furthermore, the difficulty with the general two-fluid approach is that the exchange processes coupling the phases are currently not thoroughly understood^[20]. As a result, despite the possible shortcoming of the HEM, it is not evident that a homogeneous equilibrium model is incapable of predicting adequately some parameters of interest such as vapor flow rate and fuel-clad temperature. Furthermore, in some flow regimes HEM gives surprisingly good results. Therefore, as a final conclusion, for normal operation and mild transients, COBRA-IIIC/MIT and COBRA-IIIP/MIT are still the best available tools. Several shortcomings of these codes such as the fuel-rod model and lack of a heat transfer package and the like may be overcome by implementing the state-of-the-art models used in the sophisticated codes such as THERMIT. Severe transients and accident analysis are certainly advised to be analysed by THERMIT. Upon the completion of a subchannelwise version with coolant centered control volume, THERMIT will

be the first three dimensional non-homogeneous fully thermodynamic non-equilibrium subchannel code for the nuclear reactor core thermal hydraulic analysis.

6.2.2 Loop Codes

The RELAP series as well as RETRAN and TRAC fall in this category. The old versions of RELAP such as RELAP2, and most of the RELAP3 versions are obsolete and need not be further considered. RELAP3B-MOD101 is the only updated version which uses several options for heat exchanger and steam-generator modeling, not even used in TRAC. This version is specially devised for the ATWS analysis. It uses a combination of old models of RELAP3 such as heat transfer package, and new models introduced in RELAP4, such as homologous pump, in addition to several unique features such as variable gap size during a transient and heat exchanger modeling. In light of the detailed information about RELAP4/MOD6 and MOD7 and RELAP5, the available highlights of loop modeling are presented in the related tables. From these tables it is clear that the major step toward the non-homogeneous non-equilibrium modeling of the primary loop is taken in the model making process of RELAP5. Such effort is also done in RETRAN through introducing

the unique feature, DSM, "dynamic slip model". RETRAN is specially devised to analyse the reflood phase of a LOCA. The state-of-the-art of the loop codes however is TRAC which is capable of handling all the phases of a hypothetical LOCA.

As a result the following codes are suggested for the transient loop calculations. First, RELAP3B-101 which is a one-dimensional, HEM code and it may be used for ATWS transients. Second, RETRAN which uses 1-D, DSM and improved physical models, third, TRAC which uses 1-D, DFM for the loop calculations, and realistic physical models.

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APPENDIX 1

In the following, the assumptions which have been made in derivation of the continuity equation in the COBRA codes are discussed.

The mass balance for liquid is given by:

$$\frac{\partial}{\partial t} [(1-\alpha)\rho_l] + \nabla \cdot [(1-\alpha)\rho_l \vec{V}_l] = -\Gamma \quad (1)$$

Mass balance for vapor follows

$$\frac{\partial}{\partial t} [(\alpha\rho_v)] + \nabla \cdot [(\alpha\rho_v \vec{V}_v)] = \Gamma \quad (2)$$

where Γ is the phase change rate. Now using a pseudo single phase concept by assuming:

$$\begin{aligned} \bar{\rho} &= (1-\alpha)\rho_l + \alpha\rho_v \\ \bar{\rho}\vec{V} &= (1-\alpha)\rho_l \vec{V}_l + \alpha\rho_v \vec{V}_v \end{aligned} \quad (3)$$

and adding terms (1) and (2) we will come up with

$$\frac{\partial}{\partial t} [(1-\alpha)\rho_l + (\alpha\rho_v)] + \nabla \cdot [(1-\alpha)\rho_l \vec{V}_l + \alpha\rho_v \vec{V}_v] = 0 \quad (4)$$

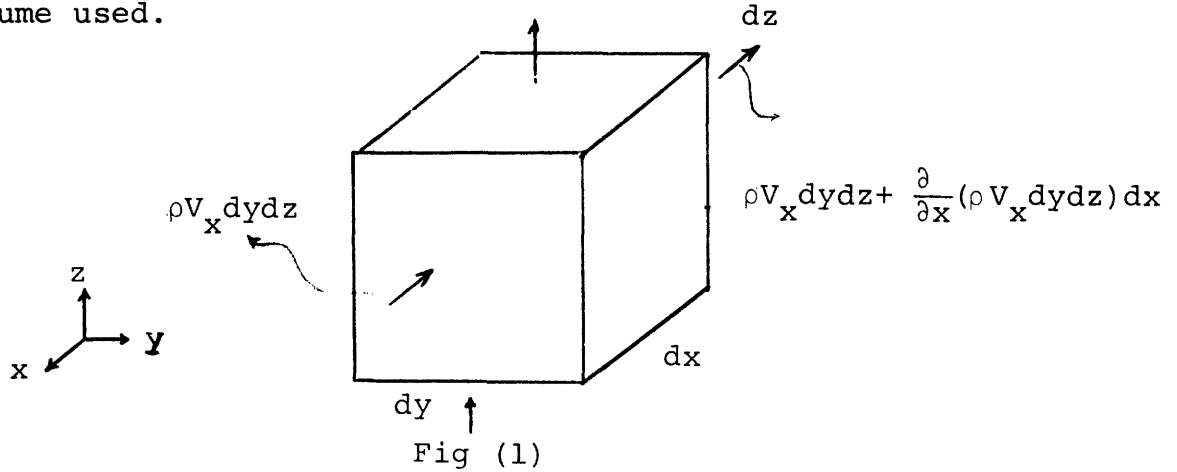
Now using the "averaged values" introduced in (3) in the equation (4), we will come up with the single phase mass balance

$$\frac{\partial}{\partial t} (\bar{\rho}) + \nabla \cdot (\bar{\rho}\vec{V}) = 0 \quad (5)$$

Equation (5) in a more involved form becomes:

$$\frac{\partial}{\partial t}(\bar{\rho}dx dy dz) + \frac{\partial}{\partial x}(\rho \bar{V} dx dz) dx + \frac{\partial}{\partial y}(\rho \bar{V} dx dz) dy + \frac{\partial}{\partial z}(\rho \bar{V} dx dy) dz = 0 \quad (6)$$

where the element of volume, $dx dy dz$, represents the control volume used.



Now in equation (6) by eliminating dz we'll have:

$$\frac{\partial}{\partial t}(\bar{\rho} dx dy) + \frac{\partial}{\partial z}(\rho \bar{V} dx dy) + \frac{\partial}{\partial y}(\rho \bar{V} dx dy) + \frac{\partial}{\partial x}(\rho \bar{V} dy dx) = 0 \quad (7)$$

In equation (7) if we assume that the control volume shown in Fig. (1) has only an infinitesimal height, dz , and its cross section normal to the z -axis "finite" instead of "infinitesimal" (i.e., $dx dy \rightarrow A$, we will come up with:

$$\frac{\partial}{\partial t}(\bar{\rho} A) + \frac{\partial}{\partial z}(\rho \bar{V} A) + \frac{\partial}{\partial y}(\rho \bar{V} A) + \frac{\partial}{\partial x}(\rho \bar{V} A) = 0 \quad (8)$$

and by substituting $\rho \bar{V} A = m$ we have:

$$\frac{\partial}{\partial t}(\bar{\rho} A) + \frac{\partial}{\partial z}(m) + \frac{\partial}{\partial y}(m) + \frac{\partial}{\partial x}(m) = 0 \quad (9)$$

A "subchannel control volume" concept used in COBRA assumes a finite cross section normal to the z -axis which

has the same area, A, equal to the physical subchannel size, and an infinitesimal height, dz.[†]

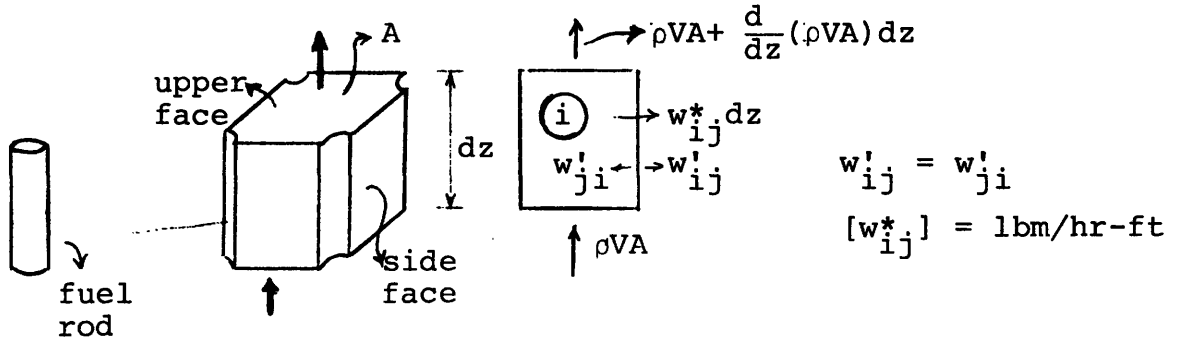


Fig (2)

In Fig. (3), w^*_{ij} represents the "cross flow" concept. Mass balance written for control volume shown in Fig. (3) becomes:

$$\rho VA + \frac{d}{dz}(\rho VA)dz + w^*_{ij}dz - \rho VA = - \frac{d}{dt}(\rho Adz)$$

$$\frac{d}{dt}(\rho A) + \frac{d}{dz}(\rho VA) + w^*_{ij} = 0 \quad m = \rho VA \quad (10)$$

Comparing equation (10), used in HEM versions of COBRA, with the pseudo single phase continuity equation, equation (9), we conclude that the cross flow term w^*_{ij} represents the two dimensional form of mass flow rate per unit length, i.e.,

$$\left(\frac{\partial}{\partial x} + \frac{\partial}{\partial y}\right)m \leftrightarrow w^*_{ij} \quad (11)$$

Flow Blockage:

This phenomena is only considered in COBRA-IV-I. None of the remaining HEM versions of COBRA account for this

[†]This infinitesimal length becomes the axial interval between the selected axial mesh points in the finite difference solution technique.

feature. Therefore in these versions, i.e. COBRA III and COBRA III-C and ..., equation (10) is simplified by assuming

$$\frac{\partial A}{\partial t} = 0$$

$$A \frac{dp}{dt} + \rho \frac{dA}{dt} + \frac{d}{dz}(m) + w^*_{ij} = 0$$

$$A_i \frac{\partial P_i}{\partial t} + \frac{\partial m_i}{\partial x} = - \sum_{j=1}^N w^*_{ij} \quad (12)$$

where the subscript i represents the subchannel under consideration.

Rewriting again equation (11):

$$A \frac{d}{dt} \bar{\rho} + \frac{d}{dz}(\bar{\rho}\bar{V})A + \sum_{j=1}^N w^*_{ij} = 0 \quad (11)$$

where in this equation, the averaged values are emphasized by using a (-) sign. This in fact follows using an integral form of mass equation which reads

$$\frac{\partial}{\partial t} \int_{\tau} \rho d\tau + \int_{s} \rho (\vec{v} \cdot \vec{n}) dA = 0 \quad (12)$$

where τ , s and \vec{v} represent volume, surface and velocity respectively.

Define the volume and surface averaged values for density and mass flux:

$$\bar{\rho} = \langle\langle \rho \rangle\rangle = \frac{1}{V} \int_{\tau} \rho d\tau \quad (13)$$

$$\bar{\rho}\bar{V} = \langle \rho V \rangle = \frac{1}{A} \int_{s} \rho (\vec{v} \cdot \vec{u}) dA$$

where $V = \int_{\tau} d\tau$ and $A = \int_{s} dA$. Also one bracket show area averaged and two bracket shows a volume averaged value.

Evaluating equation (12) for the control volume shown in Fig. 2 we have

$$V \frac{\partial}{\partial t} \bar{\rho} + \bar{\rho} \bar{V} A |_{\text{upper face}} - \bar{\rho} \bar{V} A |_{\text{lower face}} + (\sum_j w_{ij}^*) \Delta z = 0 \quad (14)$$

where $\sum_j w_{ij}^* = \frac{1}{\Delta z} \int_{\text{side faces}} \rho (\vec{v} \cdot \vec{n}) dA$ and $V = A \cdot \Delta z$. (15)

If Δz becomes small, in limit we will have equation (11).

APPENDIX 2

The details of the correlations introduced in Tables 3-1 to 3-3 are presented here.

1. Pre-CHF Heat Transfer Correlations

The heat transfer correlations used in the pre-CHF heat transfer regimes are as follows:

Dittus-Bolter (A-1) correlation:

$$h = 0.023 \left(\frac{GD_h}{\mu} \right)^{0.8} \cdot \left(\frac{C_p \mu}{k} \right)^{0.4} \cdot \left(\frac{k}{D_h} \right)$$

Data Base:

$$Re = \frac{GD_h}{\mu} \text{ greater than } 10,000$$

$$Pr = \frac{C_p \mu}{k} \text{ } 0.7 \text{ to } 100$$

$$\frac{L}{D_h} \text{ greater than } 50$$

Sieder-Tate (A-2) correlation:

$$h = 0.023 \left(\frac{GD_h}{\mu} \right)^{0.8} \cdot \left(\frac{C_p \mu}{k} \right)^{0.4} \cdot \left(\frac{k}{D_h} \right) \cdot \left(\frac{\mu}{\mu_w} \right)^{0.14}$$

Data Base:

As for Dittus-Boelter correlation.

Thom correlation (A-3):

$$T_w = T_{sat} + 0.072 e^{-\frac{P}{1260 Q_w^{\frac{1}{2}}}}$$

Data Base:

Vertical upflow of water

Round tube: 0.5-in. diameter, 60-in. length
 Annulus: 0.7-in. ID, 0.9-in. OD, 12-in. length
 Pressure: 750 to 2000 Psia.
 Mass flux: 0.77×10^6 to 2.80×10^6 lbm/hr-ft²
 Heat flux: to 0.5×10^6 Btu/hr-ft²

Schrock-Grossman correlation (A-4):

$$h = 0.023 \left[\frac{\rho V (1-x) D_h}{\mu} \right]^{0.8} \left(\frac{C_p \mu}{k} \right)^{0.4} \left[2.5 \left(\frac{1}{x_{tt}} \right)^{0.75} \right] \left(\frac{k}{D_h} \right)$$

where the inverse of the Martinelli-Lockhart-Nelson Parameter for turbulent flow is

$$\frac{1}{x_{tt}} = \left(\frac{x}{1-x} \right)^{0.9} \left(\frac{\rho_{ls}}{\rho_{gs}} \right)^{0.5} \left(\frac{\mu_g}{\mu_l} \right)^{0.1}$$

Data Base:

Water in round tubes

Diameter: 0.1162 to 0.4317 in.
 Length: 14 to 50 in.
 Pressure: 42 to 505 Psia
 Mass flux: 0.175×10^6 to 3.28×10^6 lbm/hr-ft²
 Heat flux: 0.06×10^6 to 1.45×10^6 Btu/hr-ft²
 Exit quality: 0.05 to 0.57

Chen* correlation (A-5):

$$Q_w = h_{NCB} (T_w - T_{sat}) + h_c (T_w - T_f(z))$$

where

$$h_c = F(0.023)k^{0.6}G^{0.8}(1-x)^{0.8}C_p^{0.4}\mu^{-0.4}D_h^{-0.2}$$

*The values of F and S factors should be found from the corresponding graphs (see Ref. 42 of this report). However, the relations given for them here are the curve fitting which are derived by Butterworth. More details are presented in Ref. 40 of this report.

and

$$h_{N_{CB}} = S(0.00122 \frac{k^{0.79} C^{0.45} P_1^{0.49} g_c^{0.25}}{\sigma^{0.5} \mu^{0.29} h_{fg}^{0.24} \rho_g^{0.24}})$$

where

$$F = \begin{cases} 1 & X_{tt}^{-1} \leq 1.0 \\ 2.35 (X_{tt}^{-1} + 0.213)^{0.736} & X_{tt}^{-1} > 0.1 \end{cases}$$

where X_{tt} is the same parameter as introduced in the Schrock-Grossman correlation. The value of S is as follows:

$$S = \begin{cases} [1 + 0.12 (R'_{TP})^{1.14}]^{-1} & ; R'_{TP} < 32.5 \\ [1 + 0.42 (R'_{TP})^{0.78}]^{-1} & ; 32.5 \leq R'_{TP} < 70 \\ 0.1 & ; R'_{TP} > 70 \end{cases}$$

where R'_{TP} is the effective two-phase Reynolds number

$$R'_{TP} = \frac{G(1-x)D_h}{\mu} F^{1.25} (10^{-4})$$

Data Base:

See Ref. 42 of this report.

2. CHF Correlations

The critical heat flux correlations named in Table are as follows:

Babcock and Wilcox (B&W-2) (A-6) correlation:

$$Q_{CHF} = \frac{1.155 - 0.407(D_h)}{(12.71)(3.054 \times 10^6 G)^A} [(0.3702 \times 10^8)(0.59137 \times 10^6 G)^B - 0.15208 h_{fg} \cdot G]$$

where

$$A = 0.71186 + (2.0729 \times 10^{-4}) \cdot (P - 1000)$$

and

$$B = 0.834 + (6.8479 \times 10^{-4}) \cdot (P - 2000)$$

Data Base:

Vertical upflow of water in rod bundles

Heated equivalent diameter of subchannels: 0.20 to 0.50 in.

Heated length: 72 in.

Pressure: 2000 to 2400 Psia

Mass flux: 0.75×10^6 to 4.0×10^6 lbm/hr-ft²

Thermodynamic quality: -0.03 to 0.20

Uniform axial flux distribution.

Westinghouse (W-3) (A-7) correlation:

$$Q_{CHF} = \{ [2.02 - 0.430(0.001P)] + [0.172 - 0.000]P \\ \exp[18.2x - 0.00413P \cdot x] \} (1.16 - 0.87x) \\ [(0.148 - 1.6x + 0.173x|x|) (G/10^6) + 1.04] \\ [0.266 + 0.836 \exp(-3.15D_h)] \\ [0.826 + 0.0008(h_{sat} - h_i)] 10^6 .$$

Data Base:

Diameter: 0.2 to 0.7 in.

Length: 10 to 144 in.

Pressure: 1000 to 2400 Psia

Mass flux: 1×10^6 to 5×10^6 lbm/hr-ft²

Quality: less than 0.15

Barnett Correlation (A-8):

$$Q_{CHF} = 10^6 \frac{A(h_{fg}/649) + B(h_{ls} - h_i)}{C - Z}$$

where

$$A = 67.45 D_h^{0.08} (G \times 10^{-6})^{0.192} \{1 - 0.744 \exp[-0.512 D_{hy} (G \times 10^{-6})]\}$$

$$B = 1.85 \cdot D_h^{1.261} (G \times 10^{-6})^{-0.0817}$$

$$C = 185 \cdot D_{hy}^{1.415} (G \times 10^{-6})^{0.212}$$

For Annuli the heated and wetted equivalent diameters,

D_h and D_{hy} , are given by

$$D_{hy} = (D_s - D_I)$$

and

$$D_h = (D_s^2 - D_I^2) / D_I$$

where D_s is the diameter of the shroud and D_I is the diameter of the inner rod.

Data Base:

Vertical upflow of water in annuli geometry

Diameter of inner rod: 0.375 to 3.798 in.

Diameter of shroud: 0.551 to 4.006 in.

Heated length: 24.0 to 108.0 in.

Mass flux ($G \times 10^{-6}$): 0.140 to 6.20 lbm/hr-ft²

Inlet subcooling: 0 to 412 Btu/lbm

Inlet Pressure: 1000 Psia

Uniform axial heat flux.

Biasi correlation (A-9):

The critical heat flux is given as the higher of the two values from the following equations.

For the low quality region

$$Q_{CHF} = \frac{1.883 \times 10^3}{D_{hy}^n G^{1/6}} \left[\frac{f(p)}{G^{1/6}} - x \right]$$

For the high quality region

$$Q_{CHF} = \frac{1.78 \times 10^3 S(P)}{D_{hy}^n G^{0.6}} (1-x)$$

where

$$n = 0.4 \text{ for } D_{hy} \geq 1 \text{ cm}$$

$$n = 0.6 \text{ for } D_{hy} \leq 1 \text{ cm}$$

$$f(p) = 0.7249 + 0.099P \exp(-0.032P)$$

$$s(p) = 1.155 + 0.149P \exp(0.019P) + \frac{0.889P}{10+P^2}$$

Data Base:

Diameter: 0.3 to 3.75 cm

Length: 20 to 600 cm

Pressure: 2.7 to 140 bar

Mass flux: 10 to 600 g/cm²-s

Quality: 1/(1+ρ_l/ρ_g) to 1

VOID-CHF correlation (A-10):

$$Q_{CHF} = (1-\alpha) 0.9\pi(24)^{-1} h_{fg} \rho_g^{0.5} [gg_c \sigma (\rho_l - \rho_g)]^{0.25}$$

Data Base:

See Ref. A-10.

3. Post-CHF Heat Transfer Correlations

The transition boiling, film boiling region and thermal radiation heat transfer coefficients introduced in Table 3-3 are as follows:

McDonough-Milich and King correlation (A-11):

a) As used in RELAP

$$Q = Q_{CHF} - h(T_w - T_{w,CHF})$$

where h is dependent on Pressure as follows:

P	h
2000	979.2
1200	1180.8
800	1501.2

Data Base:

Vertical upflow of water in round tubes

Diameter: 0.152 in.

Length: 12.5 in.

Mass flux: 0.2×10^6 to 1.4×10^6 lbm/ft²-hr

Wall temperature: less than 1030°F

Pressure: 800, 1200 and 2000 Psia

b) As used in RETRAN

For pressure greater than 1200 Psia,

$$h = h|_{1200} - (h|_{1200} - h|_{2000}) \left(\frac{P-1200}{800} \right)$$

and for $P \leq 1200$ Psia,

$$h = h|_{1200} + (h|_{800} - h|_{1200}) \left(\frac{1200-P}{400} \right)$$

Bjornard-Griffith correlation (A-12):

$$h = \frac{Q_{T.B.}}{T_w - T_1}$$

where

$$Q_{T.B.} = \epsilon Q_{CHF} + (1 - \epsilon) Q_{MSFB}$$

where

$$\epsilon = [(T_w - T_{MSFB}) / (T_{CHF} - T_{MSFB})]^2$$

Data Base (see Ref. A-12).

Dougall-Rohsenow correlation (A-13)

$$h = 0.023 \left[\left(\frac{DG}{\mu_g} \right) \left(\frac{\rho_g}{\rho_l} (1-x) + x \right) \right]^{0.8} [P_{rg}]^{0.4} \left(\frac{kg}{D} \right)$$

The physical properties are evaluated at saturation conditions. If $n \leq 0.0$, the term $\left[\frac{\rho_g}{\rho_l} (1-x) + x \right]$ is taken equal to 1.0 which causes the correlation reduces to the Dittus Boelter correlation.

Data Base:

See Ref. A-13 .

Groeneveld 5.7 correlation (A-14):

$$h = 0.052 \frac{k_g}{D_h} \left[\frac{GD_h}{\mu_g} \left[x + \frac{\rho_g}{\rho_l} (1-x) \right] \right]^{0.688} P_{rw}^{1.06} Y^{-1.06}$$

and the modified Groeneveld 5.7 correlations is given by

$$h = 0.052 \frac{k_g}{D_h} \left[\frac{G \cdot D_e \cdot x}{\mu_g \alpha_{dfm}} \right]^{0.688} P_{rw}^{1.06} Y^{-1.06}$$

where

$$Y = 1 - 0.1(\rho_l \rho_g^{-1} - 1)^{0.4} (1-x)^{0.4}$$

Data Base

Diameter: 0.06 - 0.25 in.

Pressure: 500 - 1400 Psia

Mass flux: $6 \times 10^6 - 3 \times 10^6$ lbm/hr-ft²

Quality: 0.1 - 0.9

Heat flux: $1.4 \times 10^5 - 7 \times 10^5$ Btu/hr-ft²

Radiation heat transfer coefficient

$$h = \sigma F (T_w^4 - T_l^4) / (T_w - T_l)$$

where

$$F = \frac{1}{\frac{1}{\epsilon} + \frac{1}{\alpha} - 1}$$

ϵ = emissivity of the wall, and

α = absorptivity of the coolant

NOMENCLATURE FOR APPENDIX 2

C_p :	specific heat
D_h :	heated equivalent diameter
D_{hy} :	wetted equivalent diameter
h :	heat transfer coefficient
h_{fg} :	heat of vaporization
k :	thermal conductivity (evaluated at bulk temp. of coolant)
L :	length
P :	pressure
Q_w :	surface heat flux
T_{sat} :	saturation temperature
T_w :	wall temperature
x :	quality

Greek

ρ_l :	density of the liquid phase
ρ_g :	density of the vapor phase
ρ_{l_s} :	density of saturated liquid
ρ_{g_s} :	density of saturated vapor
σ :	surface tension
μ :	viscosity
μ_l :	viscosity of the liquid phase
μ_g :	viscosity of the vapor phase

Appendix 2 - References

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