A new form of governing equations of fluids arising from Hamilton’s principle
Sergey Gavrilyuk, Henri Gouin

To cite this version:

HAL Id: hal-00204738
https://hal.archives-ouvertes.fr/hal-00204738
Submitted on 15 Jan 2008

HAL is a multi-disciplinary open access archive for the deposit and dissemination of scientific research documents, whether they are published or not. The documents may come from teaching and research institutions in France or abroad, or from public or private research centers.

L’archive ouverte pluridisciplinaire HAL, est destinée au dépôt et à la diffusion de documents scientifiques de niveau recherche, publiés ou non, émanant des établissements d’enseignement et de recherche français ou étrangers, des laboratoires publics ou privés.
A new form of governing equations of fluids arising from Hamilton’s principle

S. Gavrilyuk\textsuperscript{a} and H. Gouin\textsuperscript{a}

\textsuperscript{a}Laboratoire de Modélisation en Mécanique et Thermodynamique, Faculté des Sciences, Université d’Aix - Marseille III, Case 322, Avenue Escadrille Normandie-Niemen, 13397 Marseille Cedex 20, FRANCE

Abstract

A new form of governing equations is derived from Hamilton’s principle of least action for a constrained Lagrangian, depending on conserved quantities and their derivatives with respect to the time-space. This form yields conservation laws both for non-dispersive case (Lagrangian depends only on conserved quantities) and dispersive case (Lagrangian depends also on their derivatives). For non-dispersive case the set of conservation laws allows to rewrite the governing equations in the symmetric form of Godunov-Friedrichs-Lax. The linear stability of equilibrium states for potential motions is also studied. In particular, the dispersion relation is obtained in terms of Hermitian matrices both for non-dispersive and dispersive case. Some new results are extended to the two-fluid non-dispersive case.

Key words: Hamilton’s principle; Symmetric forms; Dispersion relations

1 Introduction

Hamilton’s principle of least action is frequently used in conservative fluid mechanics [1-3]. Usually, a given Lagrangian $\Lambda$ is submitted to constraints representing conservation in the time-space of collinear vectors $j_k$

$$\text{Div} j_k = 0, \quad k = 0, \ldots, m$$  \hspace{1cm} (1.1)

where Div is the divergence operator in the time-space. Equation (1.1) means the conservation of mass, entropy, concentration, etc.

Lagrangian appears as a function of $j_k$ and their derivatives. To calculate the variation of Hamilton’s action we don’t use Lagrange multipliers to take into account the constraints (1.1). We use the same method as Serrin in [2] where...
the variation of the density \( \rho \) is expressed directly in terms of the virtual displacement of the medium. This approach yields an antisymmetric form for the governing equations

\[
\sum_{k=0}^{m} j_k \left( \frac{\partial K_k}{\partial z} - \left( \frac{\partial K_k}{\partial z} \right)^* \right) = 0 \tag{1.2}
\]

where \( z \) is the time-space variable, "star" means the transposition, and \( K_k \) is the variational derivative of \( \Lambda \) with respect to \( j_k \)

\[
K_k^* = \frac{\delta \Lambda}{\delta j_k} \tag{1.3}
\]

Equations (1.1) - (1.3) admit particular class of solutions called potential flows

\[
j_k = a_k j_0, \quad a_k = \text{const}, \quad k = 1, \ldots, m
\]

\[
K_0^* = \frac{\partial \varphi_0}{\partial z}
\]

where \( \varphi_0 \) is a scalar function. We shall study the linear stability of constant solutions for potential flows.

In Section 2, we present the variations of unknown quantities in terms of virtual displacements of the continuum. In Section 3, we obtain the governing system (1.2) using Hamilton’s principle of least action. Conservation laws admitted by the system (1.1) - (1.3) are obtained in section 4. For non-dispersive case, we obtain the equations (1.1) - (1.3) in the symmetric form of Godunov-Friedrichs-Lax [4,5]. Section 5 is devoted to the linear stability of equilibrium states (constant solutions) for potential motions. We obtain dispersion relations in terms of Hermitian matrices both for dispersive and non-dispersive flows and propose simple criteria of stability. In Section 6-7, we generalize our results for two-fluid mixtures in the non-dispersive case. Usually, the theory of mixtures considers two different cases of continuum media. In homogeneous mixtures such as binary gas mixtures, each component occupies the whole volume of a physical space. In heterogeneous mixtures such as a mixture of incompressible liquid containing gas bubbles, each component occupies a part of the volume of a physical space. We do not distinguish the two cases because they have the same form for the governing equations. We obtain a simple stability criterion (criterion of hyperbolicity) for small relative velocity of phases. In Appendix, we prove non-straightforward calculations.

Recall that "star" denotes conjugate (or transpose) mapping or covectors (line vectors). For any vectors \( \mathbf{a}, \mathbf{b} \) we shall use the notation \( \mathbf{a}^* \mathbf{b} \) for their scalar product (the line vector is multiplied by the column vector) and \( \mathbf{ab}^* \) for their tensor product (the column vector is multiplied by the line vector). The product of a mapping \( A \) by a vector \( \mathbf{a} \) is denoted by \( A \mathbf{a} \), \( \mathbf{b}^* A \) means
covector $c^*$ defined by the rule $c^* = (A^*b)^*$. The divergence of a linear transformation $A$ is the covector $\text{Div} A$ such that, for any constant vector $a$,

$$\text{Div} (A)a = \text{Div} (Aa)$$

The identical transformation is denoted by $I$. For divergence and gradient operators in the time-space we use respectively symbols $\text{Div}$ and $\frac{\partial}{\partial t}$, where $z^* = (t, x^*)$, $t$ is the time and $x$ is the space. The gradient line (column) operator in the space is denoted by $\nabla (\nabla^*)$, and the divergence operator in the space by $\text{div}$. The elements of the matrix $A$ are denoted by $a^j_i$ where $i$ means lines and $j$ columns. If $f(A)$ is a scalar function of $A$, matrix $B^* \equiv \frac{\partial f}{\partial A}$ is defined by the formula

$$(B^*)_i^j = \left( \frac{\partial f}{\partial A} \right)_i^j = \frac{\partial f}{\partial a^j_i}$$

The repeated latin indices mean summation. Index $\alpha = 1, 2$ refers to the parameters of components densities $\rho_\alpha$, velocities $u_\alpha$, etc.

2 Variations of a continuum

Let $z = \begin{pmatrix} t \\ x \end{pmatrix}$ be Eulerian coordinates of a particle of a continuum and $D(t)$ a volume of the physical space occupied by a fluid at time $t$. When $t$ belongs to a finite interval $[t_0, t_1]$, $D(t)$ generates a four-dimensional domain $\Omega$ in the time-space. A particle is labelled by its position $X$ in a reference space $D_0$. For example, if $D(t)$ contains always the same particles $D_0 = D(t_0)$, and we can define the motion of a continuum as a diffeomorphism from $D(t_0)$ into $D(t)$

$$x = X(t, X) \quad (2.1)$$

We generalize (2.1) by defining the motion as the diffeomorphism from the reference space $\Omega_0$ into the time-space $\Omega$ occupied by the medium in the following parametric form

$$\begin{cases} t = g(\lambda, X) \\ x = \phi(\lambda, X) \end{cases} \quad (2.2)$$

where $Z = \begin{pmatrix} \lambda \\ X \end{pmatrix}$ belongs to a reference space $\Omega_0$. The mappings

$$\begin{cases} \lambda = h(t, x) \\ X = \psi(t, x) \end{cases} \quad (2.3)$$
are the inverse of (2.2). Definitions (2.2) imply the following expressions for the differentials $dt$ and $dx$

\[
\begin{pmatrix}
  dt \\
  dx
\end{pmatrix} = B \begin{pmatrix}
  d\lambda \\
  dX
\end{pmatrix}
\]

(2.4)

where

\[
B = \begin{pmatrix}
  \frac{\partial g}{\partial \lambda}, & \frac{\partial g}{\partial X} \\
  \frac{\partial \phi}{\partial \lambda}, & \frac{\partial \phi}{\partial X}
\end{pmatrix}
\]

Formulae (2.2), (2.4) assume the form

\[
\begin{cases}
  dt = \frac{\partial g}{\partial \lambda} d\lambda + \frac{\partial g}{\partial X} dX \\
  dx = \frac{\partial \phi}{\partial \lambda} d\lambda + \frac{\partial \phi}{\partial X} dX
\end{cases}
\]

(2.5)

From equation (2.5) we obtain

\[
dx = u \, dt + F \, dx
\]

where velocity $u$ and deformation gradient $F$ are defined by

\[
u = \frac{\partial \phi}{\partial \lambda} \left( \frac{\partial g}{\partial \lambda} \right)^{-1}, \quad F = \frac{\partial \phi}{\partial X} - \frac{\partial \phi}{\partial \lambda} \frac{\partial g}{\partial X} \left( \frac{\partial g}{\partial \lambda} \right)^{-1}
\]

(2.6)

Let

\[
\begin{cases}
  t = G(\lambda, X, \varepsilon) \\
  x = \Phi(\lambda, X, \varepsilon)
\end{cases}
\]

(2.7)

be a one-parameter family of virtual motions of the medium such that

\[
G(\lambda, X, 0) = g(\lambda, X), \quad \Phi(\lambda, X, 0) = \phi(\lambda, X)
\]

where $\varepsilon$ is a scalar defined in the vicinity of zero. We define Eulerian displacement $\zeta = (\tau, \xi)$ associated with the virtual motion (2.7)

\[
\tau = \frac{\partial G}{\partial \varepsilon}(\lambda, X, 0), \quad \xi = \frac{\partial \Phi}{\partial \varepsilon}(\lambda, X, 0)
\]

(2.8)

We note that $\zeta$ is naturally defined in Lagrangian coordinates. However, we shall suppose that $\zeta$ is represented in Eulerian coordinates by means of (2.3).

Let us now consider any tensor quantity represented by $f(t, x)$ in Eulerian coordinates and $\circ f(\lambda, X)$ in Lagrangian coordinates. Definitions (2.2), (2.3) involve

\[
\circ f(\lambda, X) = f\left( g(\lambda, X), \phi(\lambda, X) \right)
\]

(2.9)
Conversely,
\[ f(t, x) = \hat{f}\left(h(t, x), \psi(t, x)\right) \quad (2.10) \]

Let \( \tilde{f}(\lambda, X, \varepsilon) \) and \( \hat{f}(t, x, \varepsilon) \) be tensor quantities associated with the virtual motions, such that
\[ \tilde{f}(\lambda, X, \varepsilon) \equiv \hat{f}(t, x, \varepsilon) \]
where \( \lambda, X, t, x \) are connected by relations (2.7) satisfying
\[ \tilde{f}(\lambda, X, 0) = \hat{f}(\lambda, X) \]
or equivalently \( \hat{f}(t, x, 0) = f(t, x) \). We then obtain
\[ \tilde{f}(\lambda, X, \varepsilon) = \hat{f}\left(G(\lambda, X, \varepsilon), \Phi(\lambda, X, \varepsilon), \varepsilon\right) \quad (2.11) \]

Let us define Eulerian and Lagrangian variations of \( f \)
\[ \hat{\delta} f = \frac{\partial \hat{f}}{\partial \varepsilon}(t, x, 0) \quad \text{and} \quad \tilde{\delta} f = \frac{\partial \tilde{f}}{\partial \varepsilon}(\lambda, X, 0) \]

Differentiating relation (2.11) with respect to \( \varepsilon \) at \( \varepsilon = 0 \), we get
\[ \hat{\delta} f = \tilde{\delta} f - \frac{\partial f}{\partial z} \zeta \quad (2.12) \]

### 3 Governing equations

Consider a four-dimensional vector \( j_0 \) satisfying conservation law
\[ \text{Div} j_0 = 0 \quad (3.1) \]

Actually, (3.1) represents the mass conservation law, where \( j_0 = \rho v^* \), \( \rho \) is the density and \( v^* = (1, u^*) \) is the four-dimensional velocity vector. Let \( a_k \) be scalar quantities such as the specific entropy, the number of bubbles per unit mass, the mass concentration, etc., which are conserved along the trajectories. Consequently, if \( j_k = a_k j_0 \),
\[ \text{Div} j_k = \frac{\partial a_k}{\partial z} j_0 = 0, \quad k = 1, \ldots, m \quad (3.2) \]

Hence \( j_k, k = 1, \ldots, m \) form a set of solenoidal vectors collinear to \( j_0 \). Hamilton’s principle needs the knowledge of Lagrangian of the medium. We take the Lagrangian in the form
\[ L = \Lambda(j_k, \frac{\partial j_k}{\partial z}, \ldots, \frac{\partial^n j_k}{\partial z^n}, z) \quad (3.3) \]
where \( j_k \) are submitted to the constraints (3.1), (3.2) rewritten as
\[
\text{Div} j_k = 0, \quad k = 0, \ldots, m
\]  
(3.4)

Let us consider three examples.

**a – Gas dynamics** [1-3]

Lagrangian of the fluid is
\[
L = \frac{1}{2} \rho |u|^2 - \varepsilon (\rho, \eta) - \rho \Pi (z),
\]
where \( \varepsilon \) is the internal energy per unit volume, \( \eta = \rho s \) is the entropy per unit volume, \( s \) is the specific entropy and \( \Pi \) is an external potential. Hence, in variables \( j_0 = \rho v, \quad j_1 = \rho s v, \quad z \), Lagrangian takes the form
\[
L = \frac{1}{2} \left( \frac{|j_0|^2}{l^* j_0} - l^* j_0 \right) - \varepsilon (l^* j_0, l^* j_1) - l^* j_0 \Pi (z) = \Lambda (j_0, j_1, z)
\]
where \( l^* = (1, 0, 0, 0) \).

**b – Thermocapillary fluids** [6,7]

Lagrangian of the fluid is
\[
L = \frac{1}{2} \rho |u|^2 - \varepsilon (\rho, \nabla \rho, \eta, \nabla \eta) - \rho \Pi (z).
\]
Since \( \nabla \rho = \nabla (l^* j_0) \) and \( \nabla \eta = \nabla (l^* j_1) \), we obtain Lagrangian in the form (3.3)
\[
L = \Lambda \left( j_0, \frac{\partial j_0}{\partial z}, j_1, \frac{\partial j_1}{\partial z}, z \right)
\]

**c – One-velocity bubbly liquids** [8-10]

\[
L = \frac{1}{2} \rho |u|^2 - W \left( \rho, \frac{d\rho}{dt}, N, \right), \quad \text{with} \quad \frac{d\rho}{dt} = \frac{\partial \rho}{\partial z} v
\]
where \( \rho \) is now the average density of the bubbly liquid and \( N \) is the number of identical bubbles per unit volume of the mixture. We define again \( j_0 = \rho v \) and \( j_1 = N v \). By using \( \frac{d\rho}{dt} = \frac{\partial \rho}{\partial z} v = \frac{\partial (l^* j_0)}{\partial z} \frac{j_0}{l^* j_0} \) and \( N = l^* j_1 \), we obtain Lagrangian in the form
\[
L = \Lambda \left( j_0, \frac{\partial j_0}{\partial z}, j_1 \right)
\]

The Hamilton principle reads: for each field of virtual displacements \( z \in \Omega \to \zeta \) such that \( \zeta \) and its derivatives are zero on \( \partial \Omega \),
\[
\delta \int_\Omega \Lambda \ d\Omega = 0
\]  
(3.5)
Since variation $\hat{\delta}$ is independent of domain $\Omega$ and measure $d\Omega$, the variation of Hamilton action (3.5) with the zero boundary conditions for $j_k$ and its derivatives yields

$$\delta \int_\Omega \Lambda \ d\Omega = \int_\Omega \sum_{k=0}^{m} \frac{\delta \Lambda}{\delta j_k} \hat{\delta} j_k \ d\Omega = 0$$  \hspace{1cm} (3.6)$$

where $\frac{\delta}{\delta j_k}$ refers to the variational derivative with respect to $j_k$. In particular, if Lagrangian (3.3) is

$$L = \Lambda \left( j_k, \frac{\partial j_k}{\partial z}, z \right)$$

we get

$$\frac{\delta \Lambda}{\delta j_k} = \frac{\partial \Lambda}{\partial j_k} - \text{Div} \left( \frac{\partial \Lambda}{\partial \left( \frac{\partial j_k}{\partial z} \right)} \right)$$

We have to emphasize that in (3.6), the variations $\hat{\delta} j_k$ should take into account constraints (3.4). We use the same method as in [2, p. 145] for the variation of density. This method does not use Lagrange multipliers since the constraints (3.4) are satisfied automatically. The calculation of $\hat{\delta} j_k$, $k = 0, \ldots, m$ is performed in two steps. First, in Appendix A we calculate Lagrangian variations $\hat{\delta} v$, $\hat{\delta} \rho$ and $\hat{\delta} a_k$ (expressions (A.2), (A.5) and (A.6), respectively). Second, by using (2.12) we obtain in Appendix B Eulerian variations $\hat{\delta} j_k$, $k = 0, \ldots, m$ (see (B.3))

$$\hat{\delta} j_k = \left( \frac{\partial \zeta}{\partial z} - (\text{Div} \ z) I \right) j_k - \frac{\partial j_k}{\partial z} \zeta$$

Let us now define the four-dimensional covector

$$K^*_k = \frac{\delta \Lambda}{\delta j_k}$$  \hspace{1cm} (3.7)$$

Taking into account conditions (3.4) and the fact that $\zeta$ and its derivatives are zero on the boundary $\partial \Omega$, equations (3.6), (B.3) yield

$$\delta \int_\Omega \Lambda \ d\Omega = \int_\Omega \sum_{k=0}^{m} K^*_k \left( \left( \frac{\partial \zeta}{\partial z} - (\text{Div} \ z) I \right) j_k - \frac{\partial j_k}{\partial z} \zeta \right) \ d\Omega$$

$$= \int_\Omega \sum_{k=0}^{m} \left( -\text{Div} \left( j_k K^*_k \right) + j_k \frac{\partial K^*_k}{\partial z} \right) \zeta \ d\Omega$$

$$= \int_\Omega \sum_{k=0}^{m} j_k \left( \frac{\partial K^*_k}{\partial z} - \left( \frac{\partial K^*_k}{\partial z} \right)^* \right) \zeta \ d\Omega = 0$$

7
Hamilton’s principle yields the governing equations in the form
\[
\sum_{k=0}^{m} j_k^* \left( \frac{\partial K_k}{\partial z} - \left( \frac{\partial K_k}{\partial z} \right)^* \right) = 0 \tag{3.8}
\]
where \( K_k \) are given by definition (3.7). The system (3.4), (3.8) represents \( m + d + 1 \) partial differential equations for \( m + d \) unknown functions \( u, \rho, a_k, k = 1, \ldots, m \), where \( d \) is the dimension of the \( \Omega \)-space. Since the matrix
\[
R_k = \frac{\partial K_k}{\partial z} - \left( \frac{\partial K_k}{\partial z} \right)^*
\tag{3.9}
\]
is antisymmetric and all the vectors \( j_k, k = 0, \ldots, m \) are collinear, we obtain that \( j_k^* R_k j_0 \equiv 0 \).

Consequently
\[
\sum_{k=0}^{m} j_k^* R_k j_0 \equiv 0 \tag{3.10}
\]
and the overdetermined system (3.4), (3.8) is compatible.

In the case \( a_k = \text{const}, k = 1, \ldots, m \) and \( \Lambda = \Lambda \left( j_0, \frac{\partial j_0}{\partial z} \right) \), the system (3.4), (3.8) can be rewritten in a simplified form
\[
\begin{cases}
\sum_{k=0}^{m} j_0^* \left( \frac{\partial K_0}{\partial z} - \left( \frac{\partial K_0}{\partial z} \right)^* \right) = 0 \\
\text{Div} j_0 = 0
\end{cases} \tag{3.11}
\]

We call potential motion such solutions of (3.11) that
\[
K_0^* = \frac{\partial \varphi_0}{\partial z} \tag{3.12}
\]
where \( \varphi_0 \) is a scalar function. In this case, the governing system for the potential motion is in the form
\[
\begin{cases}
\frac{\delta \Lambda}{\delta j_0} = \frac{\partial \varphi_0}{\partial z} \\
\text{Div} j_0 = 0
\end{cases} \tag{3.13}
\]
For the gas dynamics model, the system (3.12), (3.13) reads
\[
\begin{align*}
\frac{\partial \varphi_0}{\partial t} &= -\left( \frac{1}{2} |u|^2 + \frac{\partial}{\partial \rho} e(\rho) + \Pi \right), \quad e(\rho) = \varepsilon (\rho, \rho s_e) \\
\nabla \varphi_0 &= u^* \\
\frac{\partial \rho}{\partial t} + \text{div}(\rho u) &= 0
\end{align*}
\]
Eliminating the derivative $\frac{\partial \varphi_0}{\partial t}$, we obtain the classical model for potential flows \[2\]
$$
\begin{align*}
\frac{\partial u^*}{\partial t} + \nabla \left( \frac{1}{2} |u|^2 + \frac{\partial}{\partial \rho} \epsilon(\rho) + \Pi \right) &= 0 \\
u^* &= \nabla \varphi_0 \\
\frac{\partial \rho}{\partial t} + \text{div}(\rho u) &= 0
\end{align*}
$$

(3.14)

4 Conservation laws

Equations (3.4), (3.8) can be rewritten in a divergence form. The demonstration is performed for Lagrangian $\Lambda$ which depends only on $j_k$, $\frac{\partial j_k}{\partial z}$ and $z$. The following result is proved in Appendix C.

Theorem 4.1

Let $L = \Lambda \left( j_k, \frac{\partial j_k}{\partial z}, z \right)$. The following vector relation is an identity

$$
\text{Div} \left( \sum_{k=0}^{m} \left( K_k^* j_k I - j_k K_k^* + A_k^* \frac{\partial j_k}{\partial z} \right) - \Lambda I \right) + \frac{\partial \Lambda}{\partial z} + \sum_{k=0}^{m} K_k^* \text{Div} j_k - \sum_{k=0}^{m} j_k^* R_k \equiv 0
$$

(4.1)

where the matrices $R_k$ are given by (3.9) and $A_k^* = \frac{\partial \Lambda}{\partial \left( \frac{\partial j_k}{\partial z} \right)}$.

In particular, we get

Theorem 4.2

The governing equations (3.4), (3.8) are equivalent to the system of conservation laws

$$
\text{Div} \left( \sum_{k=0}^{m} \left( K_k^* j_k I - j_k K_k^* + A_k^* \frac{\partial j_k}{\partial z} \right) - \Lambda I \right) + \frac{\partial \Lambda}{\partial z} = 0
$$

$$
\text{Div} j_k = 0, \quad k = 0, \ldots, m
$$
In the special case of potential flows (3.13) the governing equations admit additional conservation laws

\[
\text{Div} \left( cK_0 - (K_0^* c) / \right) = 0 \tag{4.2}
\]

where \( c \) is any constant vector. Indeed,

\[
\text{Div} (cK_0) = c^* \left( \frac{\partial K_0}{\partial z} \right)^*, \quad \text{Div} (K_0^* c) = \frac{\partial}{\partial z} (K_0^* c) = c^* \left( \frac{\partial K_0}{\partial z} \right)
\]

Since for the potential flows

\[
\frac{\partial K_0}{\partial z} = \left( \frac{\partial K_0}{\partial z} \right)^*
\]

we obtain (4.2). In particular, if we take \( c = 1 \) for the gas dynamics equations, we get conservation laws (3.14). Multiplying (4.1) by \( j_0 \) and taking into account the identity (3.10), we obtain the following theorem

**Theorem 4.3**

The following scalar relation is an algebraic identity

\[
\left( \text{Div} \left( \sum_{k=0}^{m} (K_k^* j_k I - j_k K_k^* + A_k^* \frac{\partial j_k}{\partial z} - \Lambda I) \right) \right) j_0 + \frac{\partial \Lambda}{\partial z} j_0 + \sum_{k=0}^{m} (K_k^* j_0) \text{Div} j_k \equiv 0 \tag{4.3}
\]

This theorem is a general representation of the Gibbs identity expressing that the ”energy equation” is a consequence of the conservation of ”mass”, ”momentum” and ”entropy”. Examples of this identity for thermocapillary fluids and bubbly liquids were obtained previously in [6,10].

Identity (4.3) yields an important consequence. Let us recall the Godunov-Friedrichs-Lax method of symmetrisation of quasilinear conservation laws [4-5] (see also different applications and generalizations in [11-12]). We suppose that the system of conservation laws for \( n \) variables \( q \) has the form

\[
\frac{\partial f_i}{\partial t} + \text{div} F_i = 0, \quad f_i = f_i(q), \quad F_i = F_i(q), \quad i = 1, \ldots, n \tag{4.4}
\]

Let us also assume that (4.4) admits an additional ”energy” conservation law

\[
\frac{\partial e}{\partial t} + \text{div} E = 0, \quad e = e(q), \quad E = E(q)
\]
which is obtained by multiplying each equation of (4.4) by some functions \( p^i \) and then by summing over \( i = 1, \ldots, n \)

\[
\frac{\partial e}{\partial t} + \text{div} E \equiv p^i \left( \frac{\partial f_i}{\partial t} + \text{div} F_i \right)
\]  

(4.5)

In particular, if we consider \( e, \ E \) and \( F_i \) as functions of \( f_i \), we obtain from (4.5)

\[
\frac{\partial e}{\partial f_i} = p^i, \quad \frac{\partial E}{\partial f_j} = p^i \frac{\partial F_i}{\partial f_j} = \frac{\partial e}{\partial f_i} \frac{\partial F_i}{\partial f_j}
\]  

(4.6)

Let us introduce functions \( N \) and \( M \) such that

\[
N = f_i p^i - e, \quad M = F_i p^i - E
\]  

(4.7)

Consequently, from equations (4.6) and (4.7) we find that

\[
\frac{\partial N}{\partial p^i} = f_i, \quad \frac{\partial M}{\partial p^i} = F_i
\]  

(4.8)

Hence, substituting (4.8) into (4.4), we get a symmetric system in the form

\[
\frac{\partial^2 N}{\partial p^i \partial p^j} + \text{tr} \left( \frac{\partial^2 M}{\partial p^i \partial p^j} \frac{\partial p^j}{\partial x} \right) = 0
\]  

(4.9)

If matrix \( N_{ij} = \frac{\partial^2 N}{\partial p^i \partial p^j} \) is positive definite then the symmetric system (4.9) is \( t \)-hyperbolic symmetric in the sense of Friedrichs. Obviously, matrix \( \frac{\partial^2 N}{\partial p^i \partial p^j} \) is positive definite if and only if matrix \( e^{ij} = \frac{\partial^2 e}{\partial f^i \partial f^j} \) is positive definite, since \( e^{ij} N_{jp} = \delta^i_p \), where \( \delta^i_p \) is the Kronecker symbol.

Now, let us rewrite identity (4.3) in the non-dispersive case

\[
\left( \text{Div} \left( \sum_{k=0}^m (K_k^* j_k - j_k K_k^*) - \Lambda I \right) \right) j_0 + \sum_{k=0}^m (K_k^* j_k) \text{div} j_k \equiv 0
\]  

(4.10)

Identity (4.10) is exactly of the same type as identity (4.5). It means that the system (3.4), (3.8) can be always rewritten in a Godunov-Friedrichs-Lax symmetric form. Actually, if we take

\[
e = \sum_{k=0}^m K_k^* j_k - (j_k K_k^*)^1 \Lambda \quad \text{and} \quad E = -P j_k K^*
\]

where

\[
P = \begin{pmatrix} 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{pmatrix}
\]
the conjugate variables $p$ are

$$ p = \begin{pmatrix} u \\ -\frac{1}{\rho} K_k^* j_0 \end{pmatrix}, \quad k = 0, \ldots, m $$

Consequently, the Gibbs identity (4.10) gives directly a set of conjugate variables. Therefore, we have proved the following theorem

**Theorem 4.4**

If $L = \Lambda(j_k)$, the system (3.4), (3.8) is always symmetrisable.

The property of convexity, needed for the hyperbolicity of the governing system, should be verified for each particular case. For example, in the gas dynamics, the energy of the system is given by the formula

$$ e = \sum_{k=0}^{1} K_k^* j_k - \sum_{k=0}^{1} (j_k K_k^*)_{1} - \Lambda = \varepsilon(\rho, \eta) + \rho \frac{|u|^2}{2} $$

Since

$$ K_0^* j_0 = \rho \left( \frac{|u|^2}{2} - \frac{\partial \varepsilon}{\partial \rho} \right), \quad K_1^* j_0 = -\rho \frac{\partial \varepsilon}{\partial \eta} $$

the conjugate variables are (see also [4])

$$ p = \begin{pmatrix} u \\ \frac{\partial \varepsilon}{\partial \rho} - \frac{|u|^2}{2} \\ \frac{\partial \varepsilon}{\partial \eta} \end{pmatrix} = \begin{pmatrix} u \\ \mu - \frac{|u|^2}{2} \\ T \end{pmatrix} $$

where $\mu$ is the Gibbs potential and $T$ is the temperature. Obviously, if $\varepsilon(\rho, \eta)$ is convex, the total energy $e$ is convex with respect to $\rho u$, $\rho$ and $\eta$.

5 Stability of equilibrium states for potential flows.

We assume that $L$ is a function of $j_k$ and $\frac{\partial j_k}{\partial z}$ besides it does not depend on $z$, i.e.

$$ L = \Lambda \left( j_k, \frac{\partial j_k}{\partial z} \right) $$

12
Let us give some definitions. An equilibrium state is a solution of equations (3.4), (3.8) such that \( j_k = j_{ke} = \text{const} \). Let \( \nu \) be a real unit vector. Any vector \( \beta \) can be represented in the form
\[
\beta = \omega \nu + \beta_{\sigma}, \quad \text{where} \quad \nu^* \beta_{\sigma} = 0
\]
The equilibrium state \( j_{ke} \neq 0 \) is linearly stable in the direction \( \nu \) if and only if all non-trivial solutions of the form \( J_k e^{i\beta^* z} \) of the system (3.4), (3.8) linearized at the equilibrium state \( j_{ke} \) are such that \( \omega \) is real for any real \( \beta_{\sigma} \).

5.1 Non-dispersive case

We note that in a non-dispersive case the stability means hyperbolicity of governing system [13-14]. We omit the index “0” and rewrite system (3.13) in the form
\[
\begin{cases}
K^* = \frac{\partial \Lambda}{\partial j} = \frac{\partial \varphi}{\partial z} \\
\text{Div} j = 0
\end{cases}
(5.1)
\]
where \( \Lambda = \Lambda(j) \). The Legendre transformation of \( \Lambda(j) \) is
\[
\Delta(K) = K^* j - \Lambda(j)
(5.2)
\]
If the matrix \( \Lambda''(j) = \frac{\partial}{\partial j} \left( \left( \frac{\partial \Lambda}{\partial j} \right)^* \right) \) is non-degenerate, (5.2) involves the formula
\[
J^* = \frac{\partial \Delta}{\partial K}
(5.3)
\]
Hence, relations (5.1) - (5.3) yield
\[
\frac{\partial j}{\partial z} = \frac{\partial}{\partial K} \left( \left( \frac{\partial \Delta}{\partial K} \right)^* \right) \frac{\partial K}{\partial z} = \Delta''(K) \frac{\partial K}{\partial z} = \Delta''(K) \varphi''(z)
(5.4)
\]
where
\[
\Delta''(K) = \frac{\partial}{\partial K} \left( \left( \frac{\partial \Delta}{\partial K} \right)^* \right), \quad \varphi''(z) = \frac{\partial}{\partial z} \left( \left( \frac{\partial \varphi}{\partial z} \right)^* \right)
\]
The second equation (5.1) and (5.4) involve
\[
\text{tr} \left( \Delta''(K) \varphi''(z) \right) = 0
\]
If we replace \( \varphi \) by \( i \Phi e^{i\beta^* z} \), we get immediately the dispersion relation
\[
\beta^* \Delta''(K_e) \beta = 0
(5.5)
\]
where index “e” corresponds to the equilibrium state. The matrix $\Delta''(K_e)$ can be easily calculated in terms of Lagrangian $\Lambda(j)$. Indeed,

$$I = \frac{\partial j}{\partial j} = \frac{\partial j}{\partial K} \frac{\partial K}{\partial j} = \Delta''(K) \Lambda''(j)$$

It follows that

$$\Delta''(K_e) = \left(\Lambda''(j_e)\right)^{-1} \quad (5.6)$$

Relations (5.5) - (5.6) imply the following result.

**Theorem 5.1** (criterion of stability for potential non-dispersive motions)

*If the symmetric matrix $G = \Lambda''(j_e)$ has the signature $(-,+,+,+)$, the equilibrium state $j_e$ is stable in any direction $\nu$ belonging to the intersection of two cones $C_1 = \{\nu | \nu^* G^{-1} \nu < 0\}$ and $C_2 = \{\nu | \nu^* G \nu < 0\}$.*

**Proof.** Since $C_1$ and $C_2$ contain the eigenvector corresponding to the negative eigenvalue of $G$, $C_1 \cap C_2 \neq \emptyset$. We can find orthogonal coordinates $(t, x_1, x_2, x_3)$ such that the dispersion relation (5.5) takes the form

$$-t^2 + \sum_{i=1}^{3} \lambda_i x_i^2 = 0, \quad \text{with} \quad \lambda_i > 0$$

In coordinates $(t, x_1, x_2, x_3)$ the cone $C_2$ is defined by the inequality

$$-t^2 + \sum_{i=1}^{3} \frac{x_i^2}{\lambda_i} < 0$$

Let us represent $\beta$ in the form

$$\beta = \begin{pmatrix} t \\ x_1 \\ x_2 \\ x_3 \end{pmatrix} = \omega \begin{pmatrix} 1 \\ n_1 \\ n_2 \\ n_3 \end{pmatrix} + \begin{pmatrix} -n_1 y_1 - n_2 y_2 - n_3 y_3 \\ y_1 \\ y_2 \\ y_3 \end{pmatrix}$$

The dispersion relation $\beta^* G^{-1} \beta = 0$ implies

$$-\left(\sum_{i=1}^{3} n_i y_i - \omega\right)^2 + \sum_{i=1}^{3} \lambda_i (\omega n_i + y_i)^2 = 0$$

It involves

$$\omega^2 \left(-1 + \sum_{i=1}^{3} \lambda_i n_i^2\right) + 2\omega \sum_{i=1}^{3} (1 + \lambda_i) n_i y_i + \sum_{i=1}^{3} \lambda_i y_i^2 - \left(\sum_{i=1}^{3} n_i y_i\right)^2 = 0$$
Due to the following inequality,
\[
\sum_{i=1}^{3} \lambda_i y_i^2 - \left( \sum_{i=1}^{3} n_i y_i \right)^2 = \sum_{i=1}^{3} \lambda_i y_i^2 - \left( \sum_{i=1}^{3} \frac{n_i}{\lambda_i^{3/2}} \sqrt{\lambda_i} y_i \right)^2
\geq \sum_{i=1}^{3} \lambda_i y_i^2 - \left( \sum_{i=1}^{3} \frac{n_i^2}{\lambda_i^{3/2}} \right) \left( \sum_{i=1}^{3} \lambda_i y_i^2 \right) \left( 1 - \sum_{i=1}^{3} \frac{n_i^2}{\lambda_i} \right)
\]
\(\omega\) is real if \(\nu\) belongs simultaneously to \(C_1\) and \(C_2\). The theorem is proved.

For the gas dynamics model, if the volume energy is a convex function of \(\rho\) and \(\eta = \rho s\), the matrix \(G\) satisfies the conditions of Theorem 5.1.

5.2 Dispersive case

If \(L = A \left( j, \partial_j \frac{\partial}{\partial z} \right)\), the governing system is

\[
\begin{align*}
\mathbf{K}^* &= \frac{\delta \Lambda}{\delta j} = \frac{\partial \Lambda}{\partial j} - \text{Div} \left( \frac{\partial \Lambda}{\partial (\partial_j \frac{\partial}{\partial z})} \right) = \frac{\partial \varphi}{\partial z} \\
\text{Div} \mathbf{j} &= 0
\end{align*}
\]

The linearised system in a coordinate form is

\[
\Lambda_{ks} j^s + \Gamma_{ks,p} j^s_p - D_{mp}^{mp} j^s_{mp} = \varphi, \quad \text{and} \quad j^s_{s} = 0 \tag{5.7}
\]

where the comma denotes the derivative with respect to \(z^k\) while \(\Lambda_{ks}, \Gamma_{ks}, D_{ks}^{mp}\), calculated at point \(j_e\), are defined by the relations

\[
\Lambda_{ks} = \frac{\partial^2 \Lambda}{\partial j^k \partial j^s}, \quad \Gamma_{ks} = \frac{\partial^2 \Lambda}{\partial j^k \partial j^s_p} - \frac{\partial^2 \Lambda}{\partial j^s \partial j^k_p}, \quad D_{ks}^{mp} = \frac{\partial^2 \Lambda}{\partial j^k_m \partial j^s_p} \tag{5.8}
\]

The following symmetry relations result from (5.8)

\[
\Lambda_{ks} = \Lambda_{sk}, \quad \Gamma_{ks} = -\Gamma_{sk}, \quad D_{ks}^{mp} = D_{sk}^{mp} \tag{5.9}
\]

For the solution of (5.7) in the form \(j^s = J^s e^{\beta_k z^p}, \varphi = i \Phi e^{\beta_k z^p}\), we get the following dispersion relation

\[
\det \begin{pmatrix} G & \beta \\ \beta^* & 0 \end{pmatrix} = -\beta^* \text{Adj} (G) \beta = 0 \tag{5.10}
\]

where \(\text{Adj} (G)\) is the adjoint matrix to \(G\), and the elements \(G_{ks}\) of the matrix \(G\) are defined by

\[
G_{ks} = \Lambda_{ks} + i \Gamma_{ks,p} \beta_p + D_{ks}^{mp} \beta_m \beta_p \tag{5.11}
\]
Equations (5.9) - (5.11) imply that \( G \) is hermitian matrix. In particular, if \( \det G \neq 0 \) then \( \text{Adj}(G) = \det(G) G^{-1} \) and the dispersion relation (5.10) is equivalent to \( \beta^* G^{-1} \beta = 0 \), which is a generalization of (5.5) for the dispersive case. We get the following obvious result:

**Theorem 5.2 (criterion of stability for potential dispersive motions)**

If \( \Gamma_{ks} = 0 \) and the symmetric matrix \( G \) defined by (5.8), (5.11) has the signature \( (-, +, +, +) \) for \( \beta = 0 \), the equilibrium state \( j_e \) is stable for small \( \beta \) in any direction \( \nu \) belonging to the intersection of the cones \( C_i, i = 1, 2 \) defined in theorem 5.1.

We note that \( \Gamma_{ks} \) defined by (5.8) are always zero if the expansion of Lagrangian \( \Lambda \) in Taylor series at the vicinity of equilibrium state does not contain linear terms with respect to \( \frac{\partial j_k}{\partial z} \).

Sometimes, we are able to obtain a “global” stability. For example, let us consider a particular case of bubbly liquids, for \( N/\rho = \text{const} \), defined in Section 3

\[
L = \frac{1}{2} \rho |\mathbf{u}|^2 + \frac{a}{2} \left( \frac{d\rho}{dt} \right)^2 - \varepsilon(\rho)
\]

where \( \varepsilon(\rho) \) is a convex function of the density and \( a \) is a positive function of \( \rho \). Then,

\[
L = \frac{1}{2} \left( \frac{|\mathbf{j}|^2}{\mathbf{I}^*} - \mathbf{I}^* \mathbf{j} \right) + \frac{a}{2} \left( \frac{\partial}{\partial z}(\mathbf{I}^* \mathbf{j}) \frac{\mathbf{j}}{\mathbf{I}^*} \right)^2 - \varepsilon(\mathbf{I}^* \mathbf{j})
\]

Since the governing equations are invariant with respect to the Galilean transformation

\[
t' = t, \quad x' = x + Ut, \quad u' = u + U
\]

we can always assume that \( u_e = 0 \) (it is sufficient to consider the governing system in the reference frame moving with the velocity \( U = u_e \)). Hence, \( \mathbf{J}_e = \rho_e \mathbf{I} \), where \( \mathbf{I}^* = (1, 0, 0, 0) \). Omitting index “e”, we can calculate the matrix \( G \) defined by (5.11)

\[
G = \begin{pmatrix}
-\frac{\partial^2 \varepsilon}{\partial \rho^2} + a\beta^2 & 0^* \\
0 & \frac{\mathbf{I}}{\rho}
\end{pmatrix}
\]

Hence,

\[
\text{Adj}(G) = \begin{pmatrix}
\frac{1}{\rho^3} & 0^* \\
0 & \left( a\beta^2 - \frac{\partial^2 \varepsilon}{\partial \rho^2} \right) \frac{\mathbf{I}}{\rho^2}
\end{pmatrix}
\]
and dispersion relation (5.10) reads

\[
\frac{\beta_1^2}{\rho} + \frac{a \beta_1^2}{\rho^3} \frac{\partial^2 \varepsilon}{\partial \rho^2} \left( \beta_2^2 + \beta_3^2 + \beta_4^2 \right) = 0 \tag{5.12}
\]

Let \( \nu \) be the direction of time in the time-space, then \( \nu = 1 \) and (5.12) is equivalent to

\[
\frac{\omega^2}{\partial^2 \varepsilon / \partial \rho^2} - a \omega^2 = \rho |\beta_\sigma|^2 \tag{5.13}
\]

The graph of the dispersion relation (5.13) for positive values of \( \omega \) is presented in Figure 1. We have denoted by \( \omega_* = \sqrt{\frac{\partial^2 \varepsilon}{\partial \rho^2} / a} \) the eigenfrequency of bubbles and by \( c_0 = \sqrt{\rho \frac{\partial^2 \varepsilon}{\partial \rho^2}} \) the equilibrium sound speed of bubbly liquid (see [8]).

![Fig. 1. Dispersion relation for bubbly liquids](image)

6 Governing equations for mixtures

Let us consider homogeneous binary mixtures. The mixture is described by the velocities \( u_\alpha \), the average densities \( \rho_\alpha \) and the specific entropies \( s_\alpha \) for each component \( (\alpha = 1, 2) \). We introduce two reference frames associated with each component in the form

\[
\begin{align*}
  t &= g_\alpha(\lambda_\alpha, X_\alpha) \\
  x &= \phi_\alpha(\lambda_\alpha, X_\alpha)
\end{align*}
\tag{6.1}
\]
and the inverse mappings

\[
\begin{aligned}
\lambda_\alpha &= h_\alpha(t, x) \\
X_\alpha &= \Psi_\alpha(t, x)
\end{aligned}
\]

(6.2)

The corresponding families of virtual motions generated by (6.1) are defined by

\[
\begin{aligned}
t &= G_\alpha(\lambda_\alpha, X_\alpha, \varepsilon_\alpha) \\
x &= \Phi_\alpha(\lambda_\alpha, X_\alpha, \varepsilon_\alpha)
\end{aligned}
\]

with

\[
\begin{aligned}
G_\alpha(\lambda_\alpha, X_\alpha, 0) &= g_\alpha(\lambda_\alpha, X_\alpha) \\
\Phi_\alpha(\lambda_\alpha, X_\alpha, 0) &= \phi_\alpha(\lambda_\alpha, X_\alpha)
\end{aligned}
\]

We define the two Eulerian displacement \( \zeta_\alpha = (\tau_\alpha, \xi_\alpha) \) where

\[
\begin{aligned}
\tau_\alpha &= \frac{\partial G_\alpha}{\partial \varepsilon_\alpha}(\lambda_\alpha, X_\alpha, 0), \\
\xi_\alpha &= \frac{\partial \Phi_\alpha}{\partial \varepsilon_\alpha}(\lambda_\alpha, X_\alpha, 0)
\end{aligned}
\]

As in Section 3 we define tensor quantities associated with the two virtual motions and variations \( \delta_\alpha \) and \( \bar{\delta}_\alpha \). In the general case, we have two four-dimensional solenoidal vectors \( j_{0(\alpha)} = \rho_\alpha v_\alpha \) corresponding to the \( \alpha \)th component, where \( \rho_\alpha \) is the density and \( v_\alpha^* = (1, u_\alpha^*) \) is the four-dimensional velocity vector. As in Section 3, we introduce additional physical quantities \( a_{k(\alpha)} \) associated with the four-dimensional vectors \( j_{k(\alpha)} = a_{k(\alpha)} j_{0(\alpha)} \), \( \alpha = 1, 2, k(\alpha) = 1, \ldots, m(\alpha) \) submitted to the constraints

\[
\text{Div } j_{k(\alpha)} = 0
\]

(6.3)

For Hamilton’s action in the form

\[
a = \int_{\Omega} \Lambda \left( j_{k(\alpha)}, \frac{\partial j_{k(\alpha)}}{\partial z}, \ldots, \frac{\partial^n j_{k(\alpha)}}{\partial z^n}, z \right) d\Omega
\]

the Hamilton principle reads: for each field of virtual displacements \( z \in \Omega \rightarrow \zeta_\alpha \) such that \( \zeta_\alpha \) and its derivatives are zero on \( \partial \Omega \),

\[
\delta_\alpha \int_{\Omega} \Lambda d\Omega = 0
\]

Here \( \delta_\alpha a \) is the derivative of the Hamilton action with respect to \( \varepsilon_\alpha \), and \( \zeta_\alpha \) are the virtual displacements expressed in Eulerian coordinates \( z \) by means of equations (6.2). The method developed in Section 3 yields the equations of motion in the form

\[
\sum_{k(\alpha)=0}^{m(\alpha)} j_{k(\alpha)}^\ast \left( \frac{\partial K_{k(\alpha)}}{\partial z} - \left( \frac{\partial K_{k(\alpha)}}{\partial z} \right)^\ast \right) = 0, \quad K_{k(\alpha)}^\ast = \frac{\delta \Lambda}{\delta j_{k(\alpha)}}
\]

(6.4)
\[ \text{Div} \mathbf{j}_{k(\alpha)} = 0, \quad \alpha = 1, 2 \quad \text{and} \quad k(\alpha) = 0, \ldots, m(\alpha) \]

As in Theorem 4.1, for the case \( \Lambda = \Lambda \left( \mathbf{j}_{k(\alpha)}, \frac{\partial \mathbf{j}_{k(\alpha)}}{\partial \mathbf{z}}, z \right) \) we can obtain the following identity

\[
\text{Div} \left( \sum_{\alpha=1}^{2} \sum_{k(\alpha)=0}^{m(\alpha)} \left( K_{k(\alpha)}^* \mathbf{j}_{k(\alpha)} I - j_{k(\alpha)} K_{k(\alpha)}^* + A_{k(\alpha)}^* \frac{\partial j_{k(\alpha)}}{\partial \mathbf{z}} \right) - \Lambda I \right) \\
+ \frac{\partial \Lambda}{\partial \mathbf{z}} + \sum_{\alpha=1}^{2} \sum_{k(\alpha)=0}^{m(\alpha)} K_{k(\alpha)}^* \text{Div} \mathbf{j}_{k(\alpha)} - \sum_{\alpha=1}^{2} \sum_{k(\alpha)=0}^{m(\alpha)} j_{k(\alpha)}^* R_{k(\alpha)} \equiv 0
\]

where

\[ A_{k(\alpha)}^* = \frac{\partial \Lambda}{\partial \left( \frac{\partial j_{k(\alpha)}}{\partial \mathbf{z}} \right)} \quad \text{and} \quad R_{k(\alpha)} = \frac{\partial K_{k(\alpha)}}{\partial \mathbf{z}} - \left( \frac{\partial K_{k(\alpha)}}{\partial \mathbf{z}} \right)^* \]

Hence, the governing equations (6.4) admit the conservation laws

\[
\text{Div} \left( \sum_{\alpha=1}^{2} \sum_{k(\alpha)=0}^{m(\alpha)} \left( K_{k(\alpha)}^* \mathbf{j}_{k(\alpha)} I - j_{k(\alpha)} K_{k(\alpha)}^* + A_{k(\alpha)}^* \frac{\partial j_{k(\alpha)}}{\partial \mathbf{z}} \right) - \Lambda I \right) + \frac{\partial \Lambda}{\partial \mathbf{z}} = 0
\]

\[ \text{Div} \mathbf{j}_{k(\alpha)} = 0, \quad \alpha = 1, 2 \quad \text{and} \quad k(\alpha) = 0, \ldots, m(\alpha) \]

We notice that for a one-velocity model, the number of scalar conservation laws is \( m + d + 1 \) where \( d \) is the dimension of the time-space. The number of unknown variables is \( m + d \). Due to Theorem 4.3, the \( m + d + 1 \) conservation laws are connected by the "Gibbs identity". In the case of mixtures, we obtain \( \sum_{\alpha} m(\alpha) + d + 1 \) conservation laws for \( \sum_{\alpha} m(\alpha) + 2d \) unknown variables. In the general case, the classical approach based on conservation laws, does not allow to obtain Rankine-Hugoniot conditions for this system. Nevertheless, in [15-17] we have obtained the jump conditions. The Hamilton principle provides these conditions without any ambiguity.

### 7 Linear stability of mixtures.

We consider the Lagrangian in the form [15-19]

\[ L = \frac{1}{2} \rho_1 \left| \mathbf{u}_1 \right|^2 + \frac{1}{2} \rho_2 \left| \mathbf{u}_2 \right|^2 - W \left( \rho_1, \rho_2, \eta_1, \eta_2, \mathbf{w} \right) \quad (7.1) \]

where \( \mathbf{w} = \mathbf{u}_2 - \mathbf{u}_1 \) and \( \eta_\alpha \) is the entropy per unit volume of the \( \alpha^{th} \) component. A generalisation of (7.1) for thermocapillary fluids was also proposed.
in [20]. If \( W \) is an isotropic function of \( w \), Lagrangian (7.1) can be rewritten as follows

\[
\Lambda = \frac{1}{2} \sum_{\alpha=1}^{2} \left( \frac{|j_{0(\alpha)}|^2}{l^* j_{0(\alpha)}} - l^* j_{0(\alpha)} \right) - W \left( l^* j_{0(1)}, l^* j_{0(2)}, l^* j_{1(1)}, l^* j_{1(2)}, \mu \right)
\]

(7.2)

where

\[
\mu = \frac{1}{2} \left| j_{0(2)} \right|^2 \left| j_{0(1)} \right|^2
\]

(7.3)

Here we shall also restrict our study to the case \( s_{\alpha} = \text{const} \). To avoid double indices, we will denote \( j_{0(1)}, j_{0(2)} \) by \( j_1, j_2 \). Hence, (7.2) and (7.3) involve

\[
\Lambda(j_1, j_2) = \frac{1}{2} \sum_{\alpha=1}^{2} \left( \frac{|j_{\alpha}|^2}{l^* j_{\alpha}} - l^* j_{\alpha} \right) - \tilde{W}(l^* j_1, l^* j_2, \mu)
\]

(7.4)

where \( \tilde{W}(\rho_1, \rho_2, \mu) = W(\rho_1, \rho_2, \rho_1 s_1, \rho_2 s_2, \mu) \).

Omitting the tilde symbol, we consider the potential motions

\[
\begin{align*}
K^*_{\alpha} &= \frac{\partial \Lambda}{\partial j_{\alpha}} = \frac{\partial \varphi_{\alpha}}{\partial z} \\
\text{Div} j_{\alpha} &= 0, \quad \alpha = 1, 2
\end{align*}
\]

(7.5)

Let us consider the Legendre transformation of \( \Lambda(j_1, j_2) \)

\[
\Delta(K_1, K_2) = \sum_{\alpha=1}^{2} j_{\alpha}^* K_{\alpha} - \Lambda
\]

(7.6)

If the matrices

\[
\Lambda_{(\alpha \beta)} \equiv \frac{\partial}{\partial j_{\alpha}} \left( \left( \frac{\partial \Lambda}{\partial j_{\beta}} \right)^* \right) = \Lambda_{(\beta \alpha)}^*
\]

(7.7)

are non-degenerate, relation (7.6) involves

\[
\begin{align*}
\Delta_{(\gamma \alpha)} &= \frac{\partial \Delta}{\partial K_{\alpha}} \\
&= \frac{\partial}{\partial K_{\gamma}} \left( \left( \frac{\partial \Delta}{\partial K_{\alpha}} \right)^* \right)
\end{align*}
\]

(7.8)

Substituting (7.8) in \( \text{Div} j_{\alpha} = 0 \), we get

\[
tr \left( \frac{\partial j_{\alpha}}{\partial z} \right) = tr \left( \frac{\partial}{\partial z} \left( \frac{\partial \Delta}{\partial K_{\alpha}} \right)^* \right)
\]

\[
= tr \left( \sum_{\gamma=1}^{2} \frac{\partial}{\partial K_{\gamma}} \left( \frac{\partial \Delta}{\partial K_{\alpha}} \right)^* \frac{\partial K_{\gamma}}{\partial z} \right) = tr \left( \sum_{\gamma=1}^{2} \Delta_{(\gamma \alpha)} \varphi''_{\gamma} \right) = 0
\]

with

\[
\Delta_{(\gamma \alpha)} \equiv \frac{\partial}{\partial K_{\gamma}} \left( \left( \frac{\partial \Delta}{\partial K_{\alpha}} \right)^* \right) = \Delta_{(\alpha \gamma)}^*
\]

(7.9)
Hence, the equations of potential motions are

\[
\text{tr} \left( \sum_{\gamma=1}^{2} \Delta_{(\gamma\alpha)} \varphi''_{\gamma} \right) = 0 \quad (7.10)
\]

In the equilibrium state \( \Delta_{(\gamma\alpha)} = \text{const.} \) Substituting \( \varphi_{\gamma} \) by \( i\Phi_{\gamma} e^{i\beta'_{\gamma}z} \) in (7.10), we obtain the dispersion relation in the symmetric form

\[
\det \left( \begin{pmatrix}
\beta^* \Delta_{(11)} \beta \\
\beta^* \Delta_{(12)} \beta \\
\beta^* \Delta_{(21)} \beta \\
\beta^* \Delta_{(22)} \beta
\end{pmatrix}
\right) = 0
\]

which is the generalisation of the dispersion relation (5.5). However, to calculate \( \Delta_{(\gamma\alpha)} \) in terms of \( \Lambda_{(\gamma\alpha)} \) defined by (7.7), (7.9)), we have to solve the following system of matrix equations

\[
\sum_{\gamma=1}^{2} \Delta_{(\gamma\alpha)} \Lambda_{(\beta\gamma)} = I \delta_{\alpha\beta}, \quad \delta_{\alpha\beta} = \begin{cases} 1, & \alpha = \beta \\ 0, & \alpha \neq \beta \end{cases}
\]

A simpler method is to consider the linearised system generated from (7.5)

\[
\sum_{\gamma=1}^{2} \Lambda_{(\gamma\alpha)} j_{\gamma} - \left( \frac{\partial \varphi_{\alpha}}{\partial z} \right)^* = 0, \quad \text{Div} j_{\alpha} = 0
\]

where \( \Lambda_{(\gamma\alpha)} \) are taken at the equilibrium state \( j_{ke} \). Substituting \( j_{\gamma} \) by \( J_{\gamma} e^{i\beta'_{\gamma}z} \) and \( \varphi_{\alpha} \) by \( i\Phi_{\alpha} e^{i\beta'_{\alpha}z} \), we obtain the dispersion relation in the symmetric form

\[
\det \left( \begin{pmatrix}
\Lambda_{(11)} & \Lambda_{(21)} & \beta & 0 \\
\Lambda_{(12)} & \Lambda_{(22)} & 0 & \beta \\
\beta^* & 0^* & 0 & 0 \\
0^* & \beta^* & 0 & 0
\end{pmatrix}
\right) = 0 \quad (7.11)
\]

Equation (7.11) is presented in terms of the matrices \( \Lambda_{(\gamma\alpha)} \) calculated directly from the Lagrangian. Let us consider Lagrangian (7.4). Suppose that the velocities of each component are the same at equilibrium \( \mathbf{u}_{1e} = \mathbf{u}_{2e} \). Due to the invariance of the governing equations with respect to the Galilean transformation we assume, without loss of the generality, that \( \mathbf{u}_{1e} = \mathbf{u}_{2e} = 0 \). Suppressing the index "e" to avoid double indices, we get

\[
\begin{align*}
\Lambda_{(11)} &= \left( \rho_1 - \frac{\partial W}{\partial \mu} \right) \frac{I - \Pi^*}{\rho_1^2} - \frac{\partial^2 W}{\partial \rho_1^2} \Pi^* \\
\Lambda_{(21)} &= \Lambda_{(12)} = \frac{1}{\rho_1 \rho_2} \frac{\partial W}{\partial \mu} (I - \Pi^*) - \frac{\partial^2 W}{\partial \rho_1 \partial \rho_2} \Pi^* \\
\Lambda_{(22)} &= \left( \rho_2 - \frac{\partial W}{\partial \mu} \right) \frac{I - \Pi^*}{\rho_2^2} - \frac{\partial^2 W}{\partial \rho_2^2} \Pi^*
\end{align*} \quad (7.12)
\]
All matrices $\Lambda_{(\alpha,\beta)}$ are diagonal. Now we are able to calculate the dispersion relation (7.11) which is the determinant of a square matrix of dimension 10. Let us denote

$$a = -\frac{\partial W}{\partial \mu}, \quad w_{11} = \frac{\partial^2 W}{\partial \rho_1^2}, \quad w_{12} = \frac{\partial^2 W}{\partial \rho_1 \partial \rho_2}, \quad w_{22} = \frac{\partial^2 W}{\partial \rho_2^2} \quad (7.13)$$

Suppose that

$$a > 0, \quad w_{11} > 0, \quad w_{22} > 0, \quad w_{11}w_{22} - w_{12}^2 > 0 \quad (7.14)$$

Taking into account (7.12) - (7.13) we obtain by straightforward calculations of the determinant (7.11)

$$(a(\rho_1 + \rho_2) + \rho_1 \rho_2) \beta_1^4 - \left(a(\rho_1^2 w_{11} + 2 \rho_1 \rho_2 w_{12} + \rho_2^2 w_{22}) + \rho_1 \rho_2 (\rho_2 w_{22} + \rho_1 w_{11})\right)$$

$$\times (\beta_2^2 + \beta_3^2 + \beta_4^2) \beta_1^2 + (\beta_2^2 + \beta_3^2 + \beta_4^2)^2 (w_{22} w_{11} - w_{12}^2) \rho_1^2 \rho_2^2 = 0 \quad (7.15)$$

Let $\nu = 1, \beta = \omega \nu + \beta_\sigma, \lambda = \frac{\omega}{|\beta_\sigma|}$. The dispersion relation (7.15) takes the form

$$(a(\rho_1 + \rho_2) + \rho_1 \rho_2) \lambda^4 - \left(a(\rho_1^2 w_{11} + 2 \rho_1 \rho_2 w_{12} + \rho_2^2 w_{22}) + \rho_1 \rho_2 (\rho_2 w_{22} + \rho_1 w_{11})\right) \lambda^2$$

$$+ (w_{22} w_{11} - w_{12}^2) \rho_1^2 \rho_2^2 = 0 \quad (7.16)$$

The following result is proved in [21]:

**Theorem 7.1**

*If potential $W(\rho_1, \rho_2, \mu)$ satisfies conditions (7.14), all the roots $\lambda$ of the polynomial (7.16) are real.*

Conditions (7.14) mean that internal energy $U = W - w \frac{\partial W}{\partial w}$, where $w = |w|$, is convex [15-16, 21]. Hence, if the internal energy is a convex function and the relative velocity $w$ is small enough, the equilibrium state is stable in time direction in the time-space.

It is interesting to note that the Lagrangian of a two-fluid bubbly liquid with incompressible liquid phase (heterogeneous case) has the same form as (7.1). Indeed, in [18,19] the potential $W$ for a bubbly liquid is proposed in the form

$$W = \rho_2 \varepsilon_{20}(\rho_{20}) - \frac{\rho_{10}}{2} m(c) |w|^2 \quad (7.17)$$

where $\rho_2 = c \rho_{20}, c$ is the volume concentration of gas phase and the index ”0” means the real density of gas and liquid phase. Concentration $c$ is expressed
by \( c = \frac{4}{3}\pi R^3 N \) where \( R \) is the average radius of the bubbles, \( N \) is the number of bubbles per unit volume of the mixture. The internal energy per unit mass \( \varepsilon_{20} \) of the gas phase and the virtual mass coefficient \( m \) are known functions of \( \rho_{20} \) and \( \rho_1 \). By introducing the average density of the liquid phase

\[
\rho_1 = \rho_{10} (1 - c) \quad \text{with} \quad \rho_{10} = \text{const}
\]

we can rewrite the potential (7.17) as

\[
W = \rho_2 \varepsilon_{20} \left( \frac{\rho_2 \rho_{10}}{\rho_{10} - \rho_1} \right) - \rho_{10} \frac{m}{2} \left( \frac{\rho_{10} - \rho_1}{\rho_{10}} \right) |\mathbf{w}|^2
\]

The increase of the volume fraction \( c \) producing the interaction between gas bubbles changes not only the coefficient \( m(c) \) but also the interfacial energy \( \varepsilon_{in}(c) \) so that potential \( W \) can be generalized to the form

\[
W = \rho_2 \varepsilon_{20} (\rho_{20}) - \rho_{10} \frac{m}{2} (c) |\mathbf{w}|^2 + \varepsilon_{in}(c) \quad (7.18)
\]

Quantity \( \varepsilon_{in}(c) \) produces an additional pressure term due to the interfacial effect [22]. Therefore, Lagrangian (7.1) describes not only binary molecular mixtures but also partial cases of suspensions. Case (7.17) is actually degenerated: \( w_{11}w_{22} - w_{12}^2 \equiv 0 \). The introduction of the interfacial term in (7.18) involves the convexity condition \( w_{11}w_{22} - w_{12}^2 > 0 \) provided \( \frac{\partial^2 \varepsilon_{in}}{\partial c^2} > 0 \).

A APPENDIX

For any function \( \hat{f}(\lambda, \mathbf{X}) \) (see definitions (2.9)-(2.10)) and its image \( \mathbf{f}(t, \mathbf{X}) \) in the \((t, \mathbf{X})\)-coordinates, we obtain the following relations

\[
\frac{\partial \hat{f}}{\partial \lambda} \left( \frac{\partial g}{\partial \lambda} \right)^{-1} = \frac{\partial \mathbf{f}}{\partial t} - \frac{\partial \mathbf{f}}{\partial \mathbf{X}} \frac{\partial \hat{f}}{\partial \lambda} \left( \frac{\partial g}{\partial \lambda} \right)^{-1} = \frac{\partial \mathbf{f}}{\partial \mathbf{X}} \quad (A.1)
\]

A.1 Variation of the velocity

The definition (2.6) of the velocity \( \mathbf{u} \) yields

\[
\dot{\mathbf{u}} = \frac{\partial \phi}{\partial \lambda} \left( \frac{\partial g}{\partial \lambda} \right)^{-1}
\]
In Lagrangian coordinates, perturbation of \( u \) is represented by the formula
\[
\tilde{u}(\lambda, X, \varepsilon) = \frac{\partial \Phi}{\partial \lambda}(\lambda, X, \varepsilon) \left( \frac{\partial G}{\partial \lambda}(\lambda, X, \varepsilon) \right)^{-1}
\]
Taking the derivative of \( \tilde{u} \) with respect to \( \varepsilon \) at \( \varepsilon = 0 \) and using (A.1), we get
\[
\tilde{\delta}u = \partial \xi \partial \lambda \left( \frac{\partial g}{\partial \lambda} \right)^{-1} - \tilde{u} \partial \tau \partial \lambda \left( \frac{\partial g}{\partial \lambda} \right)^{-1} = \frac{d\xi}{dt} - u \frac{d\tau}{dt}
\]
where \( \frac{d}{dt} = \frac{\partial}{\partial t} + u^* \nabla^* \) is the material derivative. We obtain then
\[
\tilde{\delta}v = (I - v^*) \frac{\partial \zeta}{\partial z} v
\]
where \( v^* = (1, u^*) \), \( I \) is the identity tensor and \( l^* = (1, 0, 0, 0) \).

A.2 Variation of the deformation gradient

In Lagrangian coordinates \((\lambda, X)\), the perturbation of the deformation gradient is given by (2.6)
\[
\tilde{F}(\lambda, X, \varepsilon) = \frac{\partial \Phi}{\partial X} - \frac{\partial \Phi}{\partial \lambda} \frac{\partial G}{\partial \lambda} \left( \frac{\partial G}{\partial \lambda} \right)^{-1}
\]
Combination of the derivative of \( \tilde{F} \) with respect to \( \varepsilon \) at \( \varepsilon = 0 \) and relation (A.1) gives
\[
\tilde{\delta}F = \frac{\partial \xi}{\partial X} - \frac{\partial \xi}{\partial \lambda} \frac{\partial g}{\partial \lambda} \left( \frac{\partial g}{\partial \lambda} \right)^{-1} - \frac{\partial \Phi}{\partial \lambda} \frac{\partial g}{\partial X} \left( \frac{\partial g}{\partial \lambda} \right)^{-1} + \frac{\partial \Phi}{\partial \lambda} \frac{\partial g}{\partial X} \left( \frac{\partial g}{\partial \lambda} \right)^{-2} \frac{\partial \tau}{\partial \lambda}
\]
\[
= \frac{\partial \xi}{\partial X} - u \frac{\partial \tau}{\partial X} = \left( \frac{\partial \xi}{\partial X} - u \frac{\partial \tau}{\partial X} \right) F \quad (A.3)
\]

A.3 Variation of the density

The mass conservation is represented by the formula
\[
\rho \det F = \rho_0(X)
\]
The perturbation of $\rho$ is in the form

$$\tilde{\rho}(\lambda, X, \varepsilon) \det \tilde{F}(\lambda, X, \varepsilon) = \rho_0(X)$$

and consequently

$$\tilde{\delta} \rho \det \tilde{F} + \tilde{\rho} \tilde{\delta} (\det F) = 0 \quad (A.4)$$

By using the Euler-Jacobi identity,

$$\delta (\det F) = \det F tr(F^{-1} \tilde{\delta} F)$$

relations (A.3) and (A.4) lead to

$$\tilde{\delta} \rho = - tr\left( \rho(I - vI^r) \frac{\partial \zeta}{\partial z} \right) \quad (A.5)$$

### A.4 Variation of specific quantities

Along each trajectory the specific quantities $a_k$ are constant. Consequently, $a_k = a_{0k}(X)$. The perturbation of $a_k$ is such that $\tilde{a}_k(\lambda, X, \varepsilon) = a_{0k}(X)$. Hence,

$$\tilde{\delta} a_k = 0 \quad (A.6)$$

## B APPENDIX

Formulae (A.2) and (A.5) yield directly the variation of $j_0$

$$\tilde{\delta} j_0 = \left( tr\left( (vI^r - I) \frac{\partial \zeta}{\partial z} \right) I - (vI^r - I) \frac{\partial \zeta}{\partial z} \right) j_0 = \left( \frac{\partial \zeta}{\partial z} - (\text{Div} \, \zeta) I \right) j_0 \quad (B.1)$$

Let us consider the variations $\tilde{\delta} j_k, k = 1, \ldots, m$ with $j_k = a_k j_0$ where the scalar fields $a_k$ are such that $a_k = a_{0k}(X)$. Due to (A.6) $\tilde{\delta} a_k = 0$ and $\tilde{\delta} j_k = a_k \tilde{\delta} j_0$. Consequently

$$\tilde{\delta} j_k = \left( \frac{\partial \zeta}{\partial z} - (\text{Div} \, \zeta) I \right) j_k, \quad k = 0, \ldots, m \quad (B.2)$$

Hence, (2.12) involves

$$\hat{\delta} j_k = \left( \frac{\partial \zeta}{\partial z} - (\text{Div} \, \zeta) I \right) j_k - \frac{\partial j_k}{\partial z} \zeta, \quad k = 0, \ldots, m \quad (B.3)$$
Let $A$, $B$ be two linear transformations depending on $z$. Let us define the linear form $\text{tr} \left( A \frac{\partial B}{\partial z} \right)$ such that for any constant vector field $a$, $\text{tr} \left( A \frac{\partial B}{\partial z} \right) a \equiv \text{tr} \left( A \frac{\partial Ba}{\partial z} \right)$. As a consequence, we get

$$\text{Div} (AB) = \text{Div} (A) B + \text{tr} \left( A \frac{\partial B}{\partial z} \right)$$  \hspace{1cm} (C.1)

Indeed,

$$\text{Div} (ABa) = \text{Div} (A) B a + \text{tr} \left( A \frac{\partial Ba}{\partial z} \right) = \left( \text{Div} A \right) B + \text{tr} \left( A \frac{\partial B}{\partial z} \right) a$$

Let us also recall the following useful formulae for any vector fields $b(z)$, $c(z)$

$$\text{Div} (b c^*) = c^* \text{Div} b + b^* \left( \frac{\partial c}{\partial z} \right)^*$$  \hspace{1cm} (C.2)

$$\text{Div} (b^* c I) = \frac{\partial}{\partial z} (b^* c) = b^* \frac{\partial c}{\partial z} + c^* \frac{\partial b}{\partial z}$$  \hspace{1cm} (C.3)

By using formulae (C.1) - (C.3), we have the following identities

$$\text{Div} (K_k j_k I) \equiv K_k^* \frac{\partial j_k}{\partial z} + j_k^* \frac{\partial K_k}{\partial z}$$  \hspace{1cm} (C.4)

$$\text{Div} (j_k K_k^*) \equiv -K_k^* \text{Div} j_k - j_k^* \left( \frac{\partial K_k}{\partial z} \right)^*$$  \hspace{1cm} (C.5)

$$\text{Div} \left( A_k^* \frac{\partial j_k}{\partial z} \right) \equiv \text{Div} \left( A_k^* \frac{\partial j_k}{\partial z \partial z} \right) + \text{tr} \left( A_k^* \frac{\partial (\frac{\partial j_k}{\partial z})}{\partial z} \right)$$  \hspace{1cm} (C.6)

$$\text{Div} (\Lambda I) \equiv -\sum_{k=0}^{m} \left( \frac{\partial \Lambda}{\partial j_k \partial z} + \text{tr} \left( \frac{\partial \Lambda}{\partial \left( \frac{\partial j_k}{\partial z} \right) \partial \left( \frac{\partial j_k}{\partial z} \right)} \right) \right) - \frac{\partial \Lambda}{\partial z}$$

$$\equiv -\sum_{k=0}^{m} \left( K_k^* \frac{\partial j_k}{\partial z} + \text{tr} \left( A_k^* \frac{\partial \left( \frac{\partial j_k}{\partial z} \right)}{\partial z} \right) \right) - \frac{\partial \Lambda}{\partial z}$$  \hspace{1cm} (C.7)

By adding (C.4) - (C.7) we obtain

$$\text{Div} \left( \sum_{k=0}^{m} \left( K_k^* j_k I - j_k K_k^* + A_k^* \frac{\partial j_k}{\partial z} \right) - \Lambda I \right) + \frac{\partial \Lambda}{\partial z}$$

$$+ \sum_{k=0}^{m} K_k^* \text{Div} j_k - \sum_{k=0}^{m} j_k^* \left( \frac{\partial K_k}{\partial z} - \left( \frac{\partial K_k}{\partial z} \right)^* \right) \equiv 0$$

which proves Theorem 4.1.
References


