



Convective Mass Flows III

- In this lecture, we shall concern ourselves once more with convective mass and heat flows, as we still have not gained a comprehensive understanding of the physics behind such phenomena.
- We shall start by looking once more at the *capacitive field*.
- We shall then study the *internal energy* of matter.
- Finally, we shall look at *general energy transport phenomena*, which by now include mass flows as an integral aspect of general energy flows.

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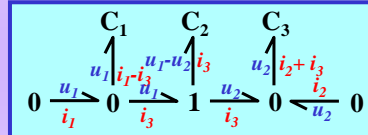
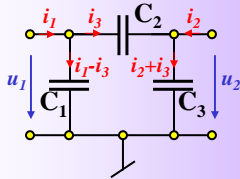
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Capacitive Fields III

- Let us briefly consider the following electrical circuit:



$$\begin{aligned} i_1 - i_3 &= C_1 \cdot du_1/dt \\ i_2 + i_3 &= C_3 \cdot du_2/dt \\ i_3 &= C_2 \cdot (du_1/dt - du_2/dt) \end{aligned}$$

$$\Rightarrow \begin{aligned} i_1 &= (C_1 + C_2) \cdot du_1/dt - C_2 \cdot du_2/dt \\ i_2 &= -C_2 \cdot du_1/dt + (C_2 + C_3) \cdot du_2/dt \end{aligned}$$



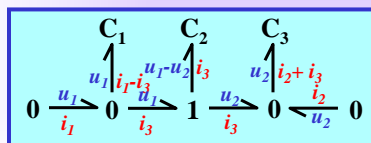
Capacitive Fields IV

$$\begin{aligned} i_1 &= (C_1 + C_2) \cdot du_1/dt - C_2 \cdot du_2/dt \\ i_2 &= -C_2 \cdot du_1/dt + (C_2 + C_3) \cdot du_2/dt \end{aligned}$$

Symmetric capacity matrix

$$\Rightarrow \begin{bmatrix} i_1 \\ i_2 \end{bmatrix} = \begin{bmatrix} (C_1 + C_2) & -C_2 \\ -C_2 & (C_2 + C_3) \end{bmatrix} \cdot \begin{bmatrix} du_1/dt \\ du_2/dt \end{bmatrix}$$

$$\Rightarrow \begin{bmatrix} du_1/dt \\ du_2/dt \end{bmatrix} = \frac{\begin{bmatrix} (C_2 + C_3) & C_2 \\ C_2 & (C_1 + C_2) \end{bmatrix}}{C_1 C_2 + C_1 C_3 + C_2 C_3} \cdot \begin{bmatrix} i_1 \\ i_2 \end{bmatrix}$$



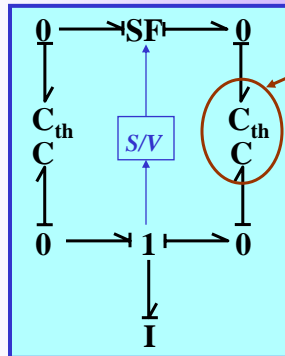
$$\Rightarrow 0 \mid \begin{matrix} u_1 \\ i_1 \end{matrix} \mid \text{CF} \mid \begin{matrix} i_2 \\ u_2 \end{matrix} \mid 0$$





Volume and Entropy Storage

- Let us consider once more the situation discussed in the previous lecture.



It was no accident that I drew the two capacitors so close to each other. In reality, the two capacitors together form a two-port capacitive field. After all, heat and volume are only two different properties of one and the same material.

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The Internal Energy of Matter I

- As we have already seen, there are three different (though inseparable) storages of matter:
 - Mass
 - Volume
 - Heat
- These three storage elements represent different storage properties of one and the same material.
- Consequently, we are dealing with a *storage field*.
- This storage field is of a capacitive nature.
- The capacitive field stores the *internal energy of matter*.

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The Internal Energy of Matter II

- Change of the internal energy in a system, i.e. the total power flow into or out of the capacitive field, can be described as follows :

$$\dot{U} = T \cdot \dot{S} - p \cdot \dot{V} + \sum_{\forall i} \mu_i \cdot \dot{N}_i$$

Flow of internal energy \rightarrow Heat flow \uparrow Volume flow \uparrow Mass flow \uparrow Molar mass flow \leftarrow Chemical potential \downarrow

- This is the *Gibbs equation*.

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The Internal Energy of Matter III

- The internal energy is proportional to the the total mass n .
- By normalizing with n , all extensive variables can be made intensive.

$$u = \frac{U}{n} \quad s = \frac{S}{n} \quad v = \frac{V}{n} \quad n_i = \frac{N_i}{n}$$

- Therefore:

$$\frac{d}{dt}(n \cdot u) = T \cdot \frac{d}{dt}(n \cdot s) - p \cdot \frac{d}{dt}(n \cdot v) + \sum_{\forall i} \mu_i \cdot \frac{d}{dt}(n \cdot n_i)$$

$$\Rightarrow \frac{d}{dt}(n \cdot u) - T \cdot \frac{d}{dt}(n \cdot s) + p \cdot \frac{d}{dt}(n \cdot v) - \sum_{\forall i} \mu_i \cdot \frac{d}{dt}(n \cdot n_i) = 0$$

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The Internal Energy of Matter IV

$$\frac{d}{dt}(n \cdot u) - T \cdot \frac{d}{dt}(n \cdot s) + p \cdot \frac{d}{dt}(n \cdot v) - \sum_i \mu_i \cdot \frac{d}{dt}(n \cdot n_i) = 0$$

$$\Rightarrow n \cdot \left[\frac{du}{dt} - T \cdot \frac{ds}{dt} + p \cdot \frac{dv}{dt} - \sum_i \mu_i \cdot \frac{dn_i}{dt} \right] + \frac{dn}{dt} \cdot \left[u - T \cdot s + p \cdot v - \sum_i \mu_i \cdot n_i \right] = 0$$

This equation must be valid independently of the amount n , therefore:

Finally, here is an explanation, why it was okay to compute with funny derivatives.

$$\frac{du}{dt} - T \cdot \frac{ds}{dt} + p \cdot \frac{dv}{dt} - \sum_i \mu_i \cdot \frac{dn_i}{dt} = 0$$

Flow of internal energy

$$u - T \cdot s + p \cdot v - \sum_i \mu_i \cdot n_i = 0$$

Internal energy

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The Internal Energy of Matter IV

$$U = T \cdot S - p \cdot V + \sum_i \mu_i \cdot N_i$$

$$\Rightarrow \dot{U} = T \cdot \dot{S} - p \cdot \dot{V} + \sum_i \mu_i \cdot \dot{N}_i + \dot{T} \cdot S - \dot{p} \cdot V + \sum_i \dot{\mu}_i \cdot N_i$$

$$= T \cdot \dot{S} - p \cdot \dot{V} + \sum_i \mu_i \cdot \dot{N}_i$$

$$\Rightarrow \dot{T} \cdot S - \dot{p} \cdot V + \sum_i \dot{\mu}_i \cdot N_i = 0$$

- This is the *Gibbs-Duhem equation*.

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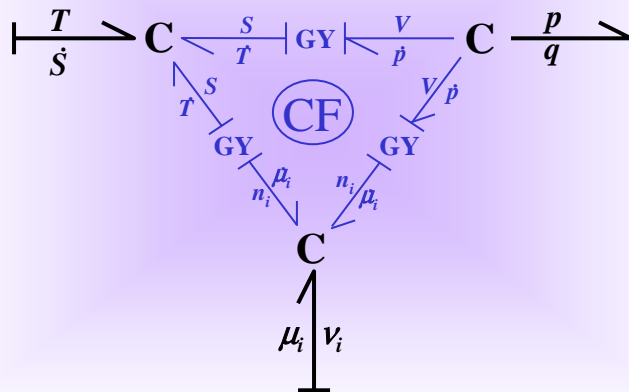
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The Capacitive Field of Matter



Simplifications

- In the case that no chemical reactions take place, it is possible to replace the *molar mass flows* by conventional *mass flows*.
- In this case, the *chemical potential* is replaced by the *Gibbs potential*.

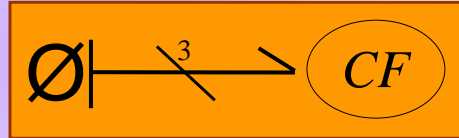
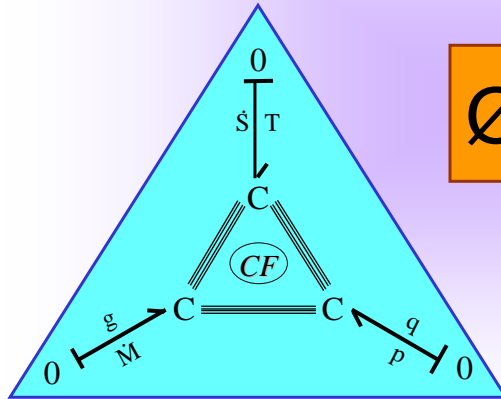
$$\frac{dU}{dt} = T \cdot \dot{S} - p \cdot \dot{V} + g \cdot \dot{M}$$





Bus-Bond and Bus-0-Junction

- The three outer legs of the CF-element can be grouped together.



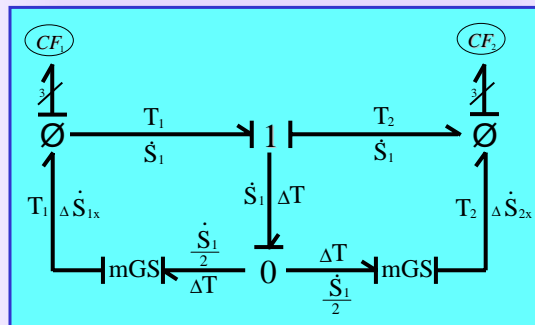
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Once Again Heat Conduction



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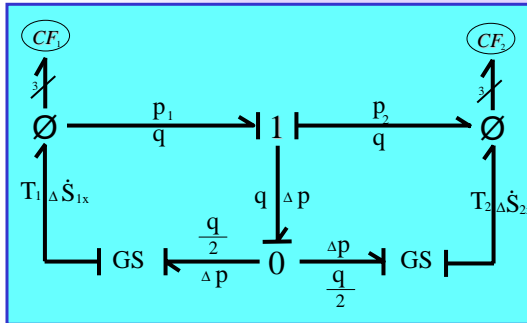
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Volume Pressure Exchange

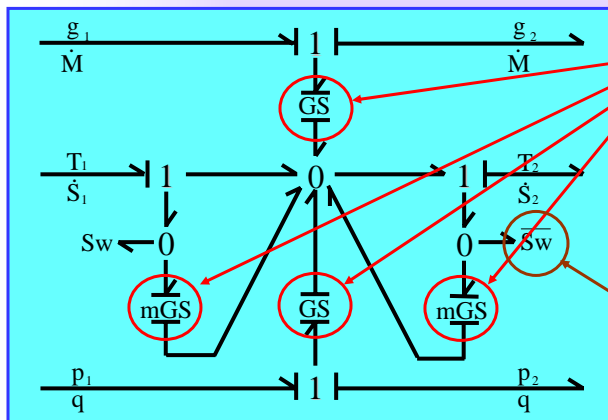


Pressure is being equilibrated just like temperature. It is assumed that the inertia of the mass may be neglected (relatively small masses and/or velocities), and that the equilibration occurs without friction.

The model makes sense if the exchange occurs locally, and if not too large masses get moved in the process.



General Exchange Element I



The three flows are coupled through RS-elements.

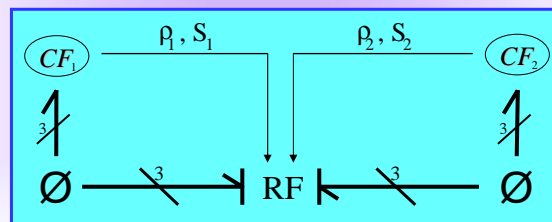
This is a switching element in bond-graph notation. This element has not yet been introduced.





General Exchange Element II

- In the general exchange element, the temperatures, the pressures, and the Gibbs potentials of neighboring media are being equilibrated.
- This process can be interpreted as a *resistive field*.



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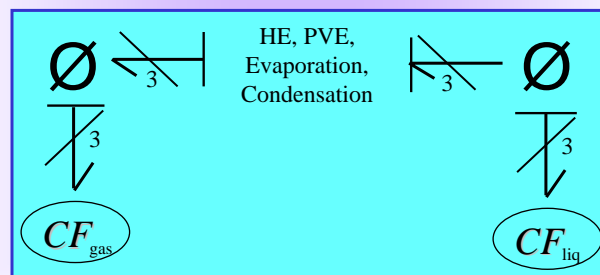
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Multi-phase Systems

- We may also wish to study phenomena such as *evaporation* and *condensation*.



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Evaporation (Boiling)

- Mass and energy exchange between capacitive storages of matter (*CF-elements*) representing different *phases* is accomplished by means of special resistive fields (*RF-elements*).
- The mass flows are calculated as functions of the pressure and the corresponding saturation pressure.
- The volume flows are computed as the product of the mass flows with the saturation volume at the given temperature.
- The entropy flows are superposed with the enthalpy of evaporation (in the process of evaporation, the thermal domain loses heat \rightarrow *latent heat*).

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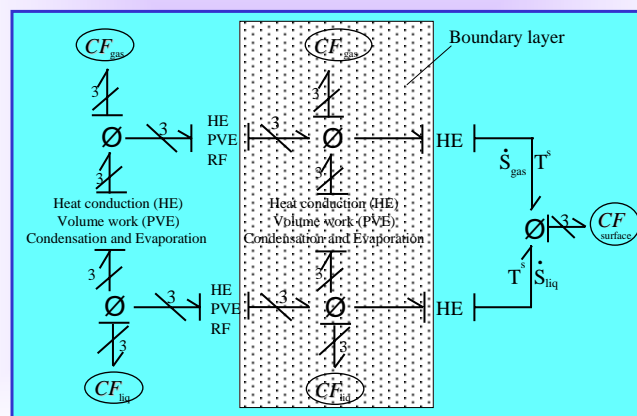
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Condensation On Cold Surfaces

- Here, a boundary layer must be introduced.



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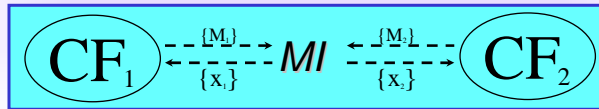
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Thermodynamics of Mixtures

- When fluids (gases or liquids) are being mixed, additional entropy is generated.
- This *mixing entropy* must be distributed among the participating component fluids.
- The distribution is a function of the *partial masses*.
- Usually, neighboring *CF-elements* are not supposed to know anything about each other. In the process of mixing, this rule cannot be maintained. The necessary information is being exchanged.



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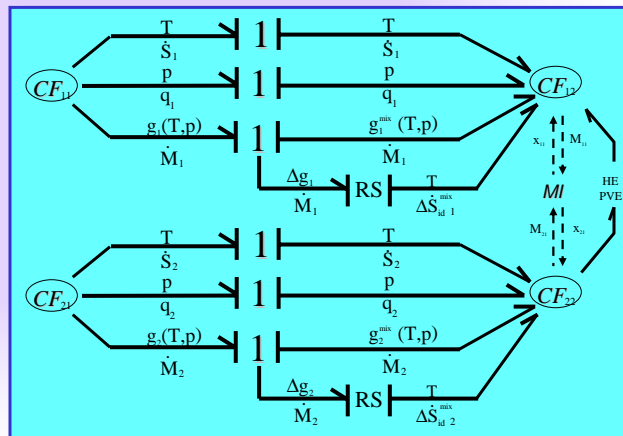
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Entropy of Mixing

- The mixing entropy is taken out of the Gibbs potential.

It was assumed here that the fluids to be mixed are at the same temperature and pressure.

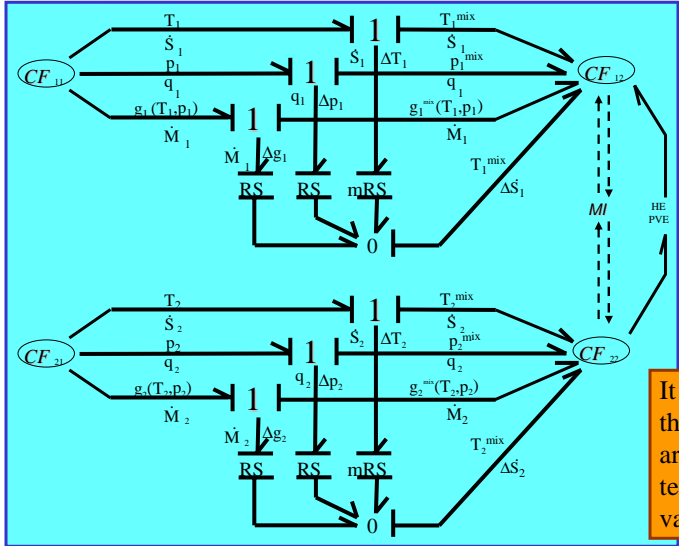


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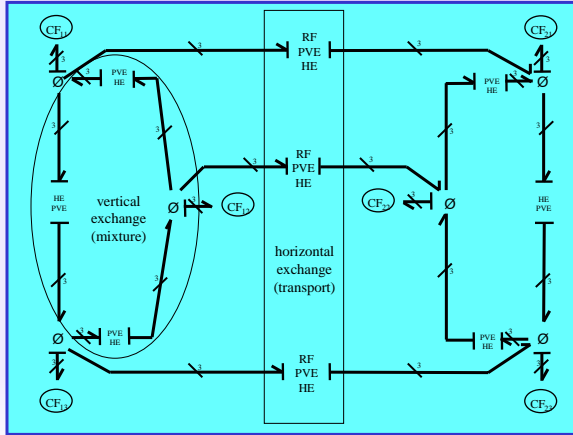




It is also possible that the fluids to be mixed are initially at different temperature or pressure values.

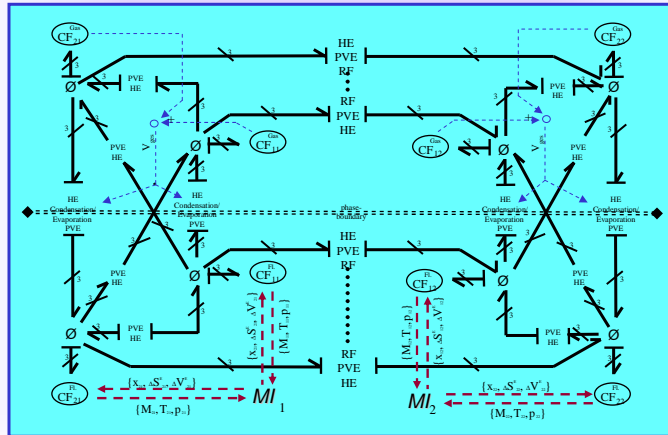


Convection in Multi-element Systems





Two-element, Two-phase, Two-compartment Convective System



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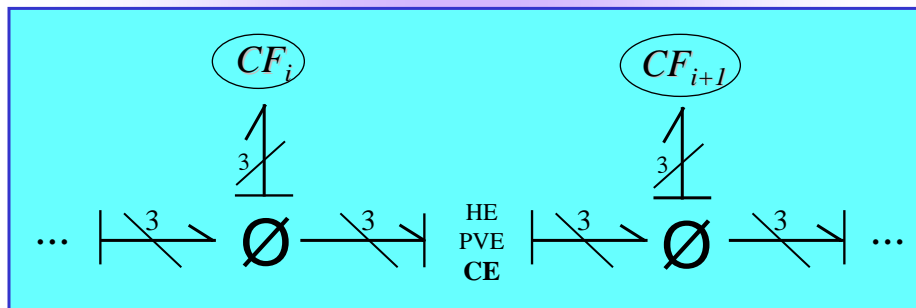
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Concentration Exchange

- It may happen that neighboring compartments are not completely homogeneous. In that case, also the concentrations must be exchanged.



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- Greifeneder, J. and F.E. Cellier (2001), “[Modeling multi-phase systems using bond graphs](#),” *Proc. ICBGM’01, Intl. Conference on Bond Graph Modeling and Simulation*, Phoenix, Arizona, pp. 285 – 291.
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