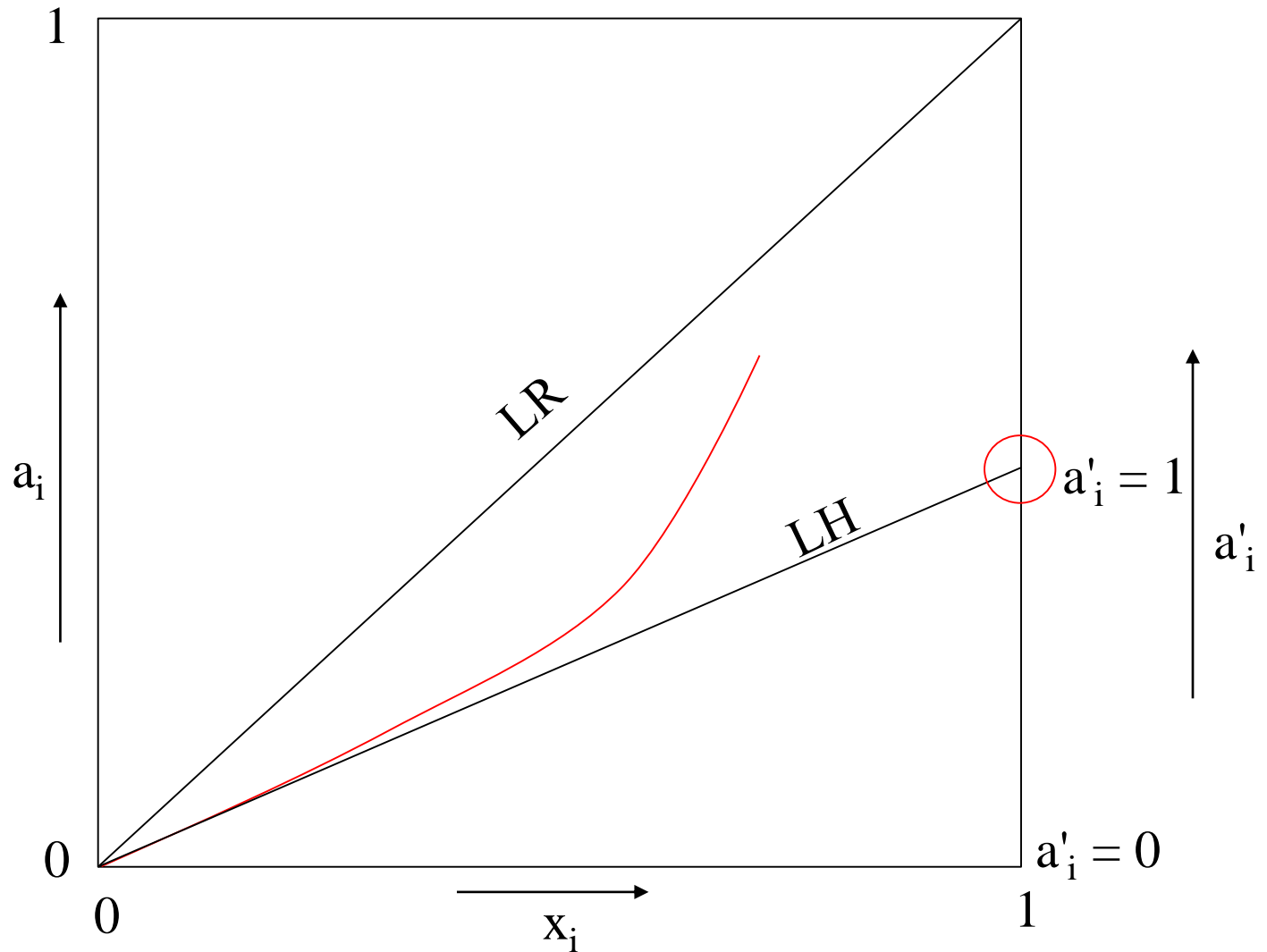




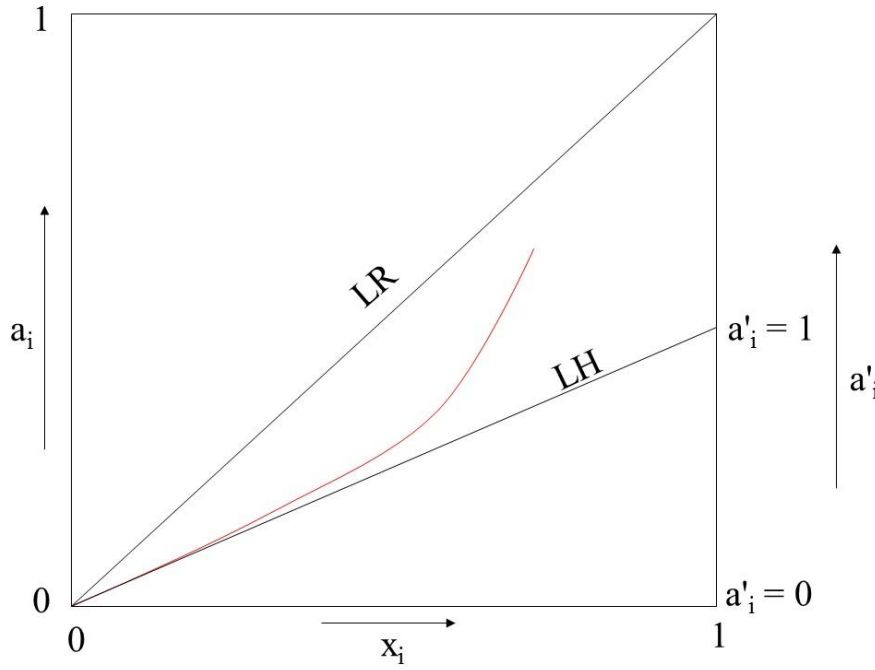
METMAT

TERMODINÂMICA DAS SOLUÇÕES

ESCALA HENRIANA- 1molar



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$$a'_i = \frac{x_i \cdot P_{\text{Total}}}{x_i^{x=1} \cdot P_{\text{Total}}} = \frac{x_i}{x_i^{x=1}} = x_i$$

$$a_i^{x=1} = \frac{p_i}{p_i^{x=1}} = a'_i$$

$$a'_i = x_i$$

Válida a LH

$$i_{\text{puro}} = i_{1\%}$$

$$K = \frac{h_i}{a_i} = \frac{\%i}{\gamma_i \cdot X_i}$$

Se válida a lei de Henry no intervalo de composição considerado:

- $X_i = \frac{\%i \cdot M_{\text{solvente}}}{100 \cdot M_i}$
- $\%i = 1$

$$K = \frac{100 \cdot M_i}{\gamma_i^0 \cdot M_{\text{solvente}}}$$

$j = \text{solvente}$



$$X_i = \frac{n_i}{n_j + n_i} = \frac{\frac{m_i}{M_i}}{\frac{m_j}{M_j} + \frac{m_i}{M_i}} = \frac{\frac{\%i}{M_i}}{\frac{\%j}{M_j} + \frac{\%i}{M_i}}$$

Para %i → 0% ⇒ %j → 100%

$$X_i = \frac{\frac{\%i}{M_i}}{\frac{100}{M_j} + \frac{0}{M_i}} = \frac{\frac{\%i}{M_i}}{\frac{100}{M_j}}$$

$$X_i = \frac{\%i \cdot M_{\text{solvente}}}{100 \cdot M_i}$$

$$i_{\text{puro}} = i_{1\%}$$

$$\Delta G_{R \rightarrow h(1\%)}^{\circ} = R.T \ln\left(\frac{\gamma_i^{\circ} \cdot M_j}{100 \cdot M_i}\right)$$

$$\text{Se } \ln \gamma^{\circ} = A + B/T$$

$$\Delta G_{R \rightarrow h(1\%)}^{\circ} = B.R + R.T \ln\left(\frac{A.M_j}{100.M_i}\right)$$

$$\Delta H_{R \rightarrow h(1\%)}^{\circ} = B.R$$

$$\Delta S_{R \rightarrow h(1\%)}^{\circ} = -R \ln\left(\frac{A.M_j}{100.M_i}\right)$$

$$i_{\text{puro}} = i_{X=1}$$

$$\Delta G_{R \rightarrow h(X=1)}^{\circ} = R.T \ln(\gamma_i^{\circ})$$

$$\text{Se } \ln \gamma^{\circ} = A + B/T$$

$$\Delta G_{R \rightarrow h(X=1)}^{\circ} = B.R + R.T.A$$

$$\Delta H_{R \rightarrow h(X=1)}^{\circ} = B.R$$

$$\Delta S_{R \rightarrow h(X=1)}^{\circ} = -A.R$$



ESCALA HENRIANA

$$i_{\text{raoultiano}} = i_{\text{henriano}}$$

Element, i^\dagger	$\gamma_i^0(1873)^*$	M (pure) = M(<i>i.d.</i> , X, liq.) $\Delta G(x)$, cal/g atom	M (pure) = M(<i>i.d.</i> , wt.-%, liq.) $\Delta G(\%)$, cal/g atom	Ref. and Notes
Ag (<i>l</i>)	200	19 700	19 700 - 10.46T	11, 2; from solubility data, regular solution is assumed.
Al (<i>l</i>)	0.029	-15 100 + 1.03T	-15 100 - 6.67T	19
B (<i>s</i>)	0.022	-15 600 + 0.71T	-15 600 - 5.15T	29; from solubility of BN, using values for nitrogen given below.
C (<i>gr</i>)	0.57	5 400 - 4.0T	5 400 - 10.1T	42, 2
Ca (<i>v</i>)	2240	-9 430 + 20.3T	-9 430 + 11.8T	20; from solubility data, regular solution is assumed.
Co (<i>l</i>)	1.07	240	240 - 9.26T	82, 2; regular solution is assumed.
Cr (<i>l</i>)	1.0	0	-9.01T	87, 86, 88, 89; liquid Fe-Cr alloys ideal at low Cr concentration.
Cr (<i>s</i>)	1.14	4 600 - 2.19T	4 600 - 11.20T	
Cu (<i>l</i>)	8.6	8 000	8 000 - 9.41T	93, 2
1/2 H ₂ (<i>g</i>)	—	—	8 720 + 7.28T	95, 96, 2; $\Delta G^\circ(\text{ppm}) = 8,720 - 11.02T$
Mn (<i>l</i>)	1.3	976	976 - 9.12T	125, 126, 2
Mo (<i>l</i>)	1	0	-10.23T	Ideal behaviour assumed.
Mo (<i>s</i>)	1.86	6 600 - 2.29T	6 600 - 12.52T	Ideal behaviour of liquid solution, transfer of standard state.
1/2 N ₂ (<i>g</i>)	—	—	860 + 5.71T	135
Nb (<i>l</i>)	1.0	0	-10.2	See Mo above.
Nb (<i>s</i>)	1.4	5 500 - 2.3T	5 500 - 12.5	
Ni (<i>l</i>)	0.66	-5 000 + 1.80T	-5 000 - 7.42T	164, 2
1/2 O ₂ (<i>g</i>)	—	—	-28 000 - 0.69T	168
1/2 P ₂ (<i>g</i>)	—	—	-29 200 - 4.6T	218
Pb (<i>l</i>)	1400	50 800 - 12.7T	50 800 - 25.4T	21; calculated from data on solubility.
1/2 S ₂ (<i>g</i>)	—	—	-32 280 + 5.6T	72
Si (<i>l</i>)	0.0013	-31 430 + 3.64T	-31 430 - 4.12T	16, 243
Sn (<i>l</i>)	2.8	3 820	3 820 - 10.62T	244; regular solution assumed.
Ti (<i>l</i>)	0.037	-11,100	-11 000 - 8.85T	19; regular solution assumed.
Ti (<i>s</i>)	0.038	-7 440 - 1.90T	-7 440 - 10.75T	19
U (<i>l</i>)	0.027	-13 400	-13 400 - 12.0T	24; regular solution assumed.
V (<i>l</i>)	0.08	-10 100 + 0.37T	-10 100 - 8.6T	28
V (<i>s</i>)	0.1	-4 950 - 1.93T	-4 950 - 10.9T	
W (<i>l</i>)	1	0	-11.5T	Assumed ideal liquid solution.
W (<i>s</i>)	1.2	+7 500 - 3.65T	+7 500 - 15.2T	
Zr (<i>l</i>)	0.037	-12 200	-12 200 - 10.13T	γ_{Zr}^0 assumed equal to γ_{Ti}^0 , and regular solution assumed.
Zr (<i>s</i>)	0.043	-8 300 - 1.82T	-8 300 - 11.95T	

$$* \gamma_i^0 = \lim_{X_i \rightarrow 0} a_i/X_i$$

† The letters in parentheses indicate the standard states used. All are of one atmosphere pressure.



Determinar a atividade raoultiana do silício numa liga líquida Fe-Si a 1600°C contendo 1%Si. Admitir válida a lei de Henry para o Si.

$$\{Si\} = \underline{Si} \quad \Delta G^\circ = -31430 - 4,12T \text{ (cal / mol)}$$

$$\Delta G^\circ = -3,91 \times 10^4 \text{ cal / mol} \quad K = \frac{h_{Si}}{a_{Si}} = \frac{1}{a_{Si}} = 3,7 \times 10^4$$

$$a_{Si} = 2,7 \times 10^{-5}$$