

Chemistry 163B

μ_i and $\Delta\mu_{\text{reaction}}$

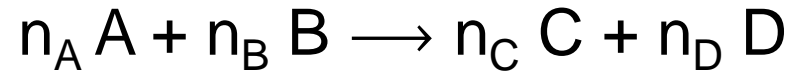
Activity

Equilibrium

goals

- Derive equilibrium and spontaneity criteria applying multicomponent thermodynamic relationships; i.e chemical potential ($\Delta\mu_{\text{reaction}}$)
- Define concentration dependence of μ in terms of **activity** (fugacity) of 'real' gases, actual solutes
- Apply **activity** to equilibrium K_{eq}
- Derive how to obtain fugacity if **REAL** gas

equilibrium in terms of $\Delta\mu$



$$v_A = -n_A \quad v_B = -n_B \quad v_C = +n_C \quad v_D = +n_D$$

$d\xi$ is extent of reaction

$d\xi > 0$ forward reaction

$d\xi < 0$ reverse reaction

$$dn_i = v_i d\xi$$

$dn_i > 0$ substance i increases

$dn_i < 0$ substance i decreases

equilibrium in terms of $\Delta\mu$

$dG_{T,P} < 0$ spontaneous

$dG_{T,P} = 0$ equilibrium

whole pot of mixed reactants and products

$$dG = -SdT + VdP + \sum_{i=1}^N \mu_i dn_i$$

$$dn_i = \nu_i d\xi$$

$$dG = -SdT + VdP + \sum_{i=1}^N \mu_i \nu_i d\xi$$

$$dG_{T,P} = \left(\sum_{i=1}^N \mu_i \nu_i \right) d\xi \leq 0$$

equilibrium in terms of $\Delta\mu$

$$dG_{T,P} = \underbrace{\left(\sum_{i=1}^N \mu_i \nu_i \right)}_{?} d\xi \leq 0$$

$$dG_{T,P} = \underbrace{\left(\sum_{i=1}^N \mu_i \nu_i \right)}_{\Delta\mu_{\text{reaction}}} d\xi \leq 0$$

$$dG_{T,P} = \Delta\mu_{\text{reaction}} d\xi \leq 0$$

$\Delta\mu_{\text{reaction}} < 0$ forward reaction spontaneous ($d\xi > 0$)

$\Delta\mu_{\text{reaction}} > 0$ reverse reaction spontaneous ($d\xi < 0$)

$\Delta\mu_{\text{reaction}} = 0$ equilibrium

just like ΔG !!!

NOTE: μ_i IS INTENSIVE (J mol^{-1})
 $\Delta\mu_{\text{reaction}}$ IS EXTENSIVE (J)

concentration dependence of μ_i

ideal gas, one component (pure substance)

$$\bar{G} = \bar{G}^\circ + RT \ln\left(\frac{P}{1 \text{ bar}}\right)$$

led to

$$\Delta G_{\text{reaction}} = \Delta G_{\text{reaction}}^\circ + RT \ln(Q_P)$$

what about if other species present?

$$\mu_i = \mu_i^\circ + RT \ln\left(\frac{P_i}{1 \text{ bar}}\right)$$

$$\Delta \mu_{\text{reaction}} = \Delta \mu_{\text{reaction}}^\circ + RT \ln Q_P$$

$$\Delta \mu_{\text{reaction}}^\circ = \sum_i \nu_i \mu_i^\circ \quad Q_P = \prod_i \left(\frac{P_i}{1 \text{ bar}}\right)^{\nu_i}$$

HANDOUT #48

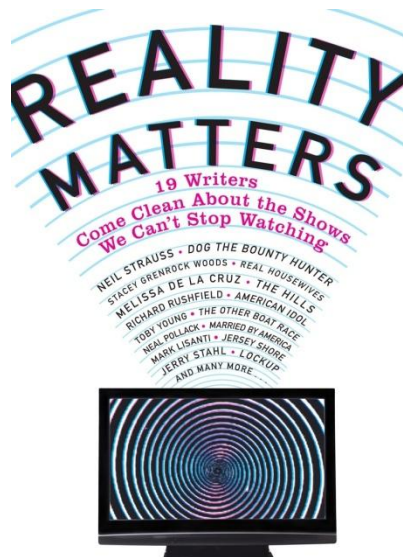
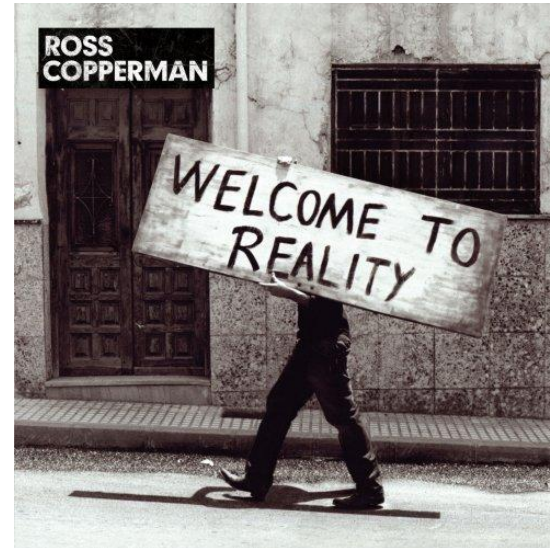
yada -yada- yada: and so forth for $\Delta\mu_{\text{reaction}}$

$$\Delta\mu_{\text{reaction}}^{\circ} = -\underline{RT} \ln K_P$$

$$\left(\frac{\partial \frac{\Delta\mu}{T}}{\partial T} \right)_P = -\frac{\Delta H}{T^2} \quad \text{where} \quad \Delta H = \sum_i \nu_i \bar{H}_i = \sum_i \nu_i \left(\frac{\partial H}{\partial n_i} \right)_{T,P,n_j \neq n_i}$$

$$\left(\frac{\partial \ln K}{\partial T} \right)_P = \frac{\Delta H^{\circ}}{\underline{RT}^2} \quad \text{where} \quad \Delta H^{\circ} = \sum_i \nu_i \bar{H}_i^{\circ}$$

correcting for REALITY (activity and fugacity)



edited by ANNA DAVID foreword by JAMES FREY

I STILL REFUSE
TO ACCEPT REALITY

a haiku



My childhood dream

was to become a mermaid
with a blue tail

correcting for REALITY (activity and fugacity)

- will define activity and fugacity coefficients γ_i 's that provides corrections for the deviation of chemical potential from the **ideal** gas and solute concentration dependence
- activity and fugacity coefficients are obtained from experimental measurements on **REAL** systems or by theory (Debye-Huckel)

$$\mu_i(T) = \mu_i^\circ(T) + RT \ln \left(\frac{a_i}{a_i^0} \right)$$

*standard
conditions*

*how μ_i 'really' changes
in going from standard
conditions to actual conditions*

a_i = activity of component i

$a_i = \gamma_i \times [\text{ideal measure of pressure, concentration, etc}]$

γ_i is activity coefficient, a correction for non-ideality

$a_i^0 = 1$ unit (bar, molar, etc)

more general μ_i and corrections for non-ideality

$$\Delta\mu_{\text{reaction}} = \Delta\mu^\circ + \underline{RT} \ln Q$$

where now Q is written in terms of activities

$$Q = \prod_i \left(\frac{a_i}{a_i^\circ} \right)^{\bar{v}_i} \quad Q = \frac{(a_C)^{\bar{n}_C} (a_D)^{\bar{n}_D}}{\underbrace{(a_A)^{\bar{n}_A} (a_B)^{\bar{n}_B}}}$$

unitless

dropped the $a^\circ=1$ 'unit'
but Q is 'unitless'

fugacity of gases

1. GASES:

$$a_i = f_i = \gamma_i P_i$$

partial pressure

fugacity of gas

fugacity coefficient

$$Q = \frac{(a_C)^{\bar{n}_C} (a_D)^{\bar{n}_D}}{(a_A)^{\bar{n}_A} (a_B)^{\bar{n}_B}}$$

$$Q = \frac{\left(\gamma_C \frac{P_C}{1\text{bar}}\right)^{\bar{n}_C} \left(\gamma_D \frac{P_D}{1\text{bar}}\right)^{\bar{n}_D}}{\left(\gamma_A \frac{P_A}{1\text{bar}}\right)^{\bar{n}_A} \left(\gamma_B \frac{P_B}{1\text{bar}}\right)^{\bar{n}_B}}$$

$$Q = \left(\frac{\gamma_C^{\bar{n}_C} \gamma_D^{\bar{n}_D}}{\gamma_A^{\bar{n}_A} \gamma_B^{\bar{n}_B}}\right) \frac{\left(\frac{P_C}{1\text{bar}}\right)^{\bar{n}_C} \left(\frac{P_D}{1\text{bar}}\right)^{\bar{n}_D}}{\left(\frac{P_A}{1\text{bar}}\right)^{\bar{n}_A} \left(\frac{P_B}{1\text{bar}}\right)^{\bar{n}_B}} = \gamma Q_P$$

correction for non-ideality

ideal gas Q_P

other conventions for activities

2. pure solids and liquids

$$\mu_i(T, P) \approx \mu_i^\circ(T, P = 1\text{bar})$$

$$\left(\frac{\partial \mu_i}{\partial P}\right)_T = \bar{V}_i \quad (\text{small for liquid or solid})$$

so $a_i \approx 1$ for pure solid or liquid
[unless extreme pressure]



other conventions for activities

3. solutes in solutions

$$a_i = \gamma_i [I]$$

concentration of I,
usually molar
but may be X_i

WHY CONCENTRATION FOR SOLUTES IN SOLUTION ?
HAVE PROVEN, $P_i = X_i P_{TOTAL} \Rightarrow Q$ FOR IDEAL GAS.
SOON WILL JUSTIFY $[I] = X_i \Rightarrow Q$ FOR ,SOLUTES IN IDEAL SOLNS.

activity coefficient γ_i corrects 'ideal' measure of 'concentration

if "activity coefficients unity"

$$a_i = [I] \quad a_i \equiv f_i = P_i \quad a_i = 1$$

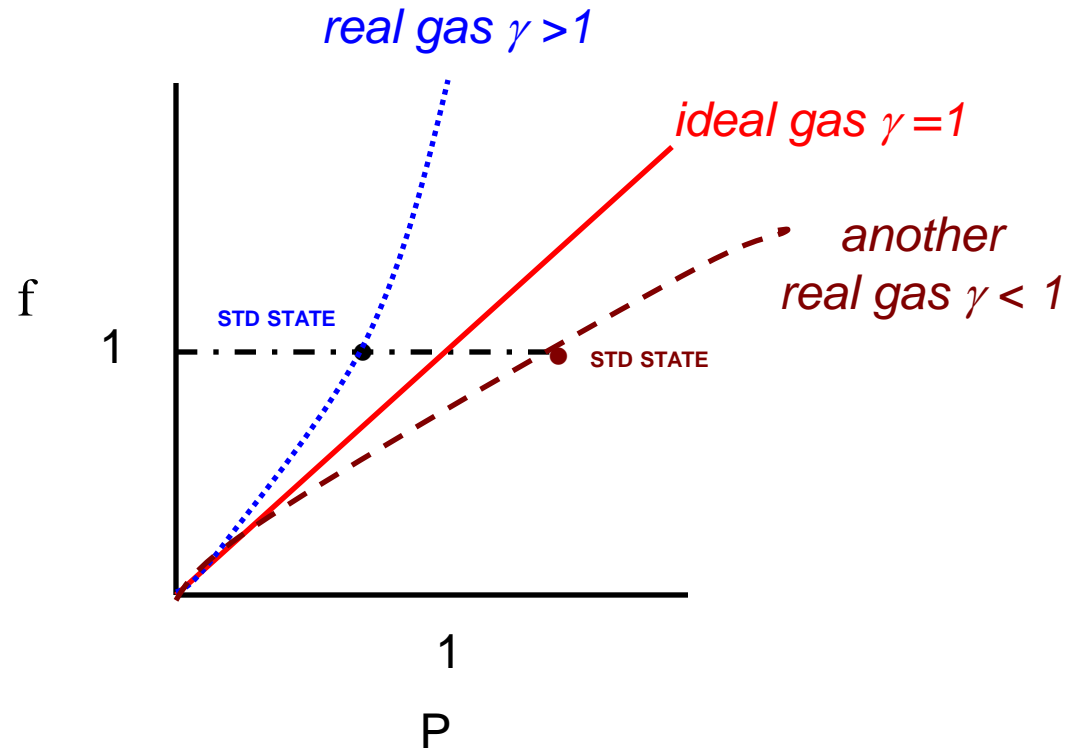
solute

gas

pure liquid or solid

HW#8 $\gamma=1$ except prob. 41* and 43.

how to evaluate activity (fugacity) coefficients for real gases (HW8 #*41)



how to evaluate activity (fugacity) coefficients for real gases

$$\left(\frac{\partial \mu}{\partial P}\right)_{T,n} = \bar{V} \quad \text{and}$$

$$\mu = \mu^\circ + RT \ln f$$

need in a moment

$$\left(\frac{\partial \mu}{\partial P}\right)_{T,n} = RT \left(\frac{\partial \ln f}{\partial P}\right)_{T,n}$$

$$\text{so } RT \left(\frac{\partial \ln f}{\partial P}\right)_{T,n} = \bar{V}$$

expression for $d \ln\left(\frac{f}{P}\right)$ will prove useful

$$\left(\frac{\partial \ln\left(\frac{f}{P}\right)}{\partial P}\right)_{T,n} = \left(\frac{\partial (\ln f - \ln P)}{\partial P}\right)_{T,n} = \left(\frac{\partial \ln f}{\partial P}\right)_{T,n} - \left(\frac{\partial \ln P}{\partial P}\right)_{T,n}$$

$$= \frac{1}{RT} \bar{V} - \frac{1}{P} = \frac{1}{RT} \left(\bar{V} - \frac{RT}{P}\right)$$

how to evaluate activity (fugacity) coefficients for real gases

$$\int_{P_1}^P d \left(\ln \frac{f}{P'} \right) = \int_{P_1}^P \frac{1}{RT} \left(\bar{V} - \frac{RT}{P'} \right) dP'$$

$$\ln \left(\frac{f(P)}{P} \right) - \ln \left(\frac{f(P_1)}{P_1} \right) = \frac{1}{RT} \int_{P_1}^P \left(\bar{V} - \frac{RT}{P'} \right) dP'$$

$$\ln \left(\frac{f(P)}{P} \right) = \ln \left(\frac{f(P_1)}{P_1} \right) + \frac{1}{RT} \int_{P_1}^P \left(\bar{V} - \frac{RT}{P'} \right) dP'$$

Diagram illustrating the limits of integration for the fugacity coefficient equation:

- Left box (limits for $\ln \left(\frac{f(P_1)}{P_1} \right)$):
 - $P_1 \rightarrow 0$
 - $\frac{f(P_1)}{P_1} \rightarrow 1$
 - $\ln \left(\frac{f(P_1)}{P_1} \right) \rightarrow 0$
- Right box (limit for the integral term):
 - $(\bar{V}_{ACTUAL} - \bar{V}_{IDEAL GAS})$

*how to evaluate activity (fugacity) coefficients for real gases
(eqn 7.20 E&R and HW8 #*41)*

$$\ln f(P) = \ln P + \frac{1}{RT} \int_{P_1 \rightarrow 0}^P \left(\bar{V} - \frac{RT}{P'} \right) dP' = \ln P + \frac{1}{RT} \int_{P_1 \rightarrow 0}^P \left(\bar{V}_{ACTUAL} - \bar{V}_{IDEAL GAS} \right) dP'$$

$$\ln \left(\frac{f(P)}{P} \right) = \ln(\gamma) = \frac{1}{RT} \int_{P_1 \rightarrow 0}^P \left(\bar{V} - \frac{RT}{P'} \right) dP' = \frac{1}{RT} \int_{P_1 \rightarrow 0}^P \left(\bar{V}_{ACTUAL} - \bar{V}_{IDEAL} \right) dP'$$

$$z = \frac{\bar{V}_{actual}}{\bar{V}_{ideal}} = \frac{P\bar{V}_{actual}}{RT} \quad (\text{compression factor E\&R eqn. 7.6})$$

$$\ln \gamma = \frac{1}{RT} \int_{P_1 \rightarrow 0}^P \bar{V}_{ideal} (z - 1) dP' = \int_{P_1 \rightarrow 0}^P \frac{(z - 1)}{P'} dP'$$

HW8 41*

$$\gamma(P, T) = \exp \left[\int_{P_1 \rightarrow 0}^P \frac{z - 1}{P'} dP' \right] \quad (\text{E \& R eqn 7.21})$$

End of Lecture

activity for solids and liquids ($P_{\text{total}} \neq 1 \text{ bar}$)

$$\mu_i(\mathbf{T}) = \mu_i^\circ(\mathbf{T}) + RT \ln \left(\frac{a_i}{a_i^0} \right)$$

$$\left(\frac{a_i}{a_i^0} \right) = e^{\frac{\mu_i(\mathbf{T}) - \mu_i^\circ(\mathbf{T})}{RT}}$$

$$\left(\frac{\partial \mu_i}{\partial P} \right)_{T,n} = \bar{V}_i$$

$$\mu_i(T, P_{\text{total}}) = \mu_i^\circ(T, 1 \text{ bar}) + \int_{1 \text{ bar}}^{P_{\text{total}}} \bar{V}_i dP \approx \mu_i^\circ(T, 1 \text{ bar}) + \bar{V}_i (P_{\text{total}} - 1 \text{ bar})$$

HW#7 37 (E&R 6.5)

$$\left(\frac{a_i}{a_i^0} \right)_{l \text{ or } s} \approx e^{\underbrace{\frac{\bar{V}_i (P_{\text{total}} - 1 \text{ bar})}{RT}}_{\approx 1}} \approx 1 \text{ for } P_{\text{total}} \text{ near } 1 \text{ bar (since } \bar{V} \text{ is small for liquids or solids)}$$

for $\Delta\mu$ at high P_{total}
would use this in Q
for liquids and solids



Midterm #2, Wednesday 26th February

- Exam WILL COVER material on lectures #8-#16 (#1-#16) (~through Friday, 21th February) and HW#5-HW#7
- Weeks 4-7 “reviews” on WWW handouts (19th February)
- HW solutions on eCommons (HW#7 Mon, 24th February)
- SAMPLE (~~practice~~) midterm on eCommons reserve Wednesday, 19th February
(see sample exam for ‘relationships provided’)
- TAs have KEY for use in sections, tutorials, and office hours (on eCommons ~23rd February)
 - REVIEW SESSION
Monday, 24th February, Classroom Unit 1, 7:00-8:00 PM
 - Midterm #2 Exam, Wednesday 26th February
done “in place”
no ear dingies (iPODs, MP3, etc)
bring charged calculators
get here ON TIME