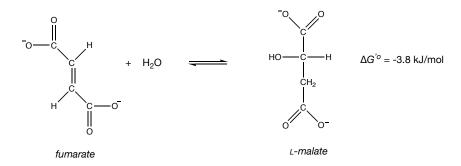
Chem 352 - Fall 2018 Quiz 1

 $R = 8.314 \ J/(mol \cdot K) = 0.08206 \ (L \cdot atm)/(mol \cdot K)$ [Ideal Gas Law Constant] $K_w = 1.00 \times 10^{-14} \text{ M}^2$ [Ionization Product for water]

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1. One of the reactions in the citric acid cycle pathway, a pathway used in the complete oxidation of glucose to CO₂ and H₂O, involves the hydration of *fumarate* to form *L-malate*.



- a. Is the reaction favorable under standard state conditions at 37°C ? (Y/N) $\underline{\text{Y}}$ $\underline{\text{Y}}$ $\underline{\text{Explain}}$: The change in the free energy for a reaction, ΔG can be used to determine if a reaction is favorable or not as written. When ΔG is < 0, the reaction is favorable, when ΔG is > 0, the reaction is unfavorable, but is favorable in the reverse direction. When $\Delta G = 0$, the reaction is at equilibrium, and is not favorable in either direction. For this reaction, you are asked if the reaction is favorable at *standard state conditions*, At standard state conditions, $\Delta G = \Delta G^{\circ}$, the standard state free energy change. It is given that $\Delta G^{\circ} = -3.8$ kJ/mol-therefore $\Delta G = -3.8$ kJ/mol and the reaction is favorable as written.
- b. What is the value of the equilibrium constant, K_{eq} , for this reaction at 37°C?

at equlibrium,
$$\Delta G^{1o} = -RT \ln \left(K_{eq}\right)$$

$$\ln \left(K_{eq}\right) = \frac{-\Delta G^{1o}}{RT}$$

$$K_{eq} = e^{\left(\frac{-\Delta G^{o}}{RT}\right)}$$

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$$K_{eq} = 4.37$$

c. If the cellular concentration of *fumarate* is 1.33 mM, and the concentration of *L-malate* is 8.55 mM, is this reaction favorable under conditions found in a cell at 37°C? (Y/N) ______
 Explain: Under the conditions given, ΔG > 0, therefore the reaction is not favorable.
 when not at standard state condition.

$$\Delta G = \Delta G^{'o} + RT \ln \left(\frac{products}{reactants} \right)$$

$$= -3.8 \frac{kJ}{mol} + \left(8.314 \times 10^{-3} \frac{kJ}{mol \cdot K} \right) \left(37 + 272 \text{ K} \right) \ln \left(\frac{8.55 \times 10^{-3} \text{ M}}{1.33 \times 10^{-3} \text{ M}} \right)$$

$$(H_{2}O \text{ is also the solvent, and so is not included as one of the reactants.})$$

$$\Delta G = 1.00 \frac{kJ}{mol}$$

2. The structural formula for the amino acid glutamic acid is shown to right. Glutamic acid contains three ionizable groups; the structure shown has all three of these groups in their deprotonated, conjugate base forms. The pK_a values for these three groups are indicated in the figure.

$pK_1 = 1.88$ $pK_1 = 9.54$ CH₂ $pK_1 = 4.27$

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a. What is the expected pH for a 150 mM solution of glutamic acid in its *fully protonated* form? (show calculations)

It is the ionizable group which is the strongest acid and has the lowest pK_a that will have the greatest influence on the pH of the fully protonated form of glutamic acid in an aqueous solution. The $[H^+]$ of a weak acid solution is approximately equal to $\sqrt{K_a C_a}$, where K_a is the acid dissociation for the weak acid, and C_a is its concentration.

$$[H^{+}] \approx \sqrt{K_a C_a}$$

$$\approx \sqrt{10^{-\rho K_1} C_a} \approx \sqrt{10^{-1.88} (150 \times 10^{-3} \text{M})}$$

$$\approx 4.44 \times 10^{-2} \text{ M}$$
(4 pts)

glutamic acid

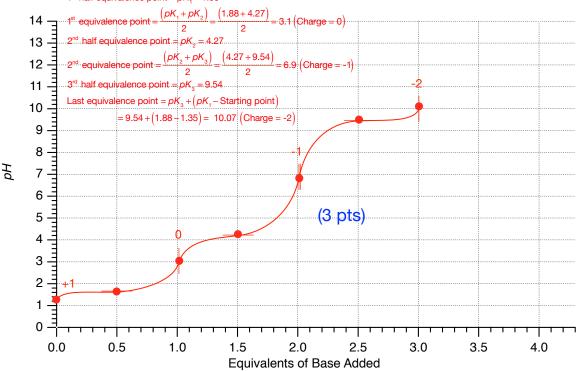
$$pH = -\log(\lfloor H^+ \rfloor)$$
$$= -\log(4.44 \times 10^{-2})$$

b. Sketch the titration curve for a 150 mM solution of glutamic acid.

Starting point = 1.35 (Charge = +1)

1st half equivalence point =
$$pK_1$$
 = 1.88

1st equivalence point = $\frac{(pK_1 + pK_2)}{(pK_1 + pK_2)}$ = $\frac{(pK_1 + pK_2)}{(pK_1 + pK_2)}$



(3 pts)

c. In solution, at what pH does the predominant (major) species of glutamic acid have a *net charge* of 0? pH =(The first equivalence or end point)

(This pH is also called the *isoelectric point*, pI, because at this pH glutamic acid does not move in an electrostatic field.)

d. Draw the structure for the predominant species of glutamic acid that is present at its isoelectric pH.

