
OpenCourseWare (2023)

CHEMISTRY II

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EVALUATION TEST 4



1. (3 points) Fill the spaces with the correct word/-s in the next statements. (Correct statements mark as +0.3 points).

A	In the following cell: $Mg(s) Mg^{2+}(0.1 M) Mg^{2+}(0.5 M) Mg(s)$ the more _____ solution is reduced in the cathode.
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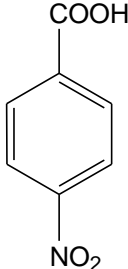
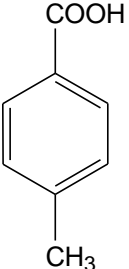
B	Given the following standard reduction potentials: $Ag^+/Ag = +0.80 V$ and $Cl_2/Cl^- = +1.36 V$. _____ is a stronger oxidizing agent than _____.
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C	_____ is a type of corrosion which occurs when a tensile stress is applied on the material under a corrosive environment.
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D	A radical is stabilized by resonance and hyperconjugation (an electron interaction between a _____ bond and a _____ orbital).
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E	A S_N2 reaction predominates in _____ solvents, when a _____ carbocation is formed, and the better the nucleophile is.
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F	<i>Trans</i> -disubstituted alkenes present a _____ melting point than the corresponding <i>cis</i> -disubstituted.
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G	<p>Given the following benzoic acids:</p> <div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;">  <p>I</p> </div> <div style="text-align: center;">  <p>II</p> </div> </div> <p>Compound I has _____ acidity than compound II, because nitro group is an _____ group.</p>
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H	In an _____ inhibition, the inhibitor can bind ONLY to enzyme-substrate complex, not to free enzyme.
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I	Fatty acids are composed of a long hydrocarbon chain (_____) ranging from 4 to 36 carbons long and a terminal _____ group (head).
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J	A nucleotide has three characteristic components: (1) a nitrogenous base, (2) a _____, and (3) one or more phosphates. The molecule without a phosphate group is called a _____.
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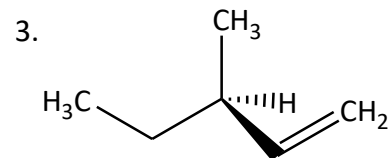
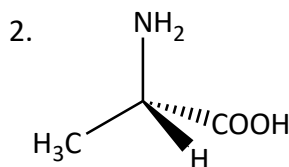
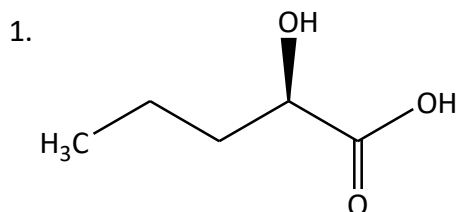
2. (1.5 points) Consider the electrolysis of an aqueous solution of KBr:

- a) (1 p) Write the half-reactions and indicate the products formed at the anode and cathode.
- b) (0.5 p) If the initial concentration of the salt (KBr in aqueous solution) is 0.5 M. Determine the pH of the medium after electrolysis.

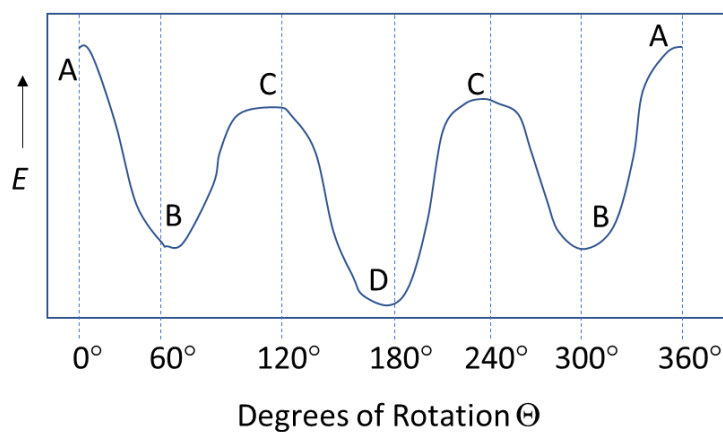
Data: $E^0(\text{K}^+/\text{K}) = -2.93 \text{ V}$; $E^0(\text{Br}_2/\text{Br}^-) = +1.07 \text{ V}$; $E^0(\text{O}_2/\text{H}_2\text{O}, \text{H}^+) = +1.23 \text{ V}$; $E^0(\text{H}_2\text{O}/\text{H}_2, \text{OH}^-) = -0.83 \text{ V}$.

3. (1.75 points) Answer the following questions:

a) (0.75 p) Write the name including the absolute configuration of the following compounds. (Hint: when alcohol group is not the main functional group, that is denominated as "hydroxy-" before the name of the main chain)

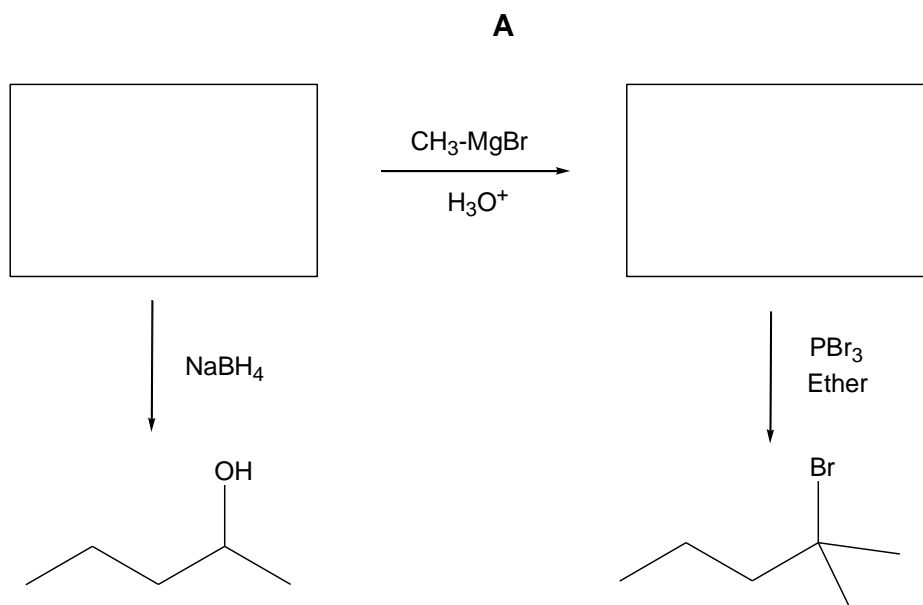


b) (1 p) Starting from the eclipsed conformation (A) for butane, draw and indicate the type of conformation using the Newman projections for B, C, D, and E according to the rotation of the C2–C3 bond indicated in the following graph. Deduce which of the structures is more stable and why.

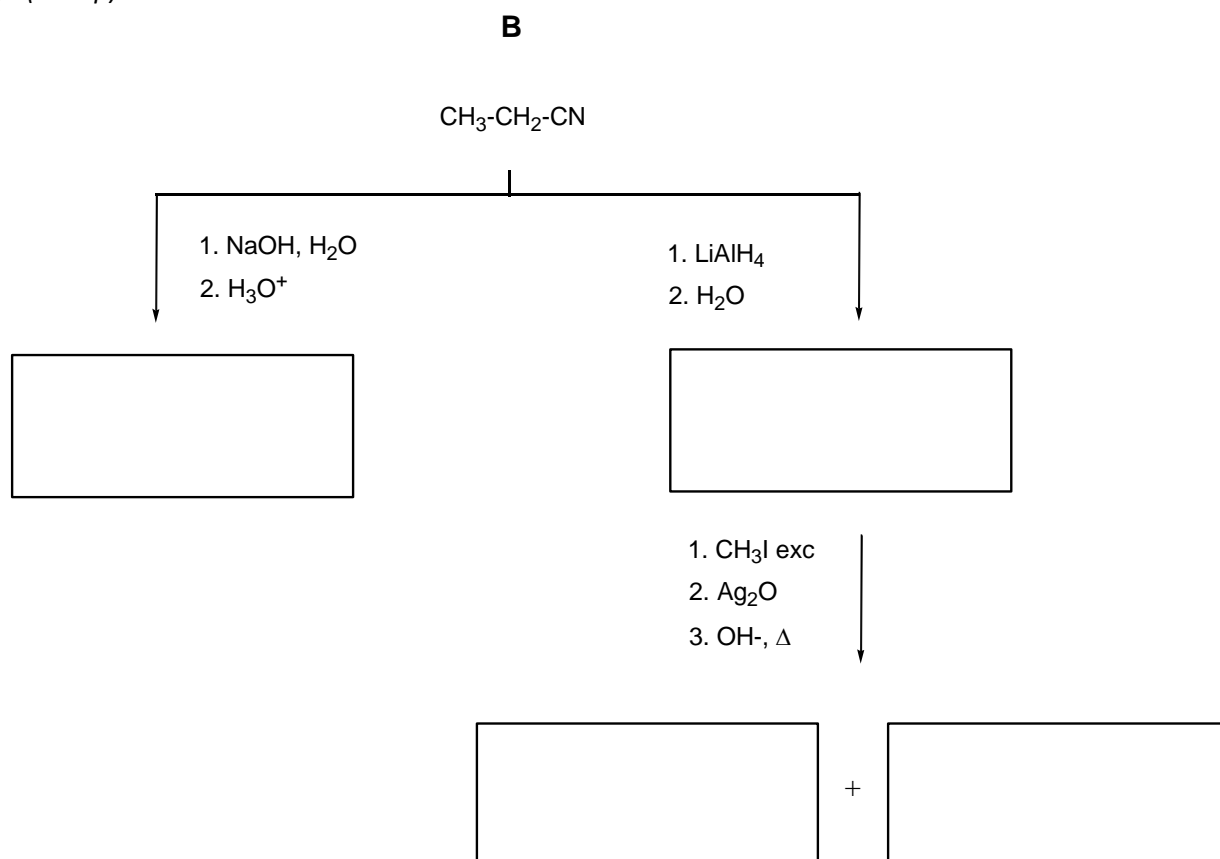


4. (2.25 points) Complete the following schemes:

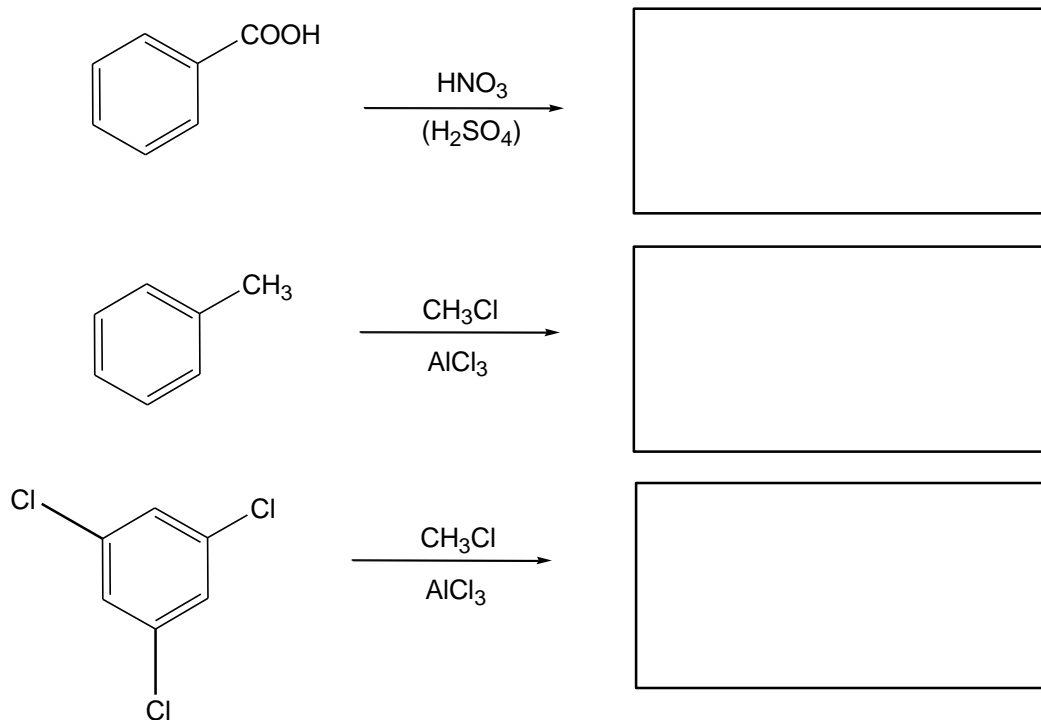
a) (0.75 p)



b) (0.75 p)



c) (0.75 p)

C

5. (1.5 points) A compound with the formula $\text{C}_3\text{H}_5\text{BrO}_2$ exhibits the following $^1\text{H-NMR}$, $^{13}\text{C-NMR}$, and IR spectra, respectively:

$^1\text{H-NMR}$: $\delta = 2.93$ (t, 2 H), 3.57 (t, 2 H), >9.0 (low field, 1 H very broad) ppm.

$^{13}\text{C-NMR}$: $\delta = 24.3, 38.6, 178.6$ ppm.

IR: characteristic bands $3067, 1717 \text{ cm}^{-1}$.

a) (0.25 p) Determine the degree of unsaturation.

b) (1.25 p) Justify its structure.

ANNEX

Chemical Shifts in ^1H NMR

Type of hydrogen		Chemical shift (δ)
Reference	$\text{Si}(\text{CH}_3)_4$	0
Alkyl (primary)	$-\text{CH}_3$	0.7–1.3
Alkyl (secondary)	$-\text{CH}_2-$	1.2–1.6
Alkyl (tertiary)	$\begin{array}{c} \\ -\text{CH}- \\ \end{array}$	1.4–1.8
Allylic	$\begin{array}{c} \text{H} \\ \\ \text{C}=\text{C}-\text{C}- \\ \end{array}$	1.6–2.2
Methyl ketone	$\begin{array}{c} \text{O} \\ \\ -\text{C}-\text{CH}_3 \end{array}$	2.0–2.4
Aromatic methyl	$\text{Ar}-\text{CH}_3$	2.4–2.7
Alkynyl	$-\text{C} \equiv \text{C}-\text{H}$	2.5–3.0
Alkyl halide	$\begin{array}{c} \text{H} \\ \\ -\text{C}-\text{Hal} \\ \end{array}$	2.5–4.0
Alcohol	$\begin{array}{c} \\ -\text{C}-\text{O}-\text{H} \\ \end{array}$	2.5–5.0
Alcohol, ether	$\begin{array}{c} \text{H} \\ \\ -\text{C}-\text{O}- \\ \end{array}$	3.3–4.5
Vinylic	$\begin{array}{c} \text{H} \\ \\ \text{C}=\text{C} \\ \end{array}$	4.5–6.5
Aryl	$\text{Ar}-\text{H}$	6.5–8.0
Aldehyde	$\begin{array}{c} \text{O} \\ \\ -\text{C}-\text{H} \end{array}$	9.7–10.0
Carboxylic acid	$\begin{array}{c} \text{O} \\ \\ -\text{C}-\text{O}-\text{H} \end{array}$	11.0–12.0

Characteristic IR bands of some common functional groups:

Functional Group		Absorption (cm^{-1})	Intensity
Alkane	C-H	2850–2960	Medium
Alkene	=C-H	3020–3100	Medium
	C=C	1640–1680	Medium
Alkyne	$\equiv\text{C-H}$	3300	Strong
	$\text{C}\equiv\text{C}$	2100–2260	Medium
Alkyl halide	C-Cl	600–800	Strong
	C-Br	500–600	Strong
Alcohol	O-H	3400–3650	Strong, broad
	C-O	1050–1150	Strong
Arene	C-H	3030	Weak
Aromatic ring		1660–2000	Weak
		1450–1600	Medium
Amine	N-H	3300–3500	Medium
	C-N	1030–1230	Medium
Carbonyl compound	$\text{C}=\text{O}$	1670–1780	Strong
	Aldehyde	1730	Strong
	Ketone	1715	Strong
	Ester	1735	Strong
	Amide	1690	Strong
	Carboxylic acid	1710	Strong

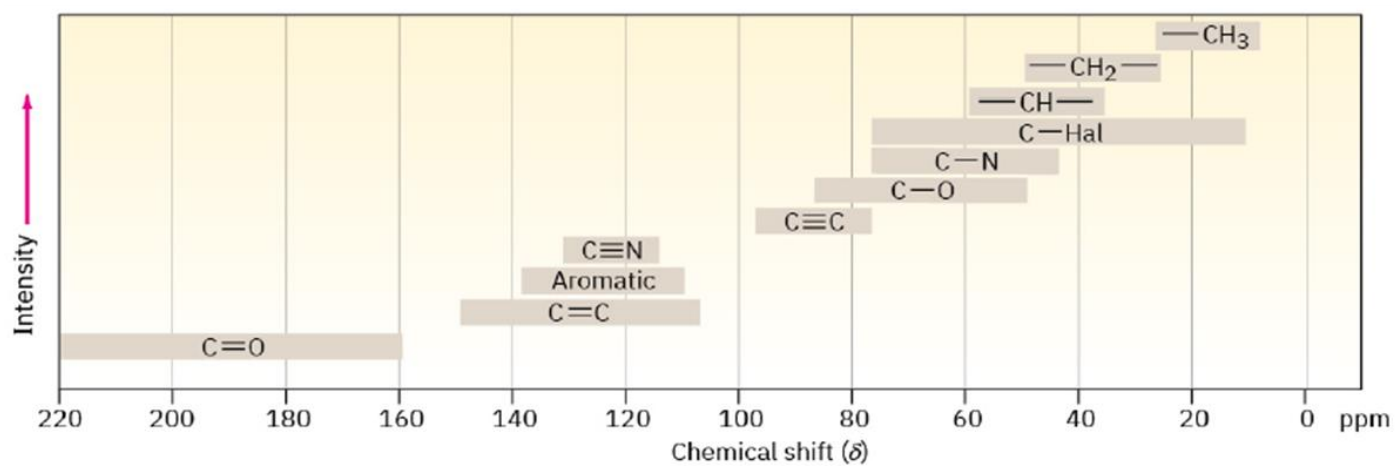
Chemical Shifts in ^{13}C NMR

IMAGE CREDITS

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