

NAME (Print): \_\_\_\_\_

Chemistry 320N  
1st Midterm Exam  
February 13, 2025

EID \_\_\_\_\_

SIGNATURE: \_\_\_\_\_

**Please print the  
first three letters  
of your last name  
in the three boxes**

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**Please Note:** Please take your time. You have three hours to take this exam. Please do not rush, we want you to show us everything you have learned this semester so far! Making careless mistakes is not good for anyone! If you find yourself getting anxious because of a problem, skip it and come back. Please do not second guess yourself! Keep track of the questions worth a lot of points. (This does not mean they are hard, it just means we think they cover important material.)

One last thing: I recommend you close your eyes for a moment, then take some nice deep breaths before you begin. **YOU GOT THIS!**

**FINALLY, DUE TO SOME UNFORTUNATE RECENT INCIDENTS YOU ARE NOT ALLOWED TO INTERACT WITH YOUR CELL PHONE IN ANY WAY. IF YOU TOUCH YOUR CELL PHONE DURING THE EXAM YOU WILL GET A "0" NO MATTER WHAT YOU ARE DOING WITH THE PHONE. PUT IT AWAY AND LEAVE IT THERE!!!**

# Student Honor Code for the University of Texas at Austin

"I pledge, as a member of The University of Texas at Austin community, to do my work honestly, respectfully, and through the intentional pursuit of learning and scholarship."

## Elaboration

1. I pledge to be honest about what I create and to acknowledge what I use that belongs to others.
2. I pledge to value the process of learning in addition to the outcome, while celebrating and learning from mistakes.
3. This code encompasses all of the academic and scholarly endeavors of the university community.

(Your signature)

## PERIODIC TABLE OF THE ELEMENTS

### ▼ Elementary Subatomic Particles

Electron	Proton	Neutron	Positron	Neutrino
Symbol $e^-$	$p^+$	$n^0$	$e^+$	$\nu$
Rest mass (kg)	$1.6726219 \times 10^{-27}$	$1.6749273 \times 10^{-27}$	$9.10938356 \times 10^{-31}$	0
Relative mass (approx)	1.836	1.839	0.511	0
Relative charge (approx)	+1	0	-1	0
Relative speed (approx)	$2.1876913 \times 10^{-8}$	$2.1876913 \times 10^{-8}$	$2.1876913 \times 10^{-8}$	0
Relative energy (approx)	$9.10938356 \times 10^{-31}$	$1.6749273 \times 10^{-27}$	$9.10938356 \times 10^{-31}$	0
Relative momentum (approx)	$9.10938356 \times 10^{-31}$	$1.6749273 \times 10^{-27}$	$9.10938356 \times 10^{-31}$	0
Relative spin (approx)	$1/2$	$1/2$	$1/2$	0
Relative magnetic moment (approx)	$9.2847637 \times 10^{-24}$	$1.8182715 \times 10^{-26}$	$9.2847637 \times 10^{-24}$	0
Relative gyromagnetic ratio (approx)	$1.7608596 \times 10^{11}$	$1.8182715 \times 10^{26}$	$1.7608596 \times 10^{11}$	0
Relative magnetic moment (approx)	$9.2847637 \times 10^{-24}$	$1.8182715 \times 10^{-26}$	$9.2847637 \times 10^{-24}$	0
Relative gyromagnetic ratio (approx)	$1.7608596 \times 10^{11}$	$1.8182715 \times 10^{26}$	$1.7608596 \times 10^{11}$	0

### ▼ Ionic Character of a Single Chemical Bond

Percent ionic character describes the nature of a bond. Bonds with  $\Delta\chi > 1.7$  are generally considered ionic. Bonds with  $\Delta\chi < 1.7$  are generally considered covalent. The curve shows that as the difference in electronegativity increases, the percent ionic character also increases.

1 IA	2 IIA	3 IIIA	4 IVA	5 VA	6 VIA	7 VIIA	8 VIIIA	9 VIIIA	10 VIIIA	11 IB	12 IIB	13 IIIB	14 IVA	15 VA	16 VIA	17 VIIA	18 VIII
1 H	2 He	3 Li	4 Be	5 B	6 C	7 N	8 O	9 F	10 Ne	11 Na	12 Mg	13 Al	14 Si	15 P	16 S	17 Cl	18 Ar
19 K	20 Ca	21 Sc	22 Ti	23 V	24 Cr	25 Mn	26 Fe	27 Co	28 Ni	29 Cu	30 Zn	31 Ga	32 Ge	33 As	34 Se	35 Br	36 Kr
37 Rb	38 Sr	39 Y	40 Zr	41 Nb	42 Mo	43 Tc	44 Ru	45 Rh	46 Pd	47 Ag	48 Cd	49 In	50 Sn	51 Sb	52 Te	53 I	54 Xe
55 Cs	56 Ba	57 La	58 Ce	59 Pr	60 Nd	61 Pm	62 Sm	63 Eu	64 Gd	65 Tb	66 Dy	67 Ho	68 Er	69 Tm	70 Yb	71 Lu	72 Hf
73 Ta	74 W	75 Re	76 Os	77 Ir	78 Pt	79 Au	80 Hg	81 Tl	82 Pb	83 Bi	84 Po	85 At	86 Rn	87 Fr	88 Ra	89 Ac	90 Th
91 Pa	92 U	93 Np	94 Pu	95 Am	96 Cm	97 Bk	98 Cf	99 Es	100 Fm	101 Md	102 No	103 Lr	104 Rf	105 Db	106 Sg	107 Bh	108 Hs
109 Mt	110 Ds	111 Nh	112 Fl	113 Lv	114 Ts	115 Og	116 Uu	117 Uub	118 Uuq	119 Uuq	120 Uuq	121 Uuq	122 Uuq	123 Uuq	124 Uuq	125 Uuq	126 Uuq

**PAPERTECH** Editors: T. K. Varga, M.A.Sc. & C. Bello, M.A.Sc.

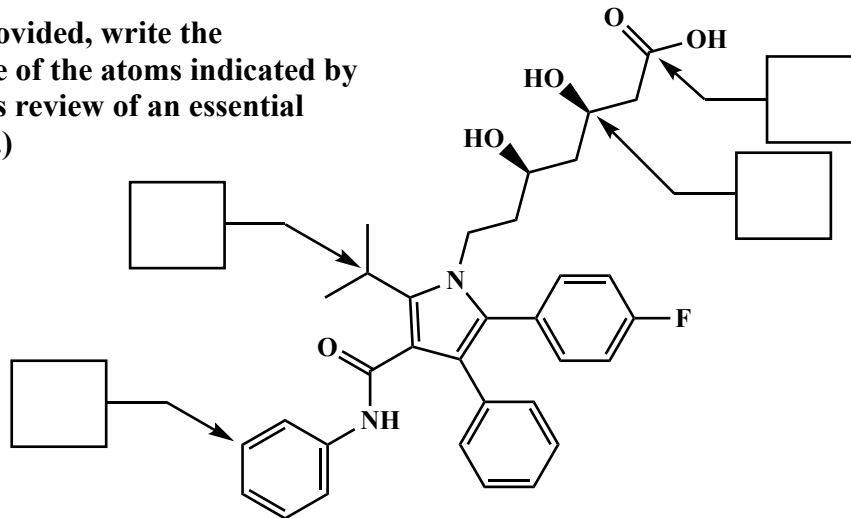
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Compound		pK <sub>a</sub>
Hydrochloric acid	$\text{H-Cl}$	-7
Protonated alcohol	$\text{RCH}_2\text{OH}_2^{\oplus}$	-2
Hydronium ion	$\text{H}_3\text{O}^{\oplus}$	-1.7
Carboxylic acids	$\text{R}-\overset{\text{O}}{\parallel}{\text{C}}-\text{H}$	3-5
Thiols	$\text{RCH}_2\text{SH}$	8-9
Ammonium ion	$\text{H}_4\text{N}^{\oplus}$	9.2
β-Dicarbonyls	$\text{RC}-\overset{\text{O}}{\parallel}{\text{C}}-\text{CH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{R}'$	10
Primary ammonium	$\text{H}_3\text{N}^{\oplus}\text{CH}_2\text{CH}_3$	10.5
β-Ketoesters	$\text{RC}-\overset{\text{O}}{\parallel}{\text{C}}-\text{CH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{OR}'$	11
β-Diesters	$\text{ROC}-\overset{\text{O}}{\parallel}{\text{C}}-\text{CH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{OR}'$	13
Water	$\text{HOH}$	15.7
Alcohols	$\text{RCH}_2\text{OH}$	15-19
Acid chlorides	$\text{RCH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{Cl}$	16
Aldehydes	$\text{RCH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{H}$	18-20
Ketones	$\text{RCH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{R}'$	18-20
Esters	$\text{RCH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{OR}'$	23-25
Terminal alkynes	$\text{RC}\equiv\text{C}-\text{H}$	25
LDA	$\text{H}-\text{N}(\text{i-C}_3\text{H}_7)_2$	40
Terminal alkenes	$\text{R}_2\text{C}=\underset{\text{H}}{\text{C}}-\text{H}$	44
Alkanes	$\text{CH}_3\text{CH}_2-\text{H}$	51



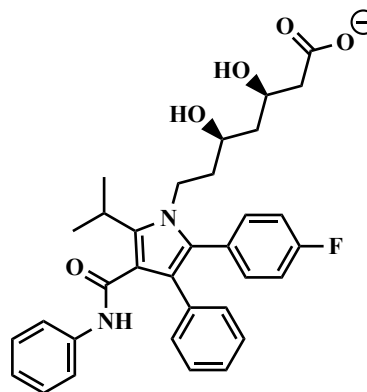
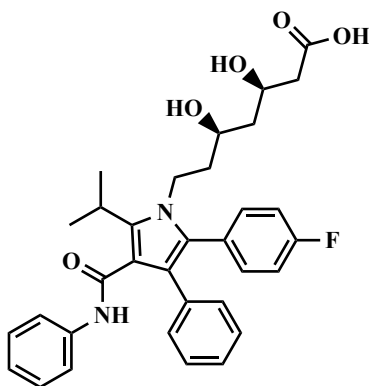
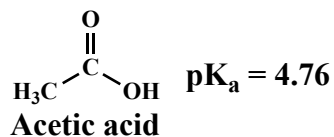
4. (8 pts) Lipitor (Atorvastatin) is a very important drug used to treat prevent atherosclerosis and heart disease. In 2023 it was 24th best selling drug, earning \$2.2 billion in sales. Answer the following questions about lipitor.

A. In the boxes provided, write the hybridization state of the atoms indicated by the arrow. (This is review of an essential OChem 1 concept.)



Lipitor (Atorvastatin)

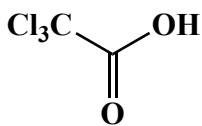
B. (6 pts) Given that the  $pK_a$  value shown for acetic acid is typical for a carboxylic acid group, fill in the circle to indicate the pH at which the protonation state of lipitor shown above would be present. You might need to fill in more than one circle, as the given structure might be present at more than one of the pH values listed.



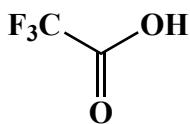
- The predominant form present at pH = 2  
 The predominant form present at pH = 7  
 The predominant form present at pH = 10

- The predominant form present at pH = 2  
 The predominant form present at pH = 7  
 The predominant form present at pH = 10

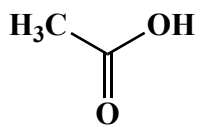
5. (4 pts) Predict the relative acidities of the following molecules. Put the number 1 under the most acidic molecule, the number 4 under the least acidic molecule, and the numbers 2 and 3 under the other two structures as appropriate.



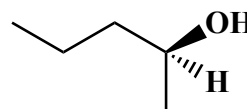
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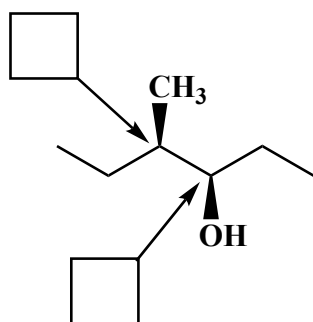


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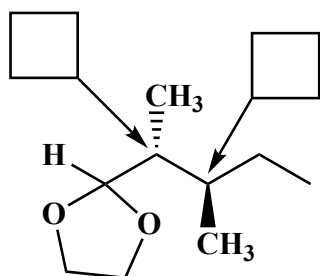
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6. (12 pts) For each pair of molecules, fill in the boxes to label each chiral center as R or S then fill in the circle that indicates the appropriate relationship between the two molecules.



Relationship:

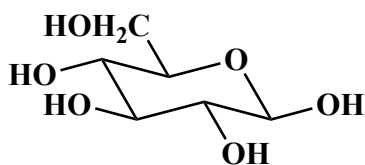
- Enantiomers  
 Diastereomers  
 Same molecule



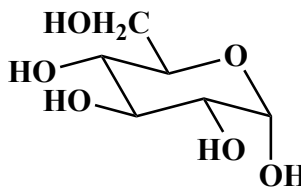
Relationship:

- Enantiomers  
 Diastereomers  
 Same molecule

7. (4 pts each) Fill in the circle to indicate the correct name for the glucose molecules shown.



- alpha( $\alpha$ ) - D - Glucose  
 beta( $\beta$ ) - D - Glucose



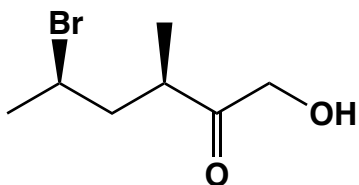
- alpha( $\alpha$ ) - D - Glucose  
 beta( $\beta$ ) - D - Glucose

Signature \_\_\_\_\_

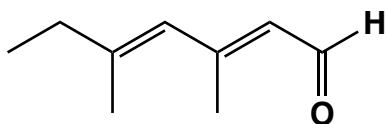
Pg 4 \_\_\_\_\_ (12)

8. (4 pts each) Write an acceptable IUPAC name or draw a structural formula for the following molecules:

A.

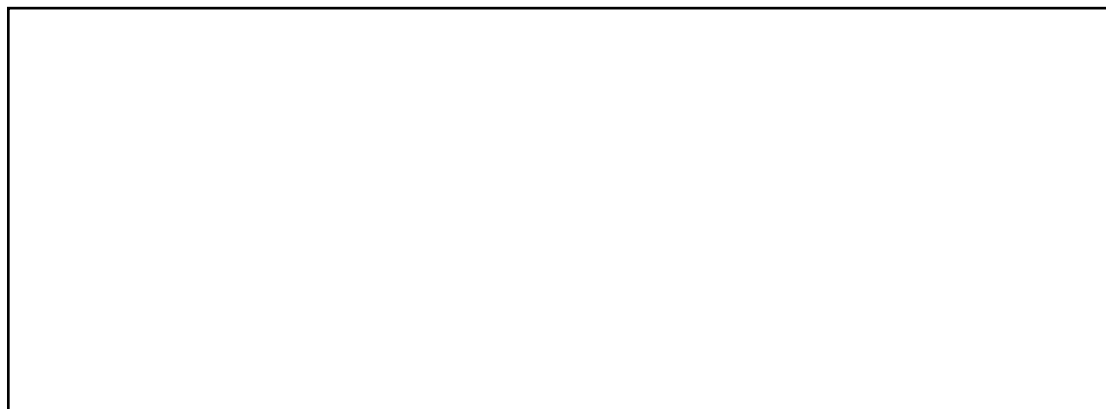


B.

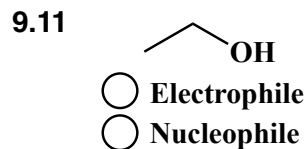
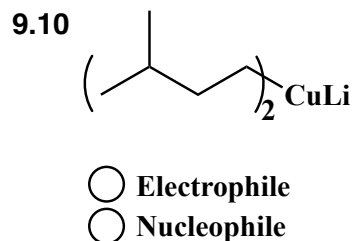
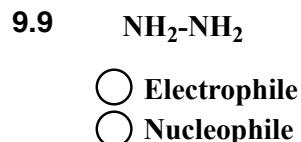
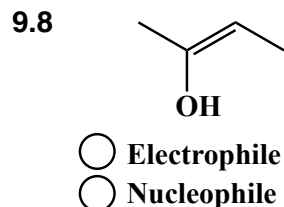
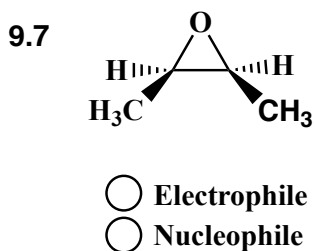
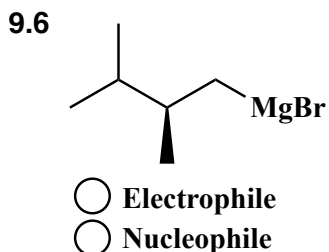
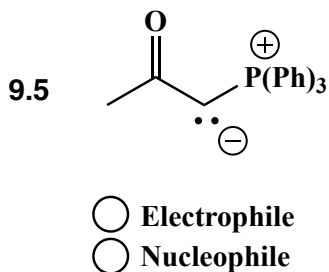
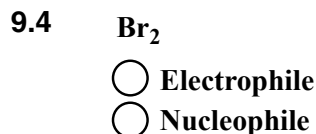
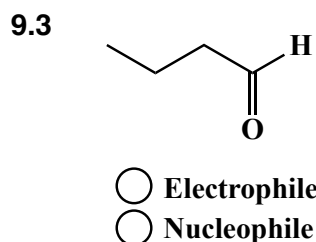
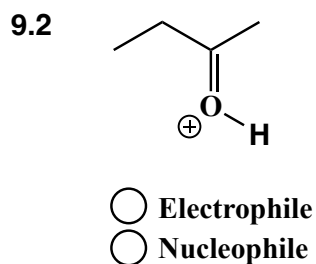
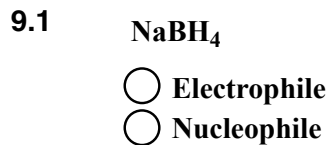


C. In the box, draw the structure corresponding to the following IUPAC name.

**(2*R*,3*R*)-2,3-dimethyl-5-oxoheptanal**



9. (11 pts) Being able to recognize the chemical personality of different species is one of the most important skills you can develop in Organic Chemistry. Fill in the appropriate circle to indicate whether each structure is a nucleophile or electrophile in the mechanisms we have seen. Note that these species might be acids or bases in certain situations, but we will ignore that for this problem.



10. (6 pts) Carboxylic acids exist as a characteristic dimer in solution, held together with hydrogen bonds. In the box provided, draw the hydrogen bonded dimer of acetic acid. Indicate the hydrogen bonds as a dashed line, and draw all lone pairs on your structures. You saw this in lecture last Thursday.



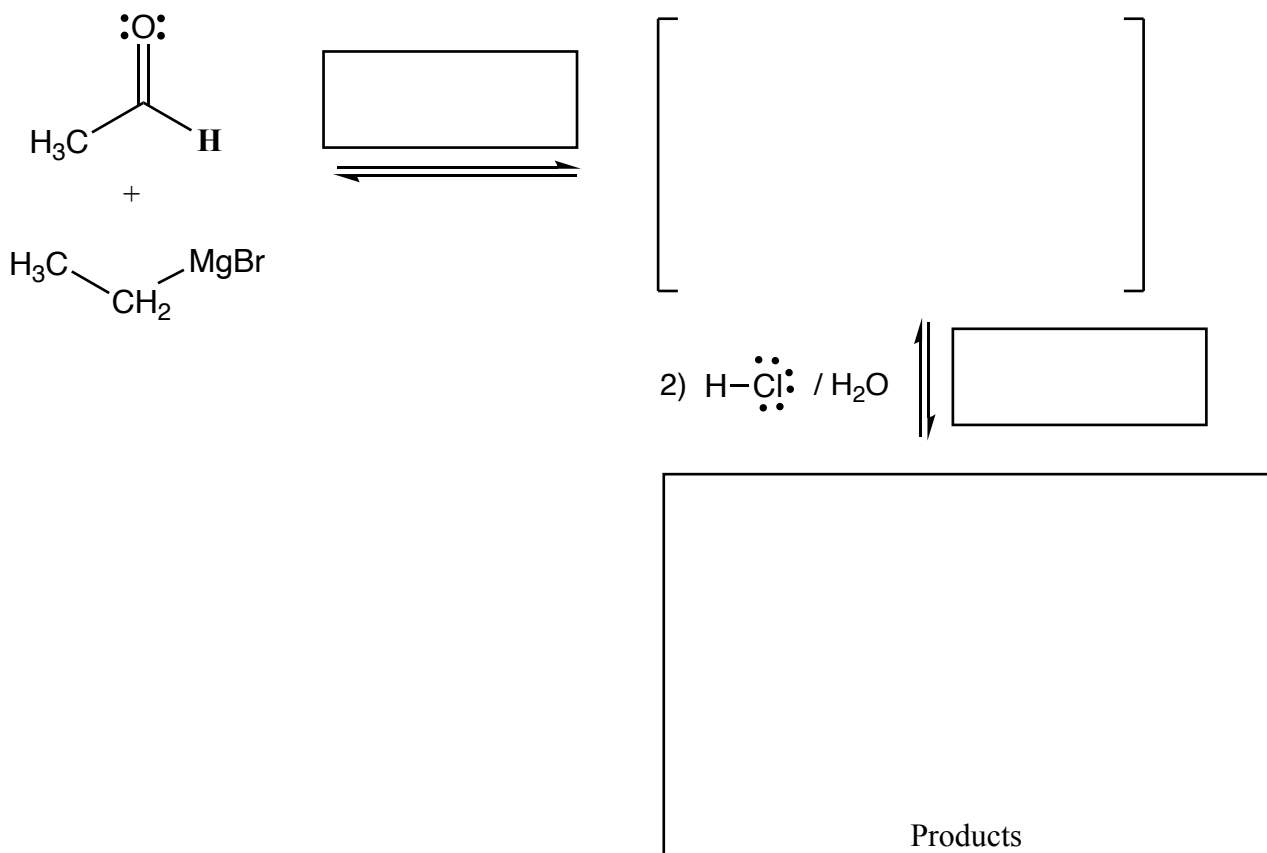


Signature \_\_\_\_\_

Pg 6 \_\_\_\_\_(12)

11. (12 pts) For this Grignard reaction, use **arrows to indicate movement of all electrons, write all lone pairs, all formal charges, and all the products for each step.** Remember, I said all the products for each step. **IF A NEW CHIRAL CENTER IS CREATED IN AN INTERMEDIATE, MARK IT WITH AN ASTERISK AND LABEL THE MOLECULE AS “RACEMIC” IF APPROPRIATE. FOR ALL CHIRAL FINAL “PRODUCTS” YOU MUST DRAW ALL ENANTIOMERS WITH WEDGES AND DASHES AND WRITE “RACEMIC” IF APPROPRIATE.** In the boxes provided by the arrows, write which of the 4 most common mechanistic elements describes each step (make a bond, break a bond, etc.).

*Grignard Reaction with an Aldehyde*

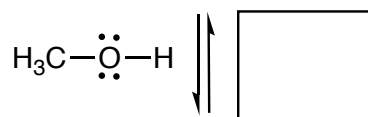
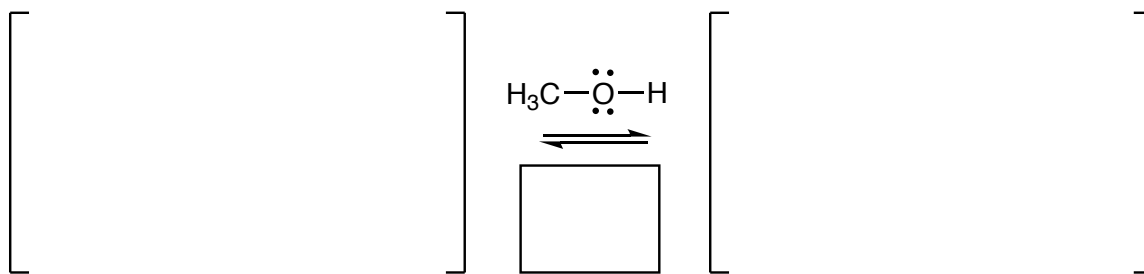
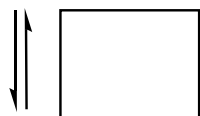
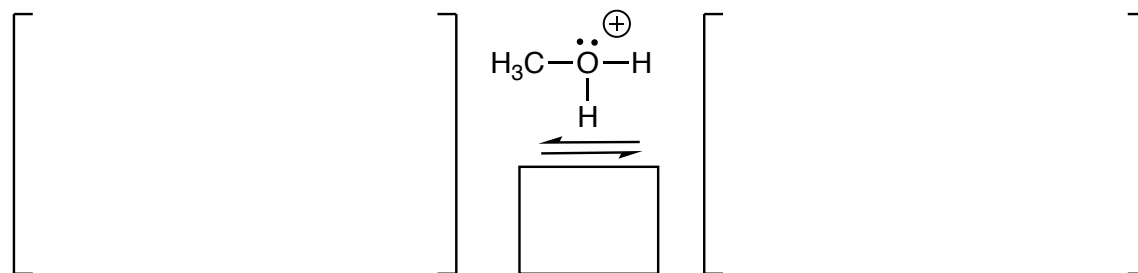
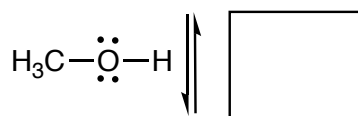
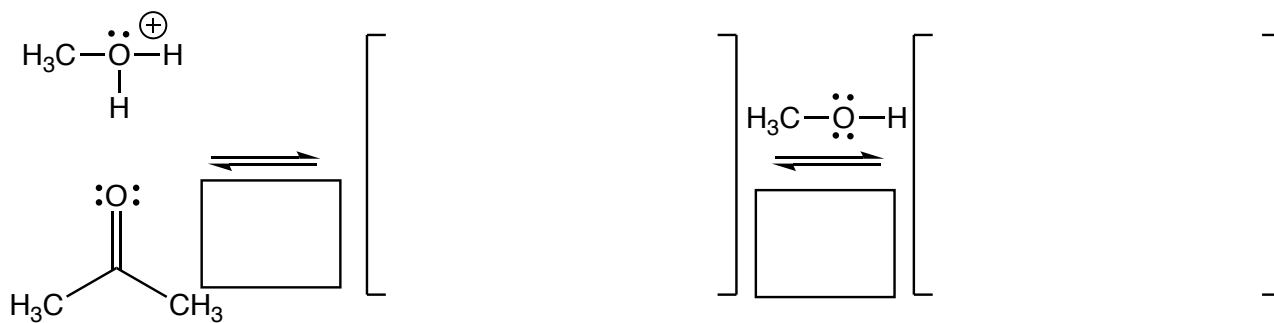


Signature \_\_\_\_\_

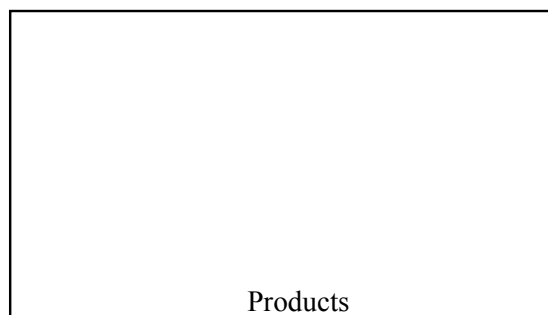
Pg 7 \_\_\_\_\_(-)

**12.** (45 pts) For the acetal reaction mechanism ON THE NEXT PAGE, use **arrows to indicate movement of all electrons, write all lone pairs, all formal charges, and all the products for each step.** Remember, I said all the products for each step. **IF A NEW CHIRAL CENTER IS CREATED IN AN INTERMEDIATE, MARK IT WITH AN ASTERISK AND LABEL THE MOLECULE AS “RACEMIC” IF APPROPRIATE. FOR ALL CHIRAL “PRODUCTS” YOU MUST DRAW ALL ENANTIOMERS WITH WEDGES AND DASHES AND WRITE “RACEMIC” IF APPROPRIATE.** In the boxes provided by the arrows, write which of the 4 most common mechanistic elements describes each step (make a bond, break a bond, etc.).

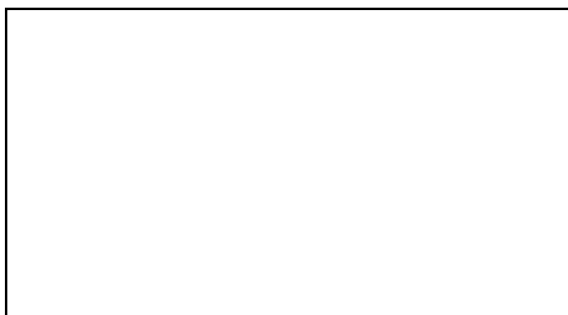
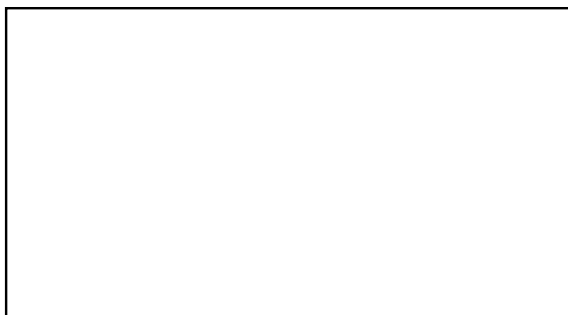
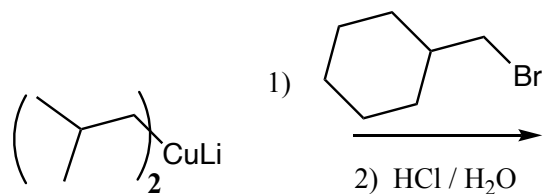
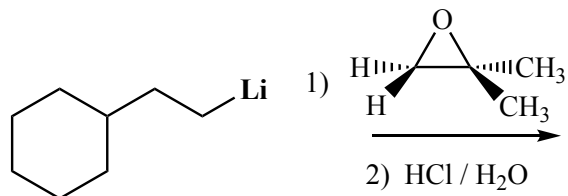
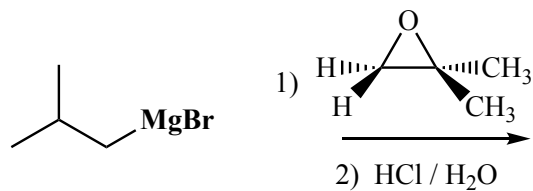
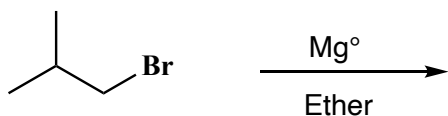
**The mechanism did not fit on the same page as the directions, so use the directions on this page to fill in the mechanism on the next page!**



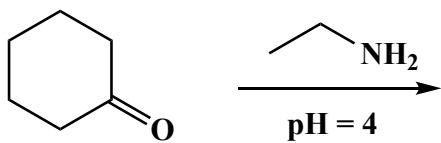
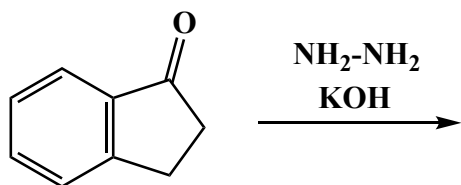
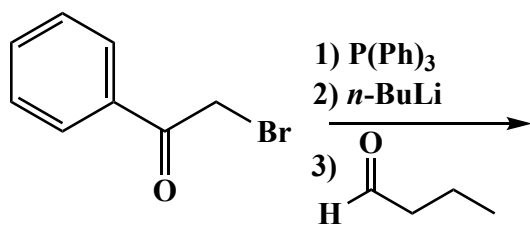
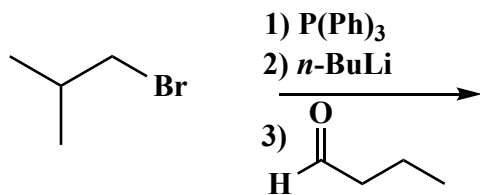
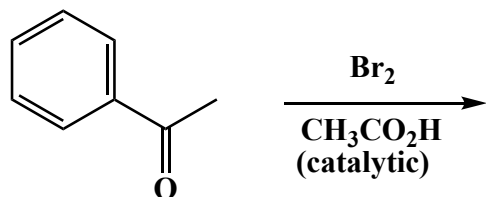
Note: For intermediates that I drew as two contributing structures in class, you only need to draw one contributing structure. Either one will be correct. Just make sure your arrows are accurate for the contributing structure you draw.



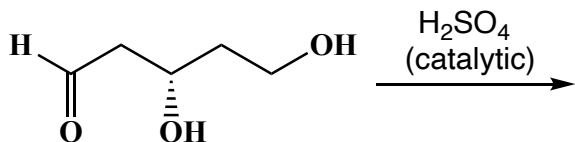
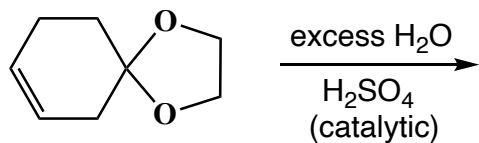
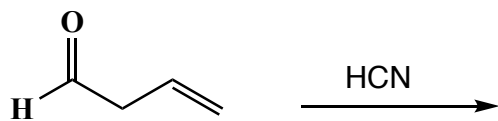
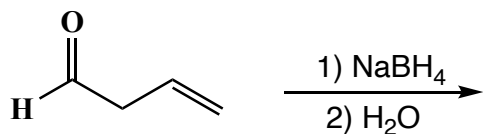
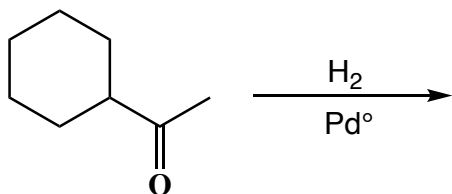
13. (3 or 5 pts.) Write the predominant product or products that will occur for each transformation. If a new chiral center is created and a racemic mixture is formed, you must draw both enantiomers and write "racemic" under the structure. Use wedges (  $\blacktriangleleft$  ) and dashes (  $\dashv$  ) to indicate stereochemistry. To get full credit, you only need to write the the major organic product for these. You do not have to worry about the other products.



13. (cont.) (3, 4 or 5 pts.) Write the predominant product or products that will occur for each transformation. If a new chiral center is created and a racemic mixture is formed, you must draw both enantiomers and write "racemic" under the structure. Use wedges (  $\blacktriangleleft$  ) and dashes (  $\cdots$  ) to indicate stereochemistry. To get full credit, you only need to write the the major organic product for these. You do not have to worry about the other products.



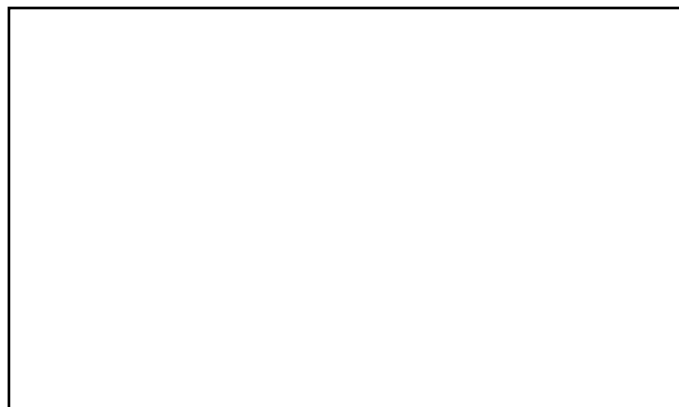
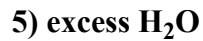
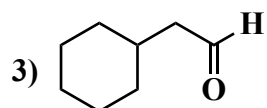
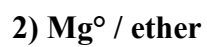
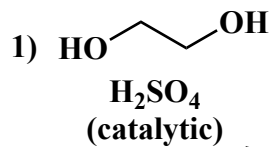
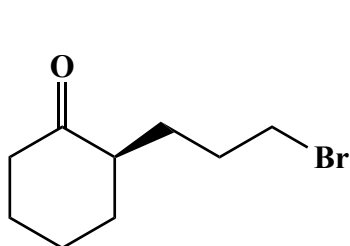
13. (cont.) (3, 4 or 5 pts.) Write the predominant product or products that will occur for each transformation. If a new chiral center is created and a racemic mixture is formed, you must draw both enantiomers and write "racemic" under the structure. Use wedges (  $\blacktriangleleft$  ) and dashes (  $\cdots$  ) to indicate stereochemistry. To get full credit, you only need to write the the major organic product for these. You do not have to worry about the other products.



Signature \_\_\_\_\_

Pg 12 \_\_\_\_\_(12)

14. (10 pts) Here is a warm-up for the synthesis problems. For the following series of reactions, write the **final** product(s) that you will see. Make sure draw all stereoisomers produced and to use wedges and dashes to indicate all stereochemistry, and you must write racemic if appropriate.

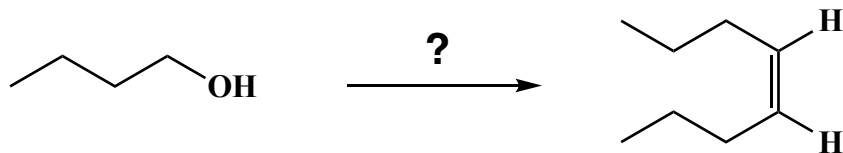


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Pg 13 \_\_\_\_\_(10)

15. These are synthesis questions. You need to show how the starting material can be converted into the product(s) shown. You may use any reactions we have learned provided that the product(s) you draw for each step is/are the predominant one(s). Show all the reagents you need. Show each molecule synthesized along the way and be sure to pay attention to the regiochemistry and stereochemistry preferences for each reaction. You must draw all stereoisomers formed, and use wedges and dashes to indicate chirality at each chiral center. Write racemic when appropriate. **All the carbons of the product must come from carbons of the starting material.**

A) (10 pts)



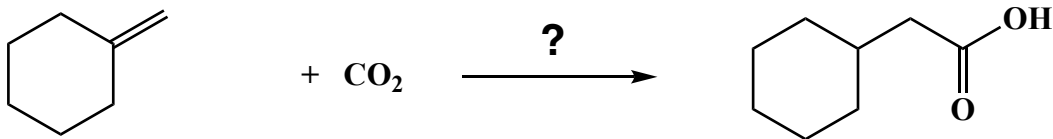


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Pg 14 \_\_\_\_\_ (10)

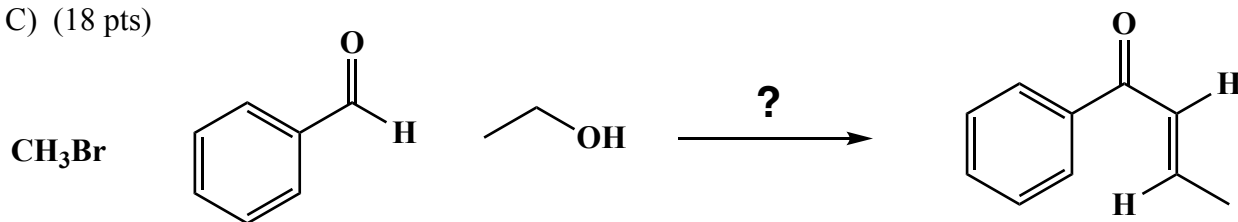
15. These are synthesis questions. You need to show how the starting material can be converted into the product(s) shown. You may use any reactions we have learned provided that the product(s) you draw for each step is/are the predominant one(s). Show all the reagents you need. Show each molecule synthesized along the way and be sure to pay attention to the regiochemistry and stereochemistry preferences for each reaction. You must draw all stereoisomers formed, and use wedges and dashes to indicate chirality at each chiral center. Write racemic when appropriate. **All the carbons of the product must come from carbons of the starting material.**

B) (10 pts)

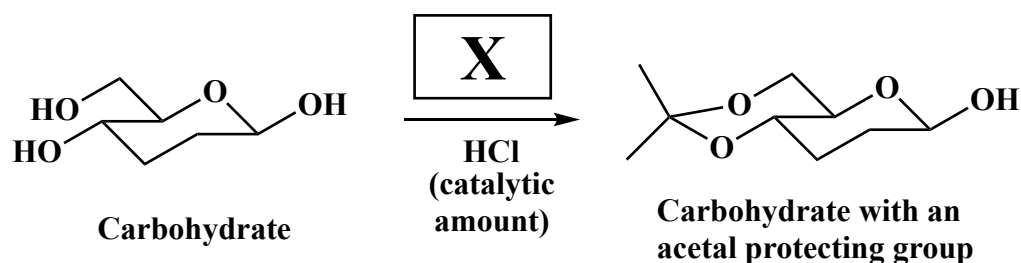


**15 (cont.)** These are synthesis questions. You need to show how the starting material can be converted into the product(s) shown. You may use any reactions we have learned provided that the product(s) you draw for each step is/are the predominant one(s). Show all the reagents you need. Show each molecule synthesized along the way and be sure to pay attention to the regiochemistry and stereochemistry preferences for each reaction. You must draw all stereoisomers formed, and use wedges and dashes to indicate chirality at each chiral center. Write racemic when appropriate. **All the carbons of the product must come from carbons of the starting material.**

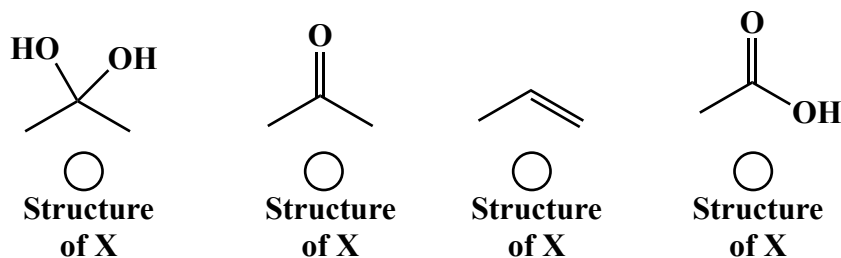
C) (18 pts)



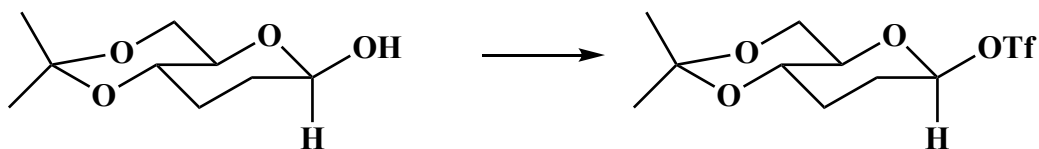
16. (12 pts) Here is an “Apply What you Know” problem. You have not seen all of this directly, but based on what you know you CAN figure it out. A major effort in modern organic chemistry involves the synthesis of carbohydrates so that we can study their properties and biochemistry. It is common to use a protecting group to prevent -OH groups on carbohydrates from taking part in unwanted reactions. The following is a common strategy to protect nearby -OH groups on a carbohydrate using an acetal.



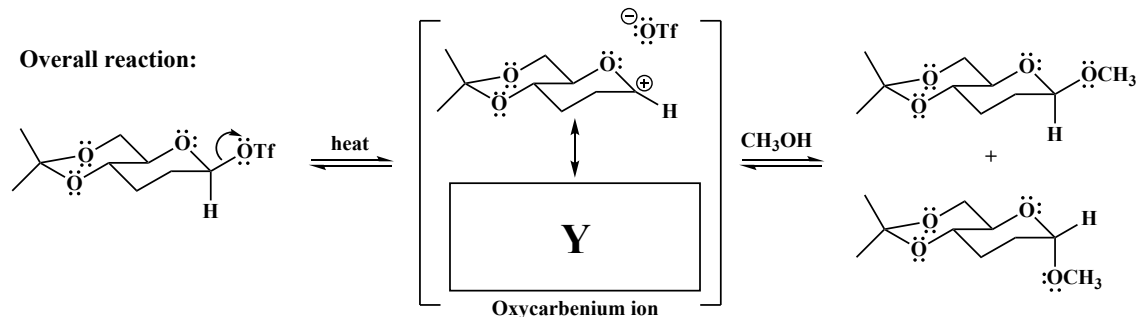
(4 pts) Look carefully at the carbohydrate with an acetal protecting group and fill in the circle beneath the molecule that corresponds to the appropriate structure “X”.



Chemists use these protected carbohydrates to make new bonds at the anomeric carbon atom. They first turn the OH on the anomeric carbon into a great leaving group. Here -OTf corresponds to something called a triflate group, the structure of which is not important. All you need to know is that the -OTf group is a great leaving group.



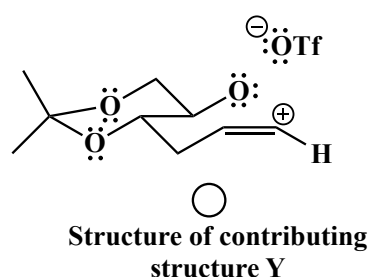
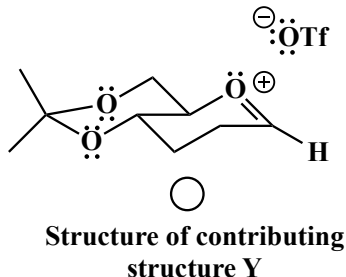
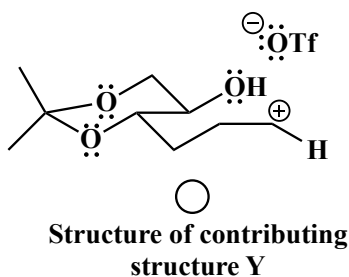
When these carbohydrates are heated, the leaving group departs to give a cation called an oxycarbenium ion intermediate that reacts with an alcohol such as methanol to make a new bond at the anomeric carbon as shown.



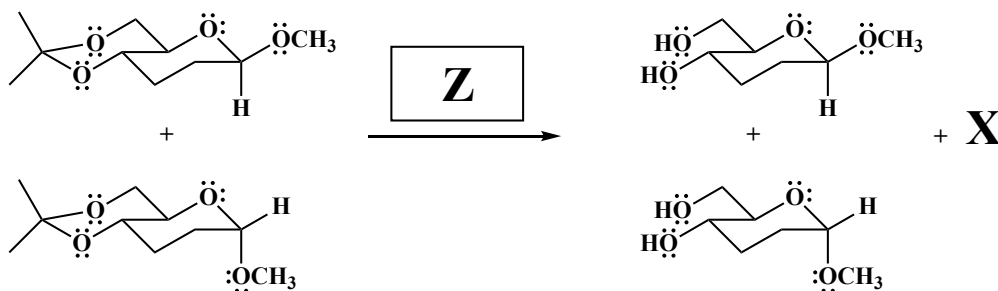
(4 pts) Fill in the appropriate circle. This overall reaction is best described as:

- An S<sub>N</sub>1 reaction     
  An S<sub>N</sub>2 reaction     
  An E1 reaction     
  An E2 reaction

(4 pts). Fill in the circle beneath the structure for Y that corresponds to the other important contributing structure of the oxycarbenium ion intermediate shown above.



To finish the reaction, the chemist must remove the acetal protecting group to regenerate the two -OH groups as well as the reagent "X" from the first part of this problem.



(4 pts) Based on what you know about acetals, fill in the circle under the reagent(s) "Z" that a chemist should use to remove the acetal protecting group.

- |                          |   |   |  |
|--------------------------|---|---|--|
| CH <sub>3</sub> OH       | NH <sub>2</sub> -NH <sub>2</sub><br>HO <sup>⊖</sup> | 1) NaBH <sub>4</sub><br>2) H <sub>2</sub> O | excess H <sub>2</sub> O<br>HCl or H <sub>2</sub> SO <sub>4</sub><br>(catalytic amount) |
| <input type="radio"/>    | <input type="radio"/>                               | <input type="radio"/>                       | <input type="radio"/>  |
| Reagents that could be Z | Reagents that could be Z                            | Reagents that could be Z                    | Reagents that could be Z   |