

CHEM 51C LEC A (40620)



Midterm 1 (Spring Qtr 2018) - LETTER SIZE

8527 (4018)

ver. D

Assigned Seat#: _____

Instructions to Instructor:

Do not alter this coversheet in ANY way. Substantial delays and additional fees may apply.

Instructions to Student:

1. Clearly print your Last Name, First Name and the Date
2. Clearly print your Student ID number in the boxes provided. Use large, dark numbers. These numbers are captured automatically during the scanning process.
3. Bubble in each number of your Student ID completely. The bubbles are used only if your written ID number is not captured.
4. Write your Name and Student ID number in the upper right corner of all following pages of your exam.

Last Name, First Name: Key

Date: ___/___/___

STUDENT ID:

For Access UCI student, leave first column blank then enter your 7-digit Student ID number.

1	0	0	0	0	0	0	0	0	1
2	0	0	0	0	0	0	0	0	2
3	0	0	0	0	0	0	0	0	3
4	0	0	0	0	0	0	0	0	4
5	0	0	0	0	0	0	0	0	5
6	0	0	0	0	0	0	0	0	6
7	0	0	0	0	0	0	0	0	7
8	0	0	0	0	0	0	0	0	8
9	0	0	0	0	0	0	0	0	9
0	0	0	0	0	0	0	0	0	0

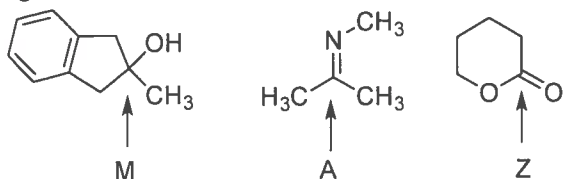
----- (This space for Instructor/TA use only) -----

Question	1	2	3	4	5	Total
Score	18	15	20	14	8	75

Do not open your exam until instructed to do so.
Answer the questions you understand best first.
 Your answers must be neat and legible.

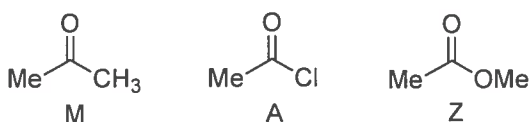
1. (22 points)

a. Rank highest to lowest oxidation state



MIFB #1d
 [Z] > [A] > [M] } 4

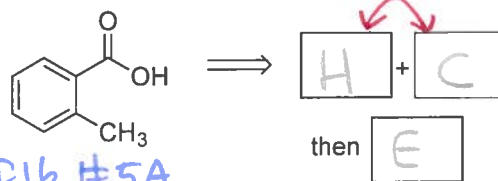
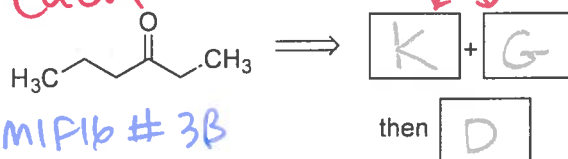
b. Rank fastest to slowest reaction with H₃CMgBr



[A] > [M] > [Z]

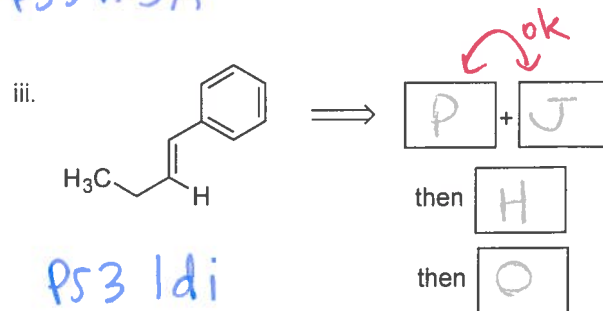
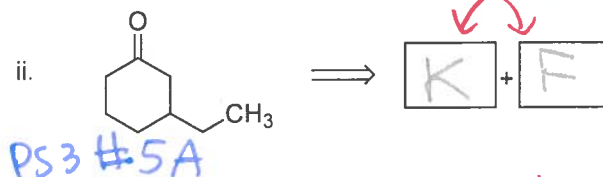
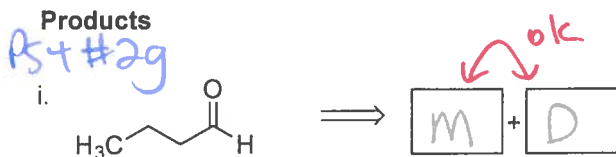
c. Fill in the correct compounds from the table to complete the retrosyntheses. You can use the same compound more than once.

Compounds			
D	PCC	C	Mg ⁰
E	CO ₂	F	NaCN
L		G	LiCu(CH ₂ CH ₃) ₂
H		K	
		M	



d. Fill in the correct compounds from the table to complete the retrosyntheses. You can use the same compound more than once.

Compounds			
C	LiAlH ₄	K	
D	Dibal-H	L	
E	BrMgCH ₂ CH ₃	M	
F	LiCu(CH ₂ CH ₃) ₂	N	
G	Mg ⁰	O	
H	n-BuLi		
J	PPh ₃		
P			



18/20

MIF16 #3b.
PS4 #6

2. 15 points

a. Match the names of the functional groups with labeled examples (5 points).

Initials: A

1 point each

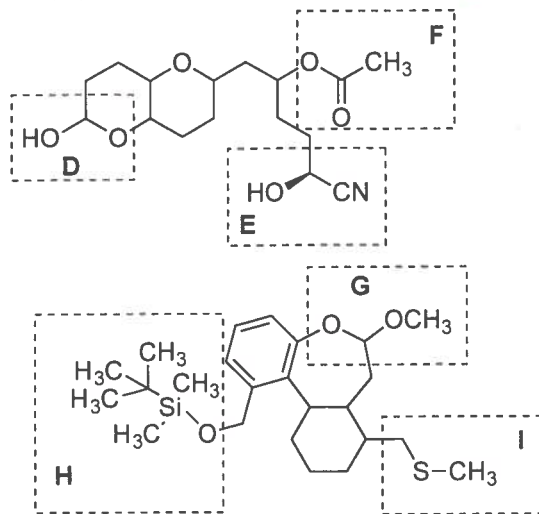
acetal B acetal should be G

1° amide A

ester F

cyanohydrin E

silyl ether H



1 pt each
5 max

b. Provide pKa's for any 5 of the following compounds (if you do them all, we will count your best 5).

MIF14 #1b
MIF16 #1c

<chem>CC(=O)C</chem>	<chem>CC(=O)O</chem>	<chem>CCO</chem>	HCl	<chem>CH4</chem>	<chem>H2SO4</chem>	<chem>CC(C)NC(C)C</chem>
20	5	16	-7	50	-3	36
18-22	4-6	14-18	-10 to -5	45-55	-11 to -7	32-38

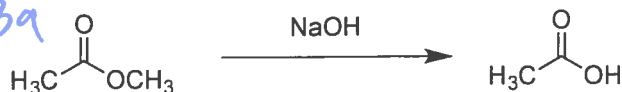
OK range

c. Provide an arrow-pushing mechanism (7 points).

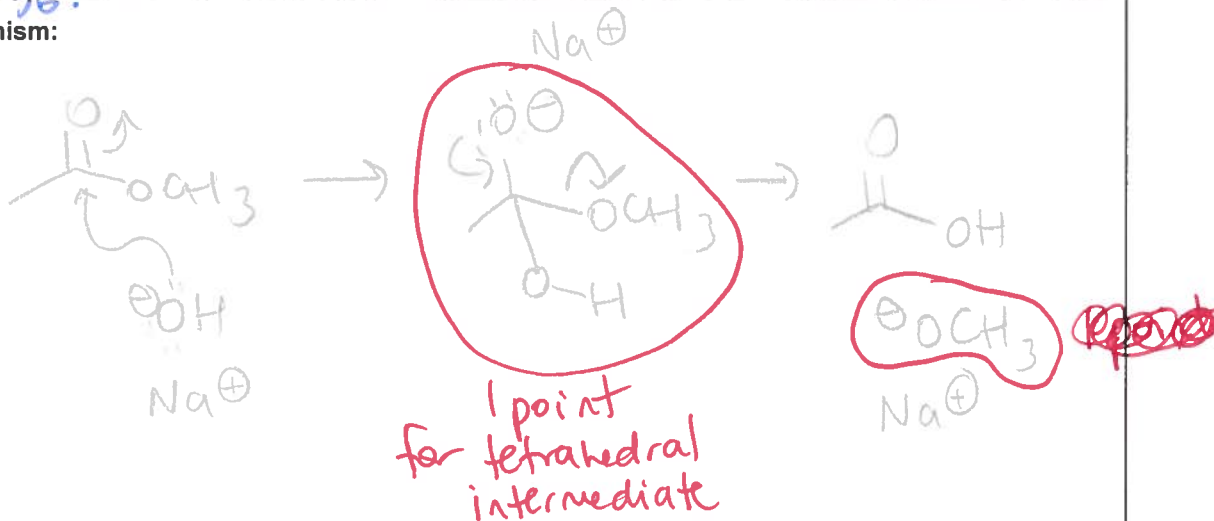
1 pt each arrow
start on electrons } -0.5
end on atom }

MIF16, 39

MIF14, #36.



Mechanism:



errors: charges -0.5

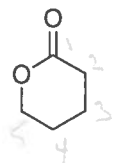
15

unnecessary proton transfers minus 1

3. (26 points) Fill in the boxes with the appropriate starting material, reagent or major product. Show stereochemistry where appropriate.

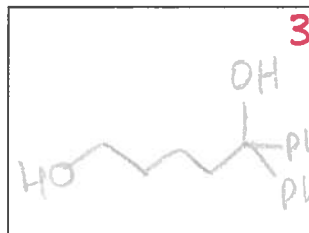
Initials: A

a.



1. PhMgBr (excess)

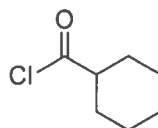
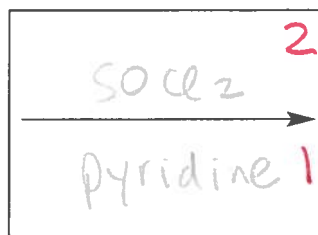
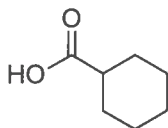
2. H₂O



What is the name for this type of reagent?

Grignard

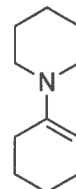
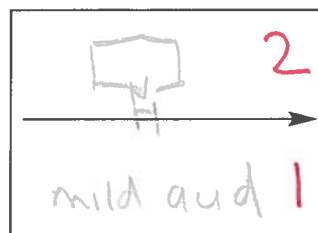
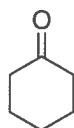
b.



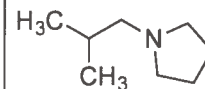
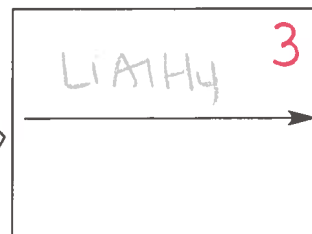
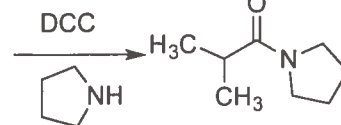
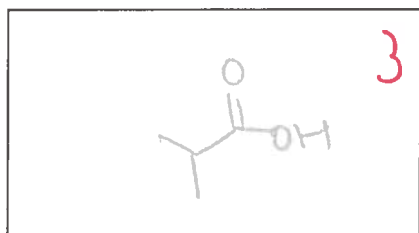
Is this reaction an oxidation, a reduction, or neither?

neither

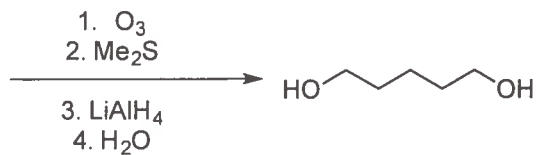
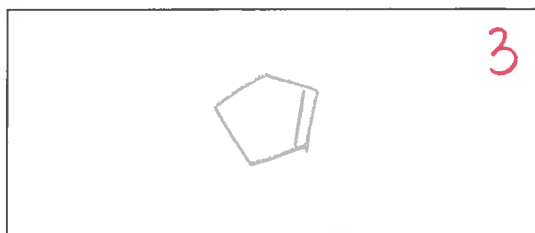
c.



d.



e.



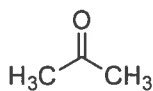
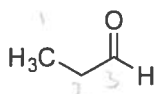
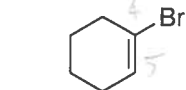
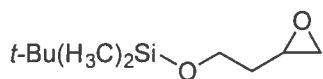
20

Initials: B

4. Propose syntheses of the target below (14 points).

All carbons in the product must come from the starting materials provided, you can use any reagent you wish.
YOU CAN IGNORE STEREOCHEMISTRY.

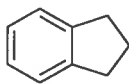
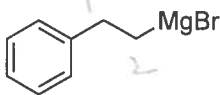
Starting Materials:



NaCN

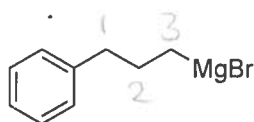
CH₃I

CO₂



ps2 #5A
ps2 #5B

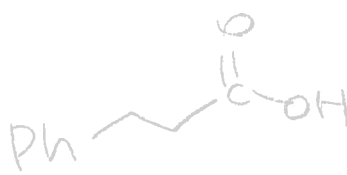
Target A.



(X)
(u)



HBr = 0

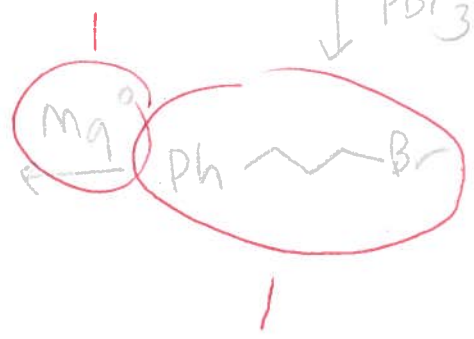


LiAlH₄

o- NaBH₄
o- B DIBAL-H



PBr₃

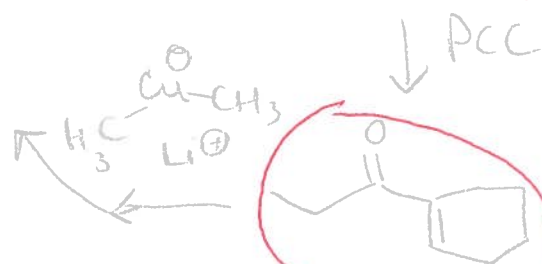
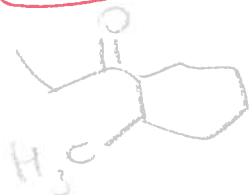
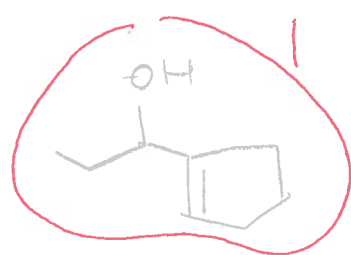
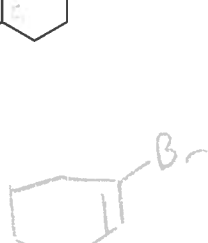
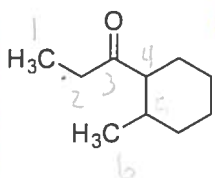


Target B.

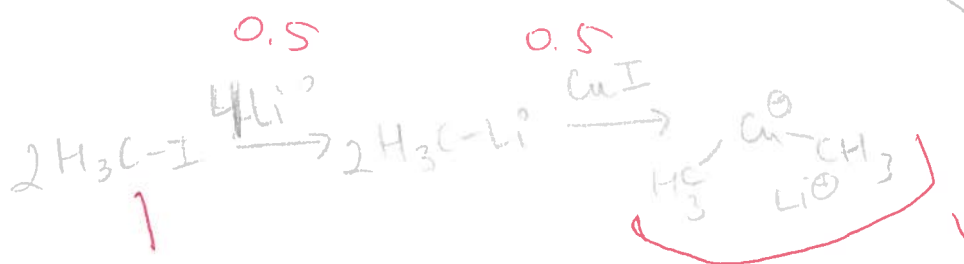
ps2 #5C

MI F14 #5A

(X)
(8)



PCC



0.5

0.5

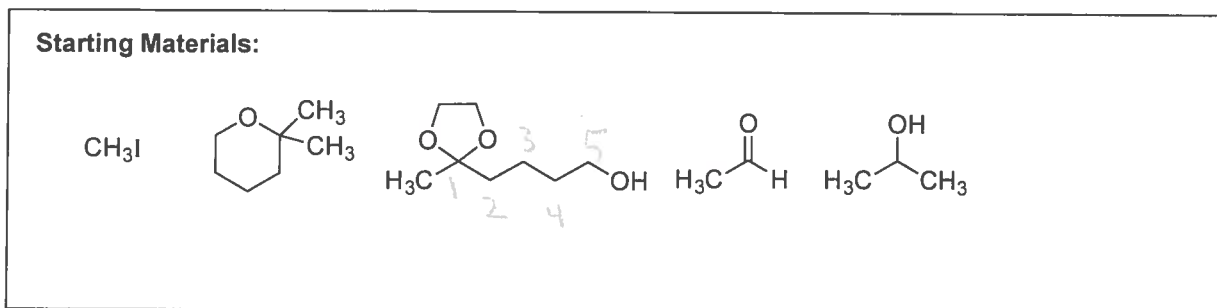
5. Propose a synthesis of the target below (8 points).

Initials: B

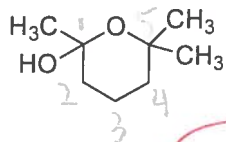
All carbons in the product must come from the starting materials provided, you can use any reagent you wish.

YOU CAN IGNORE STEREOCHEMISTRY.

mifib #5A
PS1 #5a
PS2 #2b
2C
~~PS2 #3a~~
PS2 #3c



Target.



proper Grignard use = 1

redox
correct
1 point
each
(2x)

