

EXAM 3, Sp '08

Question 1: Nomenclature (12 points)
 Give the appropriate IUPAC name for each of the following compounds. Be sure to indicate one stereochemical feature whenever

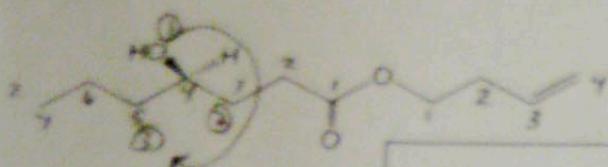


(E)-3-methyl-N-phenyl-2-hexenamide

1

FR

(2E)



3-butenyl (R)-4-hydroxyheptanoate

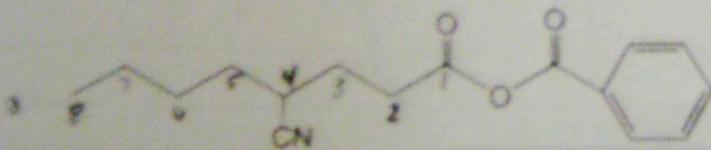
1

1

1

1

1



benzolic 4-cyanoctanoic anhydride

1

1

1

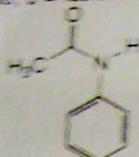
2

Final: 12 points

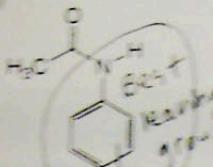
Rank the following compounds in order of increasing reactivity with H₂O. (1 = slowest)

B. Facts: 12 points

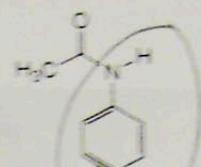
1. Rank the following compounds in order of increasing reactivity with H_2O . (1 = slowest rate, 3 = fastest rate) (3 pts)



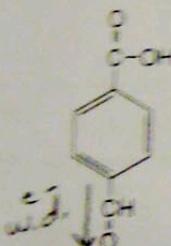
2



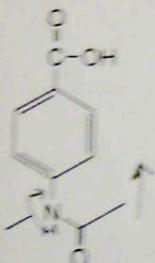
3



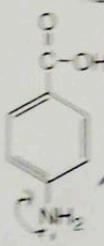
2. Rank the following compounds in order of increasing pK_a . (1=lowest pK_a , 3=highest pK_a) (3 pts)



1



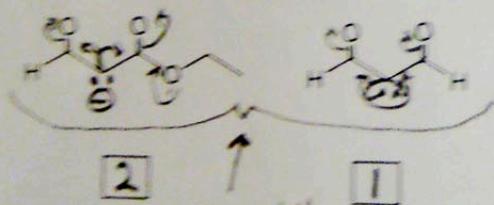
2



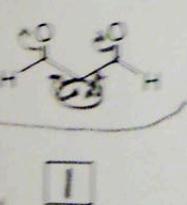
3

least acidic

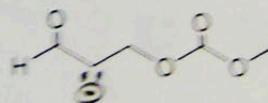
3. Rank the following compounds in order of increasing pK_a . (1=lowest pK_a , 3=highest pK_a) (3 pts)



2

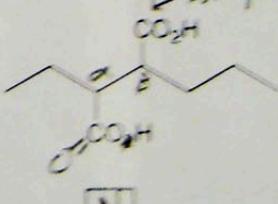
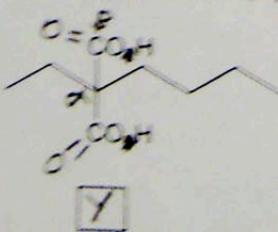
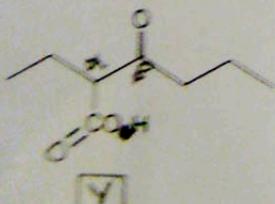


1



3

4. Consider the decarbonylation at 180 °C of the following carboxylic acids. If decarbonylation is possible place Y (for yes) in the box. If not, place N (for no) in the box. (3 pts)

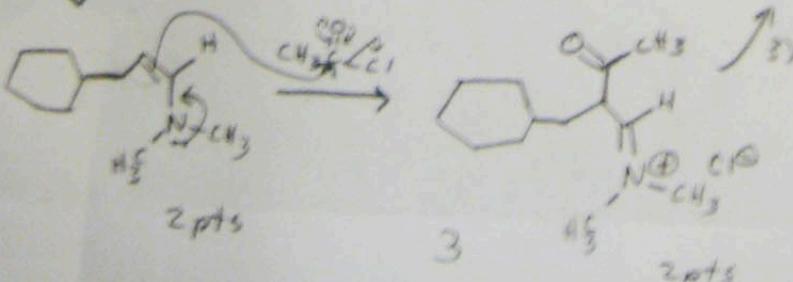
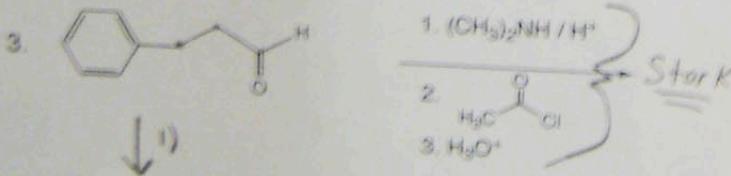
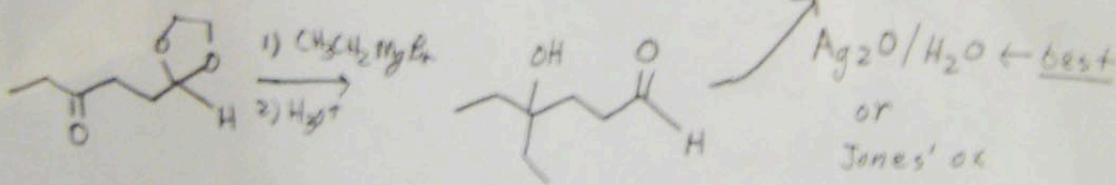
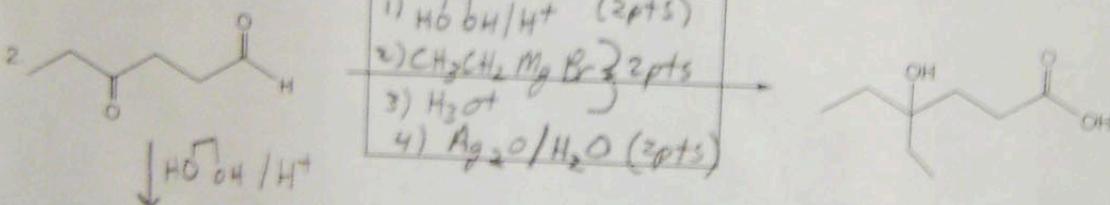
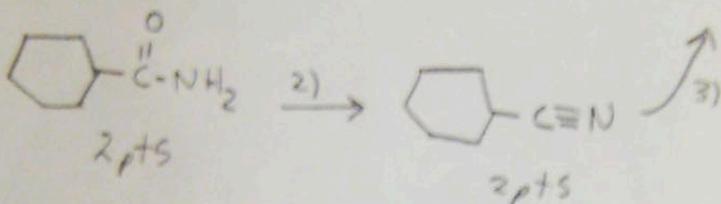
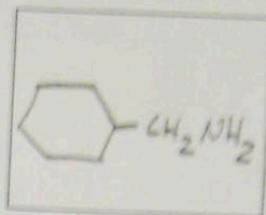
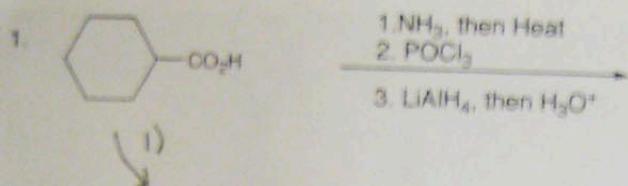


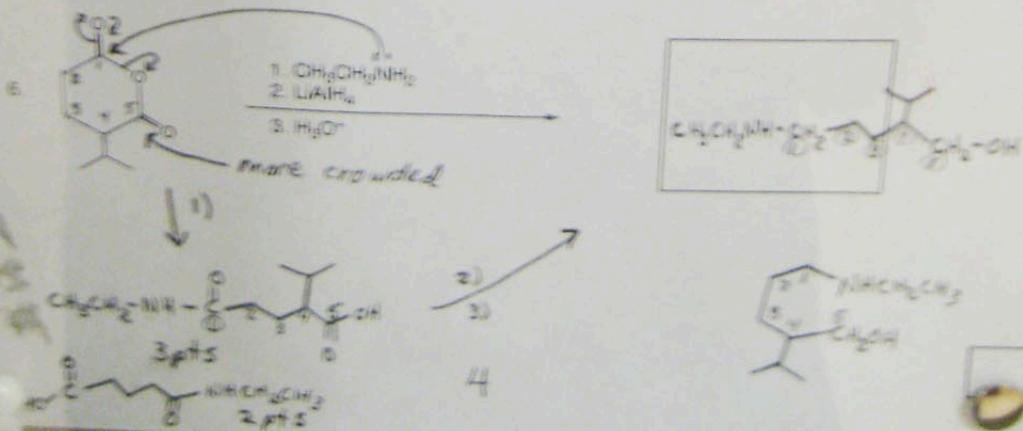
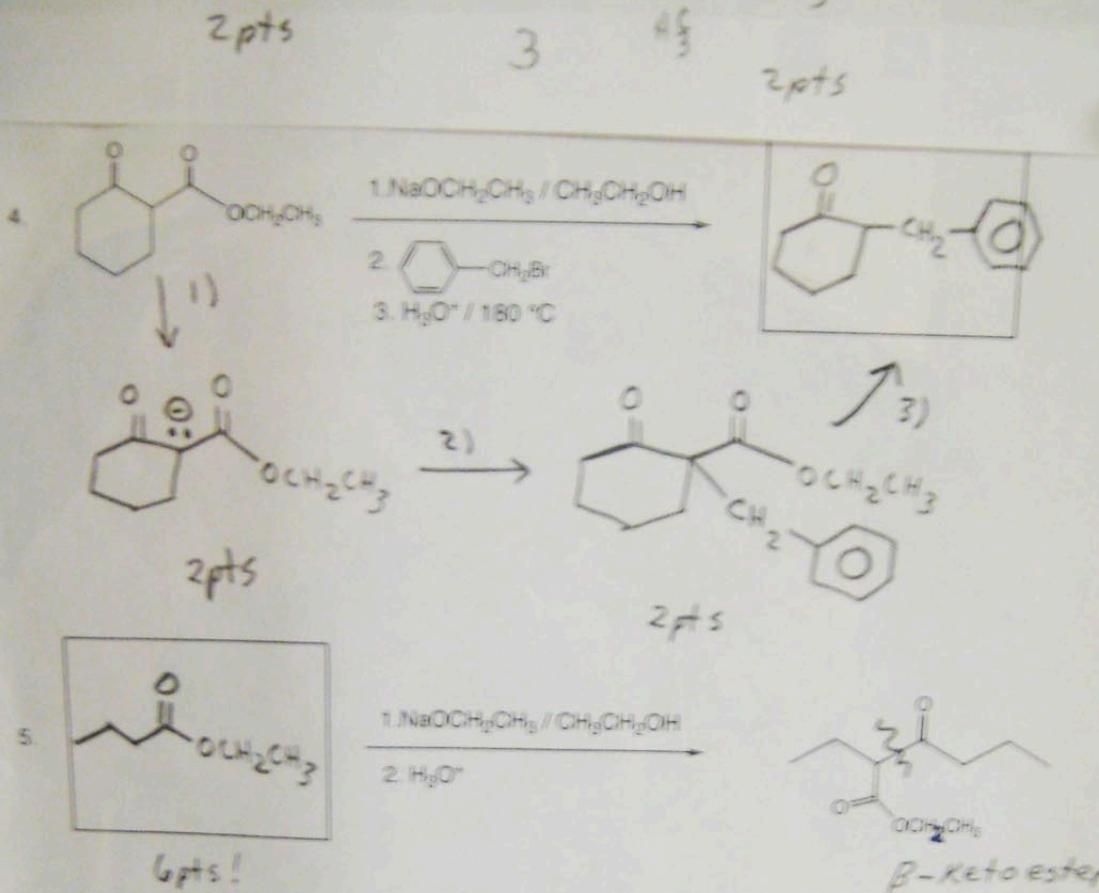
far away

β -keto acids necessary

C. Reactions Total = 36 points, 6 points each

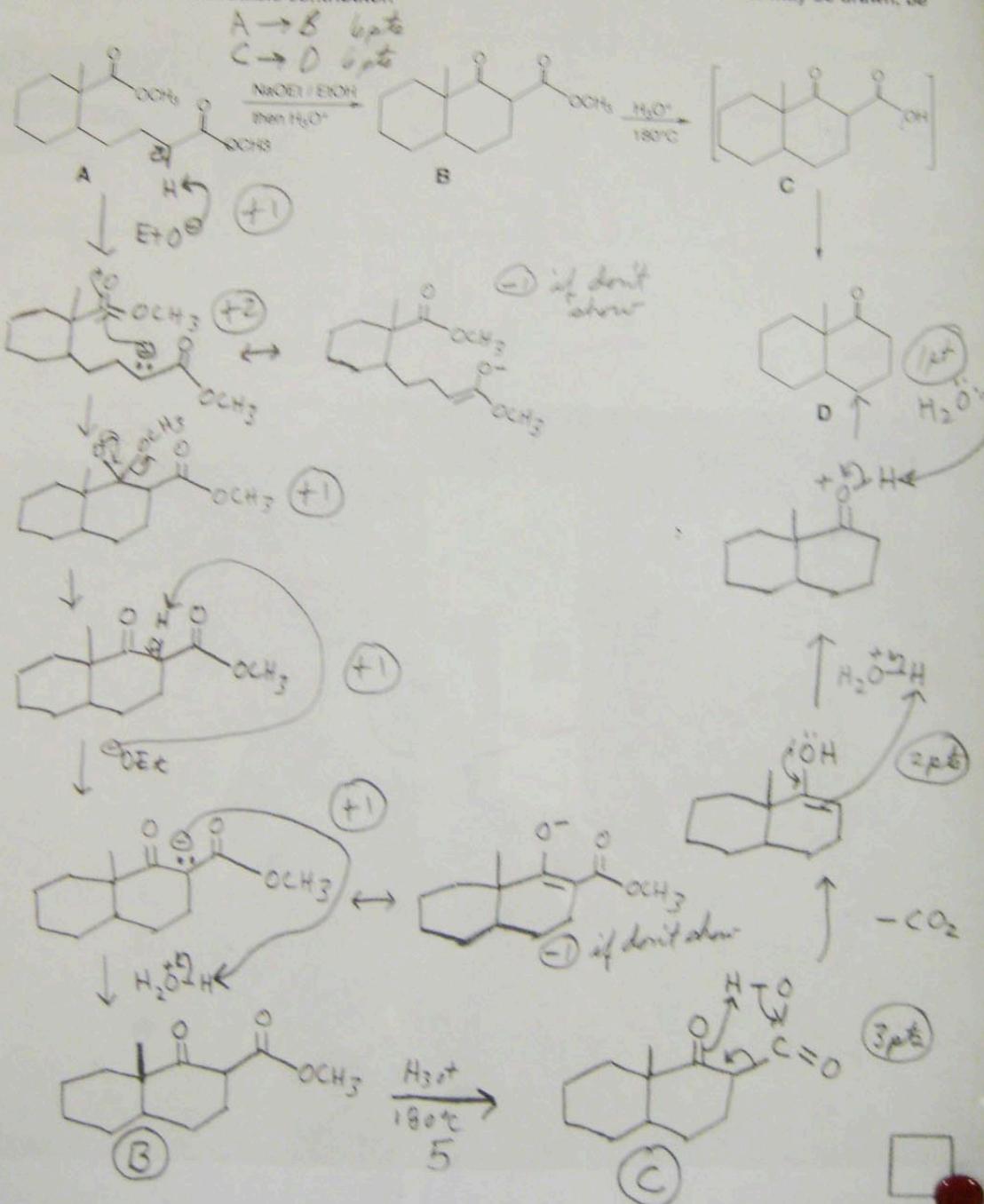
Please provide the starting material, reagents or major product in the answer box. Be sure your drawing indicates **stereochemistry** if applicable. Partial credit is awarded only when intermediate products in a multi-step reaction are shown below the reaction.





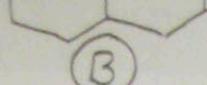
D. Mechanism: (12 points)

Provide a clear mechanism to explain the formation of the series of products below. Do not show the transformation of B to C. Remember to show only one step at a time (NO SHORTCUTS!). Show all intermediates and all formal charges. When more than one resonance contributor may be drawn, be sure to draw the most stable contributor.

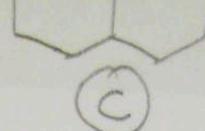


E. Synthesis: 13 Points

Synthesize the molecule below using any of the following reagents: benzene, bromobenzene; any alkenes, alcohols, or alkyl halides of three carbons or less, ethylene oxide, any inorganic reagents, oxidizing or

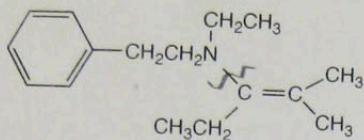


180°C
5

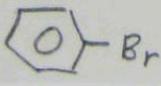
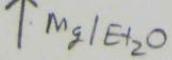
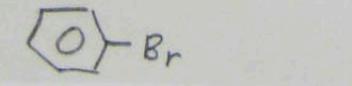
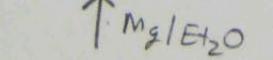
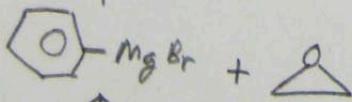
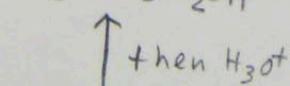
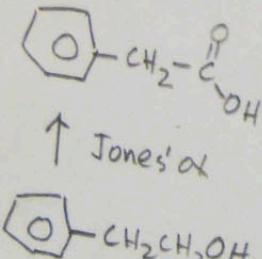
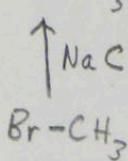
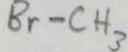
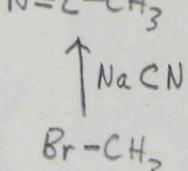
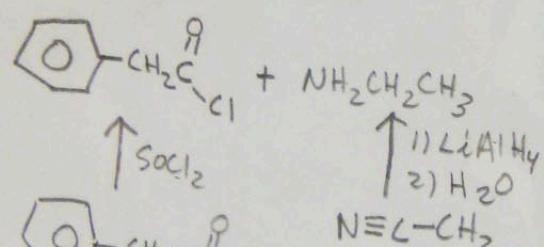
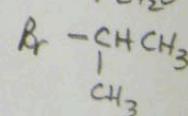
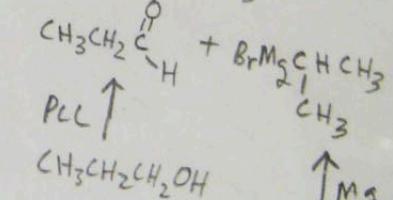
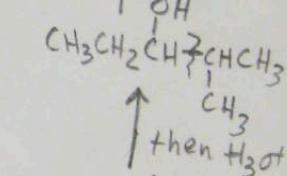
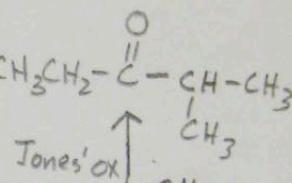
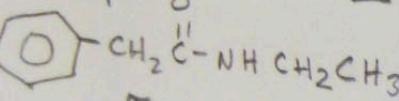


E. Synthesis: 13 Points

Synthesize the molecule below using any of the following reagents: benzene, bromobenzene; any alkenes, alcohols, or alkyl halides of **three carbons** or less; ethylene oxide; any inorganic reagents, oxidizing or reducing agents, and any peroxyacids.



1) LiAlH_4
2) H_3O^+ or H_2O

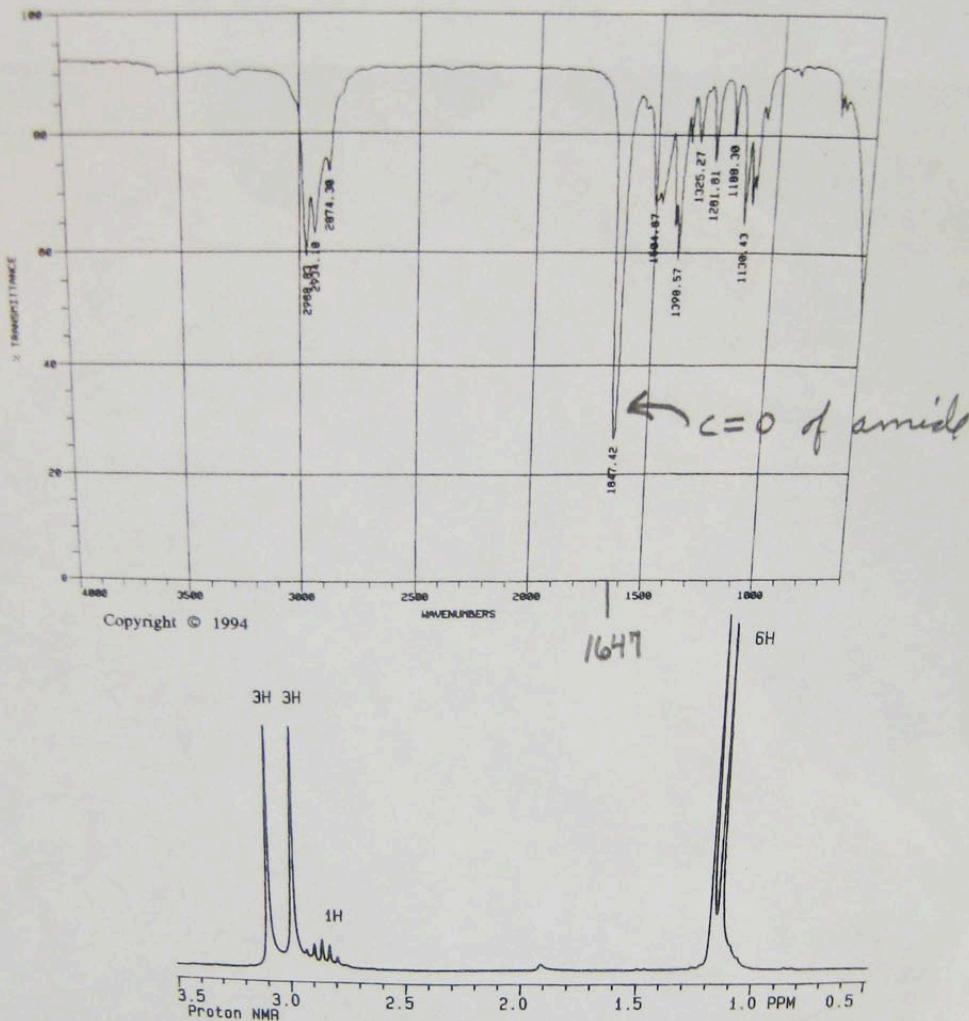


F. Spectroscopy: 12 Points

A compound with the formula $C_6H_{13}NO$ exhibits the IR, ^1H NMR and proton decoupled ^{13}C NMR spectra shown below. Please identify this compound and draw the structure in the box provided below.

F. Spectroscopy: 12 Points

A compound with the formula $C_6H_{13}NO$ exhibits the IR, 1H NMR and proton decoupled ^{13}C NMR spectra shown below. Please identify this compound and draw the structure in the box provided below.



partial credit:

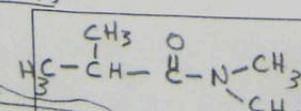
amide: 3 pts

isolated CH_3 : 1 pt each (2 max)

isopropyl: 3 pts

1H adj to max int.

CH_3 adj. to 1 H: 1 pt each (2 max)



ONLY if NO isopropyl

