Simulated Exam I (F09O1B)

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1. Name each of the compounds in the box provided using IUPAC naming including stereochemistry if applicable (16).



2. Rank the compounds in order of increasing melting point, with 1 = lowest mp, 3 = highest mp (6).



3. Rank the compounds in order of increasing solubility in water, with 1 = lowest solubility, 3 = highest solubility (6).



4. In the acid/base reaction below, provide the products that would result **AND** place an arrow to show the direction of equilibrium (8).



5. Provide the hybridization in the boxes provided (12).



Bond angle of C_AC_BO_c?

Are hydrogens "a" and "b" in the same plane (YES/NO)

6. From the monochlorination of 3,4-dimethylhexane below, draw all structural isomers (8).



7. Draw important resonance contributors for the structure below. Circle the major contributor (10).

Đ (CH₃)₂N

8. Draw the Newman projection of the most stable conformation of 3,4-dimethylheptane along the C3-C4 bond (6).



9a. Draw the most stable chair conformation for each compound in the box below (10).



b. Which is more stable (A or B)? _____

10. In the monobromination of butane below, show the propagation steps of the mechanism that explains the formation of the products (10).



a. With the bond dissociation energies below, calculate the overall ΔH° showing all work (4)

b. At equilibrium, are products or reactants favored? (2)

Bond-Dissociatio Energy		। त्र त त ।		Bond-Dissociation Energy	
kcal/mol		Bond	,	kcal/mol	
kcal/mol 104 106 38 58 46 36 136 103 88 71 119 51 104 109 84 70 56 91 98 107 81 68 53 91 98 107 81 68 53 91		Bonds to secondary c $(CH_3)_2CH - H$ $(CH_3)_2CH - F$ $(CH_3)_2CH - CI$ $(CH_3)_2CH - CI$ $(CH_3)_2CH - OH$ Boals to intitiaty carbon $(CH_3)_2CH - H$ $(CH_3)_2C - H$ $(CH_3)_2C - H$ $(CH_3)_2C - CI$ $(CH_3)_2C - Br$ $(CH_3)_2C - Br$ $(CH_3)_2C - OH$ Other C - H bonds PhCH ₂ - H CH ₂ = CHCH ₂ - H CH ₂ = CHCH ₂ - H CH ₂ = CHCH ₂ - H Ph - H C - C bonds CH ₃ - CH ₃ CH ₃ CH ₂ - CH ₃ (CH ₃) ₃ C - CH ₃ (CH ₃) ₃ C - CH ₃	arbons ons (benzylic) (allylic) (vinyl) (aromatic)	kcal/mol 95 106 80 68 53 91 91 106 79 65 50 91 85 87 108 110 88 85 82 84 81	2
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