| Name | | | | | | | | | | | |
|------|--|--|--|--|--|--|--|--|--|--|--|
| | | | | | | | | | | | |

General Instructions: Write your name in the space provided above and on the provided Scan-tron form. *Do not put your name anywhere else in this exam book.*

Make sure that you read each question carefully and provide complete answers. Time limit is 55 min.

Grading: Grading will be on the basis of a highest possible score of 100 points.

- I. Multiple Choice 2 points each, 40 points total
- II. R /S Designations 1.5 points each, 7.5 points total
- III. Structures 1.5 points each, 7.5 points total
- III. Reaction Products 5 points each, 25 points total
- IV. Spectroscopy Problem 20 points

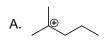
| | KJ/MOL | KCAL/MOL | | KJ/MOL | KCAL/MOL | | KJ/MOL | KCAL/MO |
|--|--------|----------|--------------------------------------|--------|----------|---------------------------------------|--------|---------|
| Bonds to H | | | H ₂ C=CH−CH ₃ | 385 | 92 | (CH ₃) ₂ CH—F | 444 | 106 |
| Н—Н | 435 | 104 | HC≡C—CH ₃ | 489 | 117 | (CH ₃)₂CH—CI | 335 | 80 |
| H—CH ₃ | 435 | 104 | | | | (CH ₃) ₂ CH—Br | 285 | 68 |
| H—CH ₂ CH ₃ | 410 | 98 | Bonds to methyl | | | (CH ₃) ₂ CH—I | 222 | 53 |
| H—CH(CH ₃) ₂ | 397 | 95 | CH ₃ —H | 435 | 104 | (CH₃)₂CH—OH | 381 | 91 |
| H—C(CH ₃) ₃ | 381 | 91 | CH ₃ —F | 456 | 109 | | | |
| H————————————————————————————————————— | 473 | 113 | CH₃—CI | 351 | 84 | ÇH₃ | | |
| | | | CH ₃ —Br | 293 | 70 | H ₃ C-C-X | | |
| | | | CH ₃ —I | 234 | 56 | CH ₃ | | |
| | 356 | 85 | СН3—ОН | 381 | 91 | (CH ₃) ₃ C—H | 381 | 91 |
| | | | //// | | | (CH ₃) ₃ C—F | 444 | 106 |
| н | 464 | 111 | H | | | (CH ₃) ₃ C—Cl | 331 | 79 |
| | 364 | 87 | H ₃ C — C — X | | | (CH ₃) ₃ C—Br | 272 | 65 |
| н 30. | | 07 | Н | | | (CH ₃) ₃ C—I | 209 | 50 |
| H—F | 569 | 136 | CH₃CH₂—H | 410 | 98 | (CH₃)₃C—OH | 381 | 91 |
| H—Cl | 431 | 103 | CH₃CH₂—F | 448 | 107 | | | |
| H—Br | 368 | 88 | CH ₃ CH ₂ —Cl | 339 | 81 | X—X bonds | | |
| H—I | 297 | 71 | CH ₃ CH ₂ —Br | 285 | 68 | F—F | 159 | 38 |
| Н—ОН | 498 | 119 | CH₃CH₂—I | 222 | 53 | CI—CI | 242 | 58 |
| H—OCH ₂ CH ₃ | 435 | 104 | CH₃CH₂—OH | 381 | 91 | Br—Br | 192 | 46 |
| | | | 19404 27 | | | — | 151 | 36 |
| C—C bonds | | | CH₃ I | | | но—он | 213 | 51 |
| CH ₃ —CH ₃ | 368 | 88 | H ₃ C— <mark>C</mark> —X | | | | | |
| CH ₃ CH ₂ —CH ₃ | 356 | 85 | Ĥ | | | | | |
| (CH ₃) ₂ CH—CH ₃ | 351 | 84 | (CH ₃) ₂ CH—H | 397 | 95 | | | |

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I. Multiple Choice

I. Each of the following multiple choice questions has one and only one correct response. Mark the letter of the correct response to each question on the Scan-Tron form provided. A #2 pencil should be used. (Make sure to write your name on the Scan-Tron form.)

Use the following choices for questions 1-3.



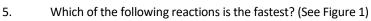
B. **

C. ••

D. \oplus

E. *

- 1. Which is the least stable carbocation?
- 2. Which is the most stable?
- 3. Which carbocation is expected to undergo a 1,2-hydride shift?
- 4. Which of these correctly describes the overall reaction $a \rightarrow e$? (See Figure 1)
 - A. non-concerted, exothermic
 - B. non-concerted, endothermic
 - C. concerted, exothermic
 - D. concerted, endothermic

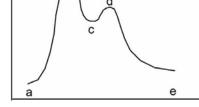




B. c **→**e

C. c →a

D. e \rightarrow c



6. What points on the graph represent transition states? (See Figure 1)

- A. c only
- B. b and d
- C. b, c, and d
- D. c and e

Figure 1 - PE Diagram for Questions 4-6

7. Use the table of BDEs on the front page to predict the ΔH of the following hypothetical reaction.

- A. +8 kcal/mol
- B. -8 kcal/mol
- C. +5 kcal/mol
- D. -5 kcal/mol
- 8. Morphine is a natural product isolated from opium poppies. It can be reacted with H_2SO_4 to form morphine sulfate which is used as an oral pain reliever. The $[\alpha]_D$ of pure morphine sulfate is -110°. If 10 morphine sulfate caplets are dissolved in 5.0 mL H_2O and the solution is measured on an automatic 1-dm polarimeter to give an observed rotation of α_{obs} = -22.0° what is the average mass (in mg) of the caplets?
 - A. 20
- B. 25
- C. 80
- D. 100
- E. 1500
- 9. The optical purity of a research compound was measured by polarimetry to be 50%. Which of these statements correctly describes the compound?
 - A. It is a mixture of enantiomers in a 4:1 ratio
 - B. It is a mixture of enantiomers in a 3:1 ratio
 - C. It is a mixture of diastereomers in a 4:1 ratio
 - D. It is a mixture of diastereomers in a 3:1 ratio.
 - E. It is a racemic mixture
- 10. Which of the following cycloalkanes is chiral?



D



C.



D.



F



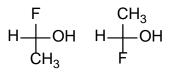
11.



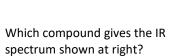
13.



15.

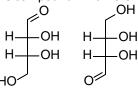


17.

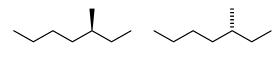


- A. 1-Octene
- B. 2-Octene
- C. 1-Octyne
- D. 2-Octyne
- E. 1-Octanol

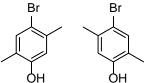
12.

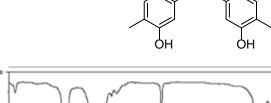


14.



16.





18.

What are the expected peak multiplicities in the ¹H NMR spectrum of

C. t, q, m, t, d D. t, q, d, m, d A. q, t, t, m, s B. s, t, t, d, m

E. q, t, t, m, q

19.

What are the expected peak multiplicities in the ¹H NMR spectrum of A. q, t, t B. s, s, s C. t, m, t C. s, s, s E. d, d, q

20. Which of the following substrates gives the *fastest* S_N2 reactions?

II. (a) Label all chirality centers in morphine as R or S as appropriate. Show priorities of all groups around each chirality center.

(b) How many stereoisomers are theoretically possible form morphine?

III. Give the structure of each of the following:

TMS (S)-3-chloro-1-hexene DMSO *meso-*3,4-dimethylhexane TsCl

IV. Predict the product of each of the following nucleophilic substitution reactions. If no reaction is expected to occur then write N.R. and explain your reasoning.

V. A compound with molecular formula $C_4H_8O_2$ displays the following IR and 1H NMR spectra. Propose a structure for this compound. Explain your reasoning by labeling all key peaks in the spectra.

