題號: 61

國立臺灣大學 106 學年度碩士班招生考試試題

科目:有機無機

行日·牙极無有 節次: 6 題號: 61 共 4 頁之第 / 頁

第一部分 有機化學 (共50分)

單選題: 第1-17題 (每題2分,共34分,請依題號順序於選擇題作答區內作答)

1. What is the index of hydrogen deficiency (or degree of unsaturation) for morphinone?

- (A) 7 (B) 8 (C) 9 (D) 10
- 2. I and II are:

I.
$$CH_3$$
 H C_2H_5 H CH_3

- (A) constitutional isomers (B) enantiomers (C) identical (D) diastereomers
- 3. How many of the following reactions is/are likely to give ketone as the major product:

I.
$$R = \frac{HgSO_4}{H_2O}$$
 II. $\frac{1. OsO_4}{2. HIO_4}$ III. $\frac{1. H^+/H_2O}{2. PCC}$

- (A) 0 (B) 1 (C) 2 (D) 3
- 4. How many of the following reactions is/are named correctly:

II.
$$CH_2I_2$$
 Siemens-Smith

Siemens-Smith

Siemens-Smith

Diels-Elder

 CO_2Me
 CO

- (A) 0 (B) 1 (C) 2 (D) 3
- 5. Based on the structure of the compound below, how many of the following states is/are correct?

- I. H_a and H_b are homotopic protons.
- II. H_f and H_e are diastereotopic protons.
- III. Ha and Hd are enantiotopic protons.
- IV. H_f and H_d are enantiotopic protons.
- (A) 1 (B) 2 (C) 3 (D) 4
- 6. About S_N2 reaction, how many of the following statements is/are correct?
 - I. Neopentyl bromide is better substrate than methyl bromide.
 - II. Anionic nucleophiles are always better than the neutral ones.
 - III. Leaving group is important and tosylates are not the good ones.
 - IV. DMSO (dimethyl sulfoxide) is a protic solvent preferred.
 - (A) 0 (B) 1 (C) 2 (D) 3

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41 日 · 月 48 無 4 節 次 : 6 題號: 61

7. Rank the chemical shifts of the protons labeled as a, b, c and d in the following compound:

- 8. Normally, the hybridization of the carbon of a methyl radical is:
 - (A) sp (B) sp^2 (C) sp^3 (D) $sp^{1.5}$
- 9. About the ¹H NMR spectrum of Q, how many of the following statements is/are correct?

$$\frac{Br_2}{\text{heat ot light}} \quad C_8H_9Br \text{ (major product)} \quad + \quad HC$$

- I. There would be only 4 aromatic protons at low field.
- II. The signal for the protons on the benzylic carbon would be a triplet.
- III. The signal for the methyl protons would be a doublet.
- IV. The signal for the benzylic protons would integrate for 2 hydrogens.
- (A) 1 (B) 2 (C) 3 (D) 4
- 10. How many of the following reactions is/are likely to proceed effectively?

I.
$$\underbrace{ \begin{array}{c} 1. \text{ BD}_3 \\ \hline 2. \text{ H}_2\text{O}_2 \text{ / OH} \end{array} }_{\text{OH}} \underbrace{ \begin{array}{c} 1. \text{ BH}_3 \\ \hline 2. \text{ CH}_3\text{COOD} \end{array} }_{\text{D}} \underbrace{ \begin{array}{c} \text{H} & \text{III.} \\ \hline \\ \text{Pd/C} \end{array} }_{\text{D}} \underbrace{ \begin{array}{c} \text{D}_2 \\ \hline \\ \text{Pd/C} \end{array} }_{\text{D}}$$

- (A) 0 (B) 1 (C) 2 (D) 3
- 11. How many of the following nomenclatures is/are correct?

- (A) 1 (B) 2 (C) 3 (D) 4
- 12. How many of the following compounds is/are aromatic?

- (A) 1 (B) 2 (C) 3 (D) 4
- 13. How many of the following reactions is/are likely to proceed effectively?

- (A) 1 (B) 2 (C) 3 (D) 4
- 14. How many of the following statements is/are correct?
 - I. [10]Annulene is aromatic.
 - II. TBAF can be used to cleave silyl ether.

題號: 61

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科目:有機無機

節次: 6

題號: 61 共 4 頁之第 3 頁

III. β -carotene absorbs orange light.

IV. Wolff-Kishner reaction proceeds under acidic condition.

(A) 1 (B) 2 (C) 3 (D) 4

15. How many of the following pairs of structures and name abbreviations is/are correct?

(A) 1 (B) 2 (C) 3 (D) 4

16. How many of the following reactions is/are likely to proceed effectively?

17. How many of the following statements is/are correct?

I. DCC is frequently used in the amide formation reactions and its full name is dicyclohexyl chromate.

II. HOOCCH₂CH₂CH₂COOH is called succinic acid.

III. McLafferty rearrangement is more likely to be observed in the mass spectra of alkenes.

IV. Generally, carbon-carbon triple bond should have IR absorption around 2150 cm⁻¹.

(A)1 (B)2 (C)3 (D)4

問答題: 第18-25題 (每題2分,共16分,請依題號順序於<u>非選擇題作答區</u>內作答)

Give the major product in line-angle structure of each of the following reactions:

節次: 6

科目:有機無機

國立臺灣大學 106 學年度碩士班招生考試試題

頁之第 4

Inorganic chemistry

- 1. Describe the point group of each following molecule: (10 pts)
- b. H₃O⁺ c. O₂F₂ d. cyclohexane (chair conformation) e. SO₄²a. Ethylene
- f. Diborane

H B H B H B

- h. Formaldehyde (H₂CO) i. Face-centered cubic j. CsCl
- 2. Explain the possible reasons for following statements:
- a. Ethylene is a planar molecule, but hydrazine is not. (3 pts)
- b. ICl_2^- is linear, but NH_2^- is bent. (3 pts)
- c. To consider two compounds, mercury (II) cyanate (Hg(OCN)₂) and mercury (II) fulminate (Hg(CNO)₂), one is highly explosive, and the other is not. (4 pts)
- 3. To consider a known molecule of NF:
- a. Construct a molecular orbital energy-level diagram for NF, showing clearly how the atomic orbitals interact to form molecular orbitals. (6 pts)
- b. What is the most likely bond order for NF. (2 pts)
- c. What are the point groups of all molecular orbitals. (3 pts)
- 4. Show that spheres occupy only 74.0 % of the total volume in face-centered cubic structure in which all the atoms are identical. (5 pts)
- 5. To consider the following half cell emf diagram for nitrogen in acidic solution:

$$HNO_2 \xrightarrow{0.996} NO \xrightarrow{1.59} N_2O \xrightarrow{1.77} N_2 \xrightarrow{-1.87} NH_3OH^+ \xrightarrow{1.41} N_2H_5^+ \xrightarrow{1.275} NH_4^+$$

- a. Construct a Frost diagram for nitrogen in acid. (4 pts)
- b. The half cell emf for $N_2O \rightarrow NH_3OH^+$. (3 pts)
- 6. [Co(H₂O)₆]³⁺ is a strong oxidizing agent, but [Co(NH₃)₆]³⁺ is stable in aqueous solution. Explain this observation by comparing the difference in LFSE (ligand field stabilization energy) for each oxidized/reduced complexes pair, in which the Co(III) complexes are low-spin and the Co(II) complexes are high-spin. (7 pts) [Orbital splitting (Δ_0) for $[Co(H_2O)_6]^{3+}$] is 16750 cm⁻¹, and that for $[Co(H_2O)_6]^{2+}$ is 8400 cm⁻¹; the value of Δ_0 for $[Co(NH_3)_6]^{3+}$ is 24000 cm⁻¹ and that for $[Co(NH_3)_6]^{2+}$ is 10200 cm⁻¹.