Chemistry 334

Examination \#1

September 29, 2014
Professor Charonnat

Name: $\qquad$

Be certain that your examination has six (6) pages including this one.
Put your name on each page of this examination booklet.
By putting your name on this examination booklet you agree to abide by California State University, Northridge policies of academic honesty and integrity.

Molecular models are allowed for this examination. All electronic devices, including calculators, are unnecessary and are not allowed.

Name: $\qquad$

1. (25 points)

Draw the structure of the expected major organic product for each of the following five (5) questions. Specify stereochemistry clearly, if relevant.
A.

B.

C.
product 1B

D.


E.


Name: $\qquad$
2. ( 25 points)

Circle the number that corresponds to the correct answer for each of the following five (5) questions.
A. The highest-occupied molecular orbital (HOMO) of the following radical is


1. $\pi_{2}$
2. $\pi_{3}$
3. $\pi_{5}$
B. For a given compound that possesses a UV chromophore, its molar absorptivity is
4. identical at all wavelengths
5. variable at different wavelengths, and always has a solitary maximum
6. variable at different wavelengths, and can have multiple maxima
C. The following compound is

7. nonaromatic
8. aromatic
9. antiaromatic
D. The reaction of an achiral, unsymmetrical sulfide with an equimolar amount of hydrogen peroxide affords
10. a sulfone
11. an optically active sulfoxide
12. a racemic sulfoxide
E. The Birch reduction of anisole derivatives occurs via a mechanism that includes
13. radical intermediates
14. carbocation intermediates
15. carbene intermediates

Name: $\qquad$
3. (10 points)

Use IUPAC nomenclature to write the systematic name of the following alkene.

4. (20 points)

Draw the specific reagents necessary to synthesize the following sulfide from benzene and any other organic compounds that contain five or fewer carbons. Distinguish individual steps clearly and draw each stable synthetic intermediate compound (i.e., the major organic product of each reaction). Do not draw mechanisms. Finally, explain why the target is racemic. In particular, state which reactions in your synthesis yield racemic products and why those outcomes are expected.

(racemic)

Name: $\qquad$
5. (20 points)

The broadband proton-decoupled ${ }^{13} \mathrm{C}$ NMR spectrum of an unknown organic compound $\left(\mathrm{C}_{9} \mathrm{H}_{12} \mathrm{O}\right)$ is shown below. The labels next to each of the resonances signify the multiplicities observed in the corresponding off-resonance proton-decoupled ${ }^{13} \mathrm{C}$ NMR spectrum ( $\mathrm{s}=$ singlet, $\mathrm{d}=$ doublet, $\mathrm{t}=$ triplet, $\mathrm{q}=$ quartet). Use the spectroscopic data to determine the structure of this compound. Make clear assignments of all resonances to explain your reasoning. (A ${ }^{13} \mathrm{C}$ NMR correlation table is included separately.)


Name: $\qquad$
5. (continued)

${ }^{13}$ C NMR assignments: chemical shift (ppm) assignment explanation of multiplicity 157.8
136.6
129.8
114.2
55.8
28.2
14.5

## Congratulations!

| 1 | $/ 25$ |
| :--- | ---: |
| 2 | $/ 25$ |
| 3 | $/ 10$ |
| 4 | $/ 20$ |
| 5 | $/ 20$ |
| Total: | $/ 100$ |

