

NAME _____

ID # _____

ORGANIC CHEMISTRY I (CHEM 2301)

9:30 – 10:20 am, July 2, 2013

Exam 1

If you want to pick your graded exam up tomorrow in class (in public), please check the box on the right:

If you do not check the box, I will not bring your exam to class on Wednesday, and you will need to pick up your exam in private from Chemistry department staff in 115 Smith beginning Monday, July 8th. Exams that are not picked up within two weeks will be disposed of.

A periodic table is attached to the back of this exam as an aid. Otherwise, you are not permitted to use any other materials (including notes, books, or electronic devices of any kind).

Right now, write your name and student ID number at the top of this page. When the exam begins, please write your name at the top of the next page.

You may use pen or pencil. However, re-grades will be considered only for exams completed in pen.

Please write your answers in the boxes/spaces provided. If your answer is not in the appropriate space (say, for example, it's on the back of the page), draw us an arrow and/or note telling us where to look.

NAME _____

Scoring: 1. _____ / 9 5. _____ / 18
2. _____ / 6 6. _____ / 16
3. _____ / 18 7. _____ / 13
4. _____ / 20

Total Score: _____ / 100

1. (9 pts) **Draw Lewis dot structures** for each of the molecules below. Draw all valence electrons as dots. If there are formal charges, draw them on the appropriate atoms.

(a) hydrazine, H_2NNH_2



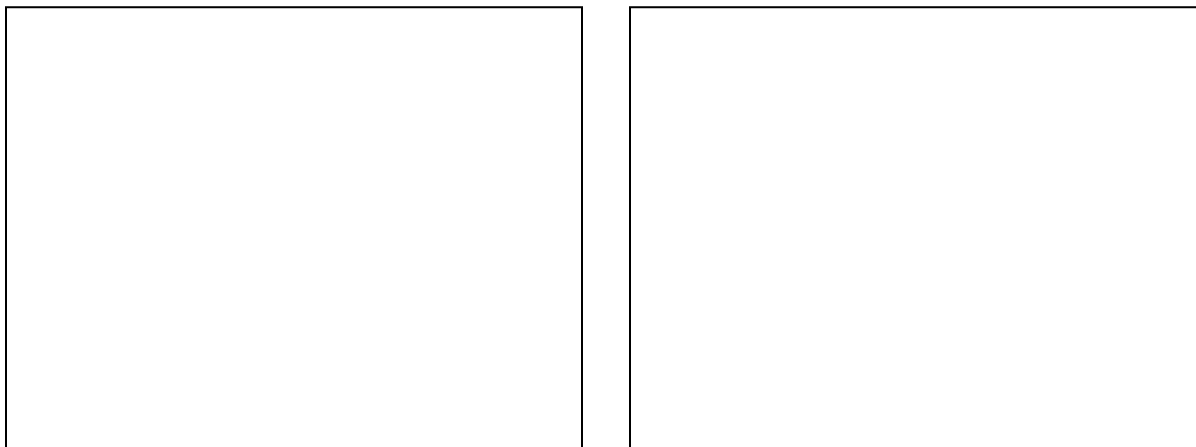
(b) sodium methoxide, CH_3ONa



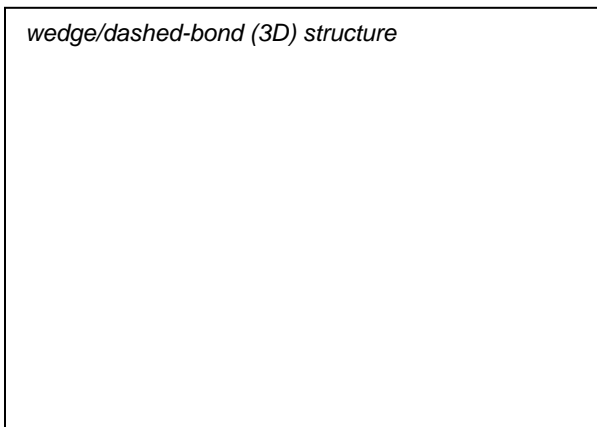
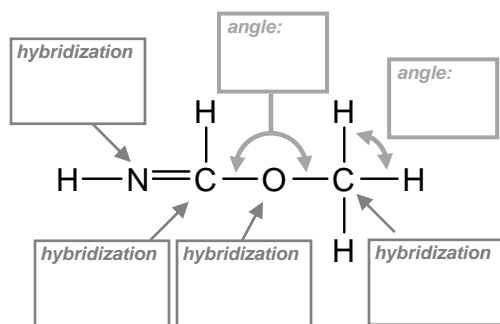
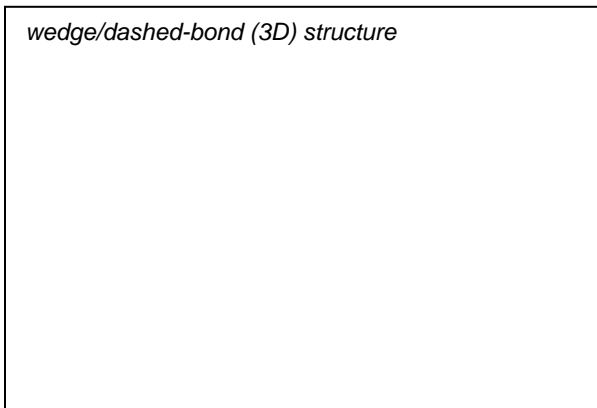
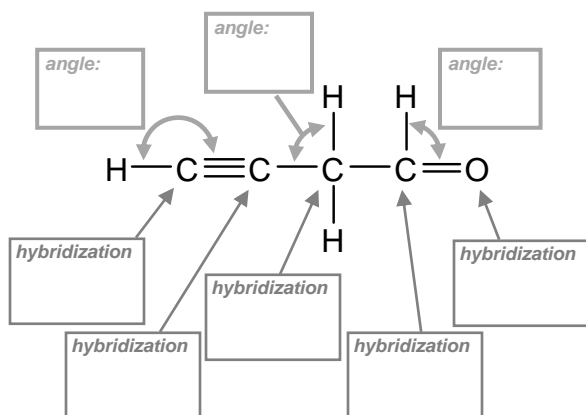
(c) propyne, HCCCH_3



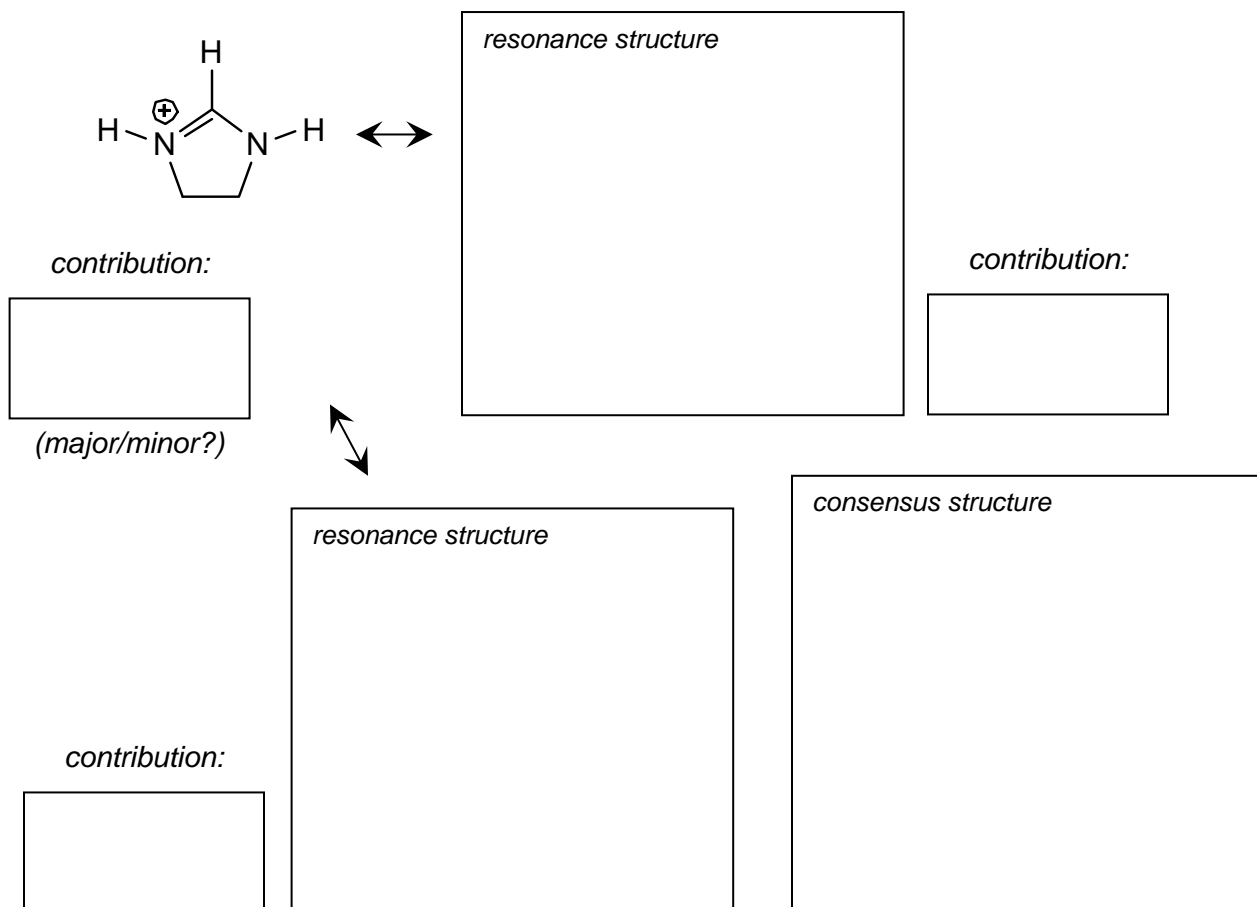
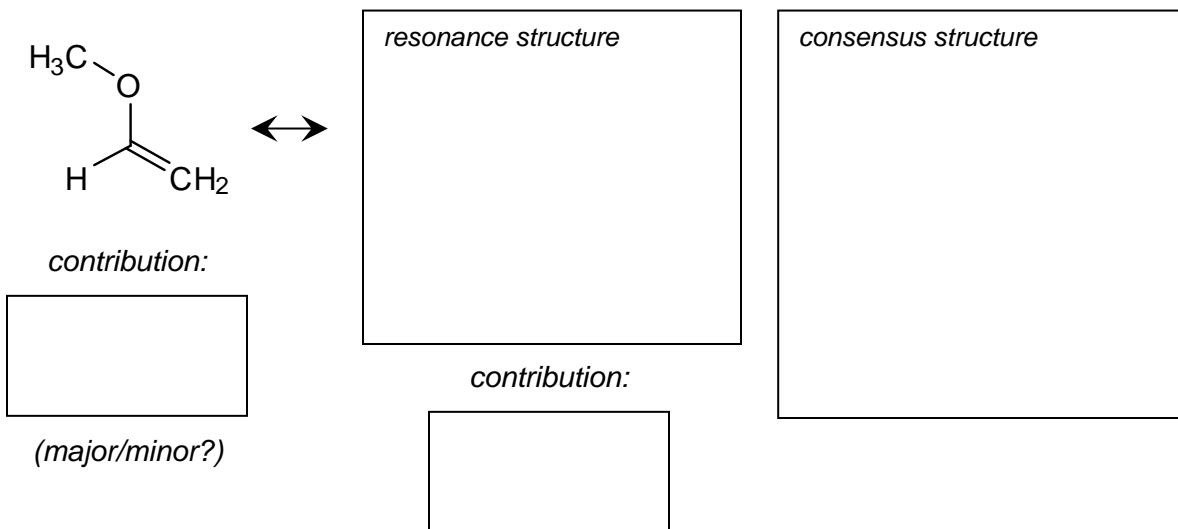
2. (6 pts) **Draw Lewis dash-bond structures** for two constitutional isomers that have molecular formula C_2H_4O , and that have **no formal charges** on any atom. Draw all atoms and lone pairs of electrons.



3. (18 pts) For each of the Lewis structures drawn below, in the boxes provided:
- Draw Lewis wedge/dashed-bond structures that illustrate the most stable three-dimensional structure of the molecule. Draw all atoms, but feel free to omit lone pairs.
 - In the boxes provided, write the hybridization state for any atom heavier than hydrogen.
 - In the boxes provided, give any bond angle indicated by curved arrows in the original Lewis structure.

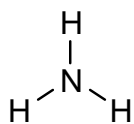


4. (20 pts) For each of the molecules on the left, draw as many Lewis dash-bond resonance structures as there are boxes. (Feel free to omit lone pairs and C-H's, or draw them—your choice.) Then, below each resonance structure, describe whether each would be a major or minor contributor. Finally, draw a consensus structure that illustrates partial charges and multiple bonds.

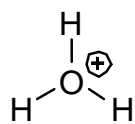


5. (18 pts) For each of the sets of molecules below:

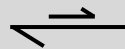
- Using “electron pushing” (with double-barbed arrows), show how the molecules on the left would react in an acid-base reaction to transfer a proton from one to the other.
- In the box on the right, draw the conjugate acid and base products of each reaction.
- In the middle, draw an equilibrium arrow that shows whether you feel the acid-base equilibrium would lie on the left or the right.



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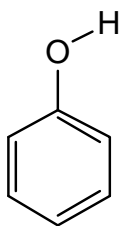


or

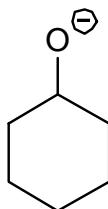


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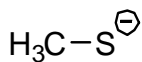
products



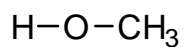
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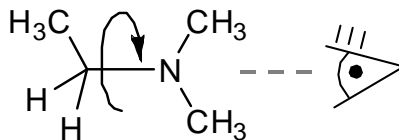


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products

6. (16 pts) For the molecule (dimethylamino)ethane (below):



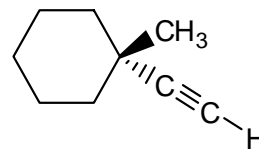
- (a) Draw Newman projections for the six conformers—three staggered and three eclipsed—accessed by rotation of the central C-N bond. Use the perspective I've indicated (with the nitrogen atom closest to the viewer).
- (b) Label which conformer you would expect to be most stable, and which you would expect to be least stable. If there are multiple conformations that are equally most or least stable, label them all.

Staggered conformers

Eclipsed conformers

7. (13 pts)

- (a) The substituted cyclohexane on the right has two chair conformers. Draw the two conformers in the boxes below. Draw all atoms except the six carbons of the cyclohexane ring—so, draw all hydrogens and both the $-\text{CH}_3$ and $-\text{C}\equiv\text{CH}$ substituents.

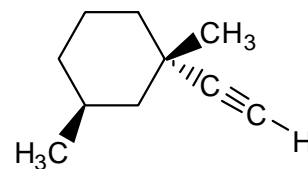


chair conformer

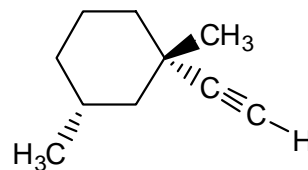
chair conformer

- (b) Circle the conformer above that you think will be more stable.

- (c) The two molecules on the right are like the molecule above, except they each have an additional methyl group. Which would be more stable? Circle one of the two molecules.



or



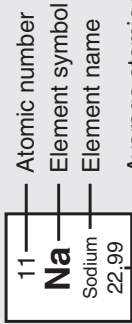
(circle one)

- (d) Interestingly, the *more* stable of the two molecules at right has the *least* stable chair conformer. For the molecule that you circled, draw the chair conformer that you think would be **least stable**. Feel free to omit the ring hydrogens—just draw the ring and substituents.

least stable chair conformer for the molecule you circled

| | | 1 | | 2 | | 3 | | 4 | | 5 | | 6 | | 7 | | 8 | | 9 | | 10 | | 11 | | 12 | | 13 | | 14 | | 15 | | 16 | | 17 | | 18 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
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| | | 1A | | 2A | | 3B | | 4B | | 5B | | 6B | | 7B | | 8B | | | | | | 1B | | 2B | | 3A | | 4A | | 5A | | 6A | | 7A | | 8A | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1 | 1 | H Hydrogen 1.01 | 2 | He Helium 4.00 | 3 | 4 | Li Lithium 6.94 | 5 | Be Beryllium 9.01 | 6 | 7 | B Boron 10.81 | 8 | C Carbon 12.01 | 9 | N Nitrogen 14.01 | 10 | O Oxygen 16.00 | 11 | F Fluorine 19.00 | 12 | Ne Neon 20.18 | 13 | Na Sodium 22.99 | 14 | Mg Magnesium 24.31 | 15 | Al Aluminum 26.98 | 16 | Si Silicon 28.09 | 17 | P Phosphorus 30.97 | 18 | S Sulfur 32.07 | 19 | Cl Chlorine 35.45 | 20 | Ar Argon 39.95 | 21 | K Potassium 39.10 | 22 | Ca Calcium 40.08 | 23 | Sc Scandium 44.96 | 24 | Ti Titanium 47.87 | 25 | V Vanadium 50.94 | 26 | Cr Chromium 52.00 | 27 | Mn Manganese 54.94 | 28 | Fe Iron 55.85 | 29 | Ni Nickel 58.69 | 30 | Cu Copper 63.55 | 31 | Zn Zinc 65.39 | 32 | Ga Gallium 69.72 | 33 | Ge Germanium 72.61 | 34 | As Arsenic 74.92 | 35 | Se Selenium 78.96 | 36 | Kr Krypton 83.80 | 37 | Rb Rubidium 85.47 | 38 | Sr Strontium 87.62 | 39 | Y Yttrium 88.91 | 40 | Zr Zirconium 91.22 | 41 | Nb Niobium 92.91 | 42 | Mo Molybdenum 95.94 | 43 | Tc Technetium (98) | 44 | Ru Ruthenium 101.07 | 45 | Rh Rhodium 102.91 | 46 | Pd Palladium 106.42 | 47 | Ag Silver 107.87 | 48 | Cd Cadmium 112.41 | 49 | In Indium 114.82 | 50 | Sn Tin 118.71 | 51 | Sb Antimony 121.76 | 52 | Te Tellurium 127.60 | 53 | I Iodine 126.90 | 54 | Xe Xenon 131.29 | 55 | Cs Cesium 132.91 | 56 | Ba Barium 137.33 | 57 | La Lanthanum 138.91 | 58 | Ce Cerium 140.12 | 59 | Pr Praseodymium 140.91 | 60 | Nd Neodymium 144.24 | 61 | Pm Promethium (145) | 62 | Sm Samarium 150.36 | 63 | Eu Europium 151.96 | 64 | Gd Gadolinium 157.25 | 65 | Tb Terbium 158.93 | 66 | Dy Dysprosium 162.50 | 67 | Ho Holmium 164.93 | 68 | Er Erbium 167.26 | 69 | Tm Thulium 168.93 | 70 | Yb Ytterbium 173.04 | 71 | Lu Lutetium 174.97 | 72 | Fr Francium (223) | 73 | Ra Radium (226) | 74 | Ac Actinium (227) | 75 | Rf Rutherfordium (261) | 76 | Hf Hafnium 178.49 | 77 | Ta Tantalum 180.95 | 78 | W Tungsten 183.84 | 79 | Re Rhenium 186.21 | 80 | Os Osmium 190.23 | 81 | Ir Iridium 192.22 | 82 | Pt Platinum 195.08 | 83 | Au Gold 196.97 | 84 | Hg Mercury 200.59 | 85 | Tl Thallium 204.38 | 86 | Pb Lead 207.2 | 87 | Bi Bismuth 208.98 | 88 | Po Polonium (209) | 89 | At Astatine (210) | 90 | Rn Radon (222) | 91 | Th Thorium 232.04 | 92 | Pa Protactinium 231.04 | 93 | U Uranium 238.03 | 94 | Np Neptunium (237) | 95 | Pu Plutonium (244) | 96 | Am Americium (243) | 97 | Cm Curium (247) | 98 | Bk Berkelium (247) | 99 | Cf Californium (251) | 100 | Fm Fermium (257) | 101 | Md Mendelevium (258) | 102 | No Nobelium (259) | 103 | Lr Lawrencium (262) |

Key



— Atomic number
— Element symbol
— Element name
— Average atomic mass*

* If this number is in parentheses, then it refers to the atomic mass of the most stable isotope.