

General Instructions:

- (i) Use scratch paper at back of exam to work out answers; final answers must be recorded at the proper place on the exam itself for credit. Models are allowed.
- (ii) Print your name on each page.
- (iii) Please keep your paper covered and your eyes on your own work. Misconduct will lead to failure in the course.

1. (10 points) Draw a structure that corresponds to each of the following names. Show all atoms in each structure, including hydrogen atoms.

(a) *E*-3-methyl-3-octene

(b) 1-bromo-3-fluorocyclohexene

Name _____

2. (16 points) $\text{H}_2\text{C}=\text{CFCF}_3$ is a relatively new refrigerant that has a smaller greenhouse effect (i.e., lower global warming effect) than currently popular refrigerants such as CHFCl_2 . Answer the following questions.

(a) Provide a drawing of $\text{H}_2\text{C}=\text{CFCF}_3$ that indicates the three-dimensional structure.

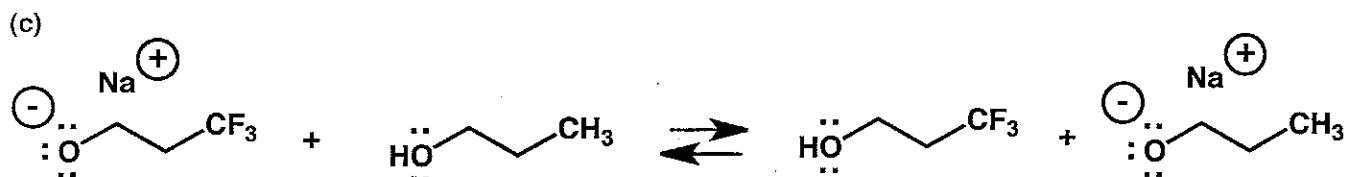
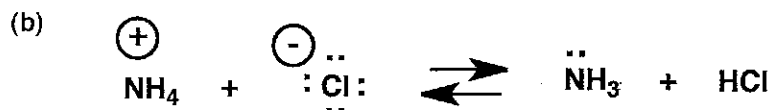
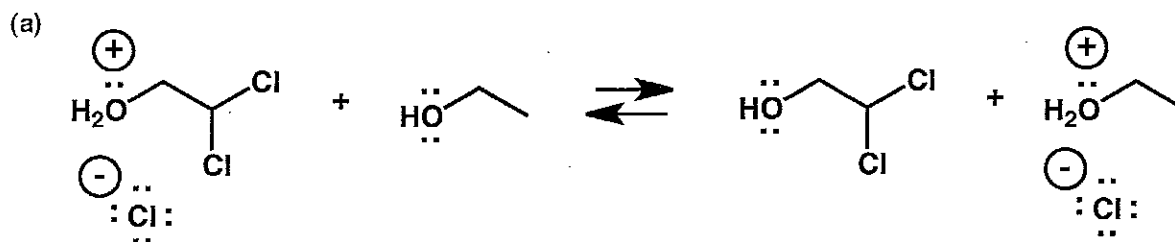
(b) Indicate the H-C-H bond angle (approximation) in the box.



(c) Provide an energy diagram that shows how the relevant atomic orbitals combine to form the molecular orbitals of the C-F bond involving the central carbon, and where the bonding electrons are expected to reside. Assume that F is sp^3 hybridized.

Name _____

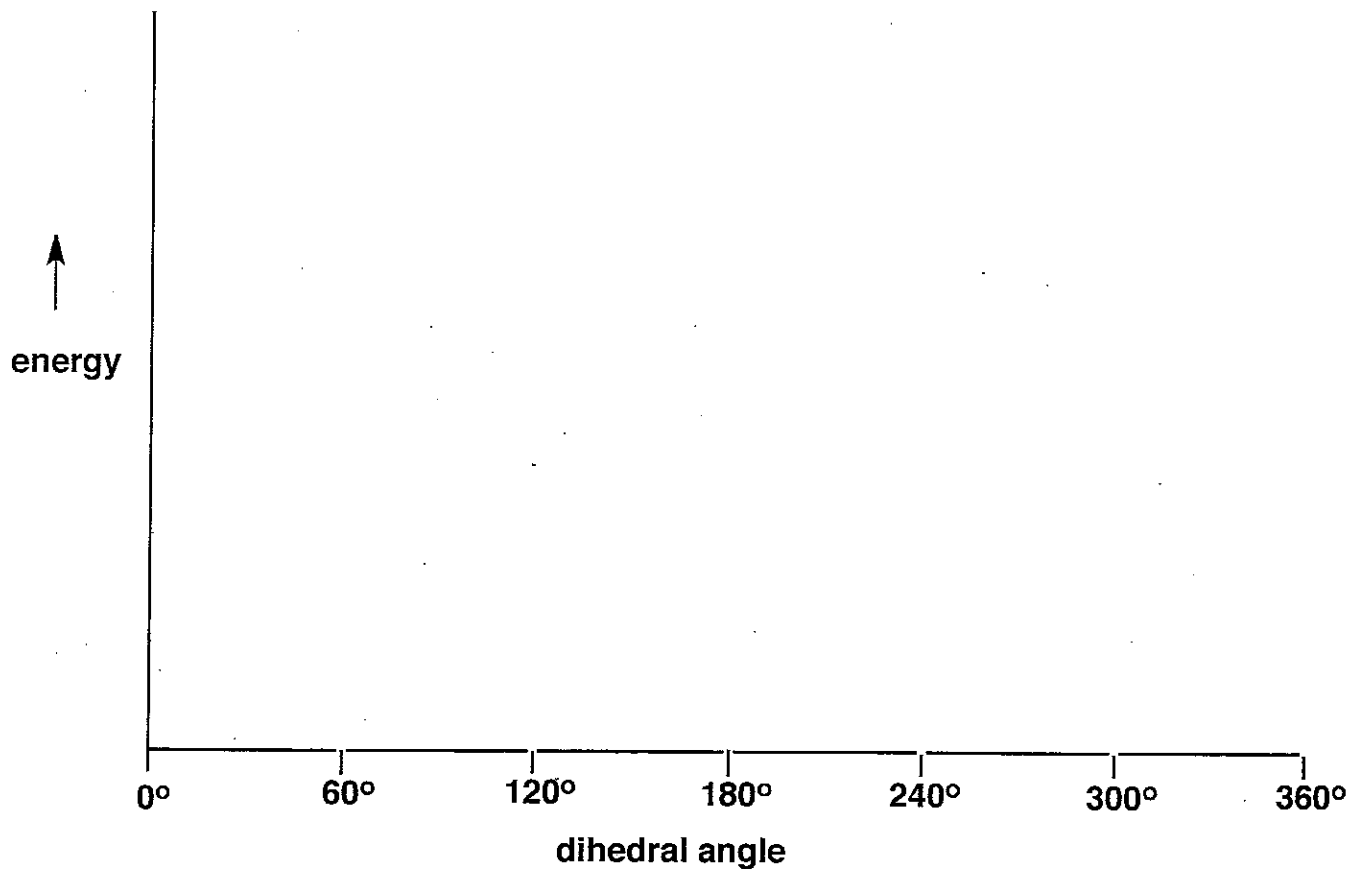
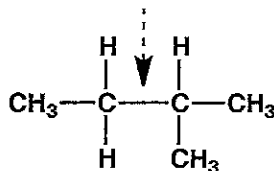
3. (12 points) For each equilibrium shown below, do two things:

(i) Put a **SQUARE** around the **WEAKER ACID**, of the two species that are serving as acids in the equilibrium.(ii) Put a **CIRCLE** around the **SIDE** of the equilibrium that you expect to be **LESS FAVORED**.

4. (25 points)

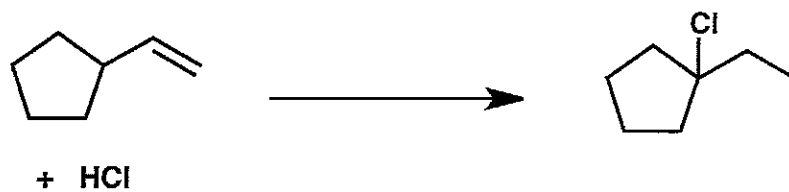
Draw an energy diagram for rotation about the indicated carbon-carbon bond (dotted arrow) of the molecule below. Show the structure for each energy minimum and each energy maximum.

CIRCLE the most stable structure(s).



Name _____

5. (12 points) Provide a mechanism ("curved arrows") for the reaction shown below. Show every atom in each structure you draw.

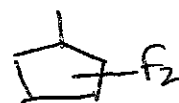


6. (25 points)

A molecule with the formula $C_6H_{10}F_2$ does not undergo any reaction when combined with H_2 in the presence of the catalyst Pd/C. Propose FIVE possible structures for this molecule (via appropriate drawings).



1 unsaturation = (ring)



| <u>Problem #</u> | <u>Score</u> |
|------------------|--------------|
| 1 | / 10 |
| 2 | / 16 |
| 3 | / 12 |
| 4 | / 25 |
| 5 | / 12 |
| 6 | / 25 |

Total: / 100

Periodic Table of the Elements

| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|---|----|---|--------|----|--------|--------|----|--------|--------|------|--------|-------|------|-----------|--------|------|--------|--------|------|-------|--------|------|--|--------|------|--------|--------|----|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|-------|--------|--------|--------|--------|----|--------|--------|----|--------|--------|----|--------|--------|----|--------|----|----|--------|----|----|--------|----|----|--------|----|----|--------|----|----|--------|--|----|----|--------|----|----|-------|----|---|--------|----|----|-------|----|----|-------|----|----|-------|----|----|-------|----|----|-------|----|----|-------|----|----|-------|-----|----|-------|-----|----|-------|-----|----|-------|-----|----|-------|
| | | <table border="1" style="margin: auto;"> <tr> <td style="padding: 2px;">1</td> <td style="padding: 2px;">H</td> <td style="padding: 2px;">1.008</td> </tr> </table> | | 1 | H | 1.008 | | | | | | | | | | | | | | | | | <table border="1" style="margin: auto;"> <tr> <td style="padding: 2px;">2</td> <td style="padding: 2px;">He</td> <td style="padding: 2px;">4.003</td> </tr> </table> | 2 | He | 4.003 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1 | H | 1.008 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 2 | He | 4.003 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 3 | Li | 6.94 | 4 | Be | 9.01 | | | | | | | | | | | | | | | | | 5 | B | 10.81 | 6 | C | 12.011 | 7 | N | 14.01 | 8 | O | 16.00 | 9 | F | 19.00 | 10 | Ne | 20.18 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 11 | Na | 22.99 | 12 | Mg | 24.31 | | | | | | | | | | | | | | | | | 13 | Al | 26.98 | 14 | Si | 28.09 | 15 | P | 30.97 | 16 | S | 32.06 | 17 | Cl | 35.45 | 18 | Ar | 39.95 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 19 | K | 39.10 | 20 | Ca | 40.08 | 21 | Sc | 44.96 | 22 | Ti | 47.90 | 23 | V | 50.94 | 24 | Cr | 52.00 | 25 | Mn | 54.94 | 26 | Fe | 55.85 | 27 | Co | 58.93 | 28 | Ni | 58.71 | 29 | Cu | 63.55 | 30 | Zn | 65.37 | 31 | Ga | 69.72 | 32 | Ge | 72.59 | 33 | As | 74.92 | 34 | Se | 78.96 | 35 | Br | 79.90 | 36 | Kr | 83.80 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 37 | Rb | 85.47 | 38 | Sr | 87.62 | 39 | Y | 88.91 | 40 | Zr | 91.22 | 41 | Nb | 92.91 | 42 | Mo | 95.94 | 43 | Tc | 98.91 | 44 | Ru | 101.07 | 45 | Rh | 102.91 | 46 | Pd | 106.4 | 47 | Ag | 107.87 | 48 | Cd | 112.40 | 49 | In | 114.82 | 50 | Sn | 118.69 | 51 | Sb | 121.75 | 52 | Te | 127.60 | 53 | I | 126.90 | 54 | Xe | 131.30 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 55 | Cs | 132.91 | 56 | Ba | 137.34 | 57 | La | 138.91 | 72 | Hf | 178.49 | 73 | Ta | 180.95 | 74 | W | 183.85 | 75 | Re | 186.2 | 76 | Os | 190.2 | 77 | Ir | 192.2 | 78 | Pt | 195.09 | 79 | Au | 196.97 | 80 | Hg | 200.59 | 81 | Tl | 204.37 | 82 | Pb | 207.19 | 83 | Bi | 208.98 | 84 | Po | (209) | 85 | At | (210) | 86 | Rn | (222) | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 87 | Fr | (223) | 88 | Ra | 226.03 | 89 | Ac | (227) | 104 | Unq* | (261) | 105 | Unp* | (262) | 106 | Unh* | (263) | 107 | Uns* | (262) | 108 | Uno* | (265) | 109 | Una* | (266) | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| <table border="1" style="width: 100%; border-collapse: collapse;"> <tr> <td style="width: 10%;"></td> <td style="width: 10%; text-align: center;">58</td> <td style="text-align: center;">Ce</td> <td style="text-align: center;">140.12</td> <td style="text-align: center;">59</td> <td style="text-align: center;">Pr</td> <td style="text-align: center;">140.91</td> <td style="text-align: center;">60</td> <td style="text-align: center;">Nd</td> <td style="text-align: center;">144.24</td> <td style="text-align: center;">61</td> <td style="text-align: center;">Pm</td> <td style="text-align: center;">(145)</td> <td style="text-align: center;">62</td> <td style="text-align: center;">Sm</td> <td style="text-align: center;">150.35</td> <td style="text-align: center;">63</td> <td style="text-align: center;">Eu</td> <td style="text-align: center;">151.96</td> <td style="text-align: center;">64</td> <td style="text-align: center;">Gd</td> <td style="text-align: center;">157.25</td> <td style="text-align: center;">65</td> <td style="text-align: center;">Tb</td> <td style="text-align: center;">158.93</td> <td style="text-align: center;">66</td> <td style="text-align: center;">Dy</td> <td style="text-align: center;">162.50</td> <td style="text-align: center;">67</td> <td style="text-align: center;">Ho</td> <td style="text-align: center;">164.93</td> <td style="text-align: center;">68</td> <td style="text-align: center;">Er</td> <td style="text-align: center;">167.26</td> <td style="text-align: center;">69</td> <td style="text-align: center;">Tm</td> <td style="text-align: center;">168.93</td> <td style="text-align: center;">70</td> <td style="text-align: center;">Yb</td> <td style="text-align: center;">173.04</td> <td style="text-align: center;">71</td> <td style="text-align: center;">Lu</td> <td style="text-align: center;">174.97</td> </tr> <tr> <td></td> <td style="text-align: center;">90</td> <td style="text-align: center;">Th</td> <td style="text-align: center;">232.04</td> <td style="text-align: center;">91</td> <td style="text-align: center;">Pa</td> <td style="text-align: center;">(231)</td> <td style="text-align: center;">92</td> <td style="text-align: center;">U</td> <td style="text-align: center;">238.03</td> <td style="text-align: center;">93</td> <td style="text-align: center;">Np</td> <td style="text-align: center;">(237)</td> <td style="text-align: center;">94</td> <td style="text-align: center;">Pu</td> <td style="text-align: center;">(244)</td> <td style="text-align: center;">95</td> <td style="text-align: center;">Am</td> <td style="text-align: center;">(243)</td> <td style="text-align: center;">96</td> <td style="text-align: center;">Cm</td> <td style="text-align: center;">(247)</td> <td style="text-align: center;">97</td> <td style="text-align: center;">Bk</td> <td style="text-align: center;">(249)</td> <td style="text-align: center;">98</td> <td style="text-align: center;">Cf</td> <td style="text-align: center;">(249)</td> <td style="text-align: center;">99</td> <td style="text-align: center;">Es</td> <td style="text-align: center;">(254)</td> <td style="text-align: center;">100</td> <td style="text-align: center;">Fm</td> <td style="text-align: center;">(257)</td> <td style="text-align: center;">101</td> <td style="text-align: center;">Md</td> <td style="text-align: center;">(258)</td> <td style="text-align: center;">102</td> <td style="text-align: center;">No</td> <td style="text-align: center;">(259)</td> <td style="text-align: center;">103</td> <td style="text-align: center;">Lr</td> <td style="text-align: center;">(260)</td> </tr> </table> | | | | | | | | | | | | | | | | | | | | | | | | | | | | | 58 | Ce | 140.12 | 59 | Pr | 140.91 | 60 | Nd | 144.24 | 61 | Pm | (145) | 62 | Sm | 150.35 | 63 | Eu | 151.96 | 64 | Gd | 157.25 | 65 | Tb | 158.93 | 66 | Dy | 162.50 | 67 | Ho | 164.93 | 68 | Er | 167.26 | 69 | Tm | 168.93 | 70 | Yb | 173.04 | 71 | Lu | 174.97 | | 90 | Th | 232.04 | 91 | Pa | (231) | 92 | U | 238.03 | 93 | Np | (237) | 94 | Pu | (244) | 95 | Am | (243) | 96 | Cm | (247) | 97 | Bk | (249) | 98 | Cf | (249) | 99 | Es | (254) | 100 | Fm | (257) | 101 | Md | (258) | 102 | No | (259) | 103 | Lr | (260) |
| | 58 | Ce | 140.12 | 59 | Pr | 140.91 | 60 | Nd | 144.24 | 61 | Pm | (145) | 62 | Sm | 150.35 | 63 | Eu | 151.96 | 64 | Gd | 157.25 | 65 | Tb | 158.93 | 66 | Dy | 162.50 | 67 | Ho | 164.93 | 68 | Er | 167.26 | 69 | Tm | 168.93 | 70 | Yb | 173.04 | 71 | Lu | 174.97 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | 90 | Th | 232.04 | 91 | Pa | (231) | 92 | U | 238.03 | 93 | Np | (237) | 94 | Pu | (244) | 95 | Am | (243) | 96 | Cm | (247) | 97 | Bk | (249) | 98 | Cf | (249) | 99 | Es | (254) | 100 | Fm | (257) | 101 | Md | (258) | 102 | No | (259) | 103 | Lr | (260) | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Lanthanides | | | | | | | | | | | | | | Actinides | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |

Numbers in parentheses: available radioactive isotope of longest half-life.

*Symbol (and name) provisional.