

1. a. (6 pts.) Determine the degree of unsaturation for the hydrocarbon with the formula C_5H_6 . 1. _____

2. _____

3. _____

b. (8 pts.) When a molecule has a degree of unsaturation of 1, what does that mean?

4. _____

5. _____

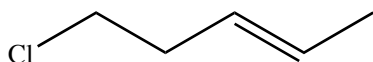
6. _____

2. (8 pts. each) Provide IUPAC names for the following alkenes. Use *Z/E* nomenclature where appropriate to specify the stereochemistry of the alkene. 7. _____

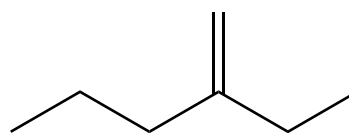
8. _____

9. _____

a.



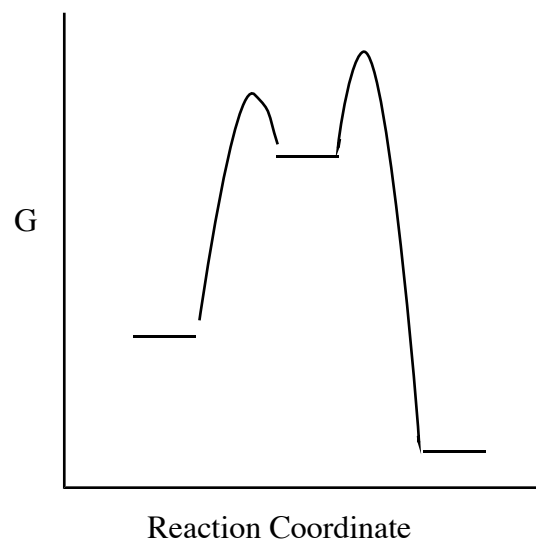
b.



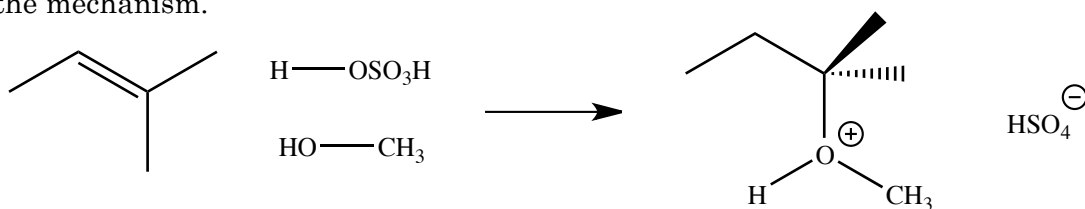
3. (8 pts.) Using valence bond theory (hybridization) explain why alkenes are nucleophilic.

4. (2 pts. ea.) The questions below refer to the reaction coordinate diagram draw to the right.

- Label the reactants with an "a".
- Label the products with a "b".
- Label the intermediates with a "c".
- Label the transition state(s) with a "d".
- Does this reaction absorb or release energy?
- Would this reaction have a positive or negative ΔG ?
- Does the equilibrium favor the reactants or products.

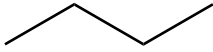
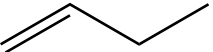
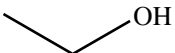


5. (12 pts.) Draw a mechanism for the reaction shown below. Include electron movement arrows with the mechanism.

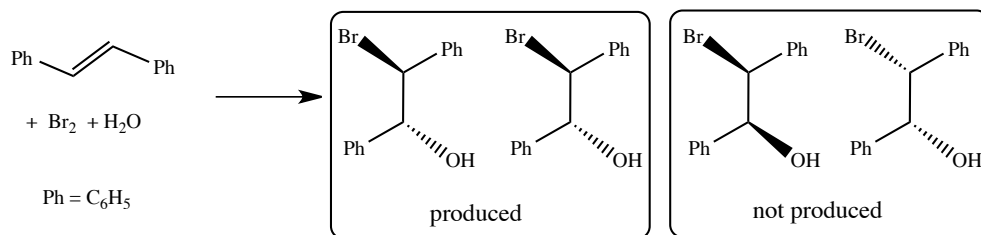


6. a. (8 pts.) Which is more stable, a primary or a tertiary carbocation. b. (12 pts.) Explain why one is more stable than the other.

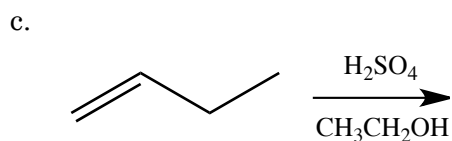
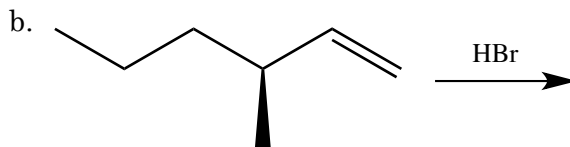
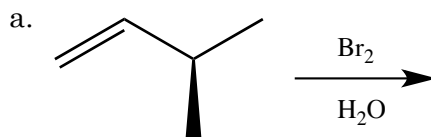
7. (16 pts.) Determine whether the following are nucleophiles, electrophiles, or neither.

H^+	CH_3OH		
Br^-	H_2SO_4		$CH_3CHCHCH_3$

8. (12 pts.) In lab, you reacted *trans*-stilbene with bromine and water. The product, 1,2-diphenyl-2-bromo-ethanol, exists as four stereoisomers, but your reaction only made two stereoisomers. Draw the intermediate that forms when the Br_2 reacts with the *trans*-stilbene and explain why only two of the four possible stereoisomers are formed.



9. (8 pts. ea.) Predict the major organic products for the following reactions. Remember to indicate the stereochemistry of the products using wedge (\blacktriangle), dashed (\cdots), or squiggly (\sim) bonds where appropriate.



1	H 1.0079																	2	He 4.0026																
3	Li 6.941	4	Be 9.012															9	F 18.998	10	Ne 20.1797														
11	Na 22.989	12	Mg 24.305															17	Cl 35.453	18	Ar 39.948														
19	K	20	Ca	21	Sc	22	Ti	23	V	24	Cr	25	Mn	26	Fe	27	Co	28	Ni	29	Cu	30	Zn	31	Ga	32	Ge	33	As	34	Se	35	Br	36	Kr
37	Cs	38	Sr	39	Y	40	Zr	41	Nb	42	Mo	43	Tc	44	Ru	45	Rh	46	Pd	47	Ag	48	Cd	49	In	50	Sn	51	Sb	52	Te	53	I	54	Xe
55	Rb	56	Ba	57	La	72	Hf	73	Ta	74	W	75	Re	76	Os	77	Ir	78	Pt	79	Au	80	Hg	81	Tl	82	Pb	83	Bi	84	Po	85	At	86	Rn
87	Fr	88	Ra	89	Ac	104	Rf	105	Db	106	Sg	107	Bh	108	Hs	109	Mt	110		111		112		114		116								118	

58	Ce	59	Pr	60	Nd	61	Pm	62	Sm	63	Eu	64	Gd	65	Tb	66	Dy	67	Ho	68	Er	69	Tm	70	Yb	71	Lu
90	Th	91	Pa	92	U	93	Np	94	Pu	95	Am	96	Cm	97	Bk	98	Cf	99	Es	100	Fm	101	Md	102	No	103	Lr