	•
Hour Exam #1 (PM) Chemistry 343	Last Name
Professor Gellman 5 October 2011	First Name
General Instructions:	
<ul><li>(i) Use scratch paper at back of example be recorded at the proper place of allowed.</li></ul>	to work out answers; final answers must n the exam itself for credit. Models are
(ii) Print your name on each page.	
(iii) Please keep your paper covered a Misconduct will lead to failure in	and your eyes on your own work. the course.
1. (15 points) Draw a structure that co Show all atoms in each structure, inclu	orresponds to each of the following names. uding hydrogen atoms.
(a) 3,3-difluoro-octane	
• •	
(b) 4-methyl-cycloheptene	
(b) 4 monty: Gyoromophone	

(c) E-3-methyl-2-hexene

Name		
Manie		

2. (14 points) For each set of three structures shown below, redraw them in the order of DECREASING  $pK_a$ , left to right. (Note: Positive charges are balanced by a chloride counterion.)

- (a) + CH<sub>3</sub>NH<sub>3</sub>
- ⊕ CH<sub>3</sub>OH<sub>2</sub>
- CH<sub>3</sub>NH<sub>2</sub>

$$(b) \qquad \begin{array}{c} O \\ \\ F \end{array} \qquad \begin{array}{c} O \\ \\ CI \end{array} \qquad \begin{array}{c} CI \end{array} \qquad$$

3. (7 points) Consider the alkane indicated below.

 $(\mathsf{CH}_3)_3\mathsf{CCH}_2\mathsf{CH}_2\mathsf{CH}_2\mathsf{CH}(\mathsf{CH}_3)\mathsf{CH}_2\mathsf{CH}_2\mathsf{CH}(\mathsf{CH}_2\mathsf{CH}_3)_2$ 

(a) Redraw the molecule so that ALL atoms are indicated by the appropriate atomic symbol (e.g, C for carbon), AND all bonds are shown as lines.

(b) Provide a SKELETAL drawing for this molecule (lines only).

$-\alpha\alpha$	170
F 35 F	1 /11
vv	

## 4. (20 points)

(a) Propane has two carbon-carbon bonds, but these bonds are equivalent in terms of their local environments (i.e., the bonding partners of the pair of carbons). Thus, one can say that propane contains only one TYPE of carbon-carbon bond.

Shown below is 3-ethyl-pentane. How many different TYPES of carbon-carbon bond are found in this molecule?

Number of TYPES of C-C bond = \_\_\_\_\_

(b) Draw Newman projections for all staggered conformations about the bond between carbon-2 and carbon-3 bond of 3-ethyl-pentane. CIRCLE the conformation(s) that you expect to be most stable.

Name	

5. (14 points) Consider the reaction below, which involve Lewis acids and bases and their complexes.

(a) Propose a mechanism (curved arrows) in which the reaction shown above occurs in a <u>single</u> step (no intermediates). [Note: Do not be confused by the sodium (Na) counterions, which are just 'bystanders' in this process.]

(b) Propose a mechanism (curved arrows) in which the reaction shown above occurs in two steps. [Note: Do not be confused by the sodium (Na) counterions, which are just 'bystanders' in this process.]

Name
. (30 points)
a) Identify (with appropriate drawings) TWO hydrocarbon molecules corresponding to the formula C <sub>6</sub> H <sub>12</sub> that would be expected to produce only ONE product upon reaction with HBr.
(b) Identify (with appropriate drawings) TWO hydrocarbon molecules corresponding to the formula $C_6H_{12}$ that would be expected to produce TWO products upon reaction with HBr.

(c) Identify (with appropriate drawings) TWO hydrocarbon molecules corresponding to the formula  $\rm C_6H_{12}$  that would NOT be expected to react with HBr.

lame ... 00170

Problem #	<u>Score</u>
1	/ 15
2	/14
3	/ 7
4	/20
5	/14
6	/30

Total:

/100

18 8A	He 4.0026	10 Ne 20.180	188	Ar	. 39,948	36	Kr	83.80	54	Xe	131.29	98	Rm	(222)	118	Uuo	(294)
<u>~</u>	17 7A	9 FF 18,998	17	۵	35.453	32.	Br	79.904	53	: ;	126.90	85	At	(210)	117	Uus	
ents t frequent	16 6A	8 O 0.	16	တ	32.06	34	Se	78.96	52	ام	127.60	84	Po	(509)	116	Uuh	(293)
The shaded elements will be encountered most frequently in the text.	15 5A	N 14,007	15	à.	30.974	33	As	74.922	51	Sb	121.76	83	Bi	208.98	115	·Uup	(288)
The sh	14 4A	6 C 11011	14	Si	28.085	32	Ge	72.61	50	Sn	118.71	82	Pb	.207.2	114	Uuq	(289)
TI.M	13 3A	. 5 B 10,811.	13	ĄĮ	26,9815	31	Ga	69.723	49	In	114.82	81	I	204.38	113	Uut	(284)
		-		12	2B	30	Zn	65.38	48	Cd	112.41	80	Hg	200.59	112	Unp	(285)
				11	1B	29	Cu	63.546	47	Ag	107.87	79	Aп	196,97	111	Rg	(280)
	n,			10		28	Z	58.693	46	Pd	106.42	8/	Pt	195.08	110	Ds	(281)
	re in greer d in	eroup pe of	\ 	6	—8B —	27	ප	58.933	45	Rh	102.91	11	I	192.22	109	Mt	(276)
	IUPAC anse are used	or me A-1	elements	œ		79	Fe	55.845	44	Ru	101.07	9/	$^{ m so}$	190,23	108	Hs	(270)
c Table ements	Group numbers recommended by the IUPAC are in green. Older group numbers are in red. (These are used in the text for calculating formal charge for the Aurona	the text for calculating formal charge for the Argroup elements.) (Atomic weights in parentheses are for the isotope of longest life.)	Transition elements	7	7B	25	Mn	54.938	43	Tc	(98)	7.5	Re	186.21	107	Bh	(272)
	commend ers are in	parenthe		9 .	6B	24	Ü	51.996	42	Mo	95.96	74	≯	183.84	106	Sg	(271)
A Periodic of the Ele	mbers rec up number	n calcular ) veights in e.)	7	5	5B	23	Λ	50.942	41	NP	92.906	73	Ta	180.95	105	Dp	(268)
A Pe of th	Group nu Older gro	elements.) (Atomic we longest life.		4	4B	22	Ξ	47.867	-40	Zr	91.224	72	Hŧ	178.49	104	Rf	(267)
					3B	21	Sc	44,956	39	>-	88.906	71	Ľu	174.97	103	Lr	(262)
	2.A	<b>Be</b>	12	-60° Z	, 24.305	2.0 2.0	లో	40.08	38	Sr	87.62	99	Ba	137,33	88	Ra	(226)
- ¥1	Z6Z001	11 6-9	#	Na	22.9898	6)	¥	39,098	37	Rb	85,468	55	౮	132.91	87	Ä	(223)
Period	1	77		eD.			4			5			9			7	

	57	58	29	9	19	62	63	64	65	99	<i>L</i> 9	89	69	. 70
Lanthanides	La	c	Pŗ	PN	Pm .	Sm	Eu	Gd	TP	Dy	Но	Hr	Tlm	Yb
	138.91	140.12	140.91	144.24	(145)	150.36	151.96	157.25	158.93	162,50	164.930	167.26	168.934	173.06
	68	06	16	92	93	94	95	96	26	86	66	100	101	102
Actinides	Ac	Th	Pa	U	ď	Pu	Am	Cm	Bk	Cť	Es	Fm	Мď	No
	(227)	232.04	231.04	238.03	(237)	(244)	(243)	(247)	(247)	(251)	(223)	(257)	(258)	(528)