NUMERICAL ANSWERS TO ASSIGNED TUTORIAL PROBLEM SETS FOR CHEM205 FROM KOTZ & TREICHEL'S CHEMISTRY & CHEMICAL REACTIVITY, **6**<sup>th</sup> **Ed.** 

NOTE: none of the answers from Ch.10 have been verified. Please report any errors.

Ch.	Q#	Comments
10	2	NF <sub>3</sub> : 26 valence ēs; electron-pair geometry = tetrahedral; molecular geometry = trigonal pyramidal. The N atom is sp³-hybridized. Three of these sp³ hybrid orbitals each overlap with a fluorine 2p orbital to form three N-F sigma (single) bonds; the N's fourth sp³ hybrid orbital contains a pair of nonbonding electrons (a lone pair).
10	4	(a) $CSe_2$ : sp; (b) $SO_2$ : sp <sup>2</sup> ; (c) $CH_2O$ : sp <sup>2</sup> ; (d) $NH_4^+$ : sp <sup>3</sup>
10	6a 6b 6c	N (both, actually): sp <sup>3</sup> ; C: sp <sup>2</sup> C of CH <sub>3</sub> : sp <sup>3</sup> ; C of C=C and C=O: sp <sup>2</sup> C of C=C: sp <sup>2</sup> ; C of C to N triple bond: sp
10	8a 8b 8c 8d	XeOF <sub>4</sub> : electron pair geometry: octahedral, molecular geometry: square pyramidal, $sp^3d^2$ . BrF <sub>5</sub> : electron pair geometry: octahedral, molecular geometry: square pyramidal, $sp^3d^2$ . SOF <sub>4</sub> : electron pair geometry: trigonal bipyramid, molecular geometry: trigonal bipyramidal, $sp^3d$ . Br <sub>3</sub> <sup>-</sup> : electron pair geometry: trigonal bipyramid, molecular geometry: linear, $sp^3d$ .
10	10a 10b	$HOSO_2F$ : 32 valence e <sup>-</sup> , "electron pair geometry" and molecular geometry at sulfur is tetrahedral: $sp^3$ . $SO_3F^-$ : same as part (a).
10	12	SO <sub>2</sub> F <sub>2</sub> : electron pair geometry: tetrahedral, molecular geometry: tetrahedral, sp <sup>3</sup> .
10	24	CIF <sub>2</sub> <sup>+</sup> : electron pair geometry: tetrahedral, molecular geometry: bent, sp <sup>3</sup> : 109°. CIF <sub>2</sub> <sup>-</sup> : electron pair geometry: trigonal bipyramid, molecular geometry: linear, sp <sup>3</sup> d: 180°.
10	26	The resonance structures are in the text. They all have the same ${\sf sp}^2$ (trigonal) hybridization. The unused 2p-orbital on N is used to make the pi-bond.
10	28	$CO_2$ : hybridization: linear - sp, bond angle: $180^\circ$ , C to O bond order: 2. $CO_3^{2-}$ : hybridization: trigonal - sp <sup>2</sup> , bond angle: $120^\circ$ , C to O bond order: 4/3.
10	32a 32b	Angle A = $120^{\circ}$ ; angle B = $109^{\circ}$ (actually will be distorted because of lone pairsto approx. $105^{\circ}$ ); angle C = $109^{\circ}$ ; angle D = $120^{\circ}$ . Carbon 1: $sp^2$ ; carbon 2: $sp^2$ ; carbon 3: $sp^3$ .
10 10 10	34a 34b 34c	1 $\pi$ -bond, 11 $\sigma$ -bonds $C(1) = sp^3$ , $C(2) = sp^2$ , $C(3) = sp^3$ The C=O bond is the shortest and strongest CO bond.
10 10	36a 36b	Structure: N (with lone pair) is bonded to 2 H's & S; S also bonded to three O's. The angles around N & S are approximately 109° (tetrahedral). The hybridization of the N atom does not change (sp $^3$ in NH $_2$ $^-$ and in H $_2$ N-SO $_3$ $^-$ ). The S atom hybridization changes from sp $^2$ in SO $_3$ to sp $^3$ in H $_2$ N-SO $_3$ $^-$ .

Ch.	Q#	Comments
10	40	XeO <sub>3</sub> : electron pair geometry: tetrahedral - sp <sup>3</sup> , molecular geometry: trigonal pyramidal.  XeO <sub>4</sub> : electron pair geometry: tetrahedral - sp <sup>3</sup> , molecular geometry: tetrahedral.
10	46a 46b 46c 46d 46e	$C_6$ ring C atoms: $sp^2$ ; side chain C atoms $sp^3$ ; N atom: $sp^3$ . angle A = $120^\circ$ ; angle B = $109^\circ$ ; angle C = $109^\circ$ . (actually will be lessbecause of lone pair occupying more space than bonding pairs) 23 sigma-bonds and 3 pi-bonds. The molecule is polar. The $H^+$ ion attaches to the most electronegative atom in the molecule, <i>i.e.</i> N.
10 10 10	52a 52b 52c	$CF_4$ is isoelectronic with $BF_4^-$ (32 valence electrons) SiF <sub>4</sub> (32 valence electrons) and SF <sub>4</sub> (34 valence electrons) are not isoelectronic. $BF_4^-$ : B is sp <sup>3</sup> ; SiF <sub>4</sub> : F is sp <sup>3</sup> ; SF <sub>4</sub> : S is sp <sup>3</sup> d.