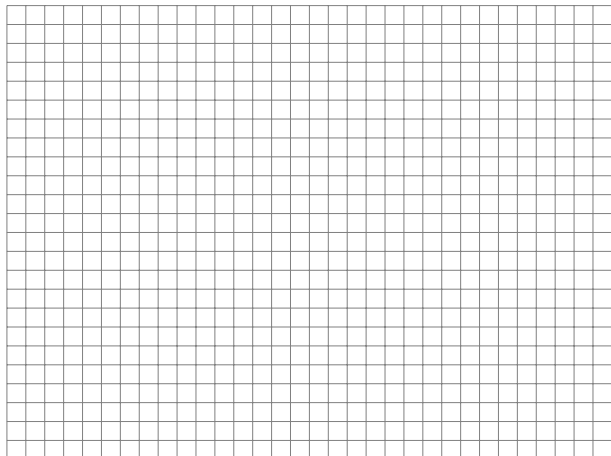


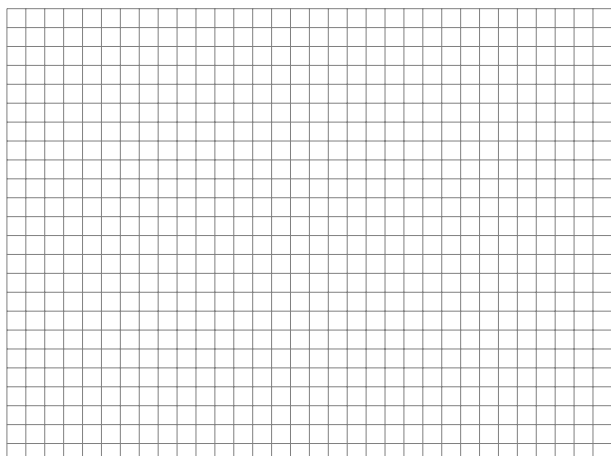
CHAPTER 0

ELECTRON-DOT STRUCTURES OF MOLECULES

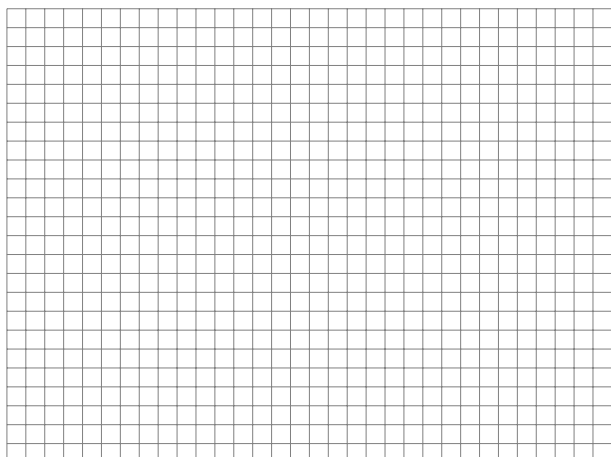
0.1 Draw electron-dot structures for the following diatomic molecules that obey the octet rule: (a) F_2 (b) Cl_2



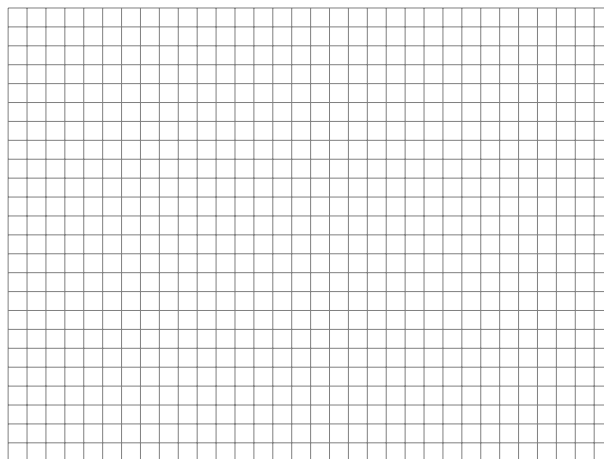
0.2 Draw electron-dot structures for the following diatomic molecules that obey the octet rule: (a) HF (b) HCl



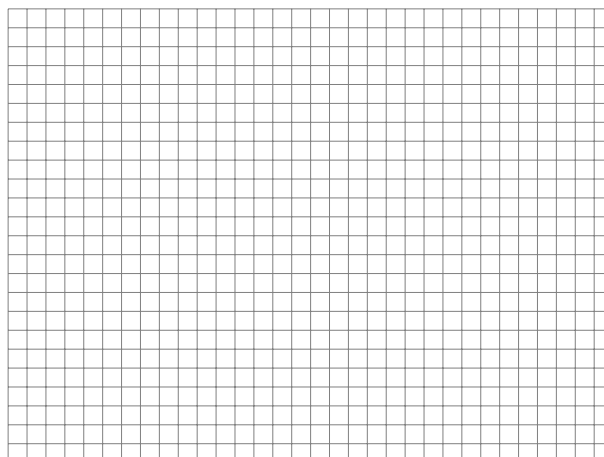
0.3 Draw electron-dot structures for the following diatomic molecules that obey the octet rule: (a) ICl (b) HI



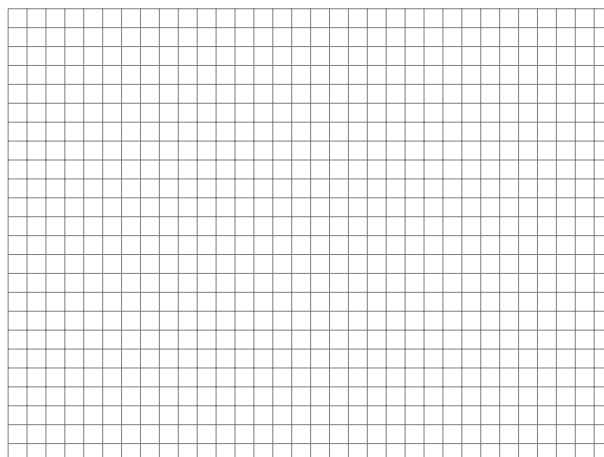
0.4 Draw electron-dot structures for the following diatomic molecules: (a) CO (b) N_2 (c) O_2



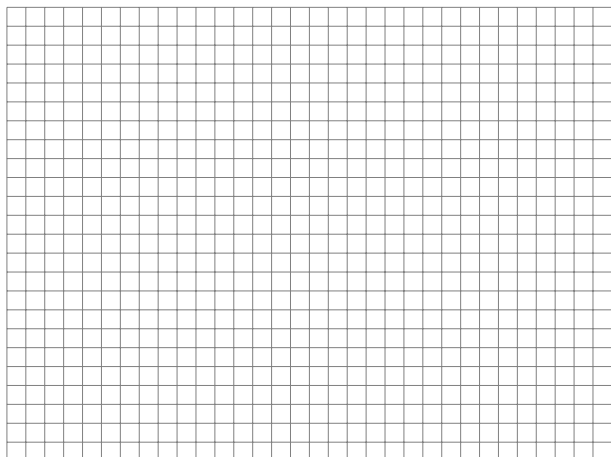
0.5 Draw electron-dot structures for the following molecules that obey the octet rule, given that the first atom listed is the central atom: (a) CHN (b) CO_2



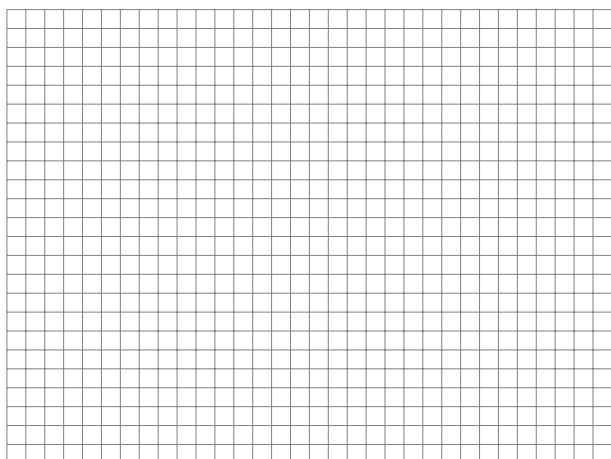
0.6 Draw electron-dot structures for the following molecules that obey the octet rule, given that the first atom listed is the central atom: (a) CH_4 (b) CH_3Cl (c) OH_2



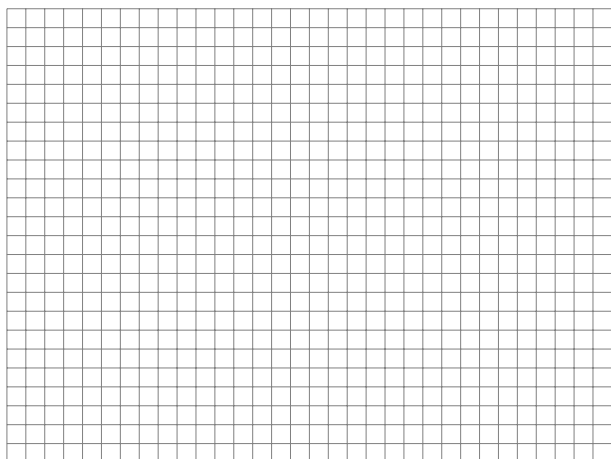
0.7 Draw electron-dot structures for the following molecules that obey the octet rule, given that the first atom listed is the central atom: (a) NH_3 (b) NCl_3



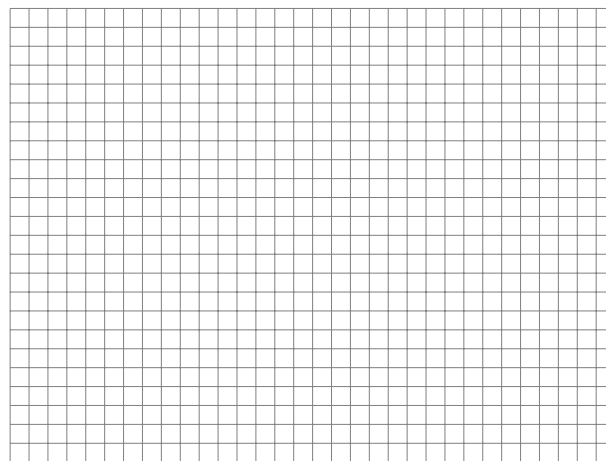
0.8 Draw electron-dot structures for the following molecules that obey the octet rule, given that the first atom listed is the central atom: (a) SeCl_2 (b) CH_2O



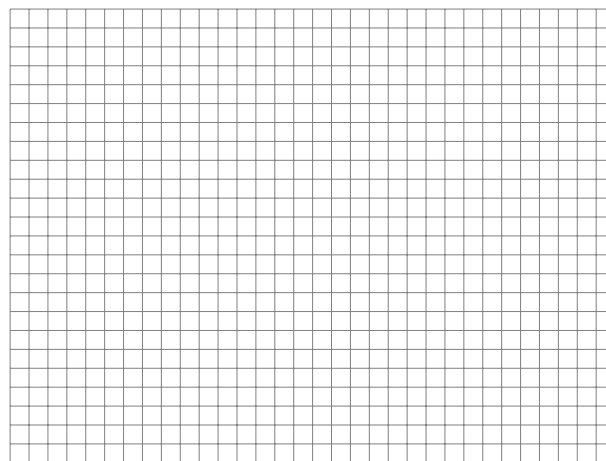
0.9 Draw electron-dot structures for the following molecules given that the first atom listed is the central atom. Some of the atoms might not obey the octet rule. If the species has a charge indicate the location of the charge: (a) BH_3 (b) BH_2F (c) POCl_3 (d) ClO_4^-



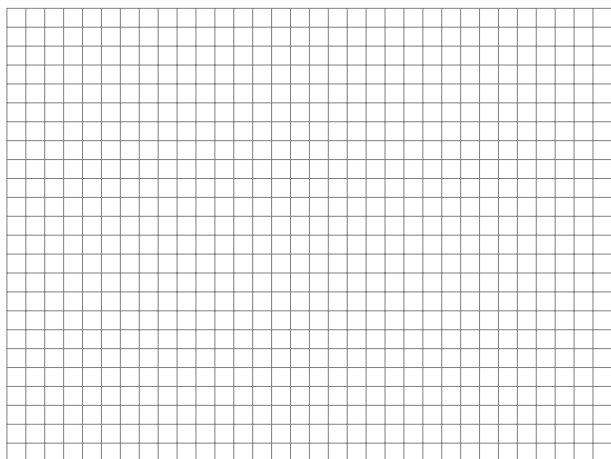
0.10 Draw electron-dot structures for the following molecules given that the first atom listed is the central atom. Some of the atoms might not obey the octet rule. If the species has a charge indicate the location of the charge: (a) BeH_2 (b) PCl_5 (c) SF_4 (d) ClF_3



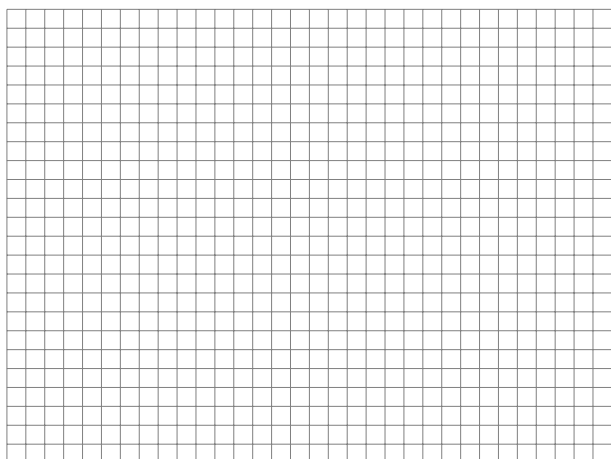
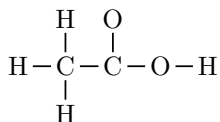
0.11 Draw electron-dot structures for the following molecules given that the first atom listed is the central atom. Some of the atoms might not obey the octet rule. If the species has a charge indicate the location of the charge: (a) I_3^- (b) Br_3^- (c) SF_6



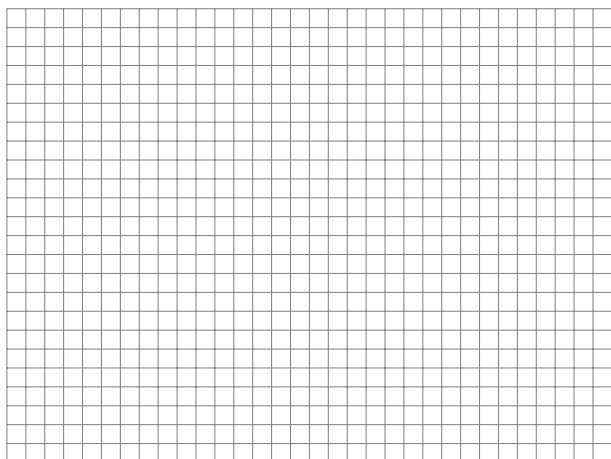
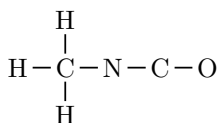
0.12 Draw electron-dot structures for the following molecules given that the first atom listed is the central atom. Some of the atoms might not obey the octet rule. If the species has a charge indicate the location of the charge: (a) ClF_5 (b) XeF_4



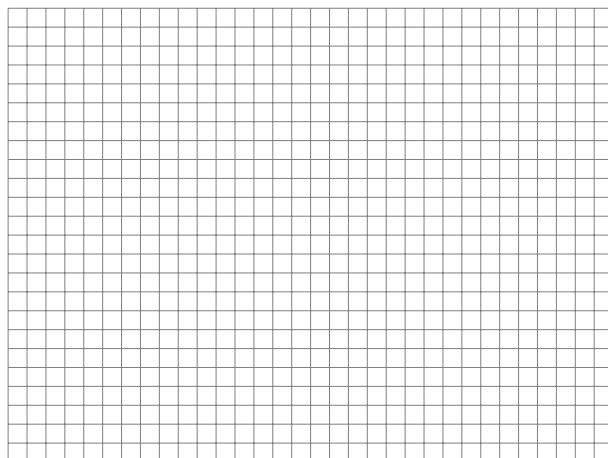
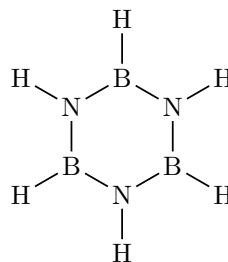
0.13 Given the skeletal structure below, draw the Lewis structure of the molecule:



0.14 Given the skeletal structure below, draw the Lewis structure of the molecule:



0.15 Given the skeletal structure below, draw the Lewis structure of the molecule:



0.16 Indicate the charge of the atom marked blue in the following electron-dot structure:



0.17 Indicate the charge of the atom marked blue in the following electron-dot structure:



0.18 Indicate the charge of the atom marked blue in the following electron-dot structure that follow the octet rule:



0.19 Indicate the charge of the atom marked blue in the following electron-dot structure that follow the octet rule:

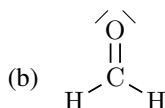
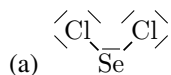


MOLECULAR SHAPE

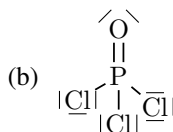
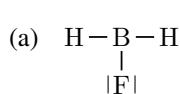
0.20 Identify the molecular shape of the molecules:
(a) NH_3 (b) CH_4

0.21 Identify the molecular shape of the molecules:
(a) H_2 (b) BeCl_2 (c) BF_3

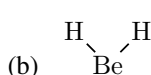
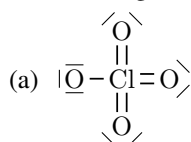
0.22 Given the following Lewis structures, predict the molecular geometry and angles:



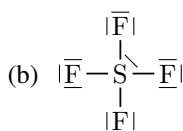
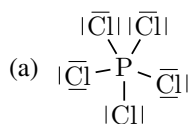
0.23 Given the following Lewis structures, predict the molecular geometry and angles:



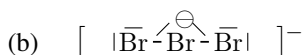
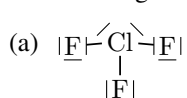
0.24 Given the following Lewis structures, predict the molecular geometry and angles:



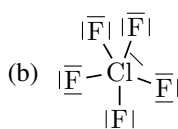
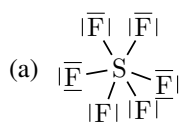
0.25 Given the following Lewis structures, predict the molecular geometry and angles:



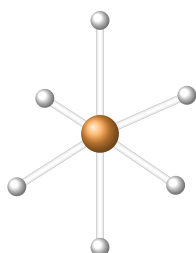
0.26 Given the following Lewis structures, predict the molecular geometry and angles:



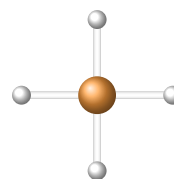
0.27 Given the following Lewis structures, predict the molecular geometry and angles:



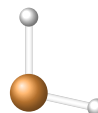
0.28 Identify the name of the following molecular structure:



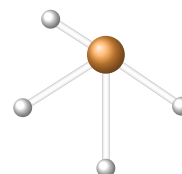
0.29 Identify the name of the following molecular structure:



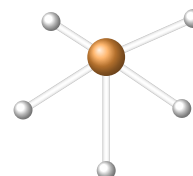
0.30 Identify the name of the following molecular structure:



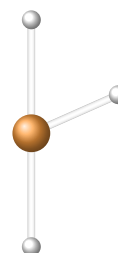
0.31 Identify the name of the following molecular structure:



0.32 Identify the name of the following molecular structure:



0.33 Identify the name of the following molecular structure:



POLARITY

0.34 Indicate the polarity or non-polarity for the following molecules: (a) H_2O (b) HCl (c) H_2

0.35 Indicate the polarity or non-polarity for the following molecules: (a) NH_3 (b) CO_2

HYBRID ORBITALS

0.36 Indicate the hybridization of: (a) NH_3 (b) CH_4
(c) H_2O

0.37 Indicate the hybridization of: (a) NH_3 (b) CH_4
(c) H_2O

MOLECULAR ORBITAL THEORY

0.38 Using the MO order provided below

$$\sigma_{2s}\sigma_{2s}^*\pi_{2p}\sigma_{2p}\pi_{2p}^*\sigma_{2p}^*$$

obtain the MO configuration for: (a) B_2 (b) C_2

0.39 Using the MO order provided below

$$\sigma_{2s}\sigma_{2s}^*\pi_{2p}\sigma_{2p}\pi_{2p}^*\sigma_{2p}^*$$

obtain the MO configuration for: (a) O_2 (b) F_2^+

0.40 Indicate the magnetic (paramagnetic or diamagnetic) configuration of the molecule with MO configuration: $\sigma_{2s}^2\sigma_{2s}^{*2}\sigma_{2p}^2\pi_{2p}^4\pi_{2p}^{*3}$

0.41 Indicate the magnetic (paramagnetic or diamagnetic) configuration of the molecule with MO configuration: $\sigma_{2s}^2\sigma_{2s}^{*2}\sigma_{2p}^2\pi_{2p}^4\pi_{2p}^{*2}$

- Answers** 0.1 (a) $|\overline{\text{F}}-\overline{\text{F}}|$ (b) $|\overline{\text{Cl}}-\overline{\text{Cl}}|$ 0.2 (a) $\text{H}-\overline{\text{F}}|$ (b) $\text{H}-\overline{\text{Cl}}|$ 0.3 (a) $|\overline{\text{I}}-\overline{\text{Cl}}|$ (b) $\text{H}-\overline{\text{I}}|$
- 0.4 (a) $|\text{C}\equiv\text{O}|$ (b) $|\text{N}\equiv\text{N}|$ (c) $\langle\text{O}=\text{O}\rangle$ 0.5 (a) $\text{H}-\text{C}\equiv\text{N}|$ (b) $\langle\text{O}=\text{C}=\text{O}\rangle$ 0.6 (a) $\begin{array}{c} \text{H} \\ | \\ \text{H}-\text{C}-\text{H} \\ | \\ \text{H} \end{array}$ (b) $\begin{array}{c} \text{H} \\ | \\ \text{H}-\text{C}-\text{H} \\ | \\ |\overline{\text{Cl}}| \end{array}$
- (c) $\begin{array}{c} \text{H} \quad \text{H} \\ \diagdown \quad \diagup \\ \text{O} \end{array}$ 0.7 (a) $\begin{array}{c} \text{H}-\overline{\text{N}}-\text{H} \\ | \\ \text{H} \end{array}$ (b) $|\overline{\text{Cl}}-\overline{\text{N}}-\overline{\text{Cl}}|$ 0.8 (a) $\begin{array}{c} \langle\overline{\text{Cl}}\rangle \quad \langle\overline{\text{Cl}}\rangle \\ \diagdown \quad \diagup \\ \text{Se} \end{array}$ (b) $\begin{array}{c} \langle\text{O}\rangle \\ || \\ \text{H}-\text{C}-\text{H} \end{array}$ 0.9 (a) $\begin{array}{c} \text{H}-\text{B}-\text{H} \\ | \\ \text{H} \end{array}$
- (b) $\begin{array}{c} \text{H}-\text{B}-\text{H} \\ | \\ |\overline{\text{F}}| \end{array}$ (c) $\begin{array}{c} \langle\text{O}\rangle \\ || \\ |\overline{\text{Cl}}|-\text{P}-|\overline{\text{Cl}}| \\ | \\ |\overline{\text{Cl}}| \end{array}$ (d) $|\overline{\text{O}}-\langle\text{O}\rangle=\text{O}\rangle$ 0.10 (a) $\begin{array}{c} \text{H} \quad \text{H} \\ \diagdown \quad \diagup \\ \text{Be} \end{array}$ (b) $|\overline{\text{Cl}}|-\text{P}-|\overline{\text{Cl}}|$ (c) $|\overline{\text{F}}|-\text{S}-|\overline{\text{F}}|$ (d) $|\overline{\text{F}}|-\text{Cl}-|\overline{\text{F}}|$
- 0.11 (a) $[\text{I}-\overline{\text{I}}-\overline{\text{I}}]^-$ (b) $[\overline{\text{Br}}-\overline{\text{Br}}-\overline{\text{Br}}]^-$ (c) $|\overline{\text{F}}|-\text{S}-|\overline{\text{F}}|$ 0.12 (a) $|\overline{\text{F}}|-\text{Cl}-|\overline{\text{F}}|$ (b) $|\overline{\text{F}}|-\text{Xe}-|\overline{\text{F}}|$
- 0.13 $\begin{array}{c} \text{H} \quad \langle\text{O}\rangle \\ | \quad || \\ \text{H}-\text{C}-\text{C}-\text{O}-\text{H} \\ | \\ \text{H} \end{array}$ 0.14 $\begin{array}{c} \text{H} \\ | \\ \text{H}-\text{C}-\overline{\text{N}}=\text{C}=\langle\text{O}\rangle \\ | \\ \text{H} \end{array}$ 0.15 $\begin{array}{c} \text{H} \\ | \\ \text{H}-\text{N}=\text{B}=\text{N}-\text{H} \\ || \quad | \\ \text{H}-\text{B}-\text{N}=\text{B}-\text{H} \\ | \\ \text{H} \end{array}$ 0.16 (a) -2 (b) +1 0.17 (a) +1 (b) -1
- 0.18 (a) +1 (b) +3 0.19 (a) +2 (b) +4 0.20 (a) NH_3 (Trigonal pyramidal) (b) CH_4 (Tetrahedral) 0.21 (a) H_2 (Linear) (b) BeCl_2 (Linear) (c) BF_3 (Trigonal planar) 0.22 (a) AB_2E_2 ; bent; 109° (b) AB_2E_2 ; bent; 109° ABE_3 ; planar trigonal; 120° 0.23 (a) ABE_3 ; planar trigonal; 120° (b) AB_4 ; tetrahedral; 109.5° 0.24 (a) AB_4 ; tetrahedral; 109.5° (b) AB_2 ; linear; 180° 0.25 (a) AB_5 ; trigonal bipyramidal; 120° and 90° (b) AB_4E ; see-saw; 120° and 90° 0.26 (a) AB_3E_2 ; T-shaped; 180° and 90° (b) AB_2E_3 ; linear; 180° 0.27 (a) AB_6 ; octahedral; 180° and 90° (b) AB_5E ; square pyramidal; 90° 0.28 Octahedral 0.29 square planar 0.30 bent 0.31 see-saw 0.32 square pyramidal 0.33 t-shaped 0.34 (a) H_2O (b) HCl (c) H_2 0.35 (a) NH_3 (b) CO_2 0.36 (a) NH_3 (sp^3) (b) CH_4 (sp^3) (c) H_2O (sp^3) 0.37 (a) NH_3 (sp^3) (b) CH_4 (sp^3) (c) H_2O (sp^3) 0.38 (a) B_2 ($\sigma_{2s}^2\sigma_{2s}^{*2}\pi_{2p}^2$) (b) C_2 ($\sigma_{2s}^2\sigma_{2s}^{*2}\pi_{2p}^4$) 0.39 (a) O_2 ($\sigma_{2s}^2\sigma_{2s}^{*2}\sigma_{2p}^2\pi_{2p}^4\pi_{2p}^{*2}$) (b) F_2^+ ($\sigma_{2s}^2\sigma_{2s}^{*2}\sigma_{2p}^2\pi_{2p}^4\pi_{2p}^{*3}$) 0.40 paramagnetic 0.41 paramagnetic