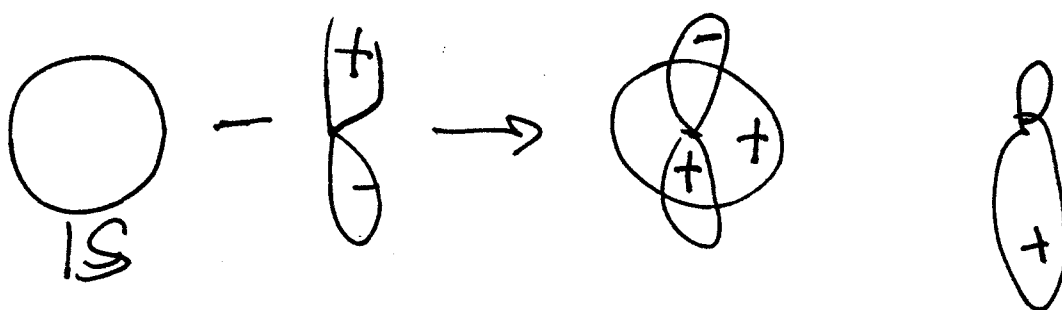
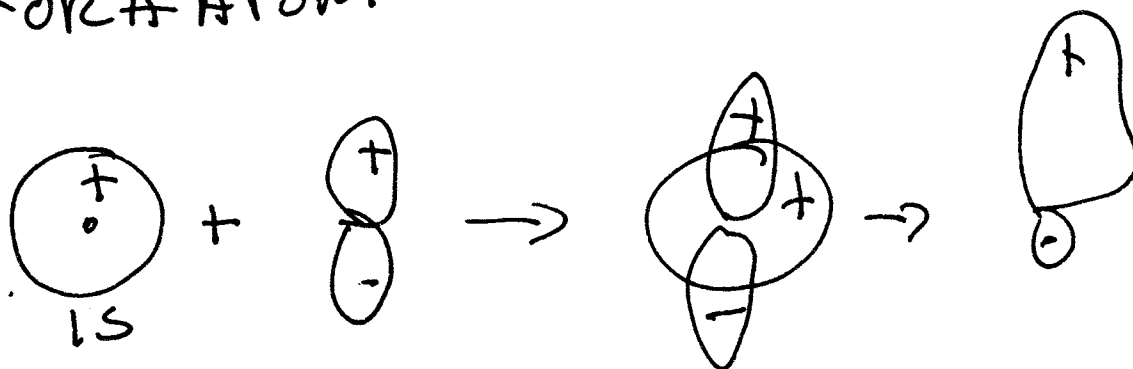


Valence Electron

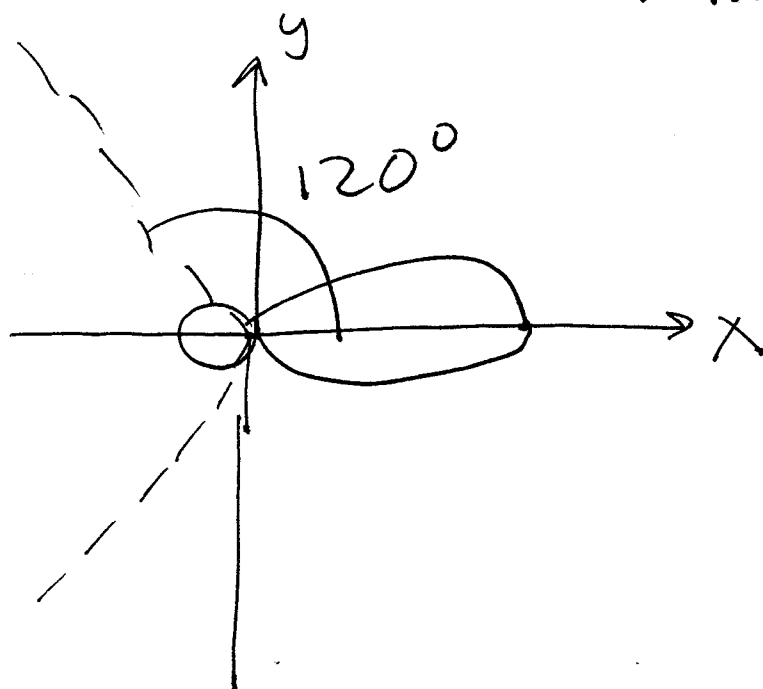
FOR A ATOM



SP \rightarrow 2 orbitals (atomic)

\Rightarrow LINEAR MOLECULES

sp^2 (S, P_x , P_y) TRIGONAL
PLANAR

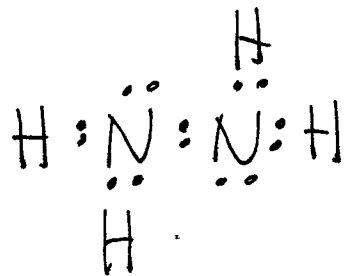
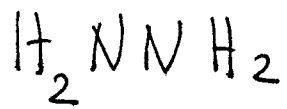


sp^3 TETRAHEDRAL

$\angle 109.5^\circ$

HYDRAZINE

(sp³)



TRIATOMIC

$\leq 16 e^- \rightarrow$ LINEAR

$17 \leq \# \leq 20 \rightarrow$ BENT

TRIATOMIC

LINEAR CO_2 , N_2O , OCS , and NO_2^+

BENT NO_2 , O_3 , NF_2 , and NO_2^-

LINEAR

A-B-C

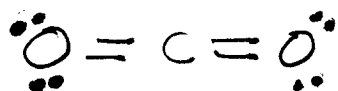
OUTER ATOMS \rightarrow OCCUPIED S ORBITAL
(A, C) ONLY P'S ORBITALS

CENTRAL ATOM \rightarrow SP ORBITALS (P_z)
(B)

$\sigma \rightarrow$ BONDS \Rightarrow SP AND P_z

EACH ATOM \Rightarrow P_x, P_y ORBITALS

$\text{CO}_2 \rightarrow 8e^-$ FOR π bonding



NONLINEAR

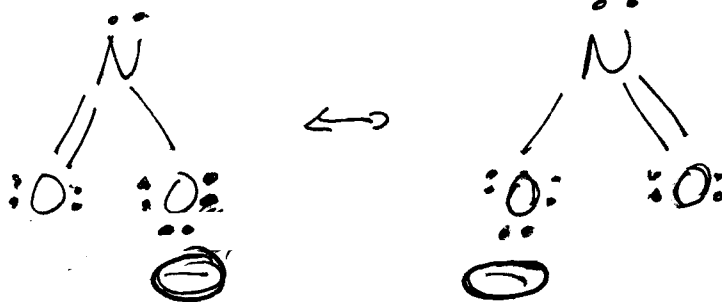
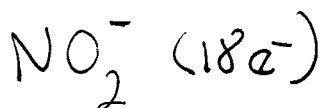
CENTRAL ATOM $\rightarrow sp^2$ ($s p_x p_y$) plane xy

$p_z \rightarrow \pi$ bonding



OUTER ATOMS $\rightarrow p_x$ or $p_y \rightarrow \sigma$ bond
(NO ORBITAL)

$p_z \rightarrow \pi$ bonding



$4e^- \rightarrow \pi$ bonding

$2e^- \rightarrow 2s$ (O) } localized
 $2e^- \rightarrow 2p$ (non bonding) } O

$2e^- \rightarrow$ lone pair in 3^{rd} sp^2

$4e^- \rightarrow \pi$ bonding $\left\{ \begin{array}{l} \text{bonding} \\ \text{nonbonding} \end{array} \right.$