

# Lewis Structures

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YALE UNIVERSITY  
CHEMISTRY 161  
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[www.mioy.org/chem161](http://www.mioy.org/chem161)

s-block												p-block							
1																			8
H																			He
←1s→	2											3	4	5	6	7	←1s→		
Li	Be											B	C	N	O	F	Ne		
←2s→												←2p→							
Na	Mg	d-block										Al	Si	P	S	Cl	Ar		
←3s→												←3p→							
K	Ca	Sc	Ti	V	Cr	Mn	Fe	Co	Ni	Cu	Zn	Ga	Ge	As	Se	Br	Kr		
←4s→		←3d→										←4p→							
Rb	Sr	Y	Zr	Nb	Mo	Tc	Ru	Rh	Pd	Ag	Cd	In	Sn	Sb	Te	I	Xe		
←5s→		←4d→										←5p→							
Cs	Ba	La	Hf	Ta	W	Re	Os	Ir	Pt	Au	Hg	Tl	Pb	Bi	Po	At	Rn		
←6s→		←5d→										←6p→							
Fr	Ra	Ac	Rf	Db	Sg	Bh	Hs	Mt	Ds	Rg	Cn	Nh	Fl	Mc	Lv	Ts	Og		
←7s→		←6d→										←7p→							

**VALENCE ELECTRONS**  
*Electrons in the highest-energy level or the outermost (largest  $n$  value) orbitals.*

N: (7 electrons,  $n = 2$ )

O: (8 electrons,  $n = 2$ )

Na: (11 electrons,  $n = 3$ )

S: (16 electrons,  $n = 3$ )

Br<sup>-</sup>: (36 electrons,  $n = 4$ )

f-block													
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←4s→		←3d→										←4p→							
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←4f→													
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Guided Example

Ammonia ( $\text{NH}_3$ )

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2. The *least* electronegative atom is usually the central atom (never Hydrogen though).

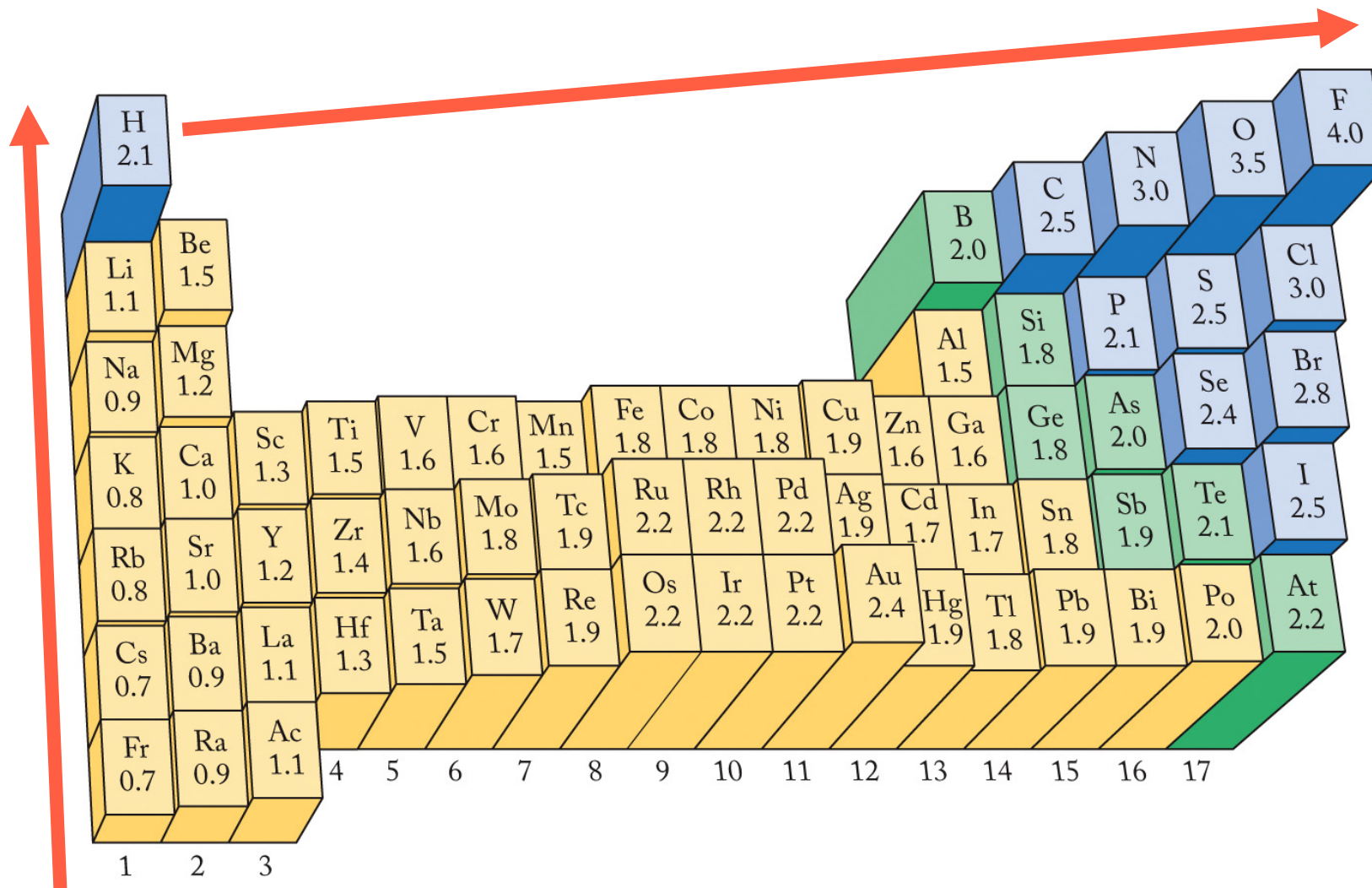
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## ELECTRONEGATIVITY

*The ability for an atom to attract electrons to itself.*



Same trend as ionization energy (IE)!

EN decreases down a column.  
EN increases across a row.

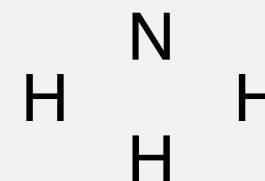
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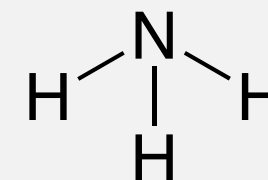
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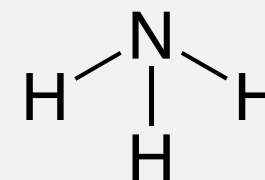
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4. Fulfill octet rule for each atom (8 electrons around each) by adding lone pairs (sets of 2 electrons).

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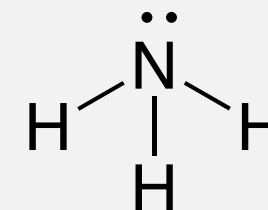
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Ammonia (NH<sub>3</sub>)

<b>Atom</b>	<b>Valence</b>
N	5 e <sup>-</sup>
H	1 e <sup>-</sup>
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<b>Total</b>	<b>8 e<sup>-</sup></b>



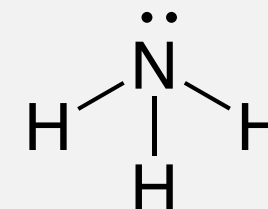
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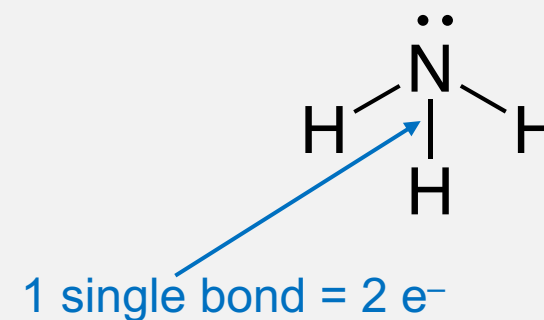
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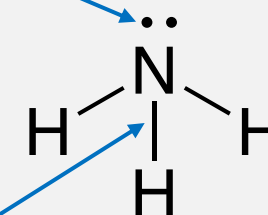
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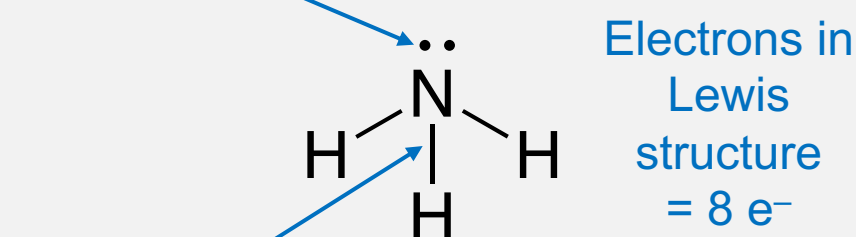
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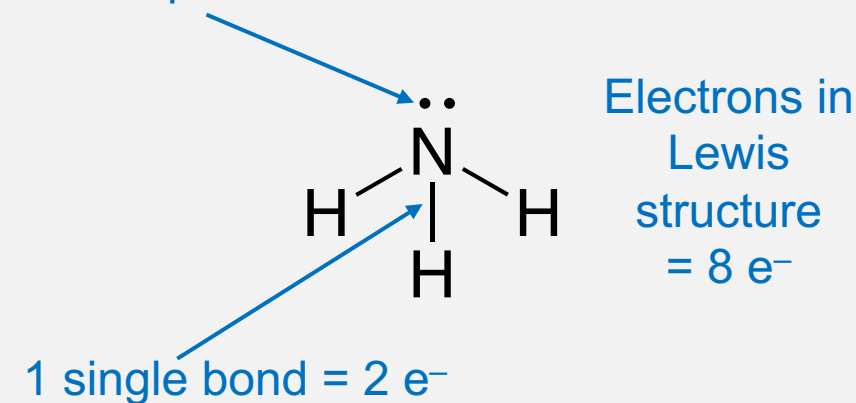
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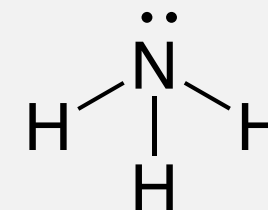
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*Make sure  
sum of FC  
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$$\text{FC} = (\text{valence } e^-) - (\# \text{ bonds}) - (\# \text{ lone } e^-)$$

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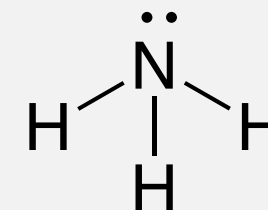
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H	1 e <sup>-</sup>
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$$\text{FC (N)} = (5 \text{ e}^-) - (3 \text{ bonds}) - (2 \text{ lone e}^-) = 0$$

$$\text{FC (H)} = (1 \text{ e}^-) - (1 \text{ bonds}) - (0 \text{ lone e}^-) = 0$$



*Make sure  
sum of FC  
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$$\text{FC} = (\text{valence e}^-) - (\# \text{ bonds}) - (\# \text{ lone e}^-)$$

## LEWIS STRUCTURES “RULES”

### Guided Example

Carbon Monoxide (CO)

## LEWIS STRUCTURES “RULES”

1. The total number of electrons in the Lewis structure must equal the total number of valence electrons.

### Guided Example

Carbon Monoxide (CO)

<b>Atom</b>	<b>Valence</b>
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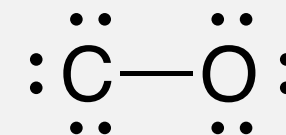
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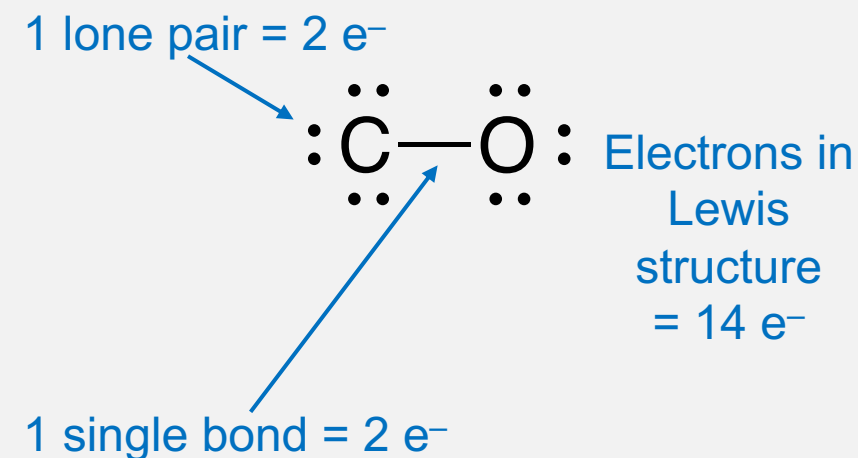
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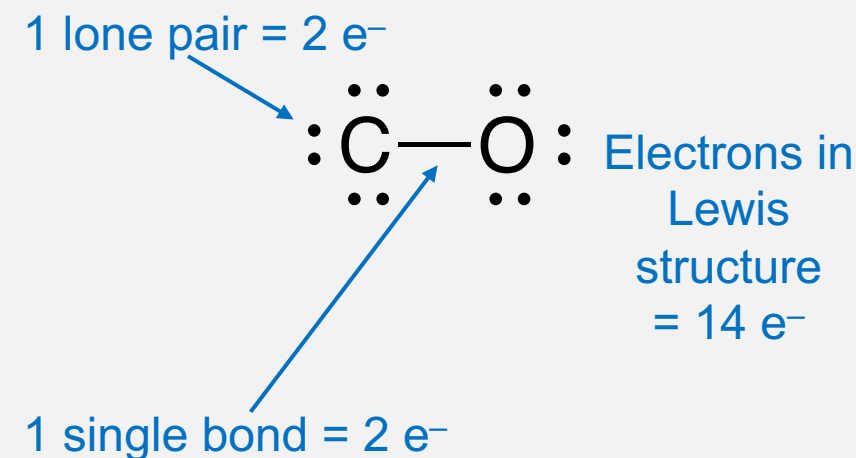
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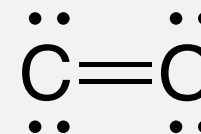
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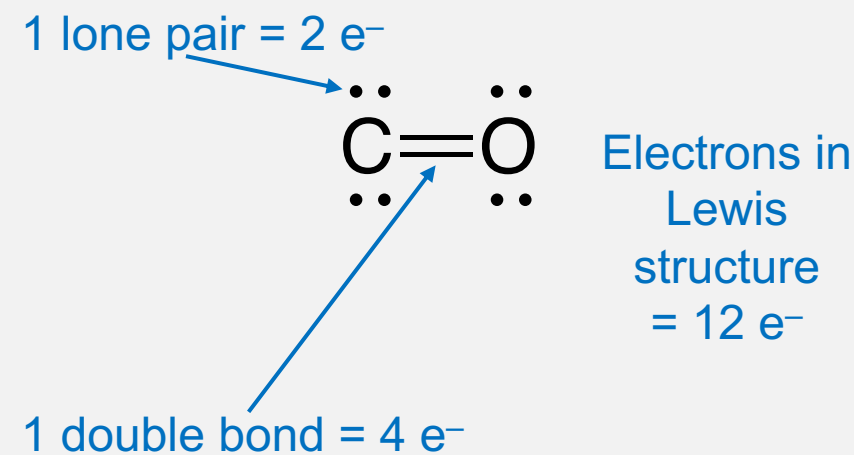
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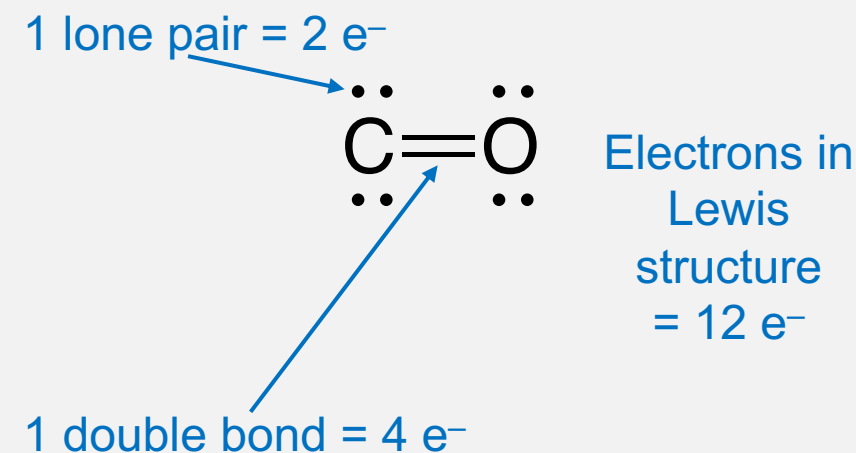
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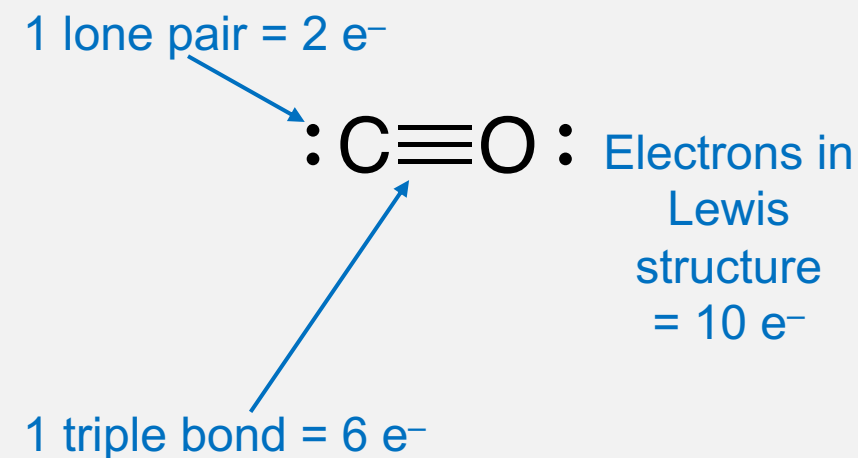
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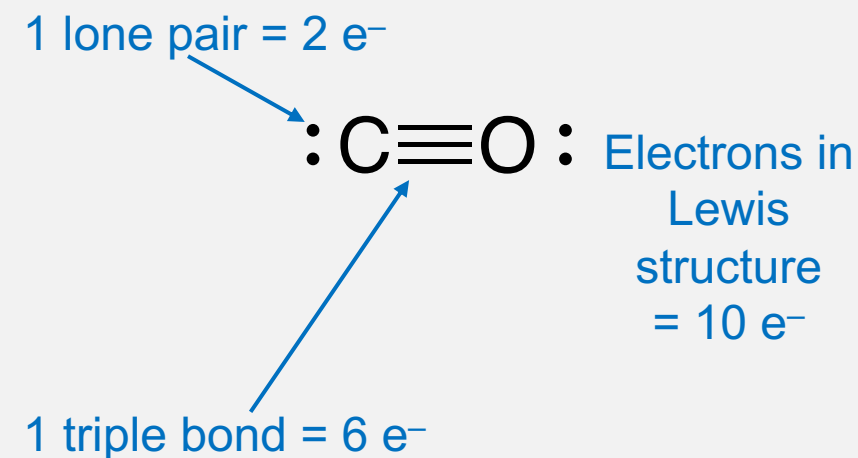
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$$\text{FC (C)} = (4 \text{ e}^-) - (3 \text{ bonds}) - (2 \text{ lone e}^-) = -1$$

$$\text{FC (O)} = (6 \text{ e}^-) - (3 \text{ bonds}) - (2 \text{ lone e}^-) = +1$$



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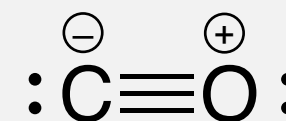
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### Guided Example

Sulfur Dioxide (SO<sub>2</sub>)

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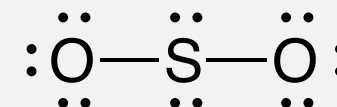
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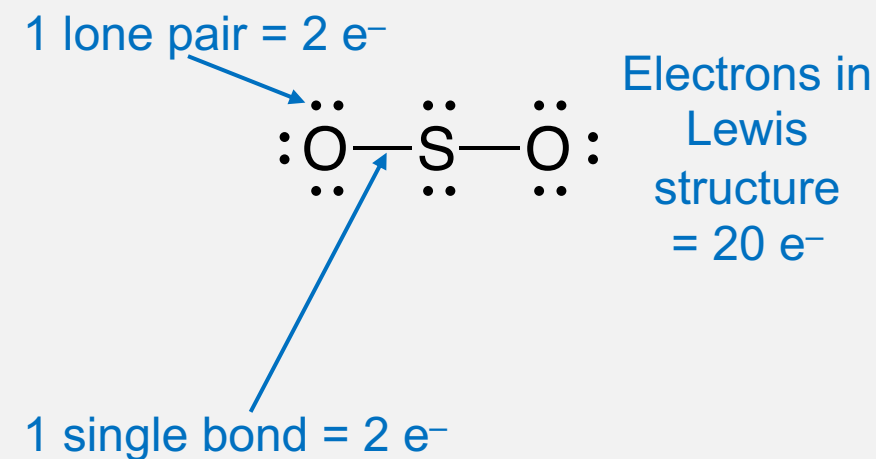
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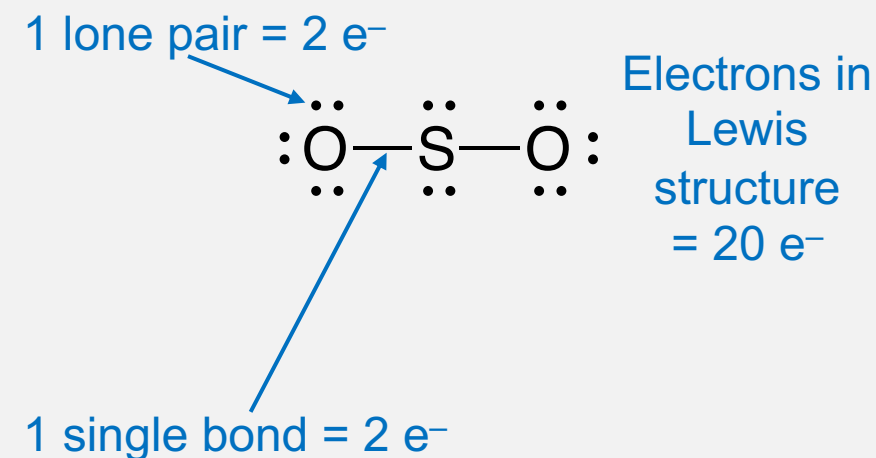
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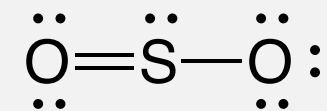
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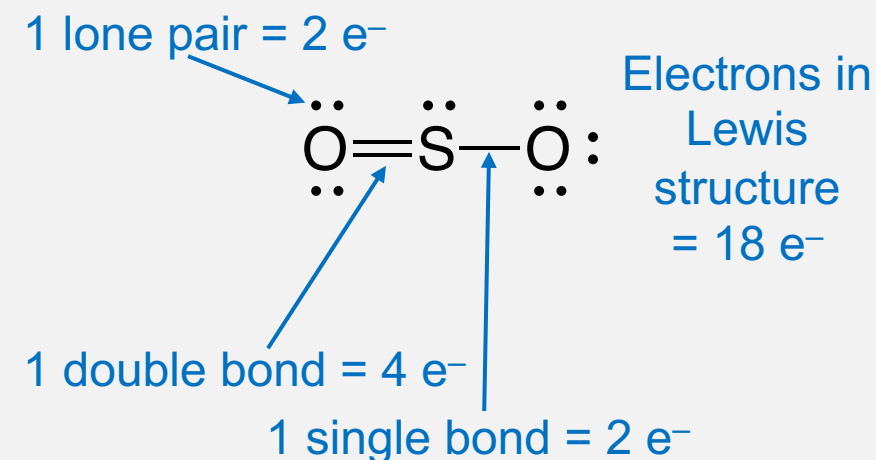
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Sulfur Dioxide (SO<sub>2</sub>)

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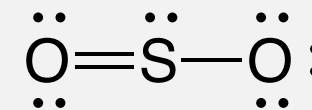
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Electrons in  
Lewis  
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= 18 e<sup>-</sup>

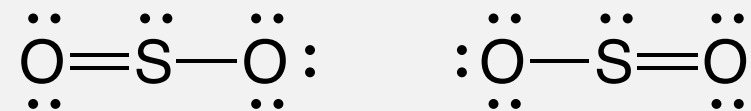
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But we arbitrarily chose the left O to be double-bonded...

We could've also chosen the right O.

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**RESONANCE:** multiple valid Lewis structures

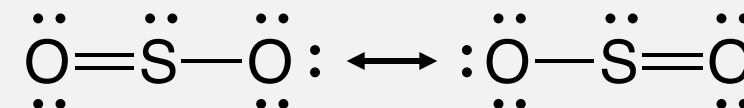
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The real structure is an average of the multiple resonance structures.

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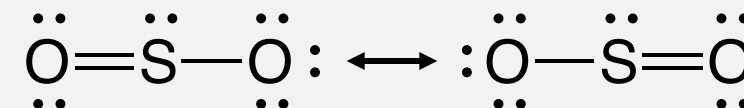
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$$\text{FC (S)} = (6 \text{ e}^-) - (3 \text{ bonds}) - (2 \text{ lone e}^-) = +1$$

$$\text{FC (O)} = (6 \text{ e}^-) - (2 \text{ bonds}) - (4 \text{ lone e}^-) = 0$$

$$\text{FC (O)} = (6 \text{ e}^-) - (1 \text{ bonds}) - (6 \text{ lone e}^-) = -1$$



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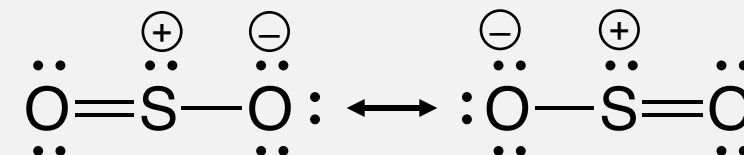
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## LEWIS STRUCTURES “RULES”

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### Guided Example

Cyanide (CN<sup>-</sup>)

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N	5 e <sup>-</sup>
charge	1 e <sup>-</sup>
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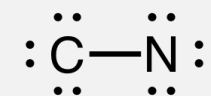
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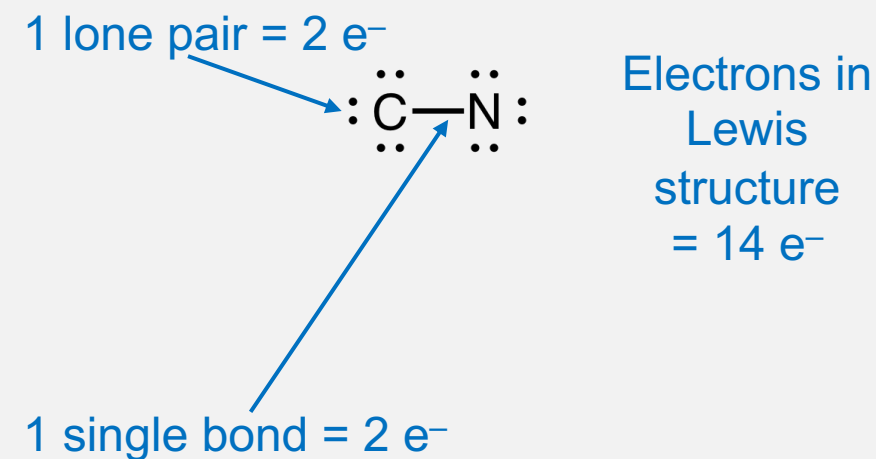
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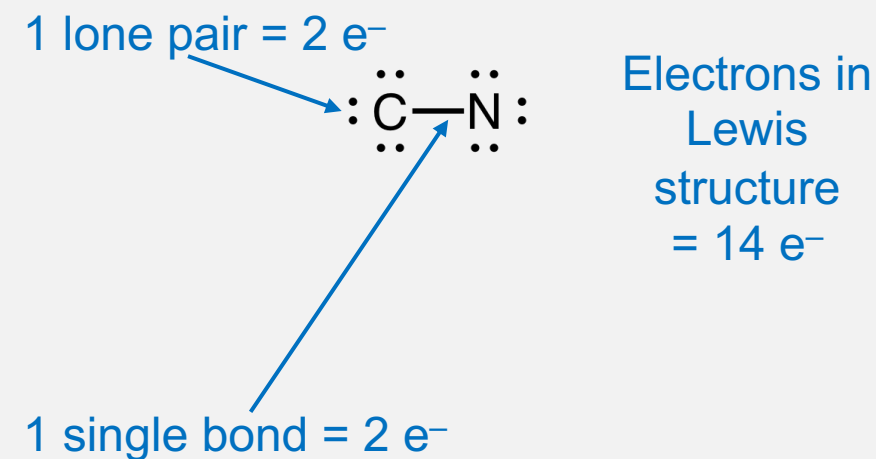
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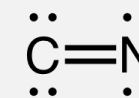
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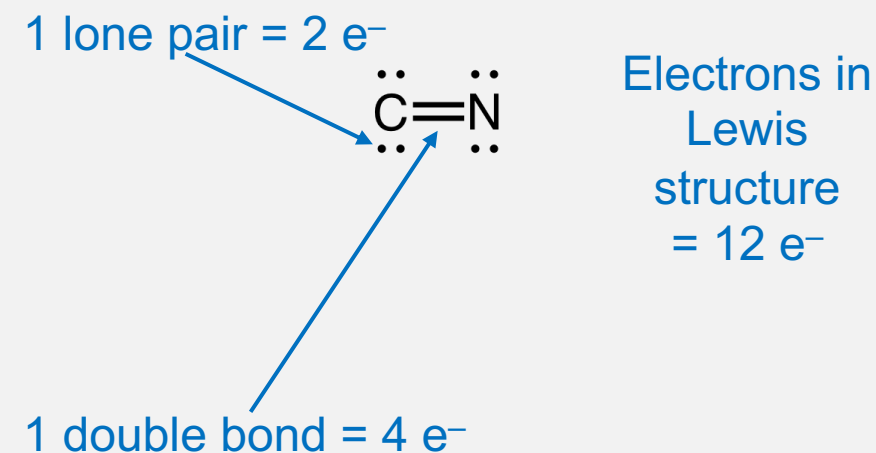
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1 lone pair = 2 e<sup>-</sup>



1 triple bond = 6 e<sup>-</sup>

Electrons in  
Lewis  
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$$\text{FC (C)} = (4 \text{ e}^-) - (3 \text{ bonds}) - (2 \text{ lone e}^-) = -1$$

$$\text{FC (N)} = (5 \text{ e}^-) - (3 \text{ bonds}) - (2 \text{ lone e}^-) = 0$$



*Make sure  
sum of FC  
equals total  
charge.*

$$\text{FC} = (\text{valence e}^-) - (\# \text{ bonds}) - (\# \text{ lone e}^-)$$

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Generally, it's best to put negative formal charges on more EN atoms and positive formal charges on less EN atoms.

*But you can't do that in this example.*