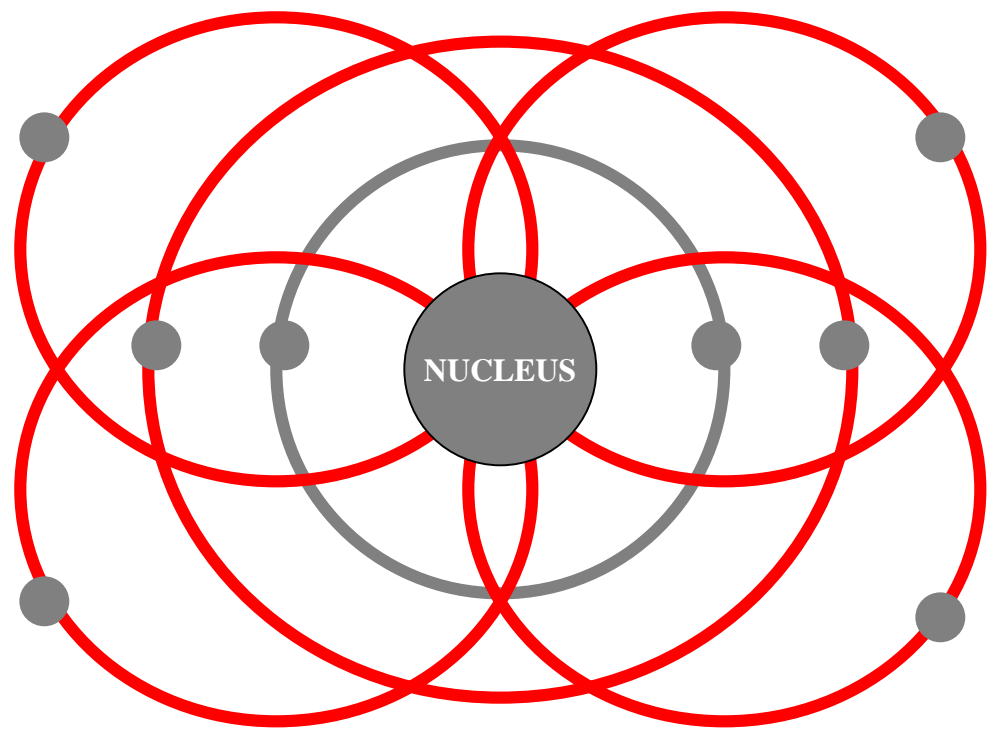


# 2<sup>ND</sup> ENERGY SHELL = 4 ORBITALS



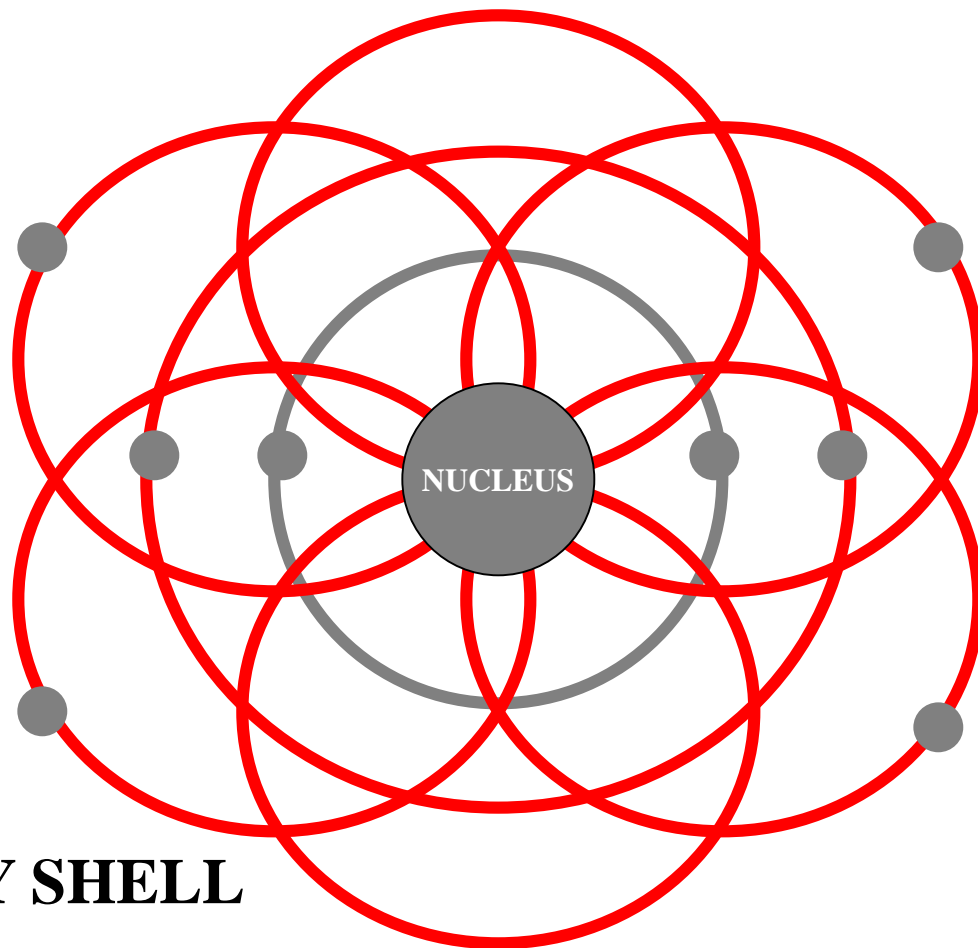
— = 1ST ENERGY SHELL

— = 2ND ENERGY SHELL

● = e-

## TYPICAL ATOM

# 2<sup>ND</sup> ENERGY SHELL = 4 ORBITALS



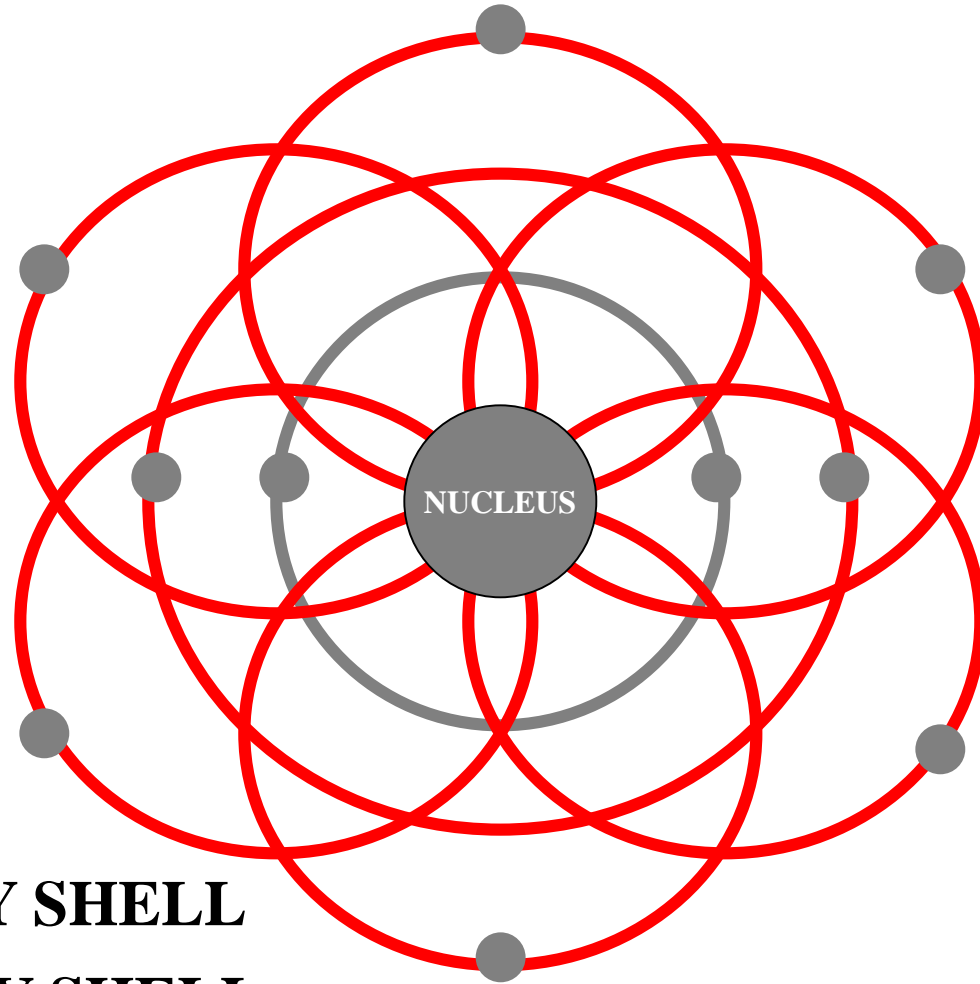
— = 1ST ENERGY SHELL

— = 2ND ENERGY SHELL

● = E-

## TYPICAL ATOM

# 2<sup>ND</sup> ENERGY SHELL = 4 ORBITALS



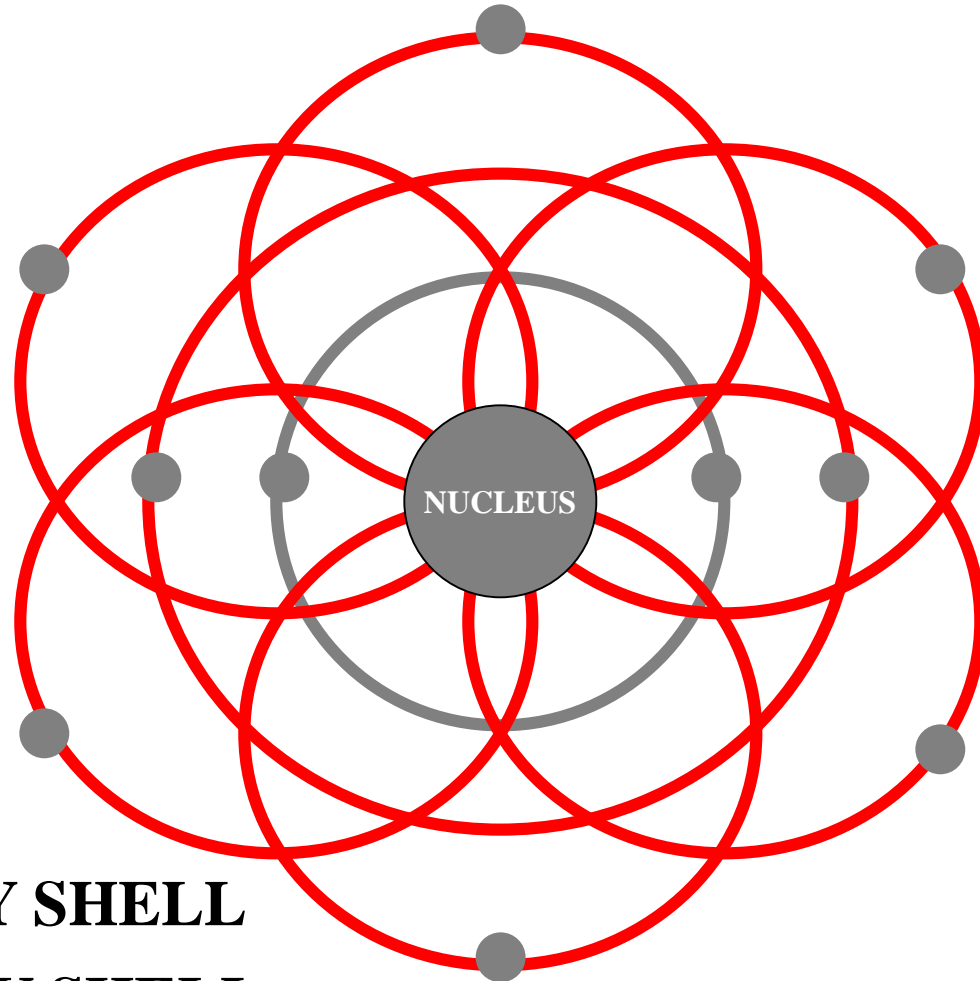
— = 1ST ENERGY SHELL

— = 2ND ENERGY SHELL

● = e-

## TYPICAL ATOM

# 2<sup>ND</sup> ENERGY SHELL = 4 ORBITALS



— = 1ST ENERGY SHELL

— = 2ND ENERGY SHELL

● = e-

2<sup>ND</sup> E.S. = 8 e- TOTAL

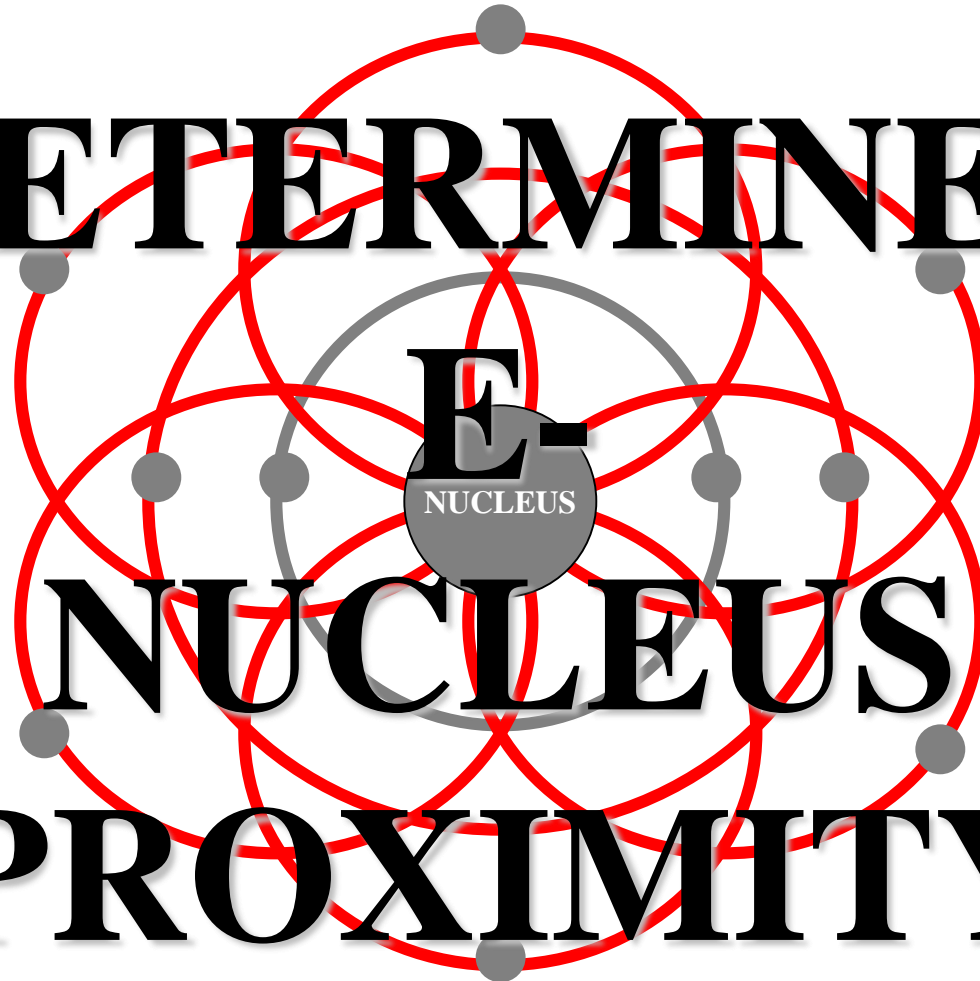
# **ELECTRON BOND POTENTIAL**

**ELECTRON  
BOND  
POTENTIAL**



**E- BOND POTENTIAL**

**DETERMINED**

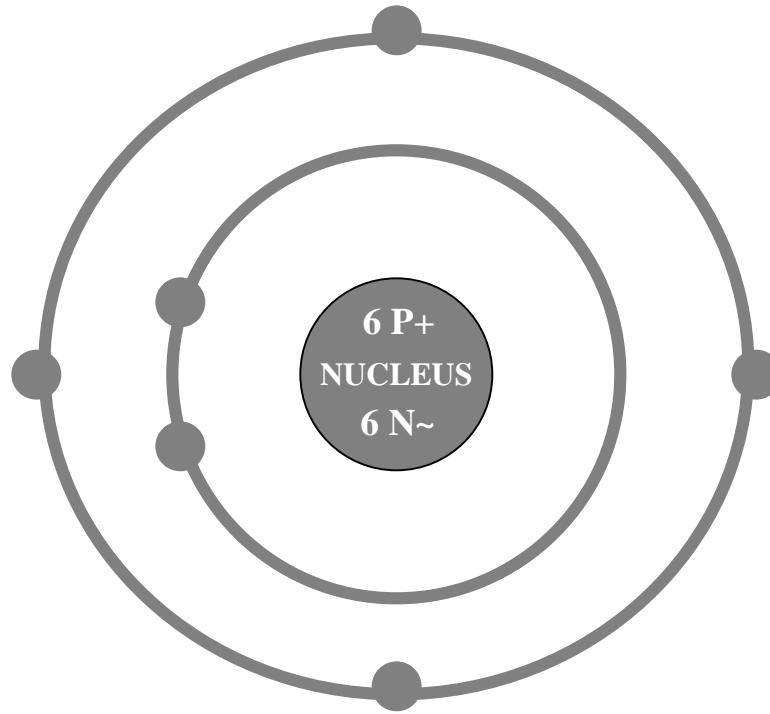


**NUCLEUS**

**PROXIMITY**

**E- BOND POTENTIAL**

# E- BOND POTENTIAL

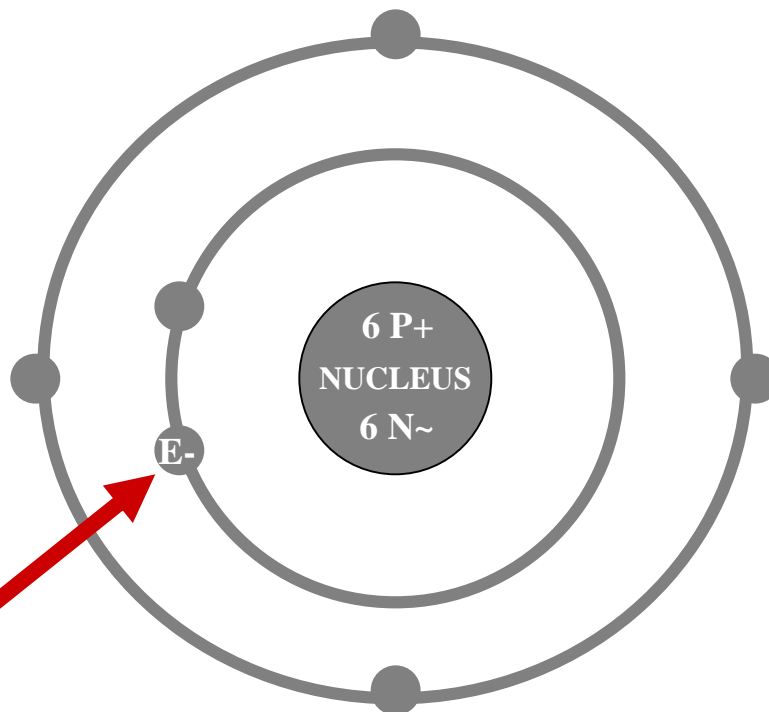


● = E-

## CARBON ATOM



# E- BOND POTENTIAL

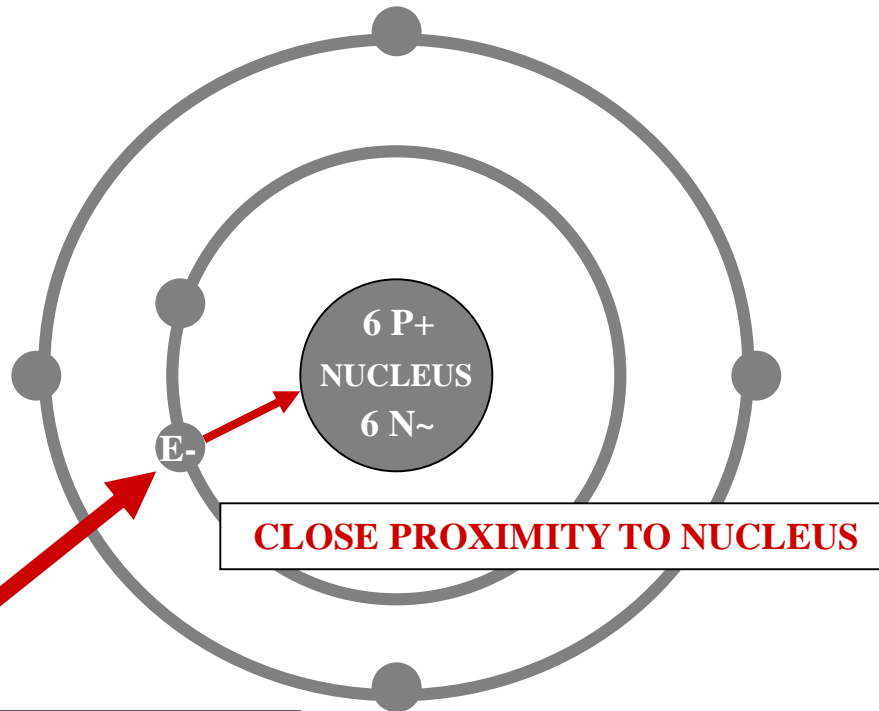


**E- W/ INNER PROXIMITY**

● = E-

**CARBON ATOM**

# E- BOND POTENTIAL



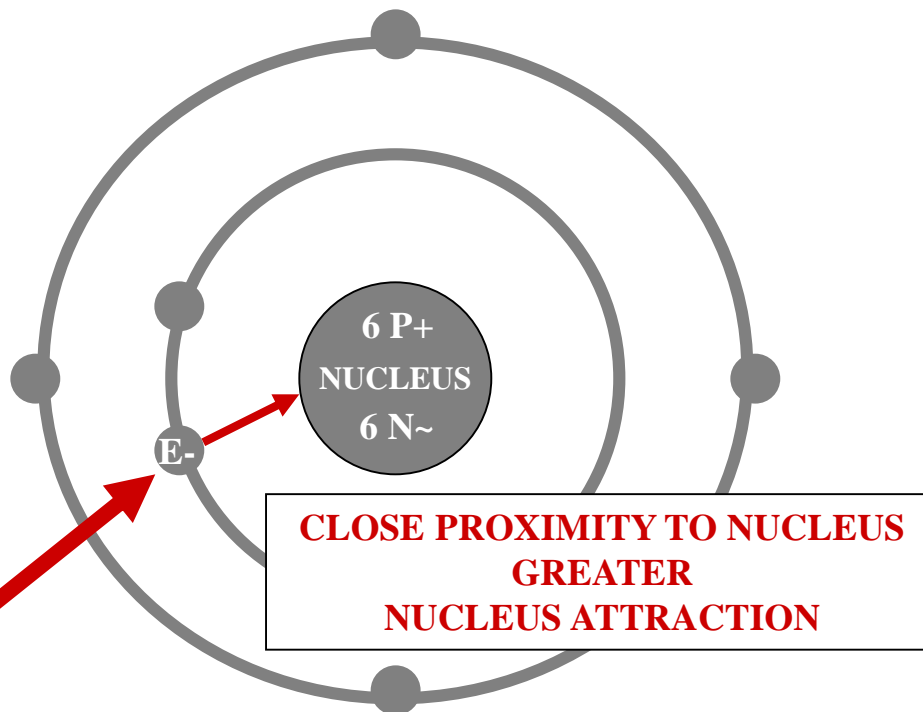
**E- W/ INNER PROXIMITY**

**CLOSE PROXIMITY TO NUCLEUS**

● = E-

**CARBON ATOM**

# E- BOND POTENTIAL

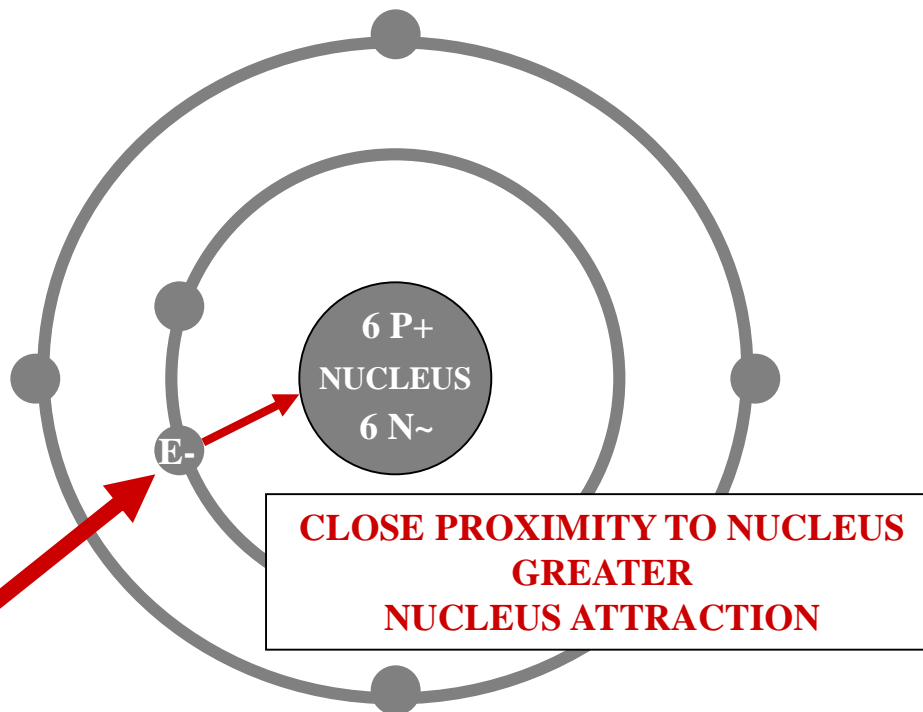


**E- W/ INNER PROXIMITY**

● = E-

**CARBON ATOM**

# E- BOND POTENTIAL



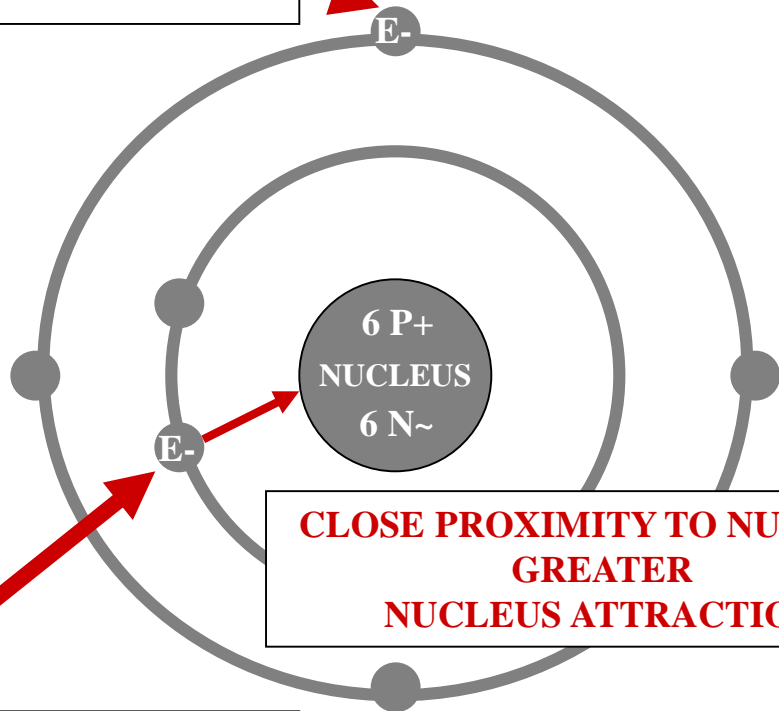
**LOW BOND POTENTIAL**

● = E-

**CARBON ATOM**

# E- BOND POTENTIAL

**E- W/ OUTER PROXIMITY**



**CLOSE PROXIMITY TO NUCLEUS  
GREATER  
NUCLEUS ATTRACTION**

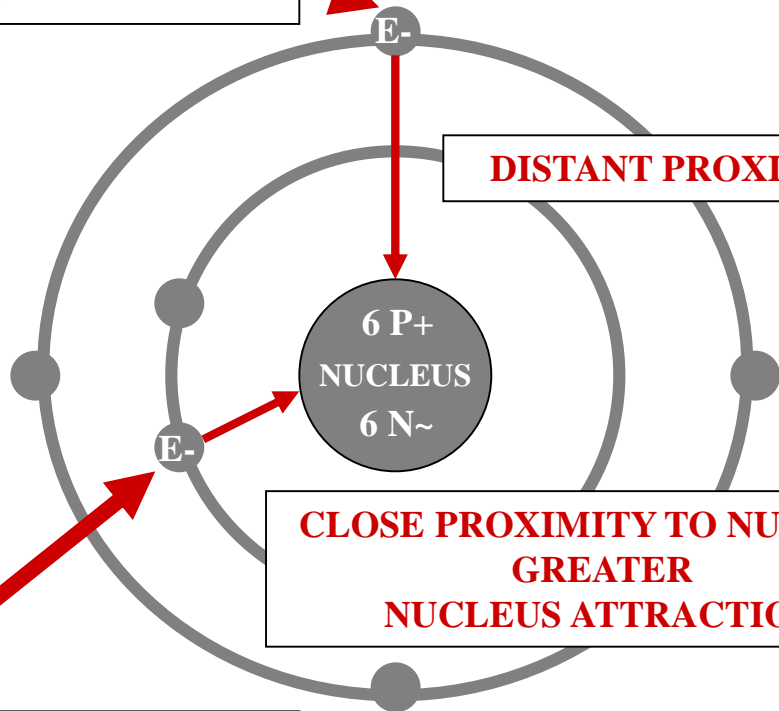
**LOW BOND POTENTIAL**

● = E-

## CARBON ATOM

# E- BOND POTENTIAL

**E- W/ OUTER PROXIMITY**



**DISTANT PROXIMITY TO NUCLEUS**

**CLOSE PROXIMITY TO NUCLEUS  
GREATER  
NUCLEUS ATTRACTION**

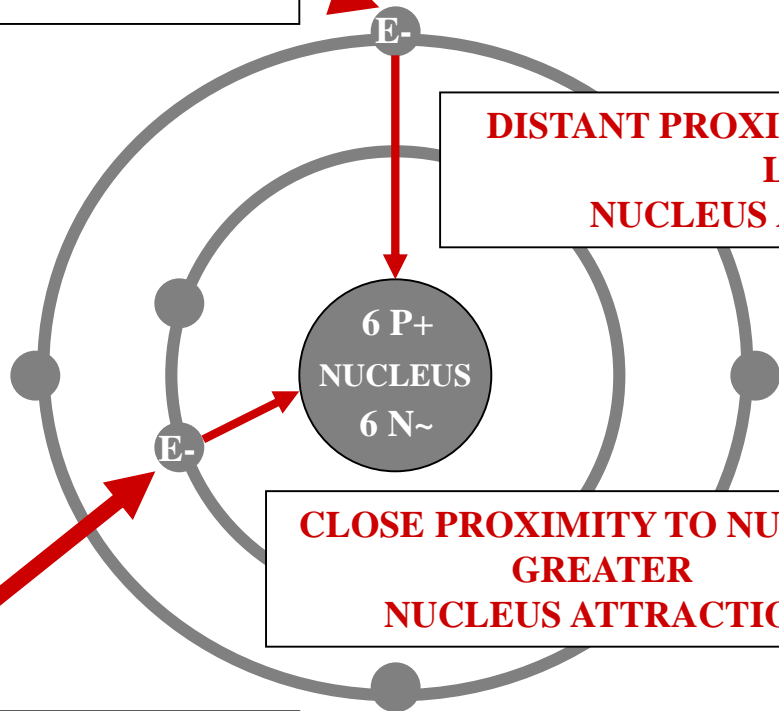
**LOW BOND POTENTIAL**

● = E-

## CARBON ATOM

# E- BOND POTENTIAL

**E- W/ OUTER PROXIMITY**



**DISTANT PROXIMITY TO NUCLEUS  
LESS  
NUCLEUS ATTRACTION**

**CLOSE PROXIMITY TO NUCLEUS  
GREATER  
NUCLEUS ATTRACTION**

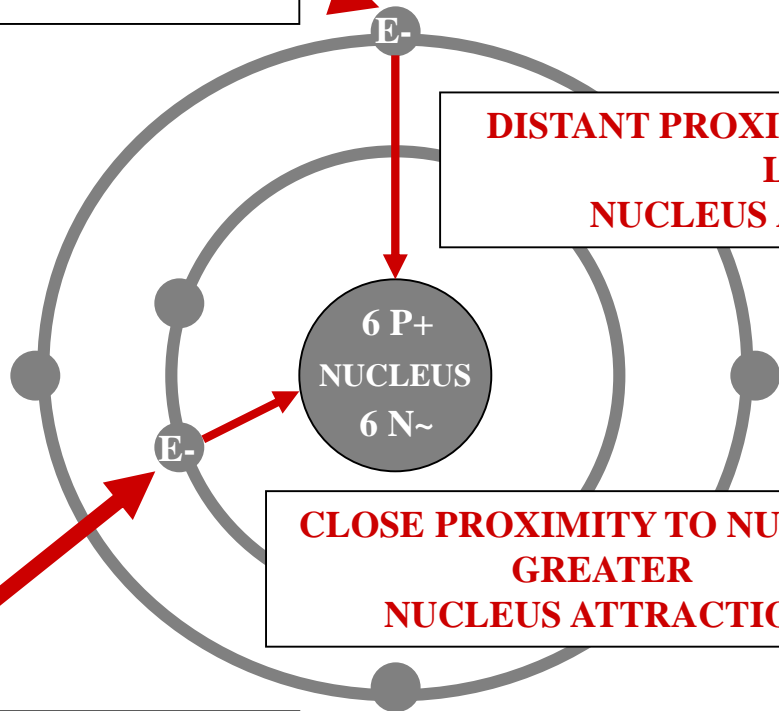
**LOW BOND POTENTIAL**

● = E-

## CARBON ATOM

# E- BOND POTENTIAL

**HIGH BOND POTENTIAL**



**DISTANT PROXIMITY TO NUCLEUS  
LESS  
NUCLEUS ATTRACTION**

**CLOSE PROXIMITY TO NUCLEUS  
GREATER  
NUCLEUS ATTRACTION**

**LOW BOND POTENTIAL**

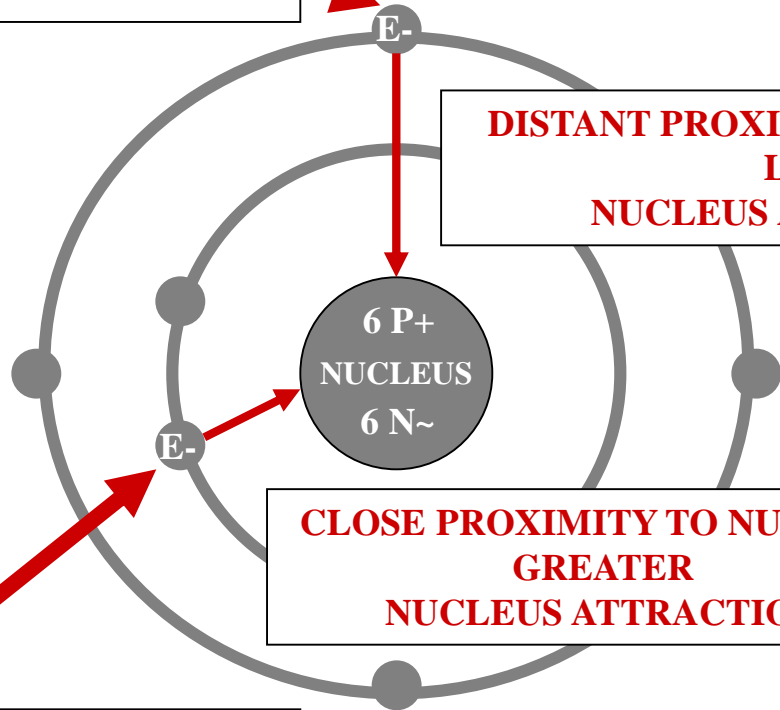
● = E-

## CARBON ATOM



# E- BOND POTENTIAL

**FORM CHEM BONDS**



**DISTANT PROXIMITY TO NUCLEUS  
LESS  
NUCLEUS ATTRACTION**

**CLOSE PROXIMITY TO NUCLEUS  
GREATER  
NUCLEUS ATTRACTION**

**LOW BOND POTENTIAL**

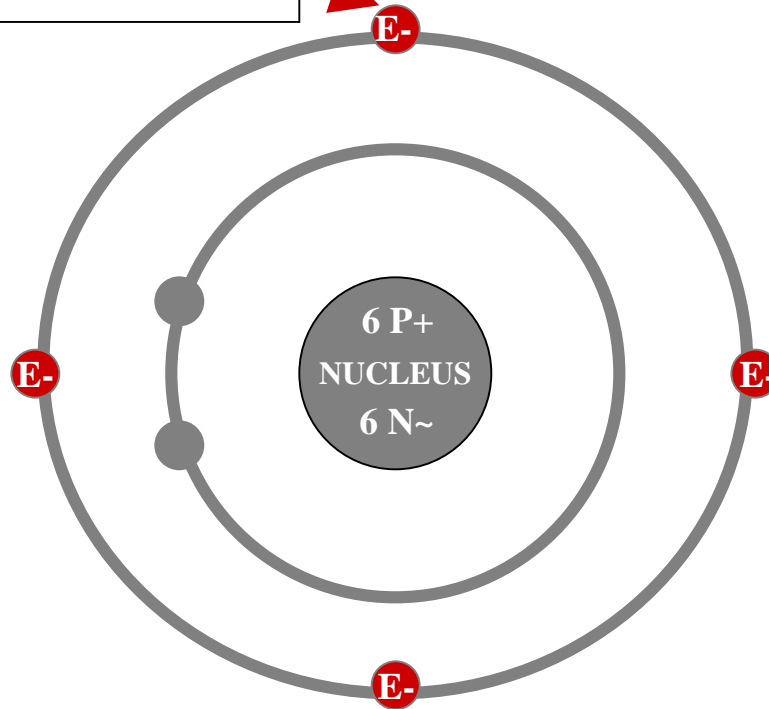
● = E-

**CARBON ATOM**



# VALENCE ELECTRONS

VALENCE ELECTRONS



● = E<sup>-</sup>

## CARBON ATOM



**CHEMICAL BONDS**

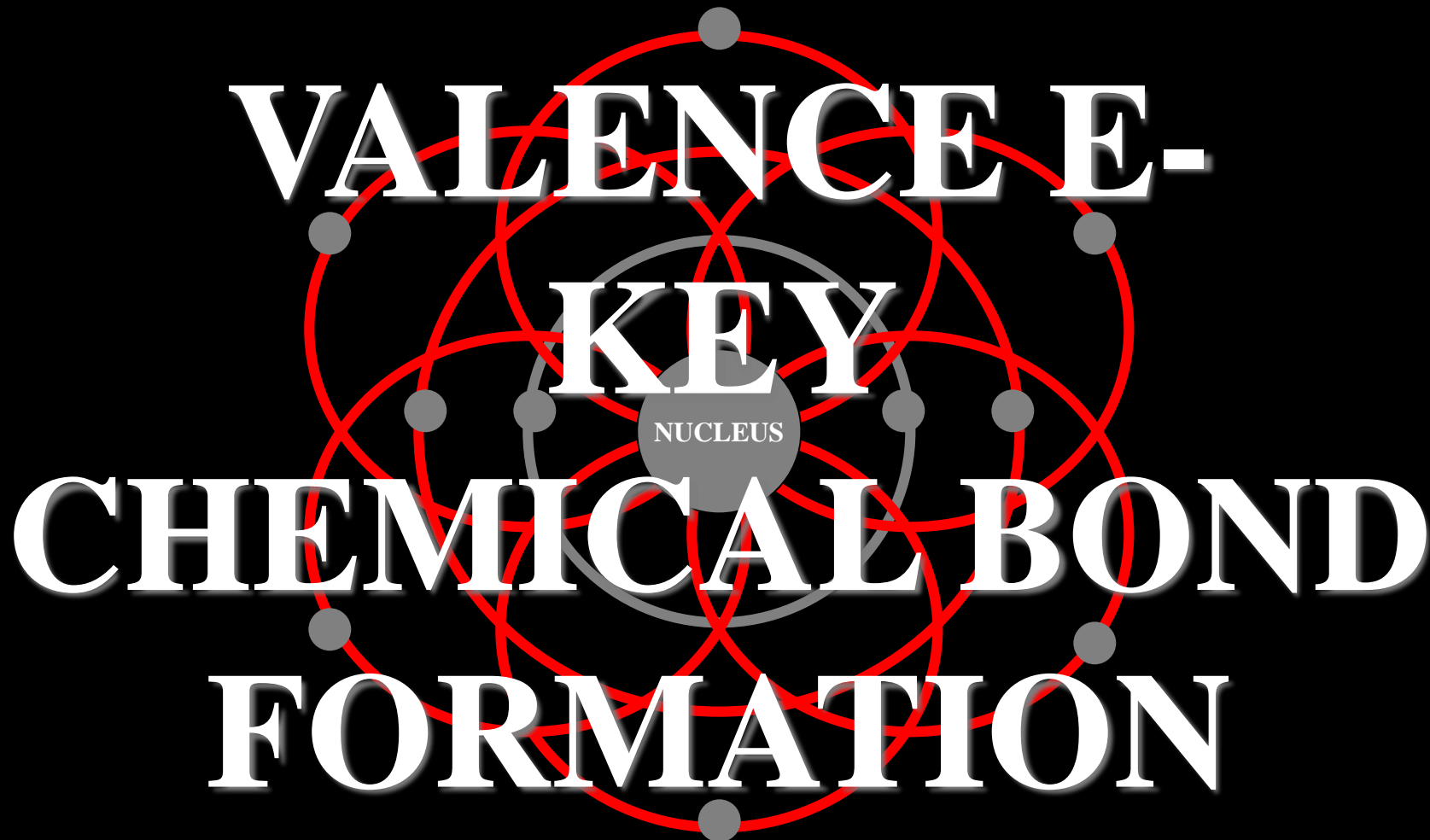
**VALENCE E-  
KEY**

NUCLEUS

**CHEMICAL BOND**

**FORMATION**

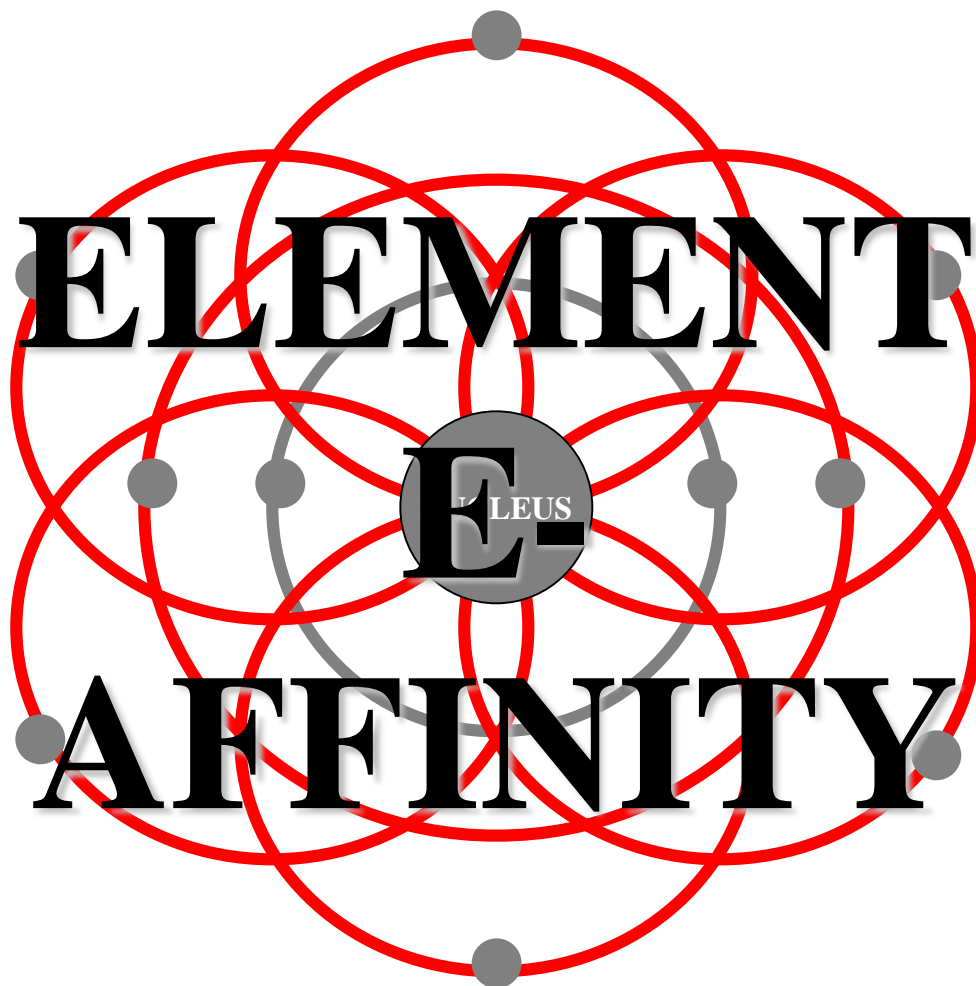
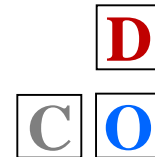
**CHEMICAL BONDS**



# VALENCE E- ELECTRO-NEGATIVITY

**ELECTRO-NEGATIVITY**

# ELECTRO-NEGATIVITY



# ELECTRO-NEGATIVITY

# ELECTRO-NEGATIVITY VALUES

ATOM: CARBON

ATOM: OXYGEN

**DIFFERENT ELEMENTS**

# ELECTRO-NEGATIVITY VALUES

**ATOM: CARBON****Symbol: C****Atomic Number: 6****Atomic Mass: 12.0107 amu****Melting Point: 3500.0 °C (3773.15 °K, 6332.0 °F)****Boiling Point: 4827.0 °C (5100.15 °K, 8720.6 °F)****Number of Protons/Electrons: 6****Number of Neutrons: 6****Electro-Negativity: 2.5****Classification: Non-metal****Crystal Structure: Hexagonal****Density @ 293 K: 2.62 g/cm<sup>3</sup>****Color: May be black****ATOM: OXYGEN****Symbol: O****Atomic Number: 8****Atomic Mass: 15.9994 amu****Melting Point: -218.4 °C (54.750008 °K, -361.12 °F)****Boiling Point: -183.0 °C (90.15 °K, -297.4 °F)****Number of Protons/Electrons: 8****Number of Neutrons: 8****Electro-Negativity: 3.5****Classification: Non-metal****Crystal Structure: Cubic****Density @ 293 K: 1.429 g/cm<sup>3</sup>****Color: colorless**

**DIFFERENT ELEMENTS**  
**DIFFERENT ELEMENTAL PROPERTIES**



# ELECTRO-NEGATIVITY VALUES



## ATOM: CARBON

Symbol: C

Atomic Number: 6

Atomic Mass: 12.0107 amu

Melting Point: 3500.0 °C (3773.15 °K, 6332.0 °F)

Boiling Point: 4827.0 °C (5100.15 °K, 8720.6 °F)

Number of Protons/Electrons: 6

Number of Neutrons: 6

Electro-Negativity: 2.5

Classification: Non-metal

Crystal Structure: Hexagonal

Density @ 293 K: 2.62 g/cm<sup>3</sup>

Color: May be black

## ATOM: OXYGEN

Symbol: O

Atomic Number: 8

Atomic Mass: 15.9994 amu

Melting Point: -218.4 °C (54.750008 °K, -361.12 °F)

Boiling Point: -183.0 °C (90.15 °K, -297.4 °F)

Number of Protons/Electrons: 8

Number of Neutrons: 8

Electro-Negativity: 3.5

Classification: Non-metal

Crystal Structure: Cubic

Density @ 293 K: 1.429 g/cm<sup>3</sup>

Color: colorless

**DIFFERENT ELEMENTS  
DIFFERENT ELECTRO-NEGATIVITY  
VALUES**

# PERIODIC TABLE

## ELECTRO-NEGATIVITY VALUES



**HIGHER E-N VALUE**

H 2.1																	He
Li 1.0	Be 1.5											B 2.0	C 2.5	N 3.0	O 3.5	F 4.0	Ne
Na 0.9	Mg 1.2											Al 1.5	Si 1.8	P 2.1	S 2.5	Cl 3.0	Ar
K 0.8	Ca 1.0	Sc 1.3	Ti 1.5	V 1.6	Cr 1.6	Mn 1.5	Fe 1.8	Co 1.8	Ni 1.8	Cu 1.9	Zn 1.6	Ga 1.6	Ge 1.8	As 2.0	Se 2.4	Br 2.8	Kr 3.0
Rb 0.8	Sr 1.0	Y 1.2	Zr 1.4	Nb 1.6	Mo 1.8	Tc 1.9	Ru 2.2	Rh 2.2	Pd 2.2	Ag 1.9	Cd 1.7	In 1.7	Sn 1.8	Sb 1.9	Te 2.1	I 2.5	Xe 2.6
Cs 0.7	Ba 0.9	La 1.1	Hf 1.3	Ta 1.5	W 1.7	Re 1.9	Os 2.2	Ir 2.2	Pt 2.2	Au 2.4	Hg 1.9	Tl 1.8	Pb 1.8	Bi 1.9	Po 2.0	At 2.2	Rn 2.4
Fr 0.7	Ra 0.7	Ac 1.1	Unq	Unp	Unh	Uns	Uno	Une									

Ce 1.1	Pr 1.1	Nd 1.1	Pm 1.1	Sm 1.1	Eu 1.1	Gd 1.1	Tb 1.1	Dy 1.1	Ho 1.1	Er 1.1	Tm 1.1	Yb 1.1	Lu 1.2
Th 1.3	Pa 1.5	U 1.7	Np 1.3	Pu 1.3	Am 1.3	Cm 1.3	Bk 1.3	Cf 1.3	Es 1.3	Fm 1.3	Md 1.3	No 1.3	Lr

# PERIODIC TABLE

## ELECTRO-NEGATIVITY VALUES

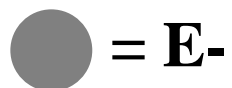
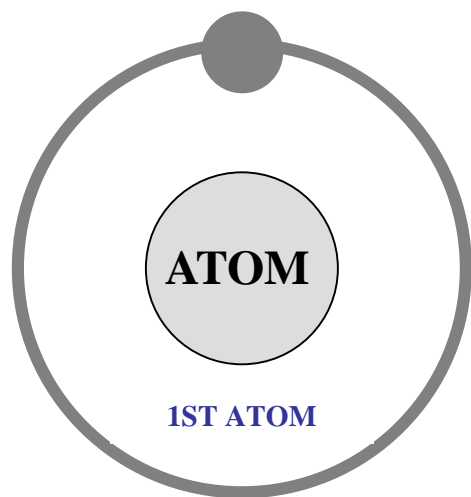
**HIGHER E-N VALUE  
THE GREATER THE  
E- AFFINITY**

H 2.1																	He
Li 1.0	Be 1.5											B 2.0	C 2.5	N 3.0	O 3.5	F 4.0	Ne
Na 0.9	Mg 1.2											Al 1.5	Si 1.8	P 2.1	S 2.5	Cl 3.0	Ar
K 0.8	Ca 1.0	Sc 1.3	Ti 1.5	V 1.6	Cr 1.6	Mn 1.5	Fe 1.8	Co 1.8	Ni 1.8	Cu 1.9	Zn 1.6	Ga 1.6	Ge 1.8	As 2.0	Se 2.4	Br 2.8	Kr 3.0
Rb 0.8	Sr 1.0	Y 1.2	Zr 1.4	Nb 1.6	Mo 1.8	Tc 1.9	Ru 2.2	Rh 2.2	Pd 2.2	Ag 1.9	Cd 1.7	In 1.7	Sn 1.8	Sb 1.9	Te 2.1	I 2.5	Xe 2.6
Cs 0.7	Ba 0.9	La 1.1	Hf 1.3	Ta 1.5	W 1.7	Re 1.9	Os 2.2	Ir 2.2	Pt 2.2	Au 2.4	Hg 1.9	Tl 1.8	Pb 1.8	Bi 1.9	Po 2.0	At 2.2	Rn 2.4
Fr 0.7	Ra 0.7	Ac 1.1	Unq	Unp	Unh	Uns	Uno	Une									

Ce 1.1	Pr 1.1	Nd 1.1	Pm 1.1	Sm 1.1	Eu 1.1	Gd 1.1	Tb 1.1	Dy 1.1	Ho 1.1	Er 1.1	Tm 1.1	Yb 1.1	Lu 1.2
Th 1.3	Pa 1.5	U 1.7	Np 1.3	Pu 1.3	Am 1.3	Cm 1.3	Bk 1.3	Cf 1.3	Es 1.3	Fm 1.3	Md 1.3	No 1.3	Lr

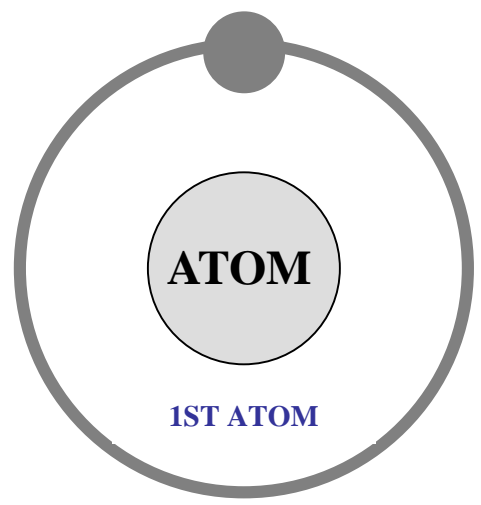
# **ELECTRO-NEGATIVITY APPLIED**

# ELECTRO-NEGATIVITY VALUES



**CHEMICAL BOND**

# ELECTRO-NEGATIVITY VALUES

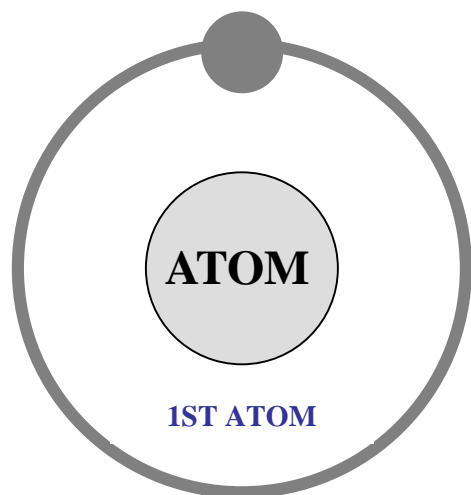


**LOW E- NEG  
VALUE**

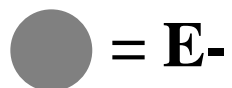
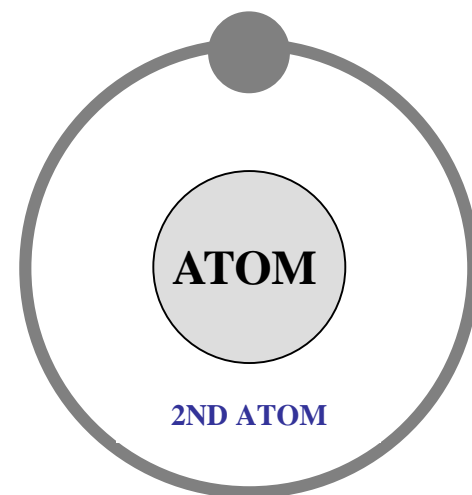
● = E-

## CHEMICAL BOND

# ELECTRO-NEGATIVITY VALUES

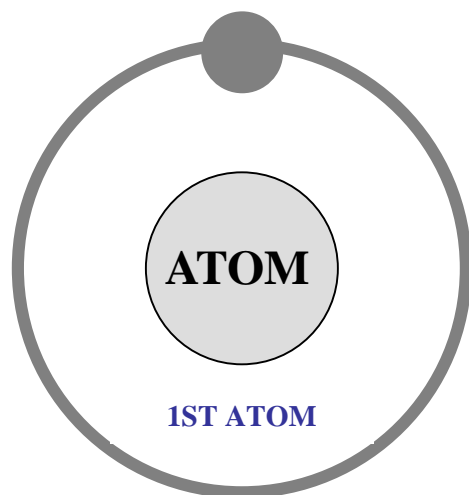


**LOW E- NEG  
VALUE**

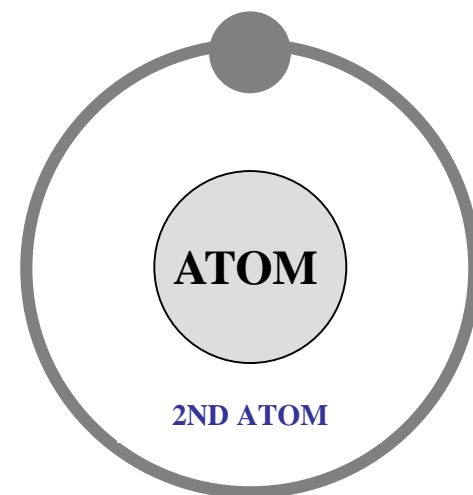


**CHEMICAL BOND**

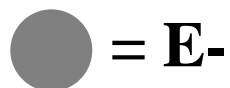
# ELECTRO-NEGATIVITY VALUES



**LOW E- NEG  
VALUE**



**HIGH E- NEG  
VALUE**

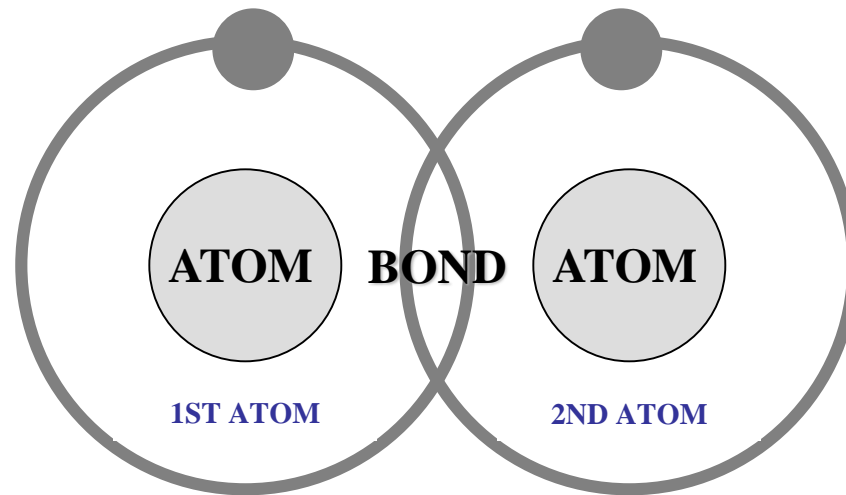


**CHEMICAL BOND**





# ELECTRO-NEGATIVITY VALUES



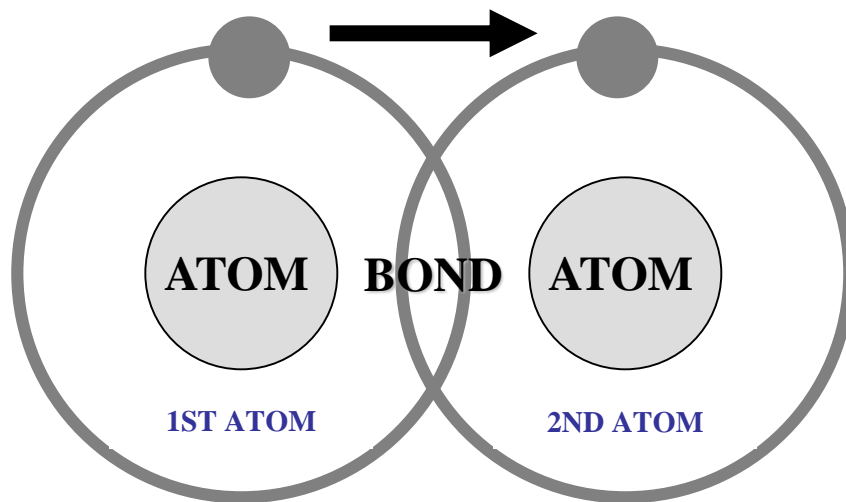
**LOW E- NEG  
VALUE**

**HIGH E- NEG  
VALUE**

● = E-

## CHEMICAL BOND

# ELECTRO-NEGATIVITY VALUES



**LOW E- NEG  
VALUE**

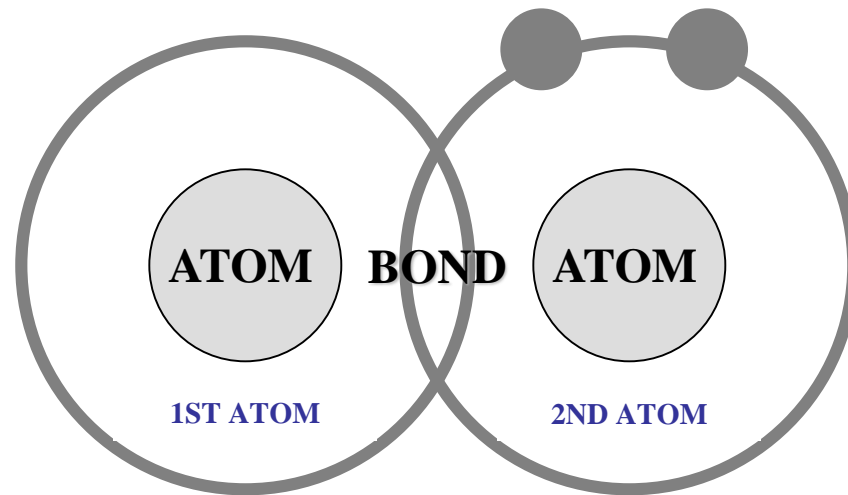
**HIGH E- NEG  
VALUE**

● = E-

## CHEMICAL BOND



# ELECTRO-NEGATIVITY VALUES



**LOW E- NEG  
VALUE**

**HIGH E- NEG  
VALUE**

● = E-

## CHEMICAL BOND

# CHEMICAL BONDS



E- MOVE FROM

ATOMS W/ LOW E- NEG

NUCLEUS



ATOMS W/ HIGH E- NEG

CHEMICAL BONDS

CHEMICAL BONDS



ELECTRO

NEGATIVITY

KEY

NUCLEUS

CHEMICAL BOND

FORMATION

CHEMICAL BONDS

# CHEMICAL BOND TYPES

# **CHEMICAL BOND TYPES**

**COVALENT BONDS**

# **CHEMICAL BOND TYPES**

# **CHEMICAL BOND TYPES**

**COVALENT BONDS**

**IONIC BONDS**

# **CHEMICAL BOND TYPES**



# CHEMICAL BOND TYPES



COVALENT BONDS  
IONIC BONDS  
HYDROGEN BONDS

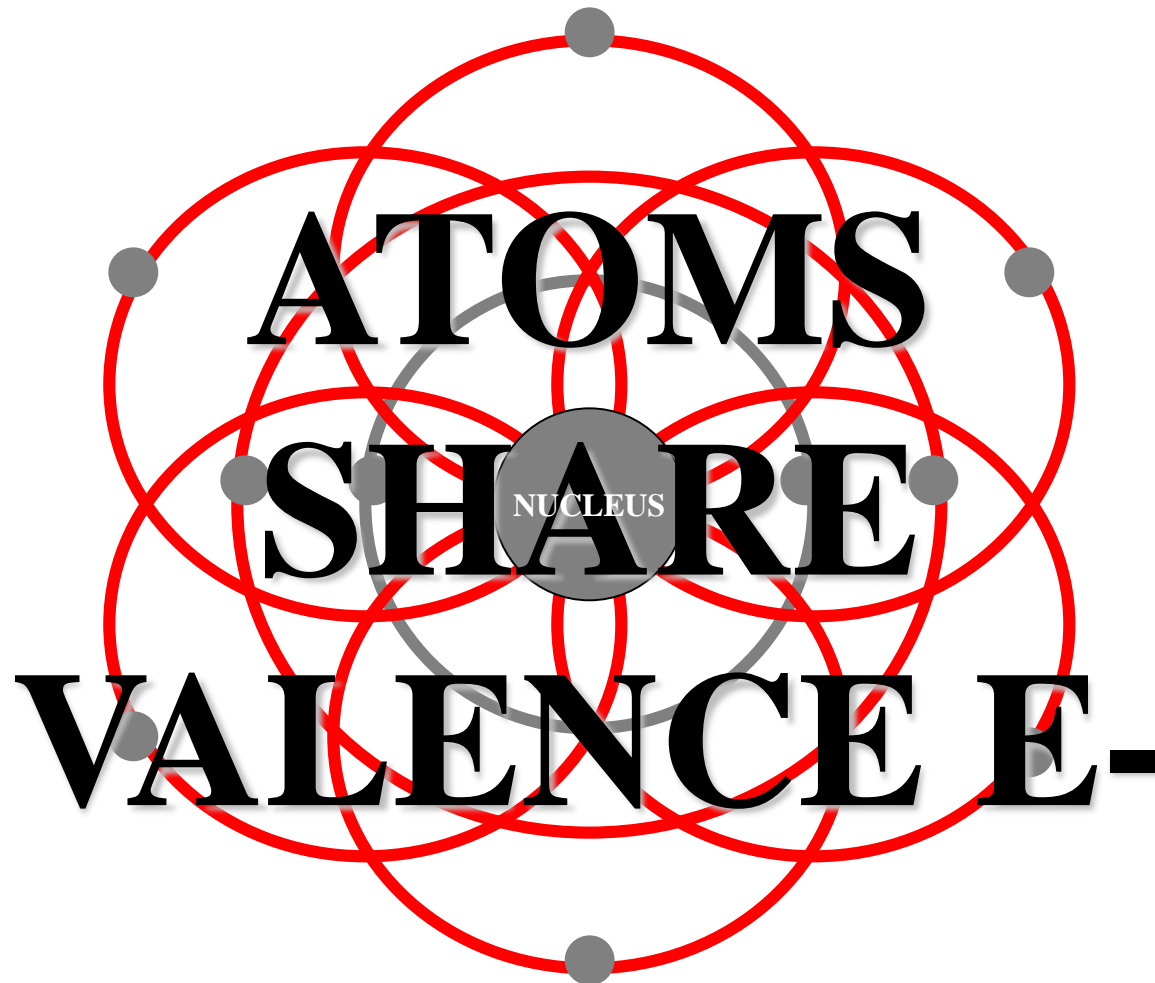
CHEMICAL BOND  
TYPES

# COVALENT BONDS

# COVALENT BONDS



# COVALENT BONDS



# COVALENT BONDS

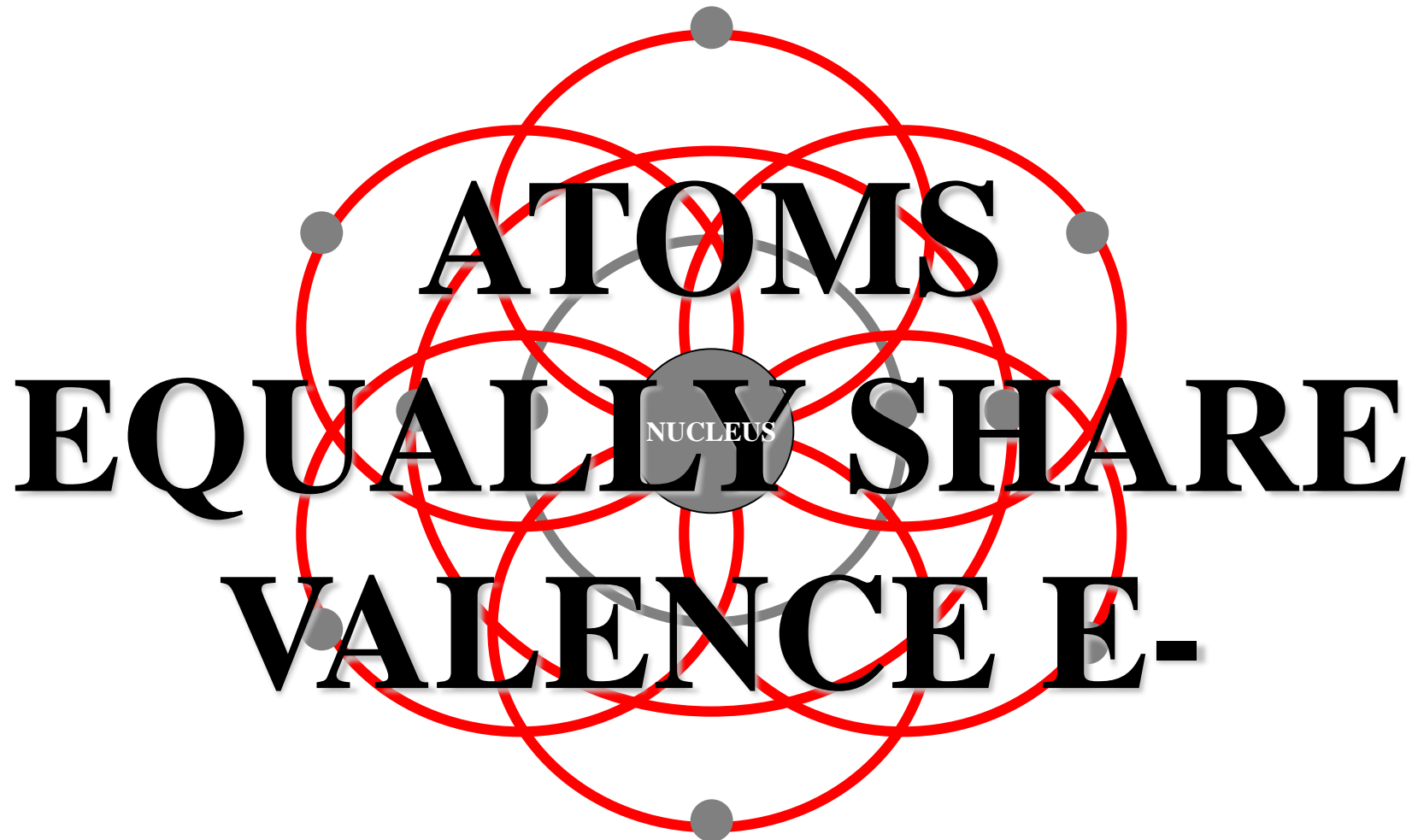


**NON-POLAR  
COVALENT  
BONDS  
&  
POLAR  
COVALENT  
BONDS**

# NON-POLAR COVALENT BOND

**NON-POLAR  
COVALENT  
BOND**

# NON-POLAR COVALENT BOND



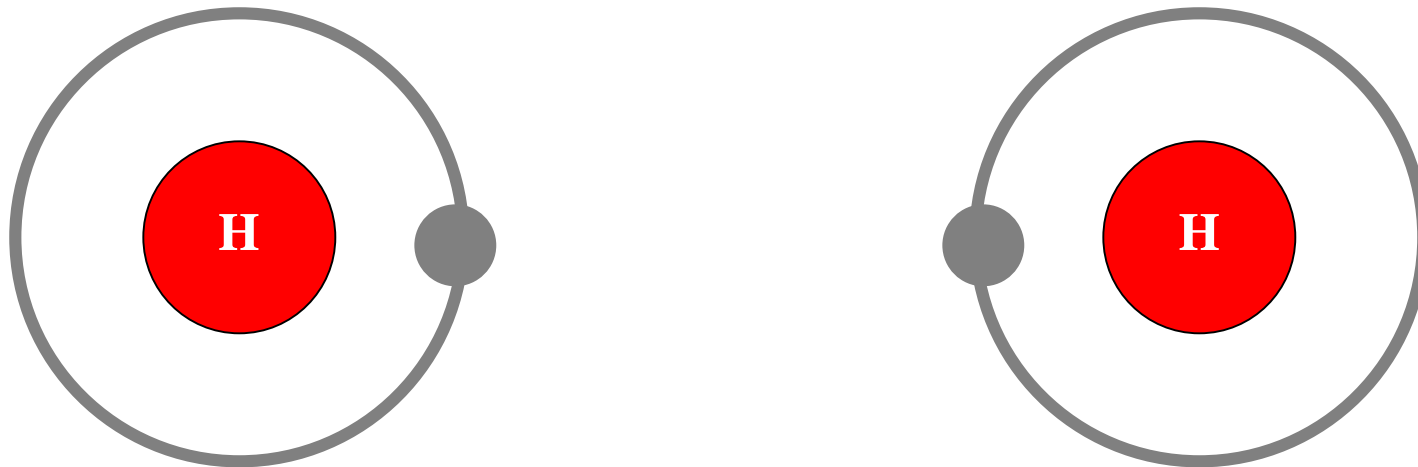
NON-POLAR COVALENT BOND





**NON-POLAR  
COVALENT BOND  
EXAMPLE  
HYDROGEN  
MOLECULE**

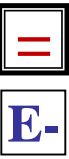
# NON-POLAR COVALENT BOND



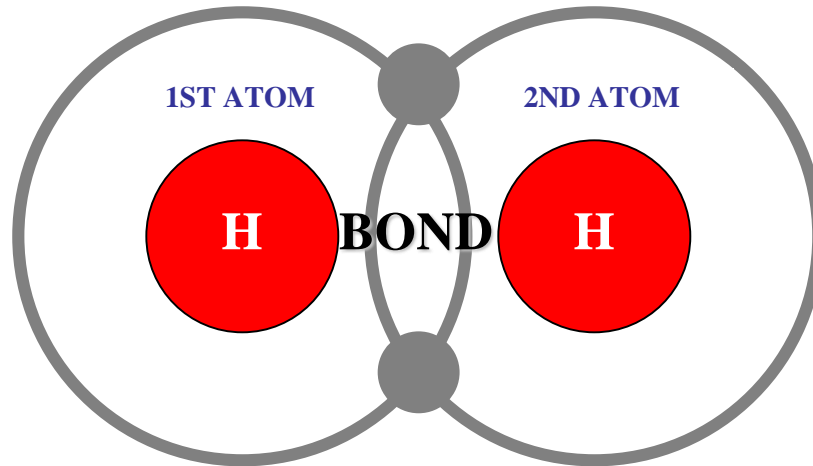
**HYDROGEN ATOMS**

● = e<sup>-</sup>

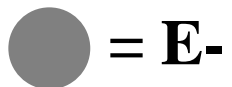
# NON-POLAR COVALENT BOND



**HYDROGEN ATOMS**



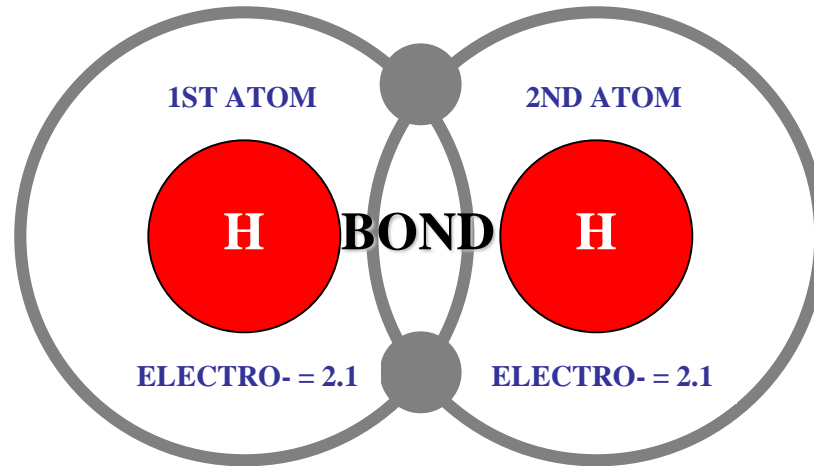
**HYDROGEN MOLECULE**




# NON-POLAR COVALENT BOND



**HYDROGEN MOLECULE**



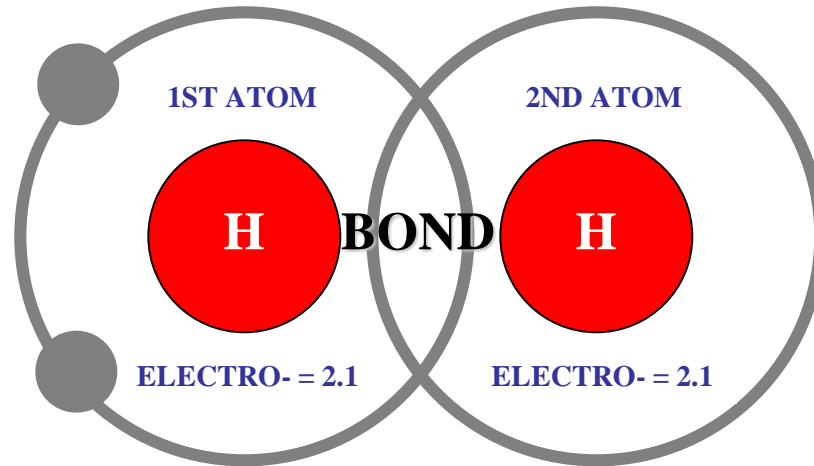
 = **E-**

**H EQUAL ELECTRO-**

# NON-POLAR COVALENT BOND



**HYDROGEN MOLECULE**



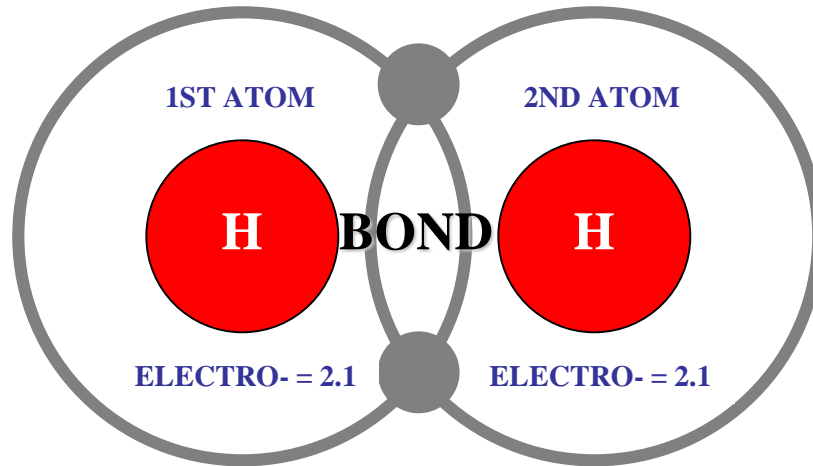
● = E-


**H EQUAL ELECTRO-**

# NON-POLAR COVALENT BOND



**HYDROGEN MOLECULE**



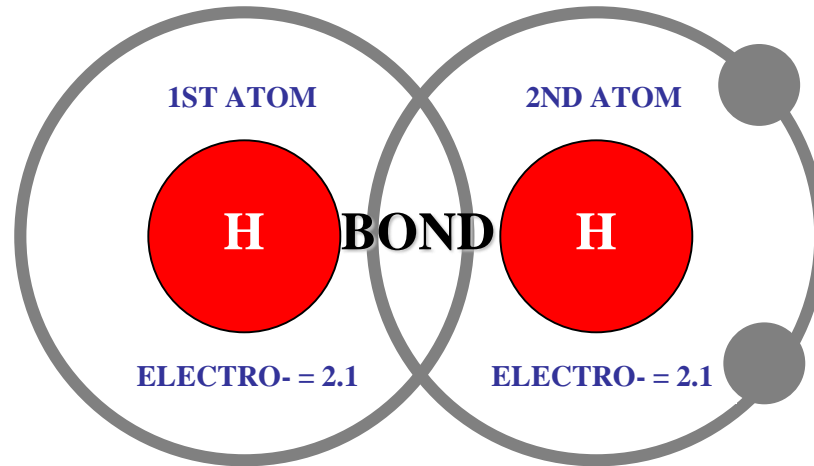
 = **E-**

**H EQUAL ELECTRO-**

# NON-POLAR COVALENT BOND



**HYDROGEN MOLECULE**



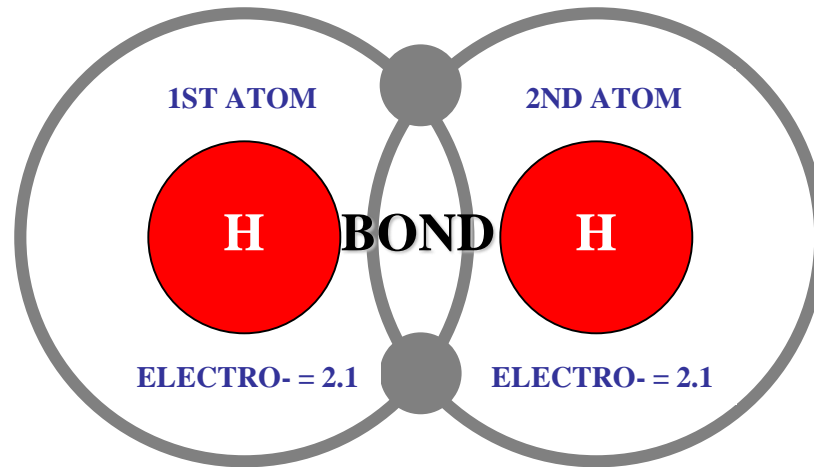
● = E-

**H EQUAL ELECTRO-**

# NON-POLAR COVALENT BOND



**HYDROGEN MOLECULE**



● = E-

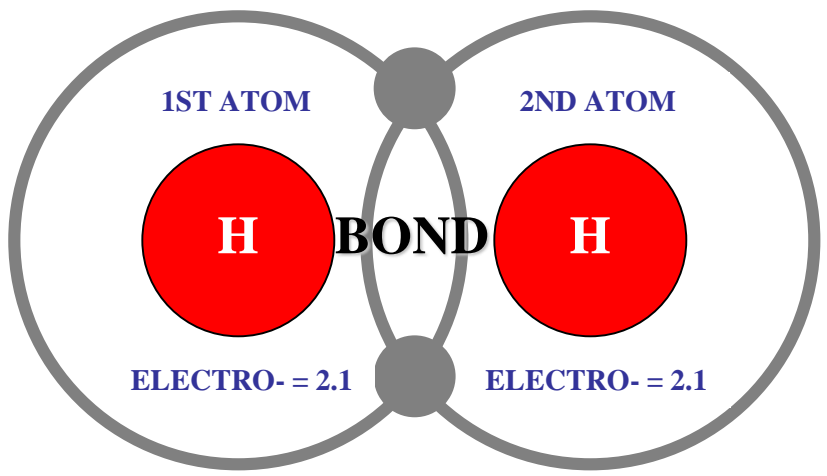
**H EQUAL ELECTRO-**



# NON-POLAR COVALENT BOND



**HYDROGEN MOLECULE**



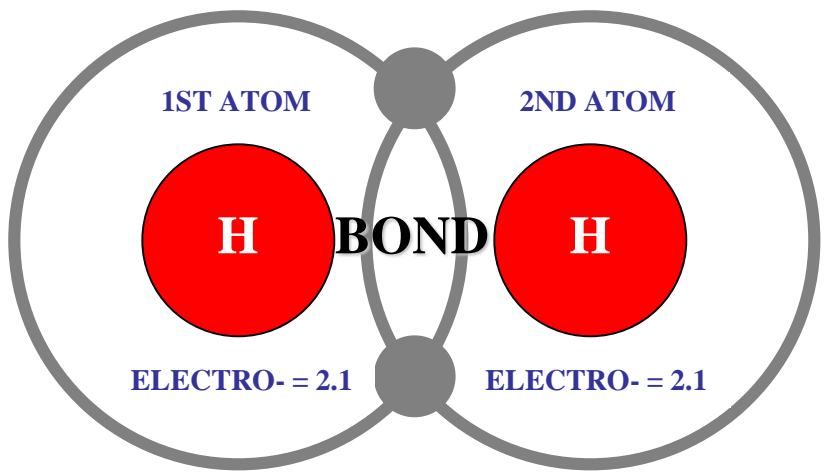
● = E-

**EQUAL E- SHARING: OUTCOME**

# NON-POLAR COVALENT BOND



**HYDROGEN MOLECULE**



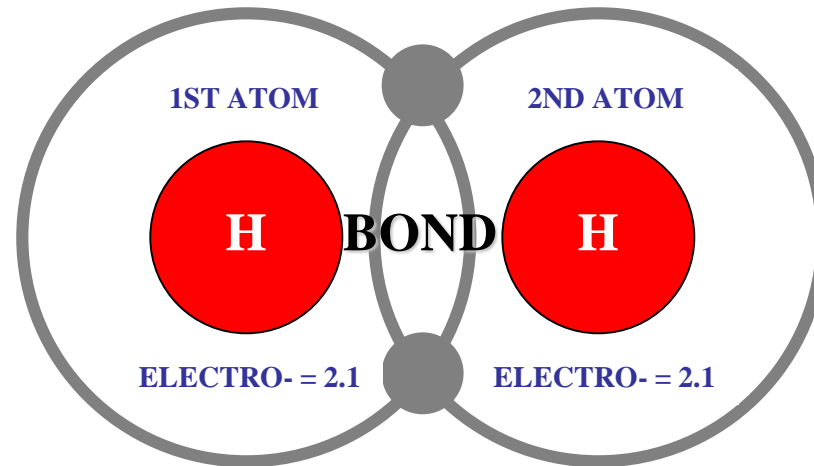
● = E-

**NON-POLAR COVALENT BOND**

# NON-POLAR COVALENT BOND



## HYDROGEN MOLECULE



● = E-

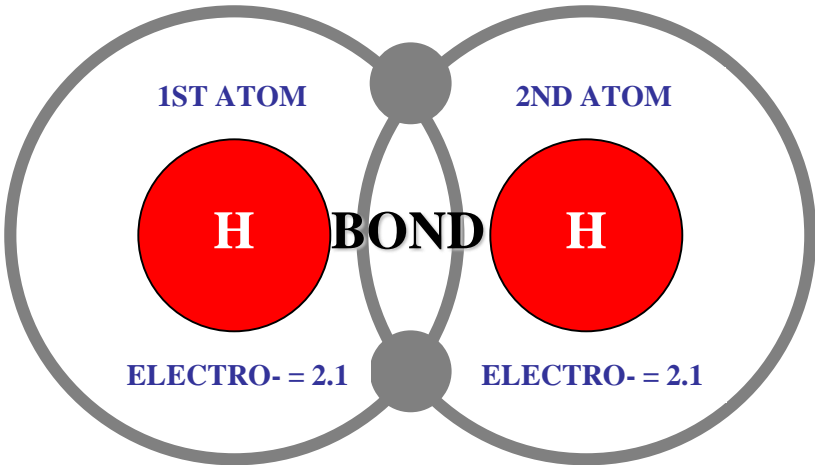
**HIGH E- SHARING: OUTCOME**



# NON-POLAR COVALENT BOND



**HYDROGEN MOLECULE**

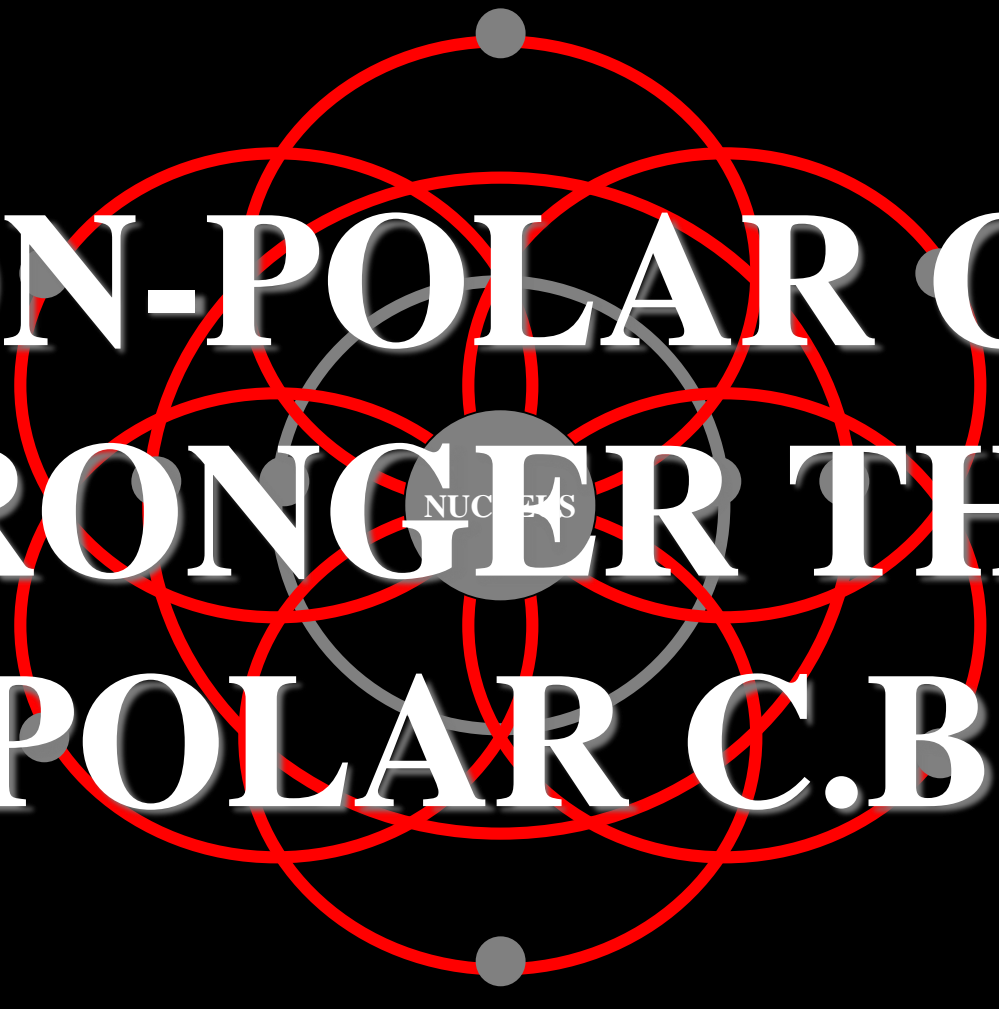


● = E-

**STRONG COVALENT BOND**



# CHEMICAL BONDS



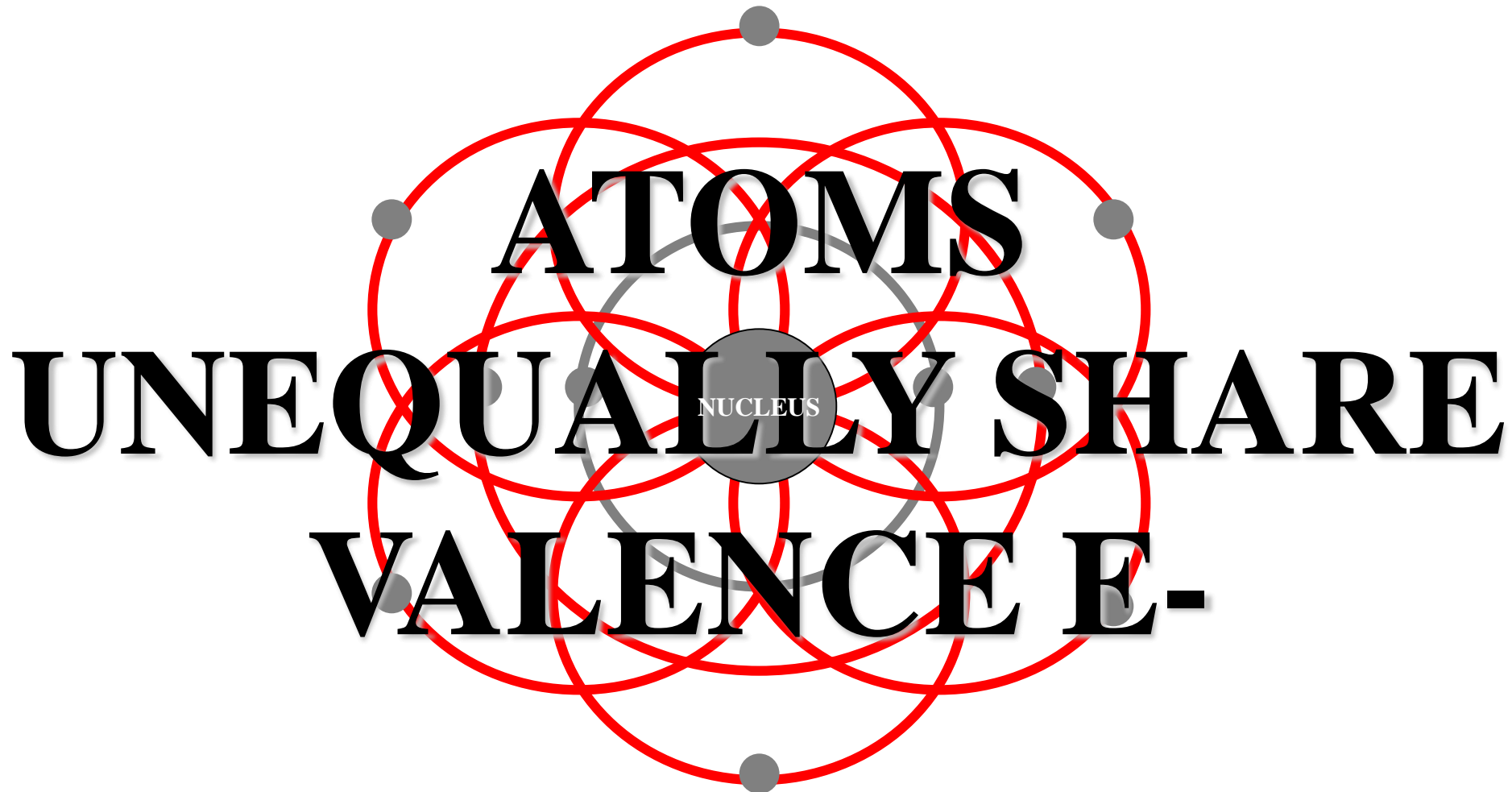
**NON-POLAR C.B.  
STRONGER THAN  
POLAR C.B.**

# CHEMICAL BONDS

# POLAR COVALENT BOND

**POLAR  
COVALENT  
BOND**

# POLAR COVALENT BOND



POLAR COVALENT BOND

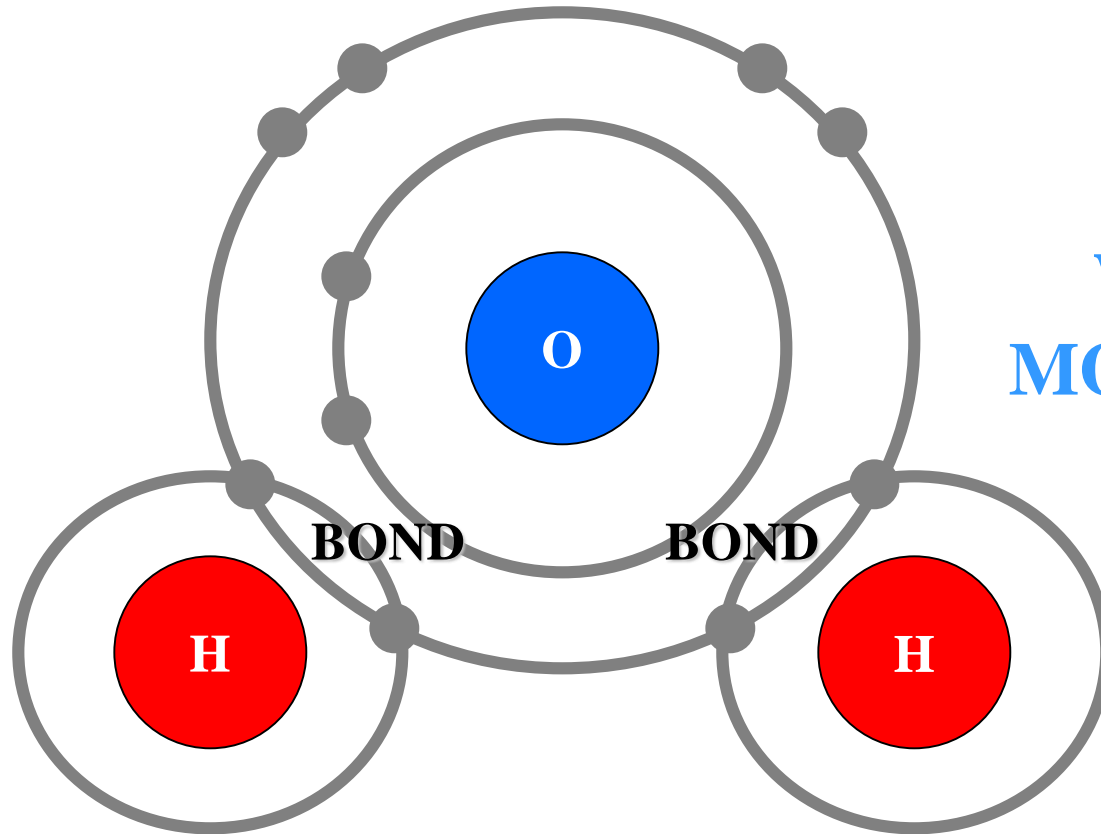


**POLAR  
COVALENT BOND  
EXAMPLE  
WATER  
MOLECULE**

# POLAR COVALENT BOND

OXYGEN

HYDROGEN



● = E-

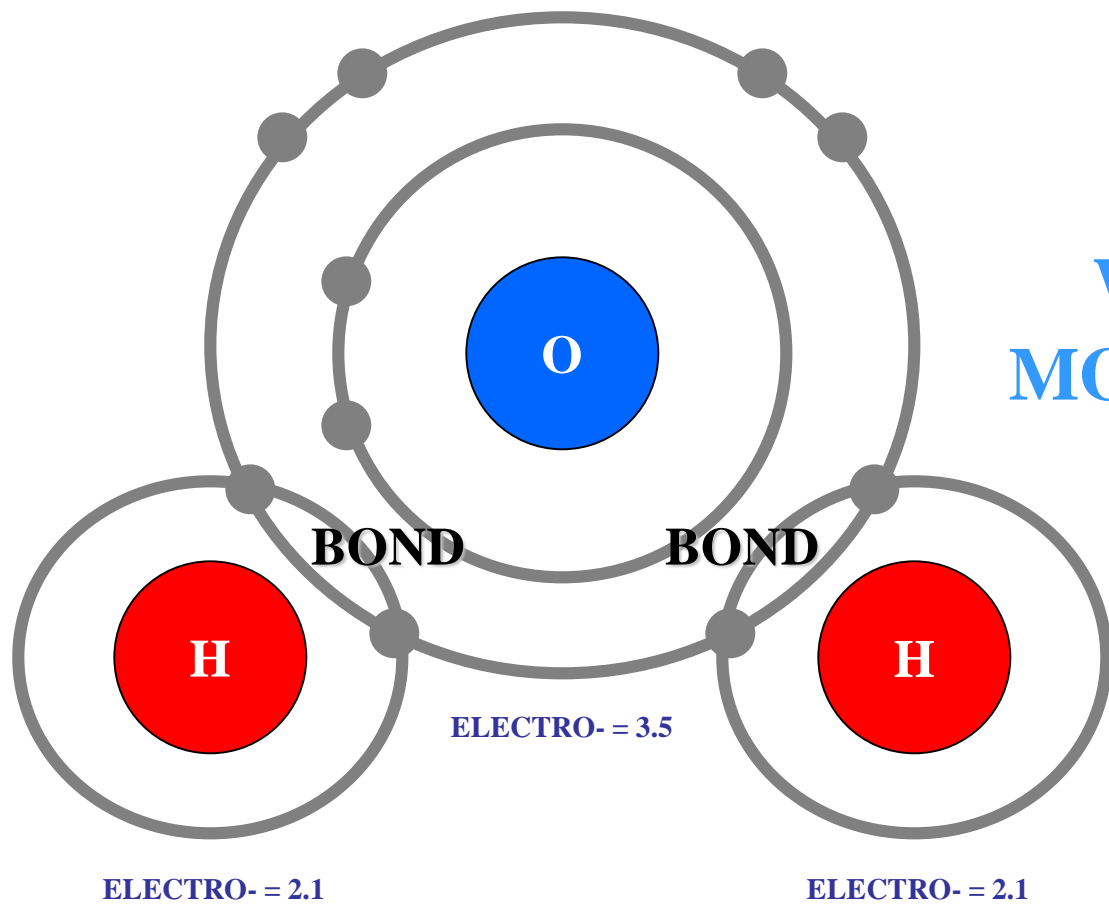
DIFF ELEMENTS

# POLAR COVALENT BOND



**OXYGEN**  
**HIGHER**  
**ELECTRO-**

**HYDROGEN**  
**LOWER**  
**ELECTRO-**



● = **E-**

**O & H DIFF ELECTRO-**

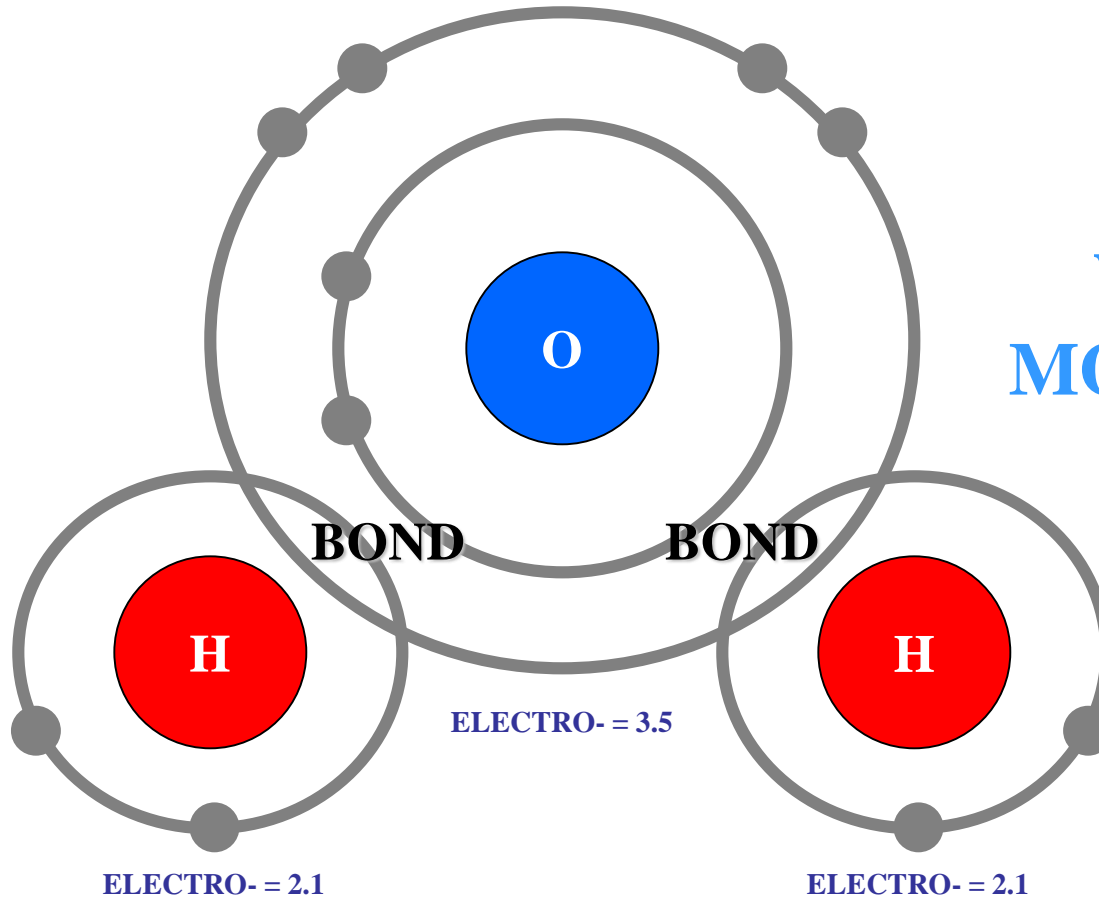
# POLAR COVALENT BOND



**OXYGEN**  
**HIGHER**  
**ELECTRO-**

**HYDROGEN**  
**LOWER**  
**ELECTRO-**

**WATER**  
**MOLECULE**



● = E-

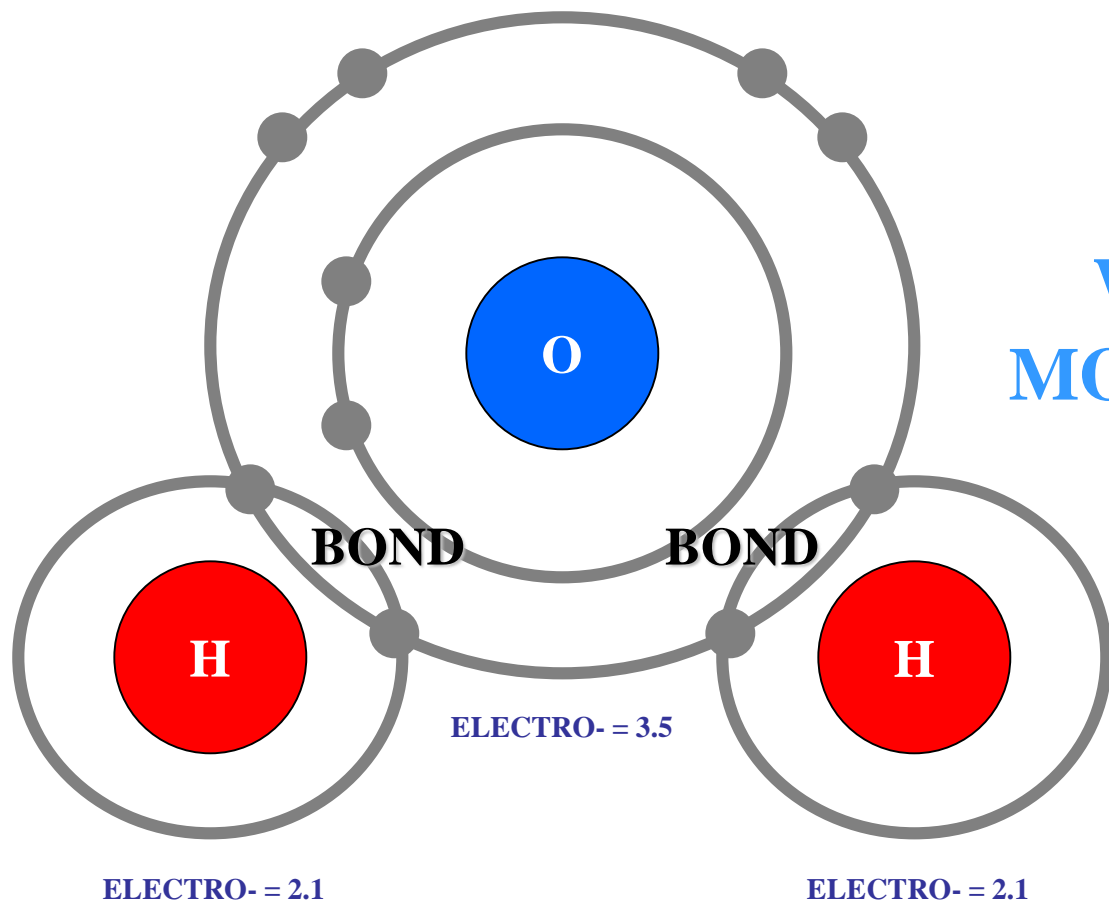
**O & H DIFF ELECTRO-**

# POLAR COVALENT BOND



**OXYGEN**  
**HIGHER**  
**ELECTRO-**

**HYDROGEN**  
**LOWER**  
**ELECTRO-**



● = E-

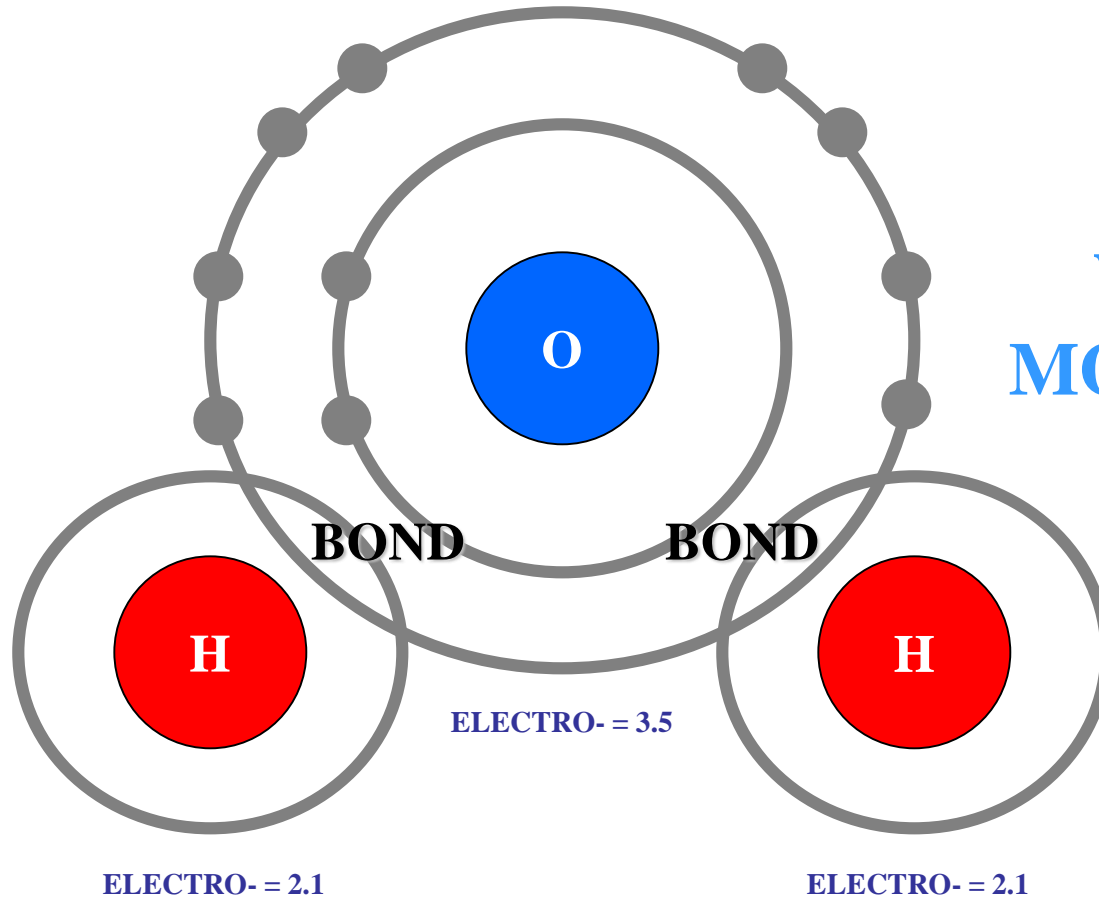
**O & H DIFF ELECTRO-**

# POLAR COVALENT BOND



**OXYGEN**  
**HIGHER**  
**ELECTRO-**

**HYDROGEN**  
**LOWER**  
**ELECTRO-**



**WATER**  
**MOLECULE**

● = E-

**O & H DIFF ELECTRO-**

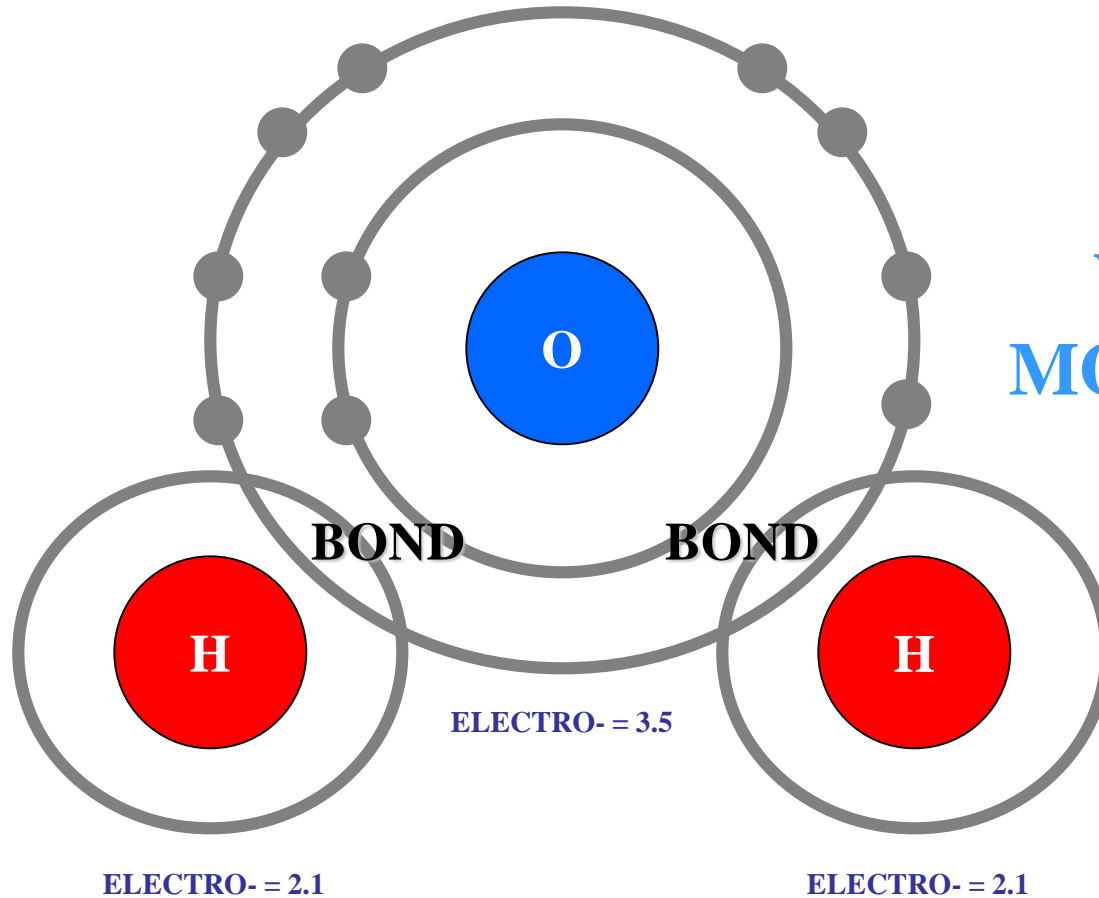
# POLAR COVALENT BOND

E- SPEND MORE TIME ABOUT **OXYGEN**

**OXYGEN**  
HIGHER  
ELECTRO-

**HYDROGEN**  
LOWER  
ELECTRO-

**WATER**  
MOLECULE



● = E-

**O & H DIFF ELECTRO-**

# POLAR COVALENT BOND

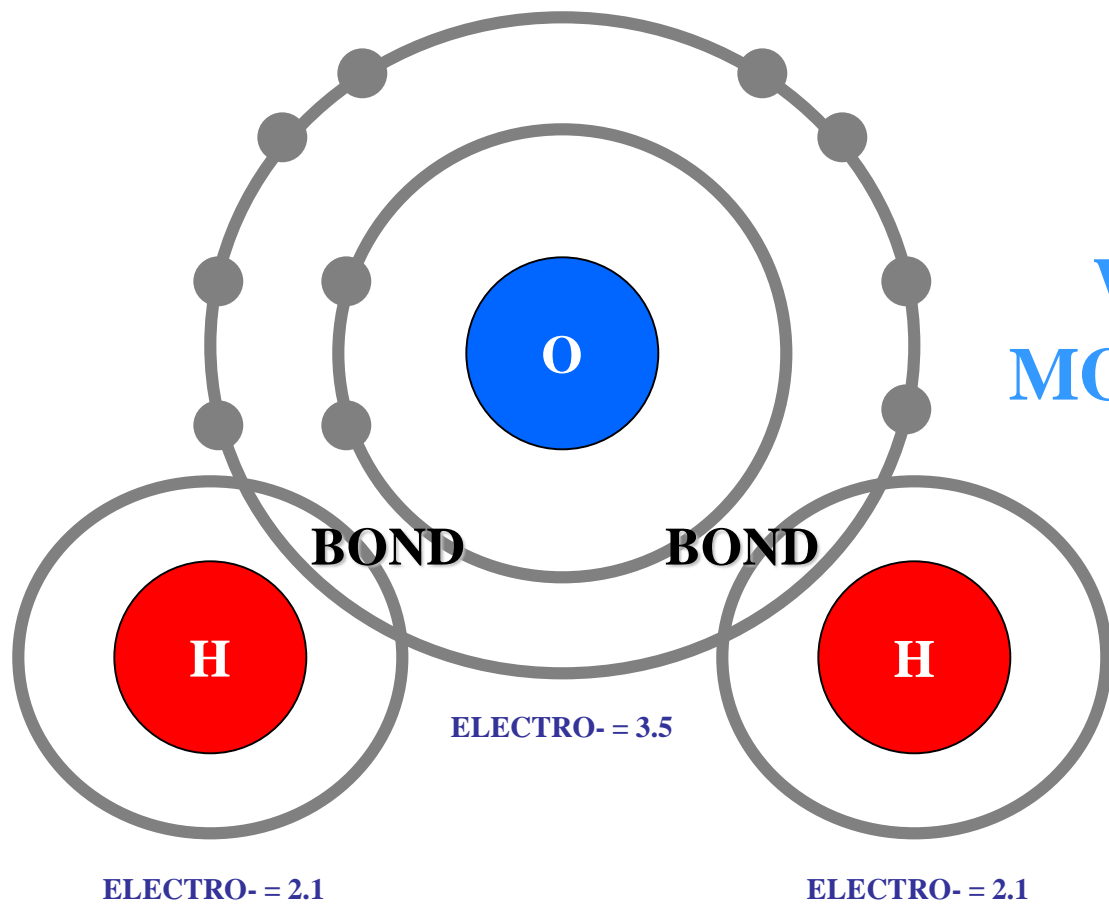


E- SPEND MORE TIME ABOUT **OXYGEN**

**OXYGEN**  
HIGHER  
ELECTRO-

**HYDROGEN**  
LOWER  
ELECTRO-

**WATER**  
MOLECULE



● = E-

**UNEQUAL E- SHARING**



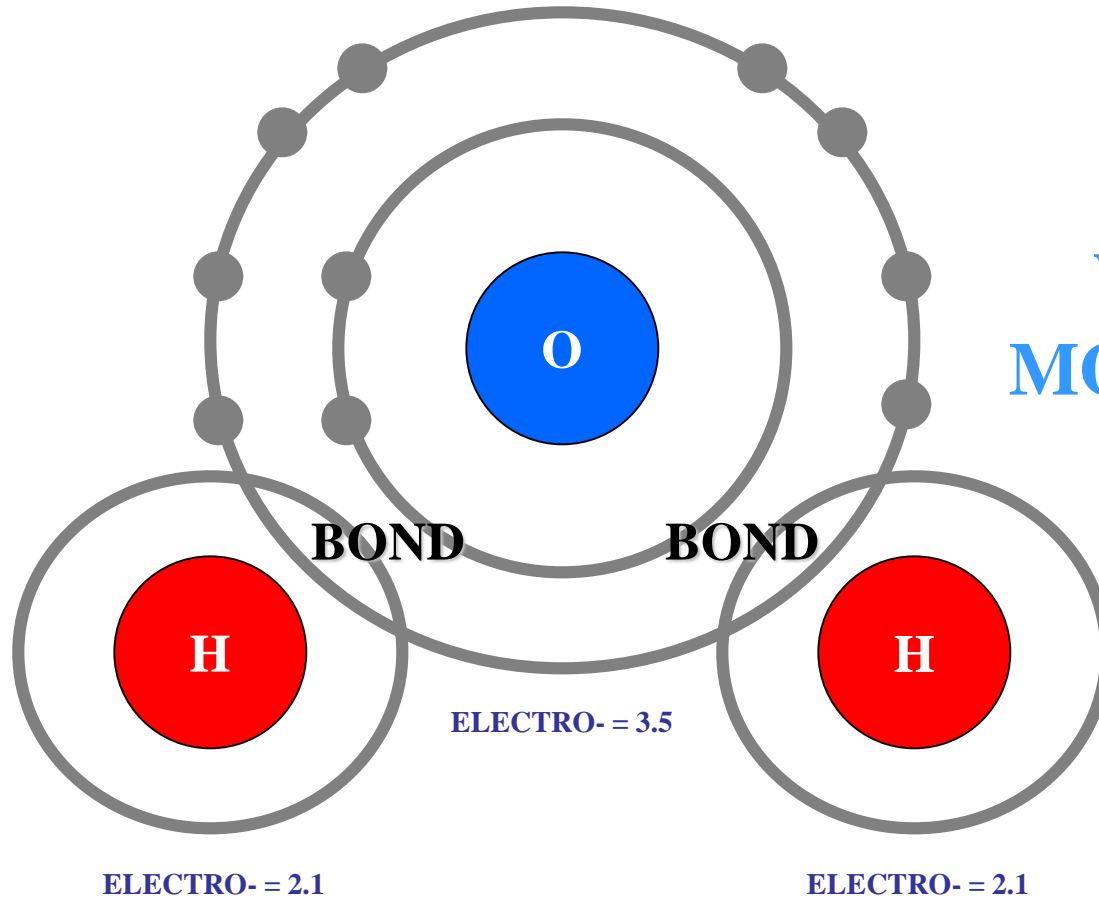
# POLAR COVALENT BOND

? CHARGE

**OXYGEN**  
HIGHER  
ELECTRO-

**HYDROGEN**  
LOWER  
ELECTRO-

**WATER**  
MOLECULE



● = E-

**UNEQUAL E- SHARING**

# POLAR COVALENT BOND

?

H

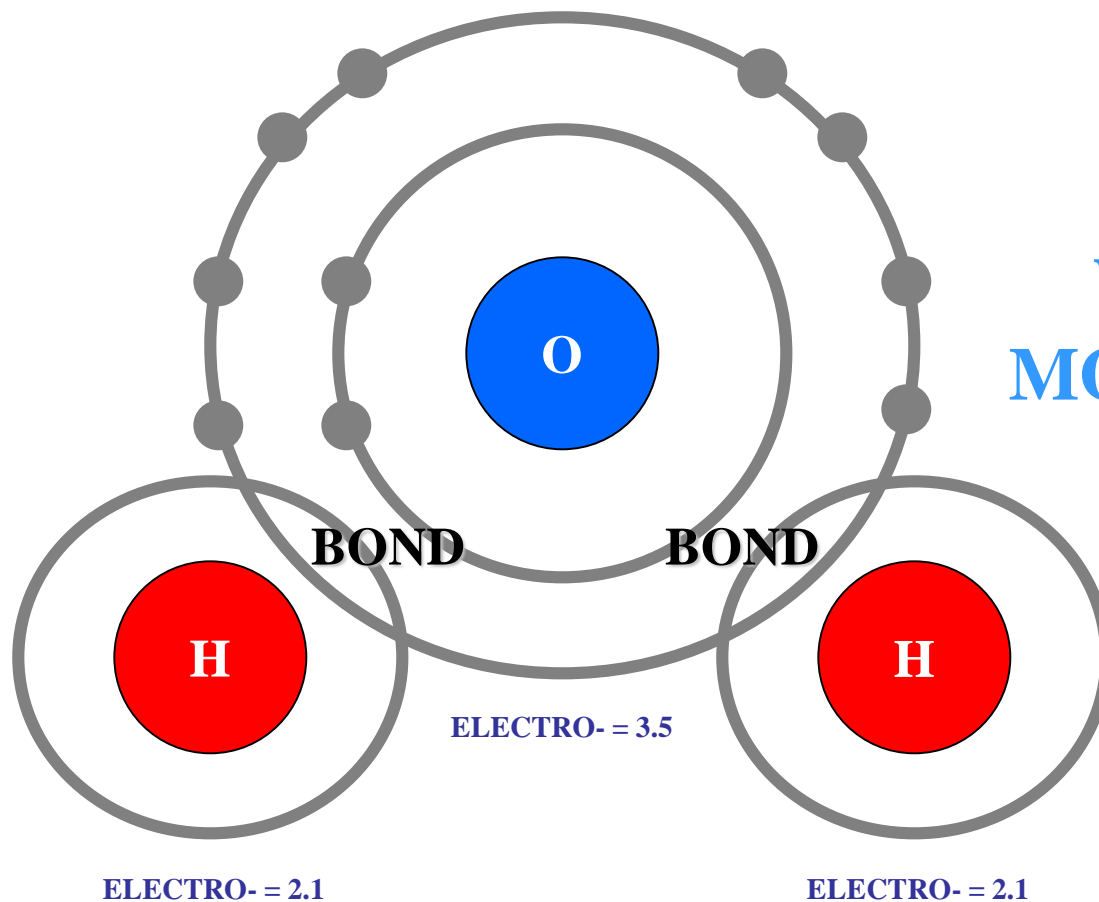
+

- CHARGE

**OXYGEN**  
HIGHER  
ELECTRO-

**HYDROGEN**  
LOWER  
ELECTRO-

**WATER**  
MOLECULE



● = E-

**UNEQUAL E- SHARING**

# POLAR COVALENT BOND

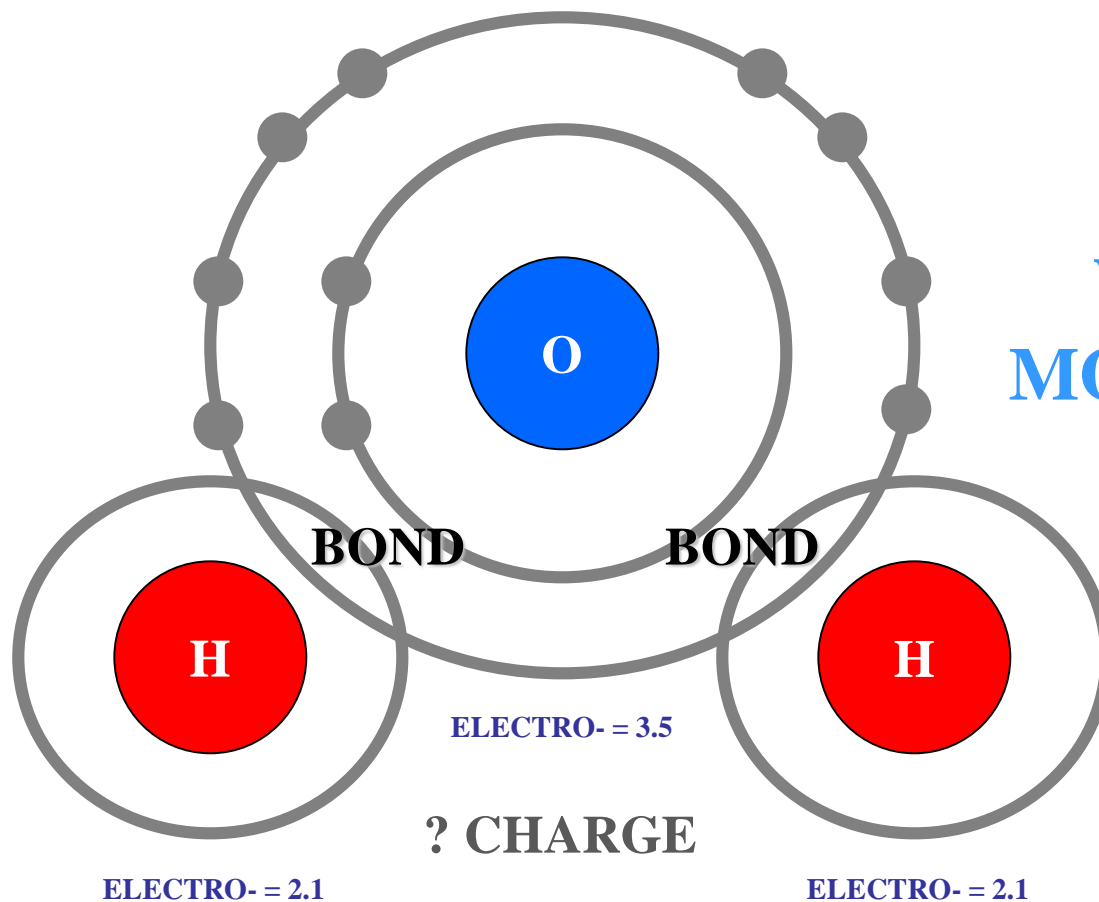
+

- CHARGE

**OXYGEN**  
HIGHER  
ELECTRO-

**HYDROGEN**  
LOWER  
ELECTRO-

**WATER**  
MOLECULE



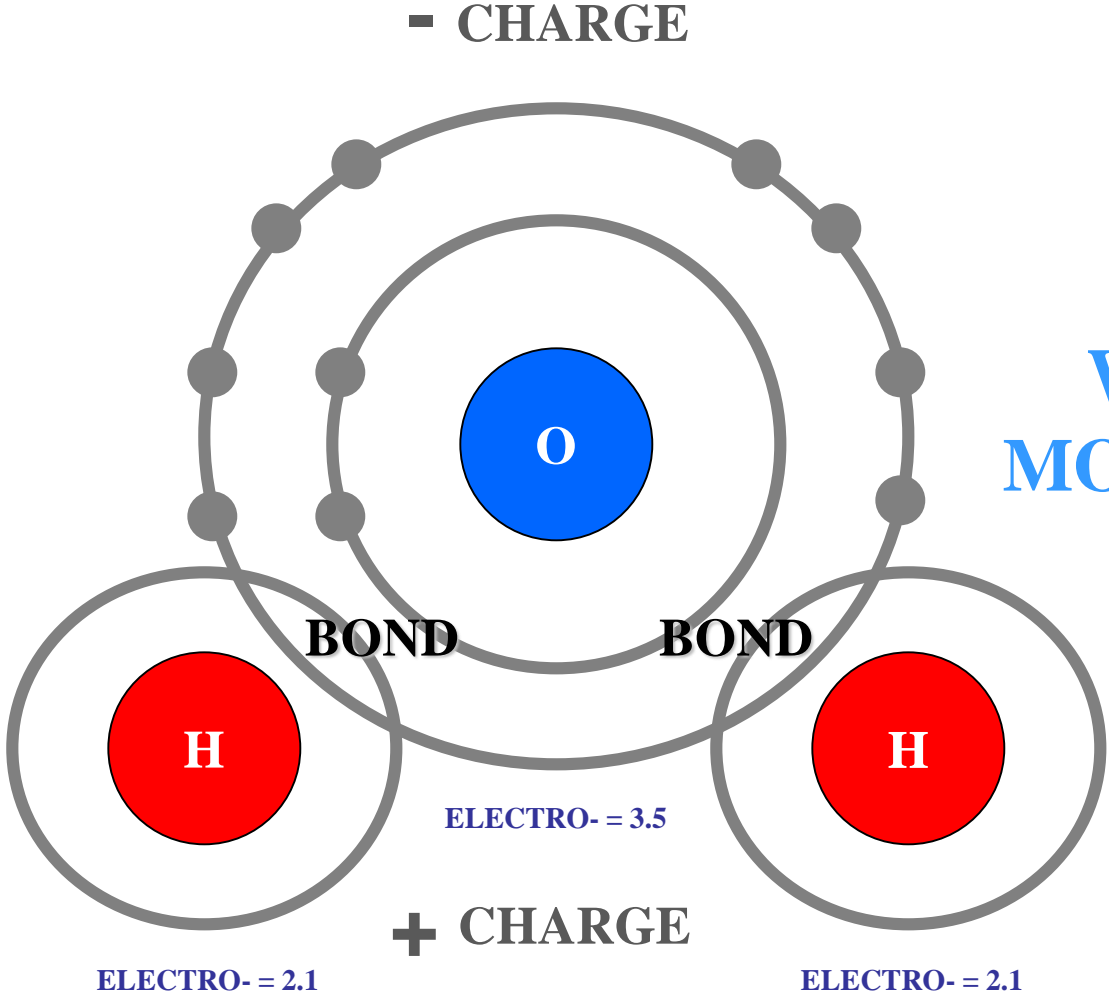
● = E-

**UNEQUAL E- SHARING**

# POLAR COVALENT BOND

**OXYGEN**  
HIGHER  
ELECTRO-

**HYDROGEN**  
LOWER  
ELECTRO-



● = E-

**UNEQUAL E- SHARING: OUTCOME**

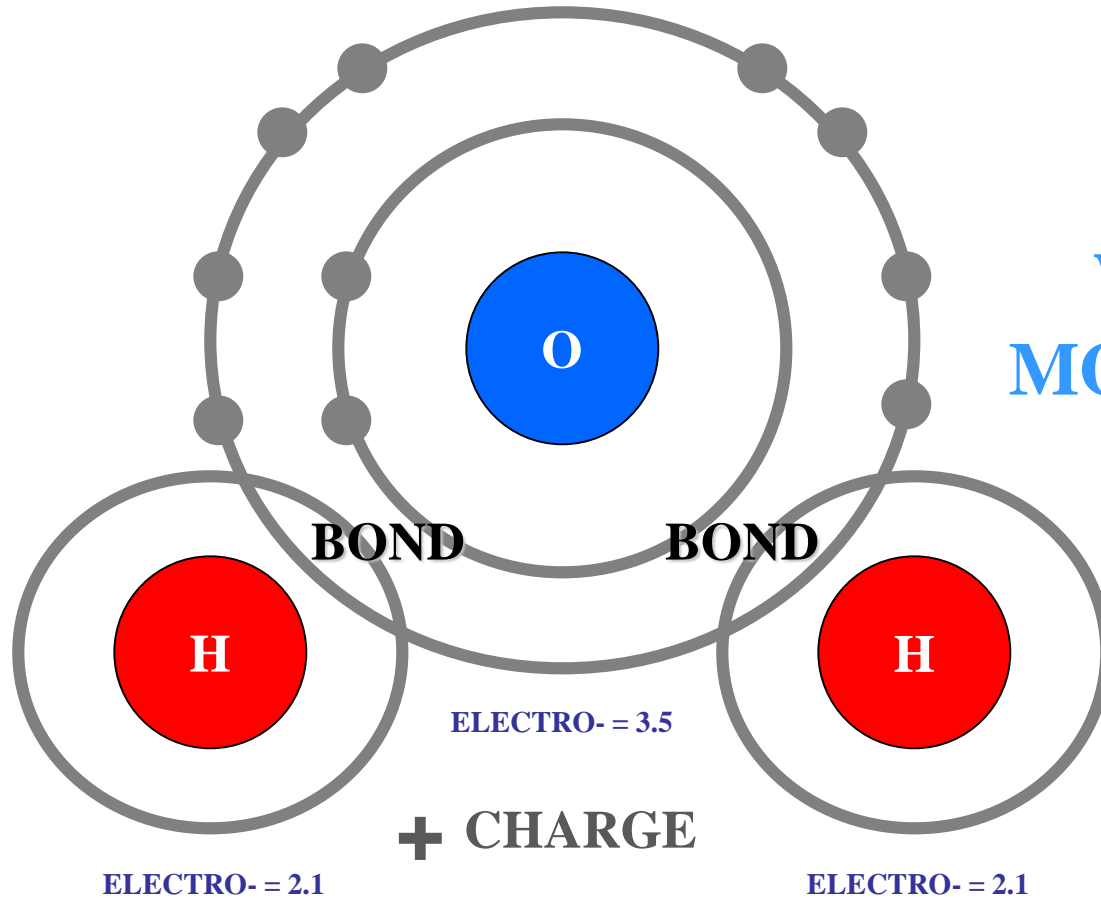
# POLAR COVALENT BOND

- CHARGE

**OXYGEN**  
HIGHER  
ELECTRO-

**HYDROGEN**  
LOWER  
ELECTRO-

**WATER**  
MOLECULE



● = E-

**POLAR COVALENT BOND**

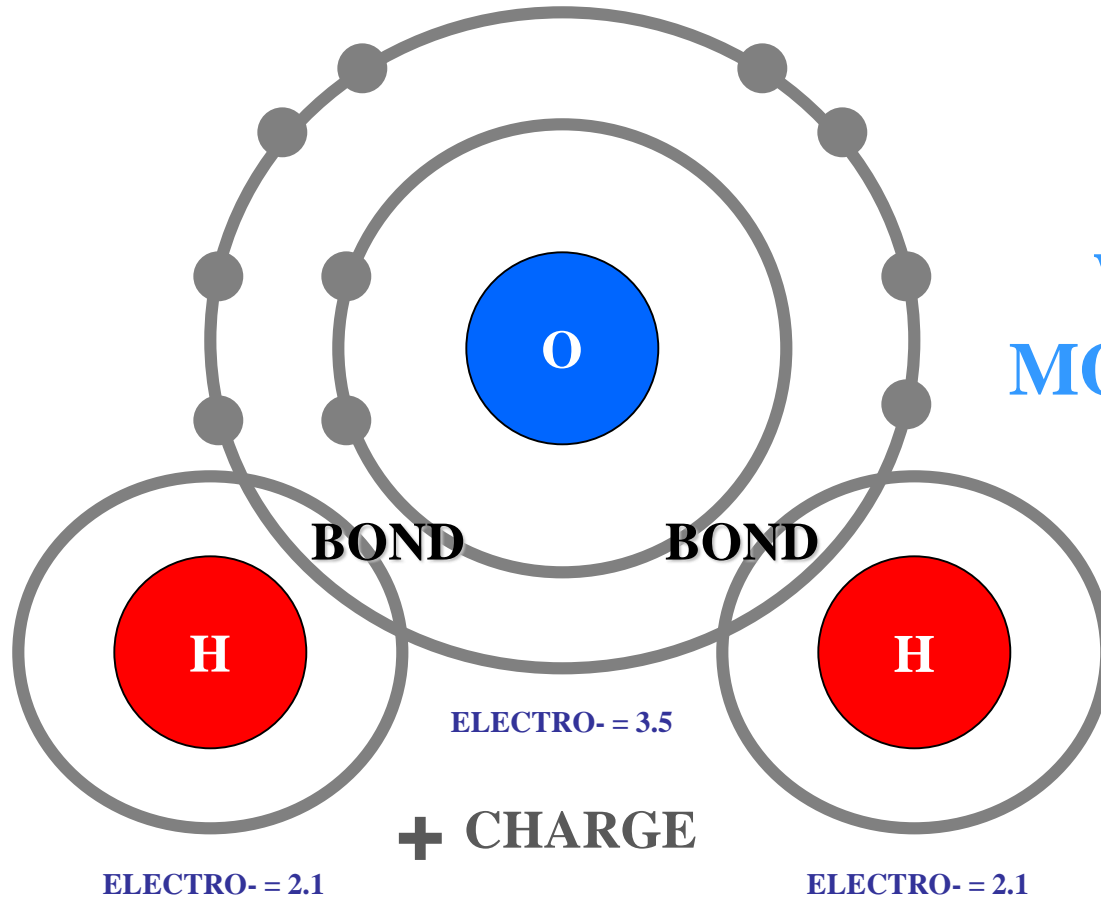
# POLAR COVALENT BOND

- CHARGE

**OXYGEN**  
HIGHER  
ELECTRO-

**HYDROGEN**  
LOWER  
ELECTRO-

**WATER  
MOLECULE**



ELECTRO- = 2.1

ELECTRO- = 3.5

ELECTRO- = 2.1

+ CHARGE

● = E-

**MODERATE E- SHARING: OUTCOME**



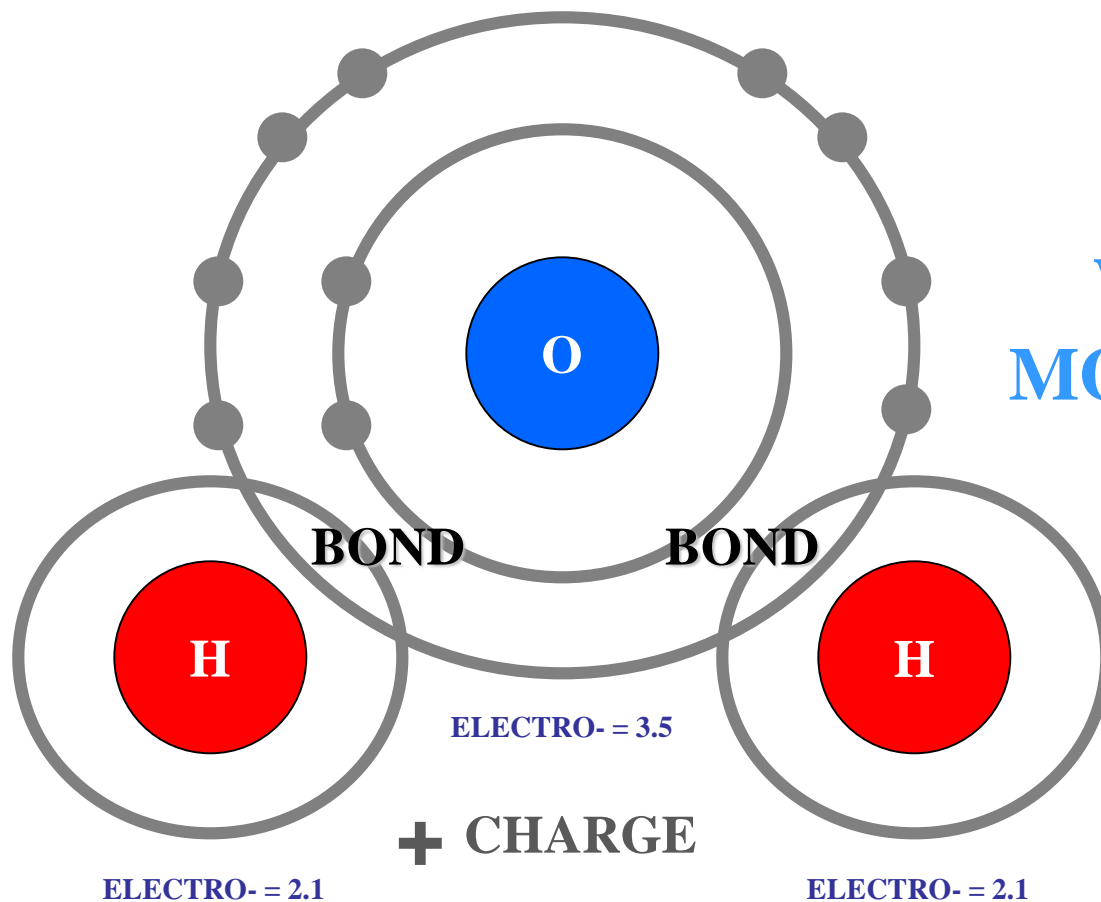
# POLAR COVALENT BOND

- CHARGE

**OXYGEN**  
HIGHER  
ELECTRO-

**HYDROGEN**  
LOWER  
ELECTRO-

**WATER**  
MOLECULE



● = E-

**WEAK COVALENT BOND**



# CHEMICAL BONDS



**POLAR C.B.  
WEAKER THAN  
NON-POLAR C.B.**

# CHEMICAL BONDS





# CHEMICAL BONDS



**COVALENT BONDS  
STRONGER THAN  
IONIC BONDS**

# CHEMICAL BONDS



# IONIC BONDS

# ION

**ION**



**ION**



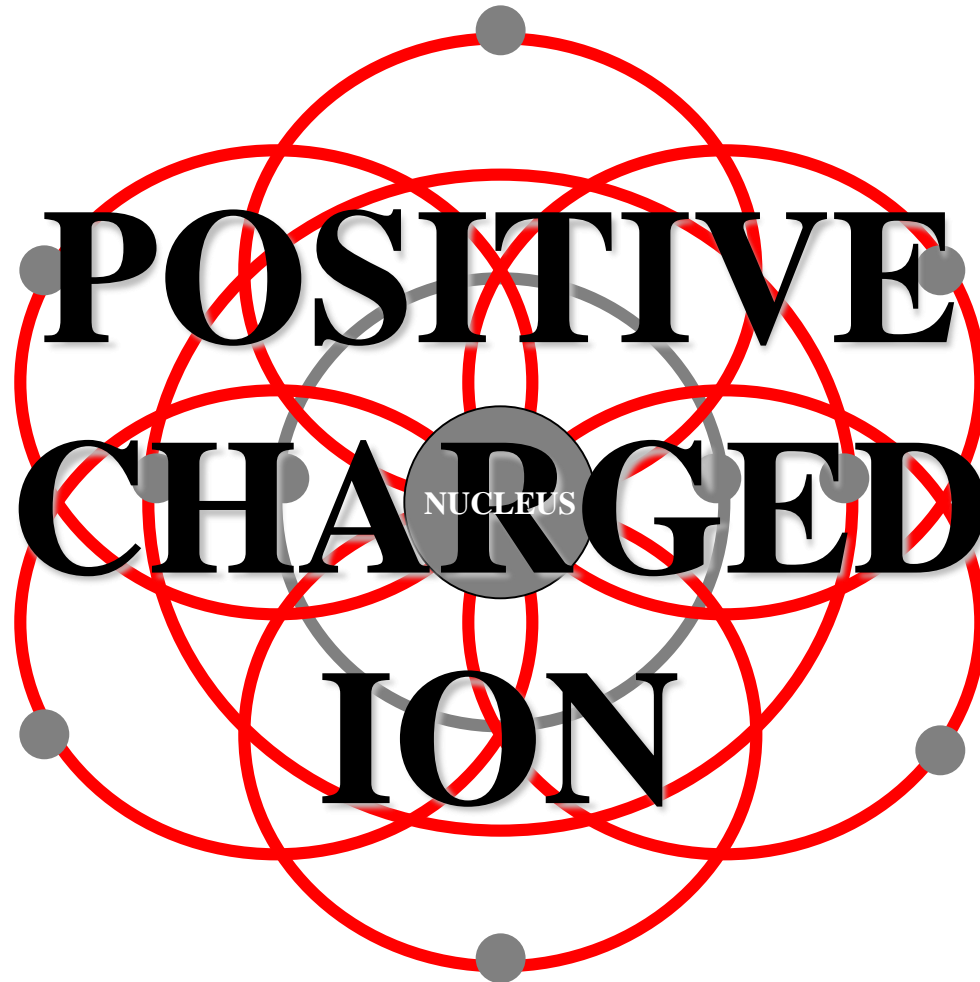
**ION**

# CATION VS ANION

**CATION**



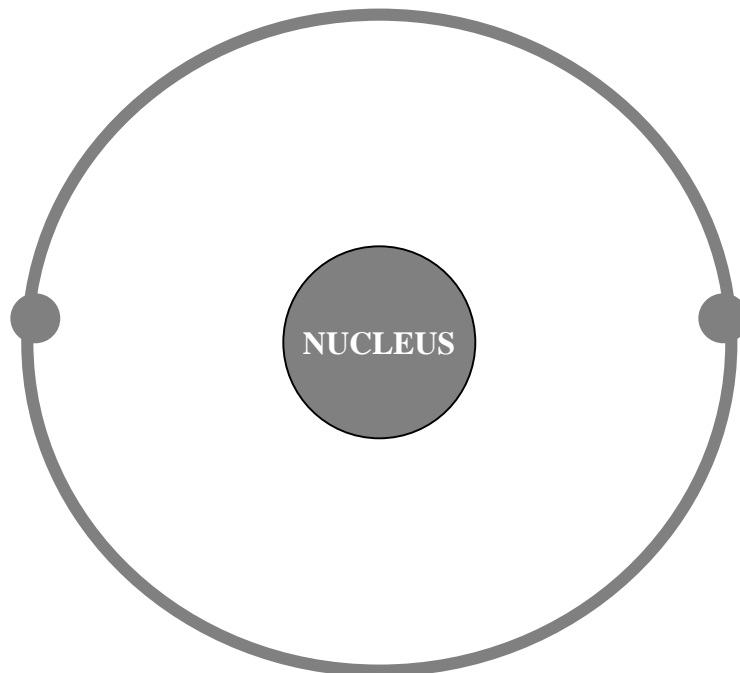
**CATION**



**CATION**



# ATOM

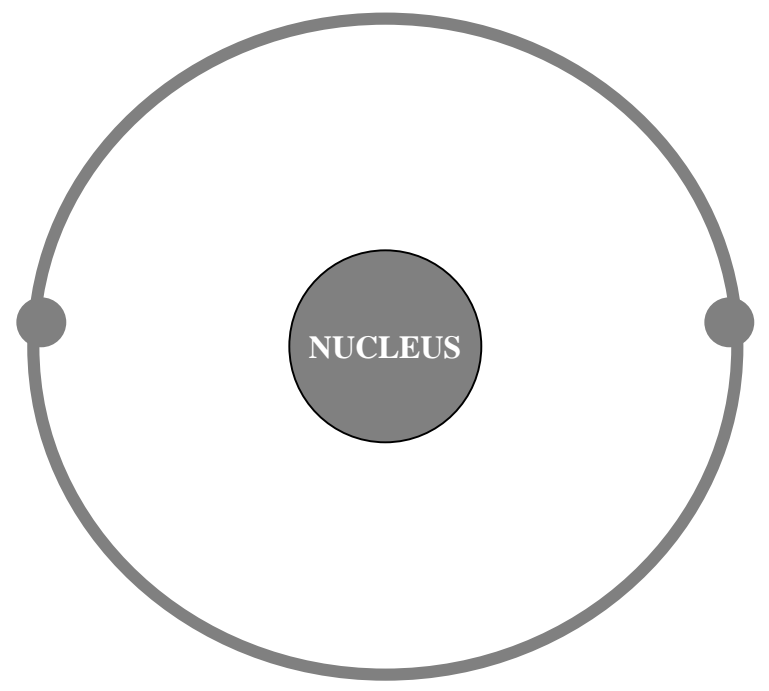


**LOW ELECTRO-NEGATIVITY**

● = E-



# ATOM

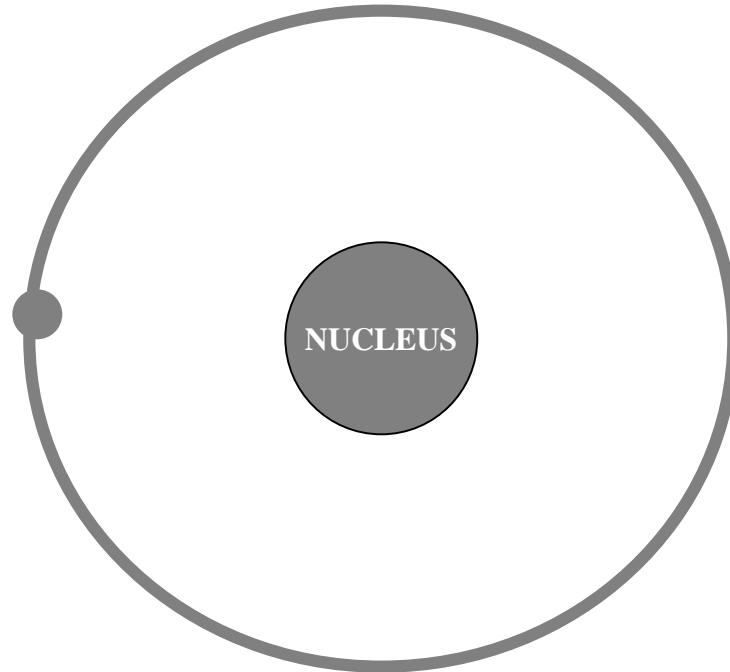


● = E-

**LOSES E-  
BOND FORMATION**



# ATOM

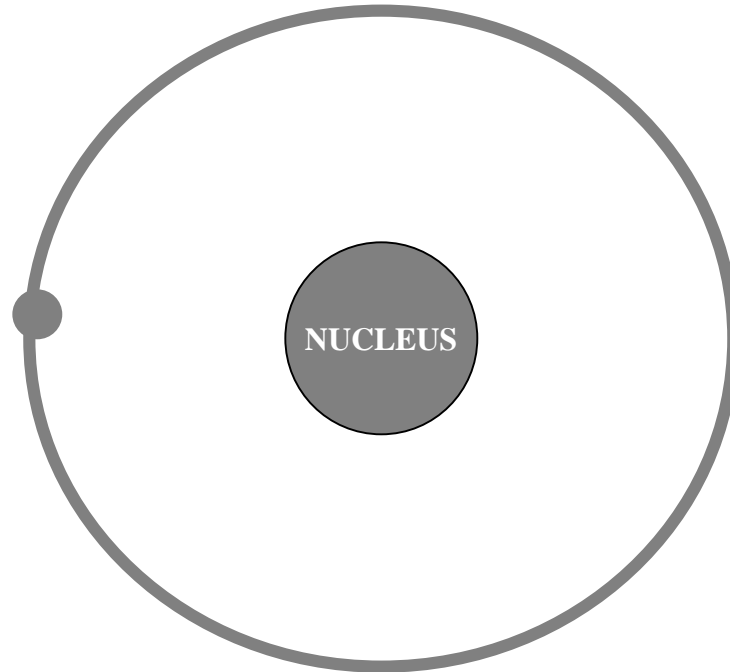


**LOSES E-**

**BOND FORMATION**

● = E-

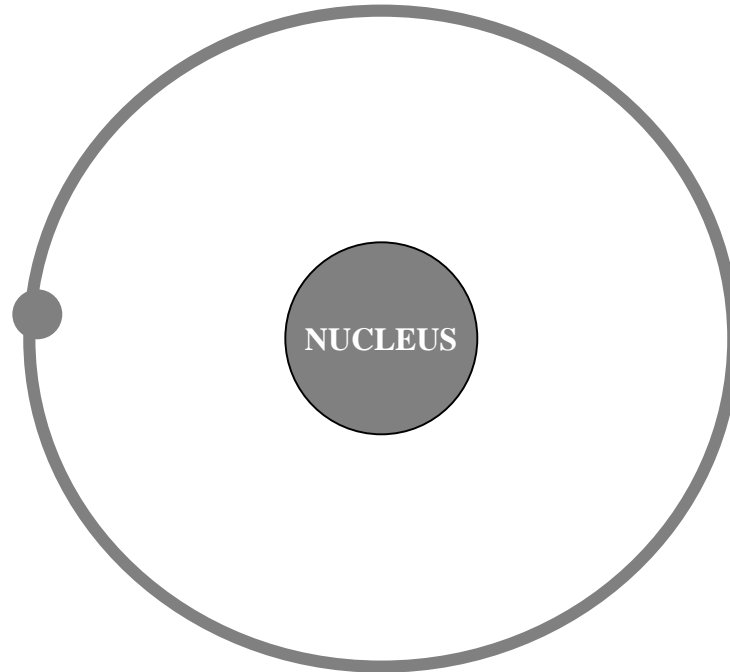
# ATOM



**IMPARTS ? CHARGE**

● = e-

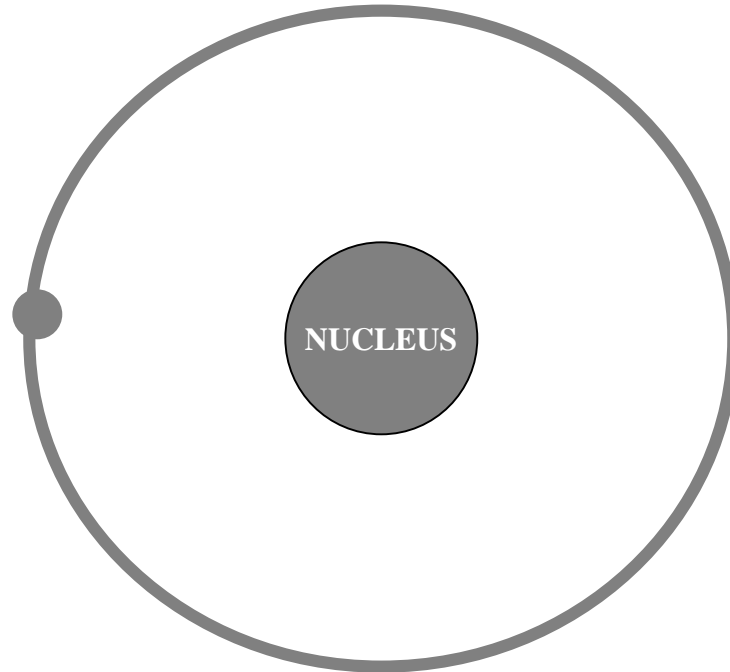
# ATOM



+ CHARGE

● = E-

# CATION



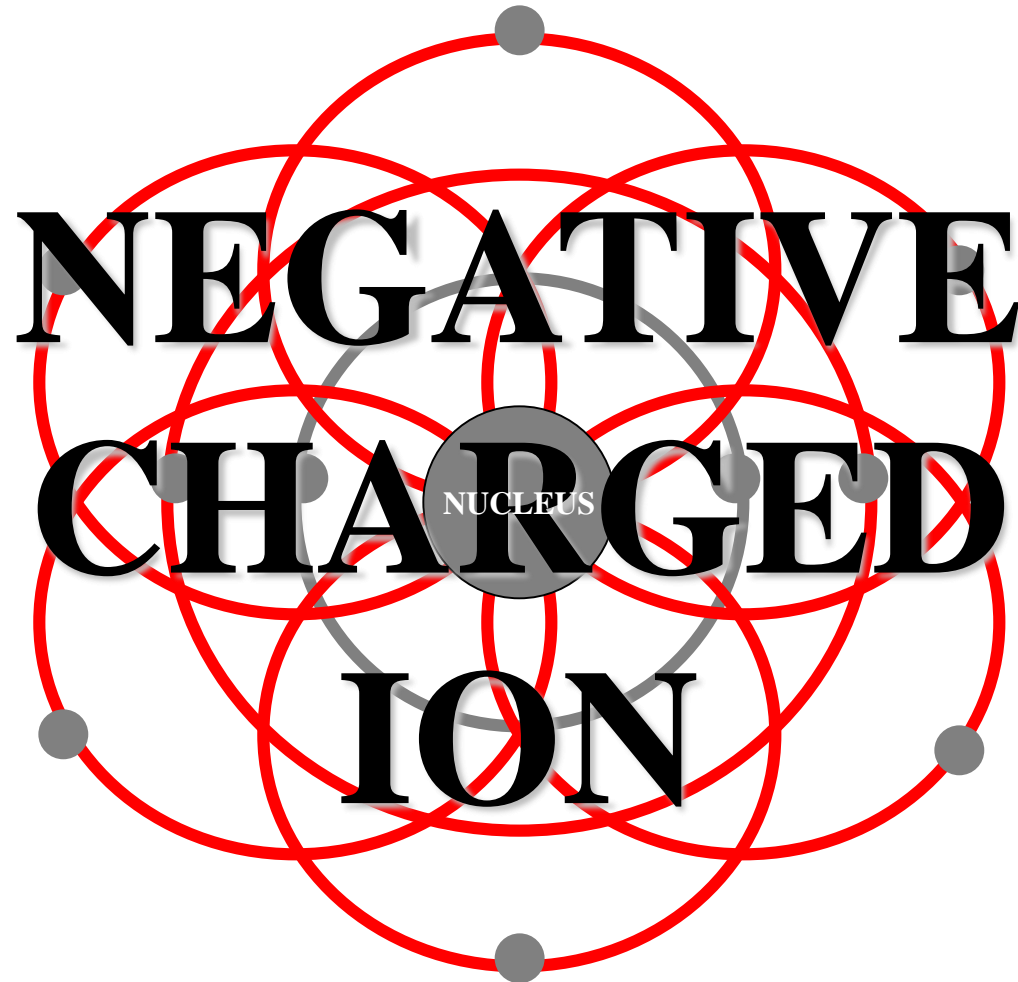
+ CHARGE

● = e-

**ANION**



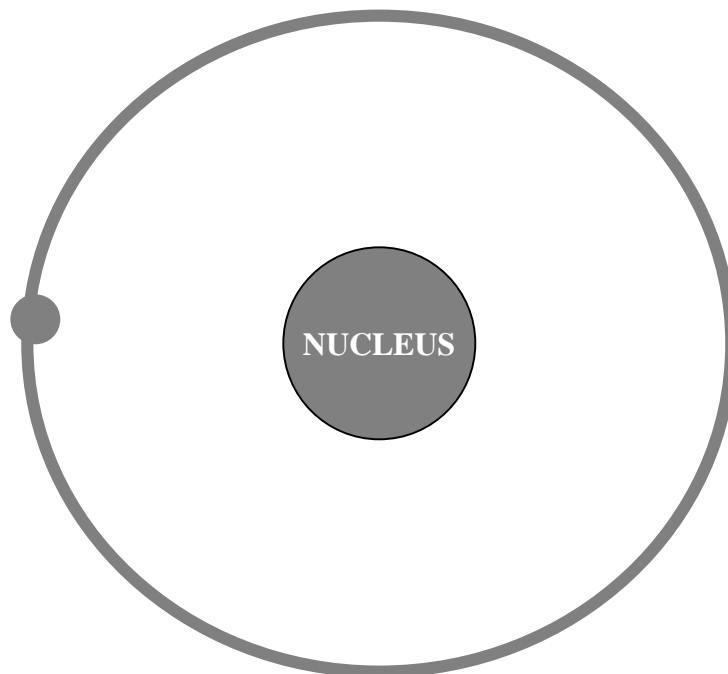
**ANION**



**ANION**



# ATOM

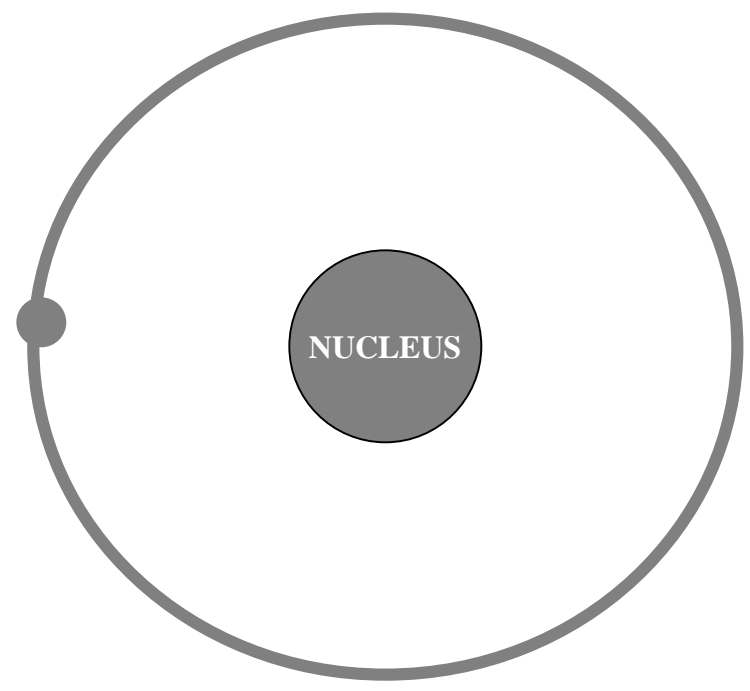


**HIGH ELECTRO-NEGATIVITY**

● = E-



# ATOM

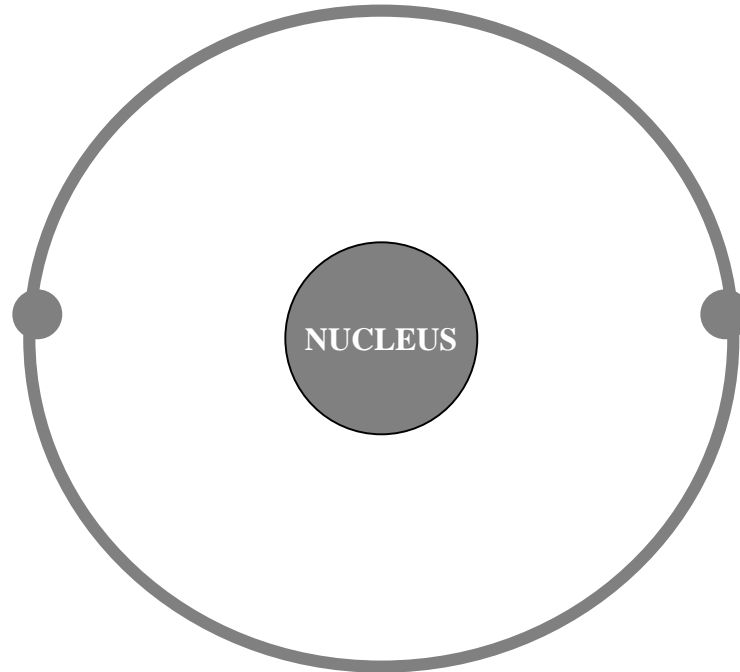


● = E-

**GAINS E-**  
**BOND FORMATION**



# ATOM

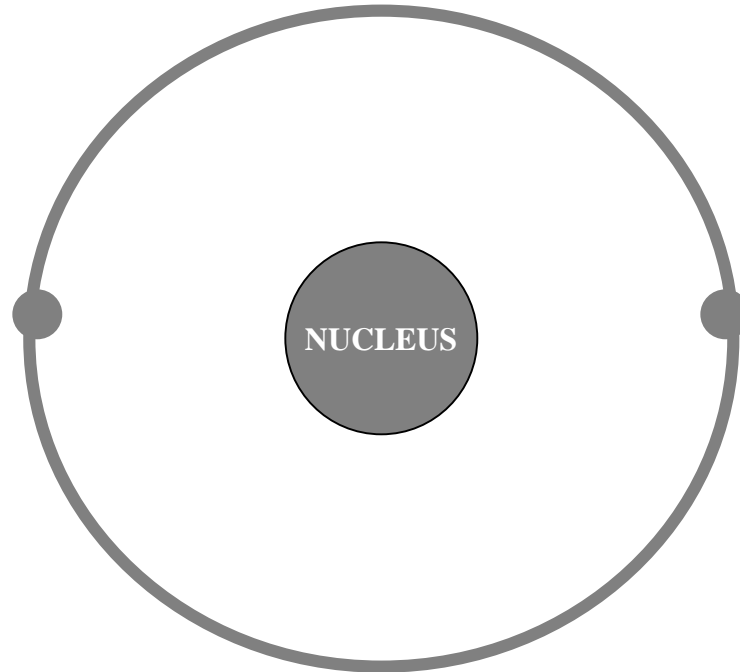


**GAINS E-**

**BOND FORMATION**

● = E-

# ATOM



**IMPARTS ? CHARGE**

● = e<sup>-</sup>