

Atomic Structure Review:

3 Parts

$p \Rightarrow$ never change

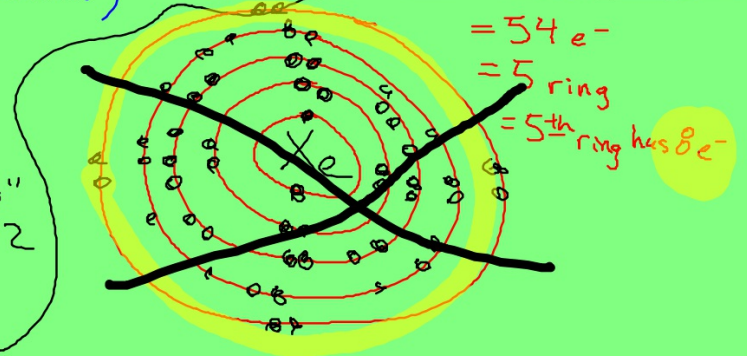
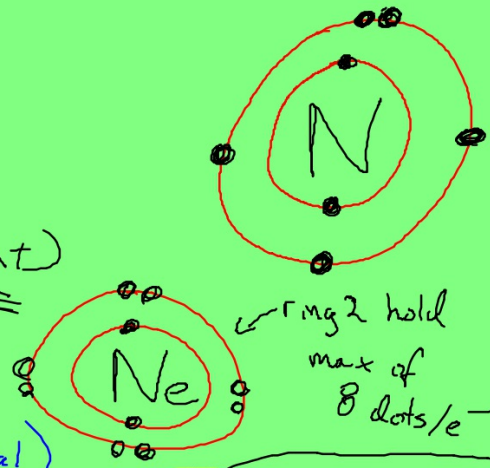
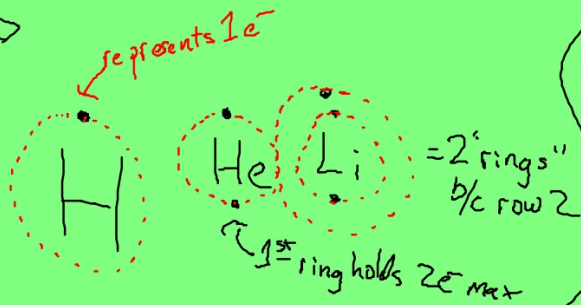
$n \Rightarrow$ isotopes

★ $e^- \Rightarrow$ change easily
to get Energy (heat & light)

How many e^- are in H?

$1e^-$ b/c $1p$ b/c $A\# = 1$ (neutral)

\Rightarrow



THE ELECTRON

- the more e^- = the greater the (-) charge = more Energy
- found in e^- cloud; spin around nucleus (positive core attracts negative cloud)
- move around in orbital \rightarrow set path & set # of e^- on each path

~~movement creates (electricity) ENERGY~~



Electron Orientation in an Atom

Quantum Mechanical Model:

- atom has no definite shape
- increase in # of e^- = increase in size = increase in ENERGY

Orbital Notation

- states location of e^- around the nucleus
- every atom has a different # of e^-

HISTORY~ 4 Models of Atoms

1) Thomson: ball of (+) charges in center with e- outside
--NOT # of p & n, or arrangements in atom

2) Rutherford: showed nucleus had most of atom's mass
--NOT nucleus contains p & n, or (+) attract (-)

3) Bohr: e- travel around nucleus in a definite orbital

-e- do not fall into the nucleus

*fixed paths are named Energy Levels

->increase level = increase E

->to inc level, you move to the next level = Quantum

->all e- not equally spaced apart or spaced from nucleus ...

think escalator

4) **Quantum Mechanical Model**: *includes models 1, 2, 3*

-math equation to tell the location & E of e-

-estimates probability of e- in certain position

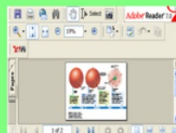
-portrays a "fuzzy cloud" model (when moving)

-> most dense where e- most likely found

***denser around smaller atoms

-specific levels tell location of e-

=> inc level = inc E



Lewis Dot refresher

-> for Group A / Representative Elements (*the peaks*)

Levels & Orientation

E levels (n) = 1, 2, 3, 4, 5, 6, 7 (rows on PT)

-> sublevels = s, p, d, f (specific location & # of e- on PT)

E Levels

Sublevels

Drawing it

add new sublevel
 1 s = hold $2e^-$ max
 add new sublevel
 2 s p = hold $6e^-$ max
 add new sublevel
 3 s p d = hold $10e^-$ max
 4 s p d f = holds $14e^-$ max

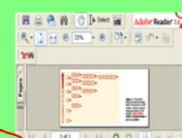
s (1 circle)

p (3 circles)

d (5 circles)

f (7 circles)

each circle can hold 2 arrows max



of circles tell you the sublevel
 # of 1s tell you how far in to go

*when you inc in E level, you add 1 more sublevel

*E levels 5, 6, 7 can have all spdf sublevels

3 Rules to using e- configuration / orbital notation:

1) Aufbau Principle

-e- enter at lowest E level first

* always start with H

2) Pauli Exclusion Principle

-an atomic orbital may have at most 2 e-

* each circle holds 2 e- (↑↓)

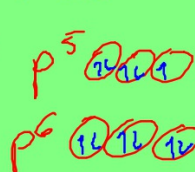
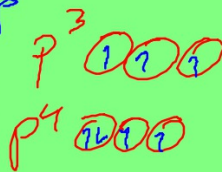
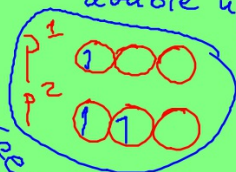
3) Hund's Rule

-when e- occupy orbitals of equal E, 1e- enters each orbital

until all orbitals contain 1e-

* put 1 arrow in each orbital then double up

ex
must keep empty circles



PRACTICE PRACTICE PRACTICE!!!

* f & d region

• always write in order of increasing Atomic Number

If you stop in 4f[#] or 5f[#] region,

you write: ... 5d¹ 4f[#]
(b/c 57 58-71)

... 6d¹ 5f[#]
(b/c 89 90-103)

If you stop in 5d²⁻¹⁰ or 6d²⁻¹⁰,

you write: ... 4f¹⁴ 5d²⁻¹⁰
(b/c 71 72-80)

... 5f¹⁴ 6d²⁻¹⁰
(b/c 103 104-112)

PRACTICE PRACTICE PRACTICE!!!

Exceptions (to Aufbau Rule)
Cu, Cr, and bunch others

DO BY NORMAL RULES!

