

QUANTUM

$E = h\nu$
 $c = \lambda\nu$ } E of photon

VIS λ 400-780nm

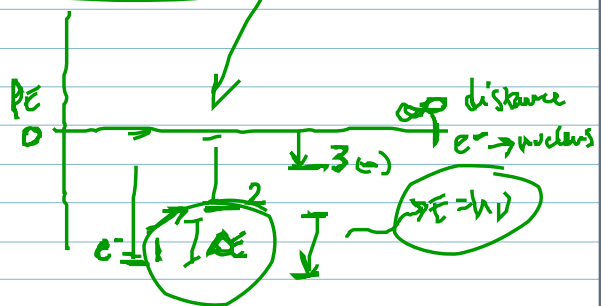
$c = 2.998 \times 10^8 \text{ m/s} = 2.998 \times 10^{17} \text{ nm/s}$
 $1 \times 10^9 \text{ nm} = 1 \text{ m}$

$F = \frac{kQ_1Q_2}{r^2}$

e^- 's closer to the nucleus experience a greater attractive force

Planck
Bohr

$E_n = \frac{-2.178 \times 10^{-18} \text{ J}}{n^2}$



e^- 's further from the nucleus have a less (-) E "higher E"

it takes less E to remove an e^- further from nucleus (IE, AES)

e^- configurations

Heisenberg \rightarrow uncertainty

Pauli \Rightarrow max 2 e^- /orbital ; opposite spin $\rightarrow \begin{matrix} +\frac{1}{2} \\ \uparrow \\ -\frac{1}{2} \\ \downarrow \end{matrix}$

Hund's Rule \Rightarrow 1 e^- /orbital in a set before second minimize e^-/e^- repulsion \downarrow E

Spectroscopy \Rightarrow orbitals / probabilities
 e^- E levels \rightarrow sublevels (orbitals)

Aufbau \rightarrow "grows up" lowest E fills first

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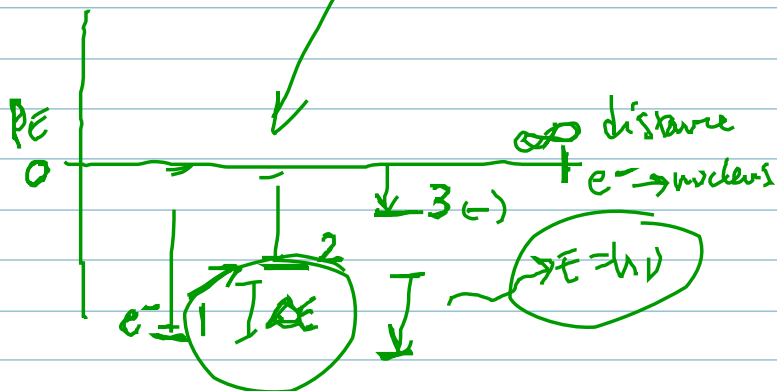
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→ Planck
→ Bohr

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e^- 's further from the nucleus have a less (-) E "higher E"

⇒ it takes less E to remove an e^- further from nucleus (IE, PES)

e^- configurations

Heisenberg → Uncertainty

Pauli ⇒ max 2 e^- /orbital ; opposite spin → $+\frac{1}{2}, -\frac{1}{2}$
↑ ↓

Mund's Rule ⇒ 1 e^- /orbital in a set before second
minimize e^-/e^- repulsion ↓ E

Schrodinger ⇒ orbitals / probabilities

e^- E levels → sublevels (orbitals)

Aufbau → "grows up" lowest E fills first

Fe [Ar] $(4s^2) 3d^6$ ^{valence} ground state

[Ar] $(4s^1) 3d^7$ "excited state"

Fe²⁺ [Ar] $3d^6$ always lose (s) before d 's } f. metals

Fe³⁺ [Ar] $(3d^5)$

paramagnetic / diamagnetic
 ✓ / ~~⊗~~

unpaired e's $3d^6$ $\uparrow\downarrow$ $(\uparrow \uparrow \uparrow \uparrow)$ $3d$

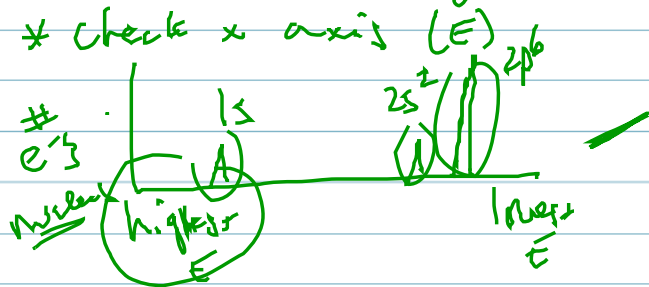
SPECTROSCOPY (V's)

$\uparrow V \rightarrow \downarrow V$

X-rays
 $\uparrow\downarrow$

P.E.S. \Rightarrow

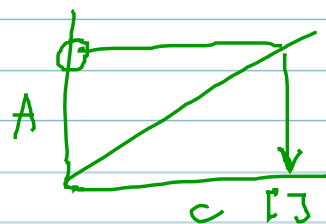
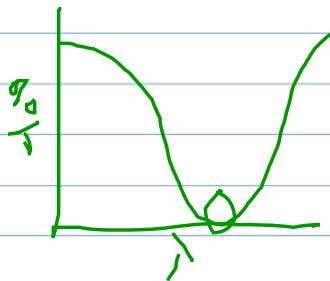
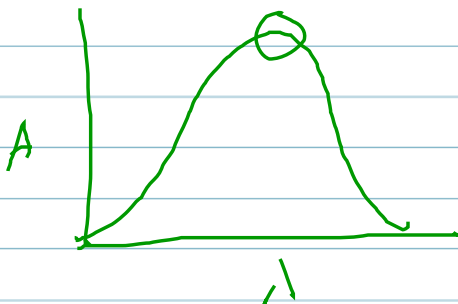
electronic arrangement transitions



vis \Rightarrow colorimetry
 Beer's Law

$A = abc$ $A \propto c$

COLORED SOLUTIONS
 + transition metals

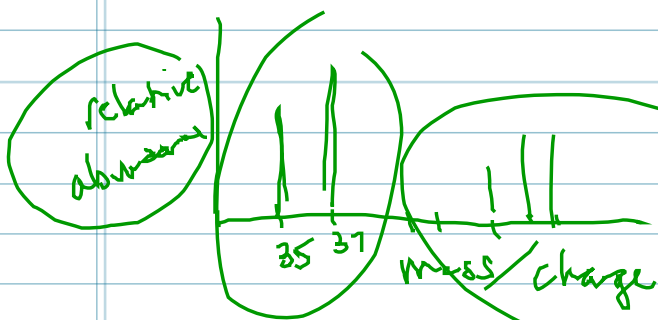


~~Radio~~ microwave \rightarrow rotations
 IR \rightarrow vibrations

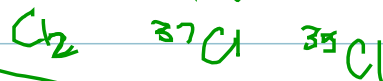
} Structure

MASS SPECTROSCOPY

\bar{x} atomic mass of element



* Watch for diatomics



$35+2 \rightarrow 70$

$35+37 \rightarrow 72$

~~37~~ $37+2 \rightarrow 74$

$\begin{matrix} \text{isotope} \\ \text{mass} \end{matrix} \times \begin{matrix} \% \text{ abundance} \end{matrix} =$

" " =

" " =

+ average atomic mass
amu