

ELECTRONIC CONFIGURATIONS

ELECTRONIC CONFIGURATIONS

CONTENTS

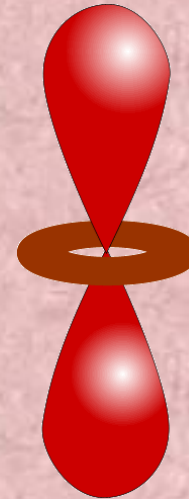
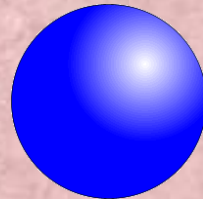
- **The Bohr Atom**
- **Levels and sub-levels**
- **Rules and principles**
- **Orbitals**
- **Rules for filling orbitals.**
- **The Aufbau principle**
- **Electronic configurations of elements 1 to 36**
- **Electronic configurations of ions**



ELECTRONIC CONFIGURATIONS

Before you start it would be helpful to...

- Know that electrons can be found outside the nucleus in energy levels (shells)
- Know the electronic configurations of the first 20 elements in 2,8,1 notation



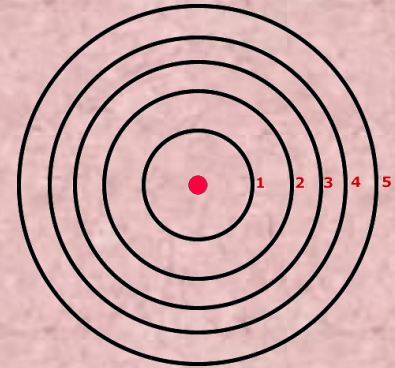
THE BOHR ATOM

Ideas about the structure of the atom have changed over the years. The **Bohr** theory thought of it as a small nucleus of protons and neutrons surrounded by circulating electrons.

Each shell or energy level could hold a maximum number of electrons.

The energy of levels became greater as they got further from the nucleus and electrons filled energy levels in order.

The theory couldn't explain certain aspects of chemistry.



Maximum electrons
per shell

1st shell **2**

2nd shell **8**

3rd shell **18**

4th shell **32**

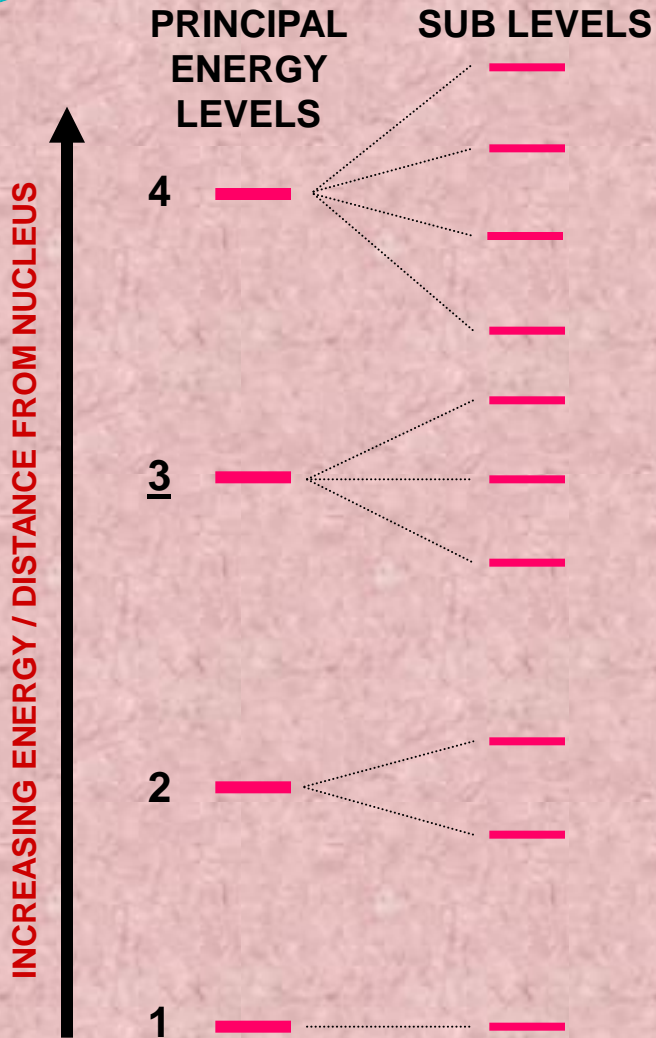
5th shell **50**

LEVELS AND SUB-LEVELS



The **energy gap between successive levels** got increasingly smaller as the **levels got further from the nucleus**. The importance of this is discussed later.

LEVELS AND SUB-LEVELS



A study of Ionisation Energies and the periodic properties of elements suggested that the **main energy levels were split into sub levels.**

Level 1 was split into 1 sub level

Level 2 was split into 2 sub levels

Level 3 was split into 3 sub levels

Level 4 was split into 4 sub levels

RULES AND PRINCIPLES

HEISENBERG'S UNCERTAINTY PRINCIPLE

“You cannot determine the position and momentum of an electron at the same time.”

This means that you cannot say *exactly* where an electron is.

THE AUFBAU PRINCIPLE

“Electrons enter the lowest available energy level.”

PAULI'S EXCLUSION PRINCIPLE

“No two electrons can have the same four quantum numbers.”

Two electrons can go in each orbital, providing they are of opposite spin.

HUND'S RULE OF MAXIMUM MULTIPLICITY

“When in orbitals of equal energy, electrons will try to remain unpaired.”

Placing two electrons in one orbital means that, as they are both negatively charged, there will be some electrostatic repulsion between them. Placing each electron in a separate orbital reduces the repulsion and the system is more stable. It can be described as the “SITTING ON A BUS RULE”!

ORBITALS

An orbital is... **a region in space where one is likely to find an electron.**

Orbitals **can hold up to two electrons as long as they have opposite spin**; this is known as PAULI'S EXCLUSION PRINCIPAL.

Orbitals have different shapes...



ORBITALS

An orbital is... **a region in space where one is likely to find an electron.**

Orbitals can hold up to two electrons as long as they have opposite spin; this is known as PAULI'S EXCLUSION PRINCIPAL.

Orbitals have different shapes...

ORBITAL	SHAPE	OCCURRENCE
s	spherical	one in every principal level
p	dumb-bell	three in levels from 2 upwards
d	various	five in levels from 3 upwards
f	various	seven in levels from 4 upwards



ORBITALS

An orbital is... **a region in space where one is likely to find an electron.**

Orbitals can hold up to two electrons as long as they have opposite spin; this is known as PAULI'S EXCLUSION PRINCIPAL.

Orbitals have different shapes...

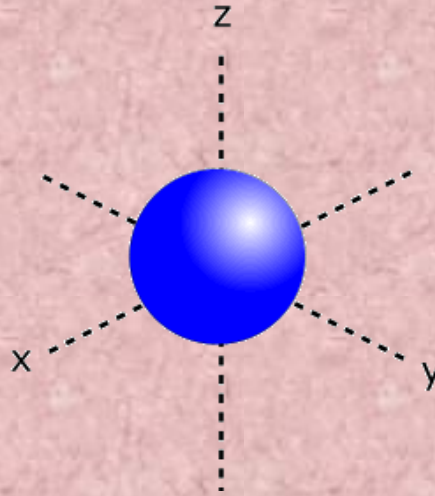
ORBITAL	SHAPE	OCCURRENCE
s	spherical	one in every principal level
p	dumb-bell	three in levels from 2 upwards
d	various	five in levels from 3 upwards
f	various	seven in levels from 4 upwards

An orbital is a 3-dimensional statistical shape showing where one is most likely to find an electron. Because, according to Heisenberg, you cannot say exactly where an electron is you are only able to say where it might be found.

DO NOT CONFUSE AN ORBITAL WITH AN ORBIT

SHAPES OF ORBITALS

s orbitals

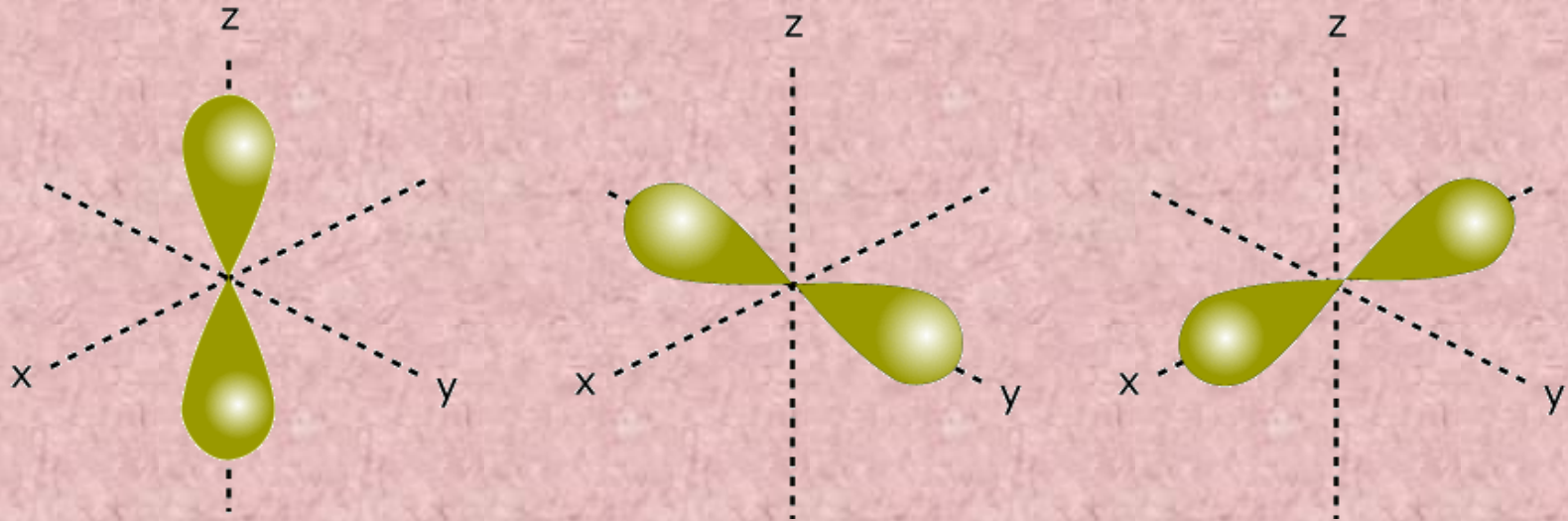


- spherical
- one occurs in every principal energy level



SHAPES OF ORBITALS

p orbitals

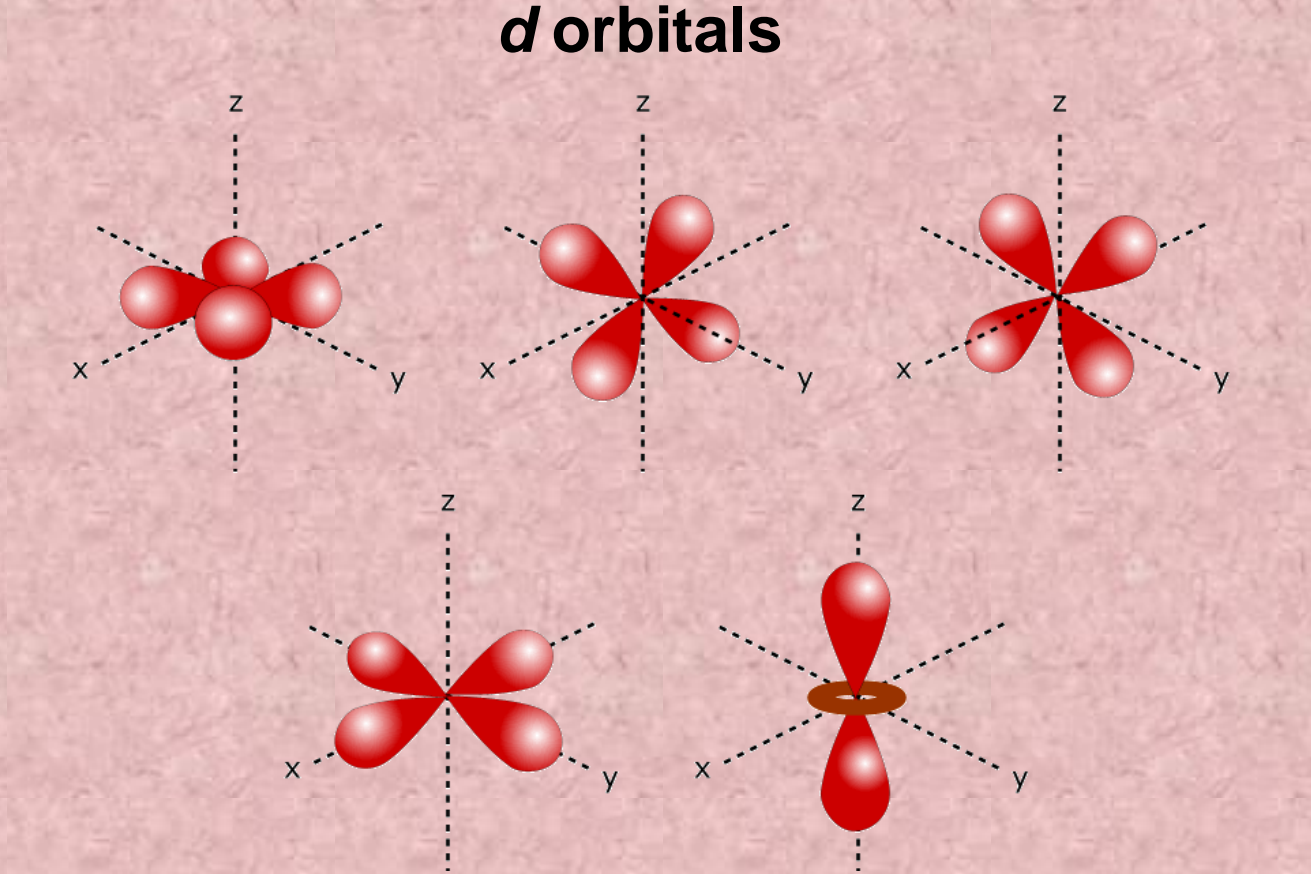


- dumb-bell shaped
- three occur in energy levels except the first



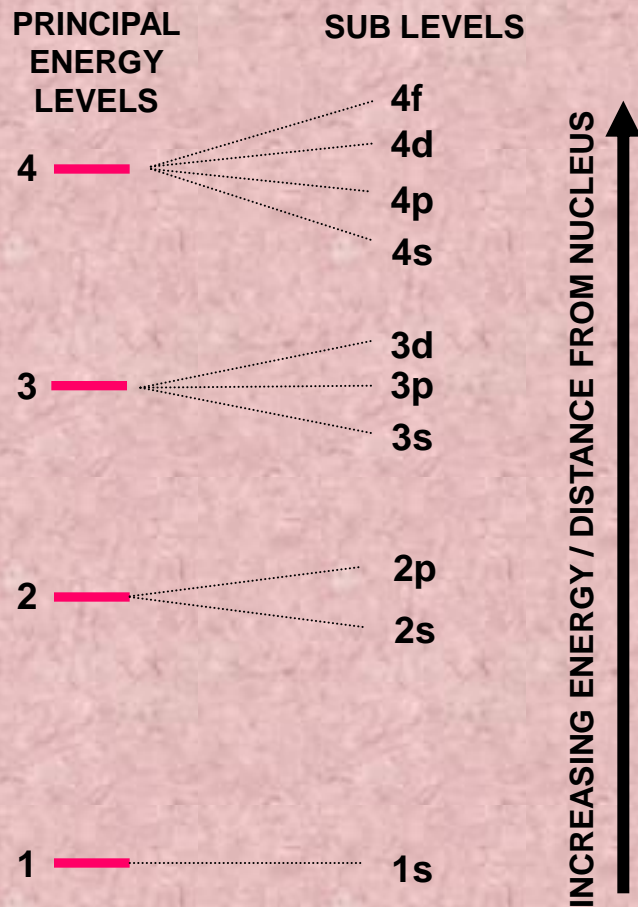
SHAPES OF ORBITALS

d orbitals



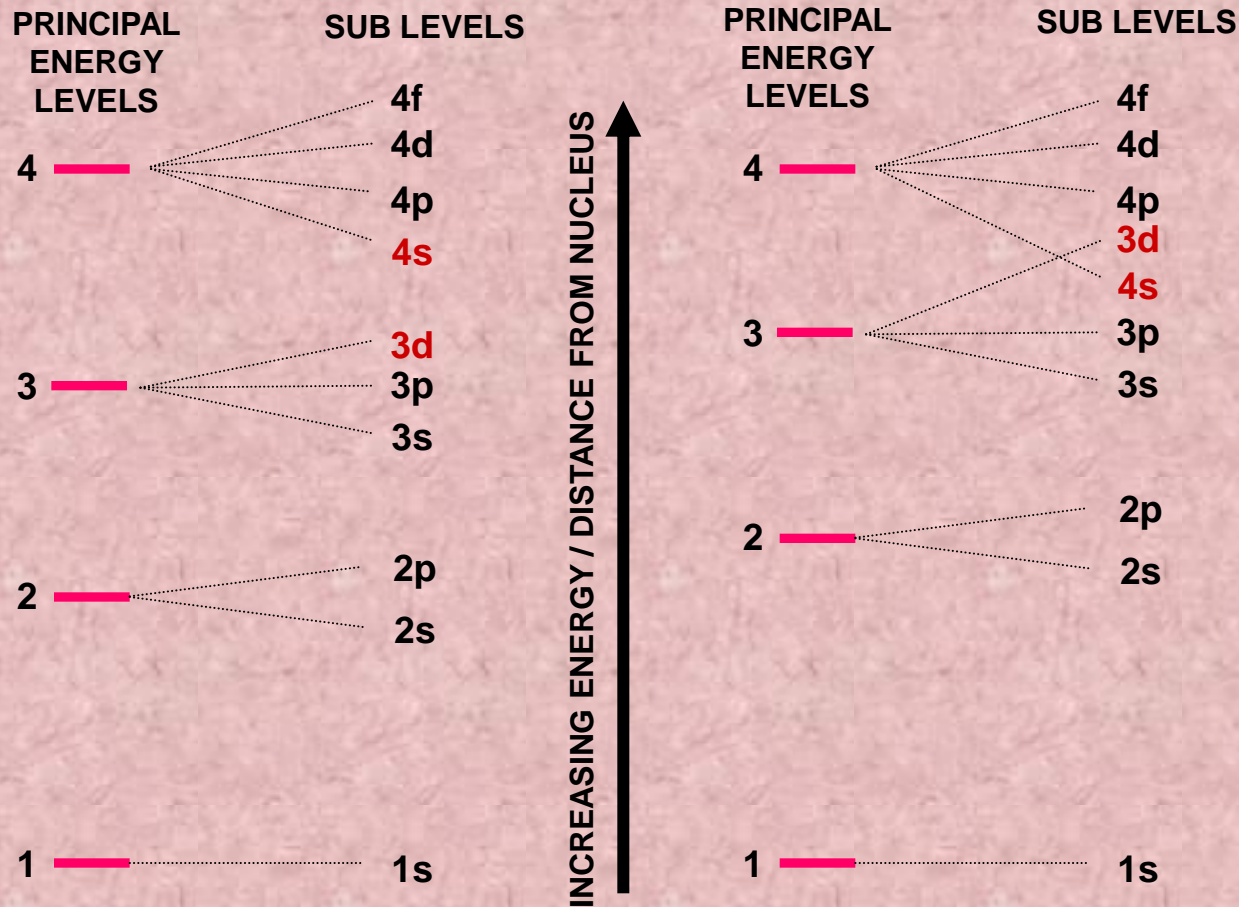
- various shapes
- five occur in energy levels except the first and second

ORDER OF FILLING ORBITALS



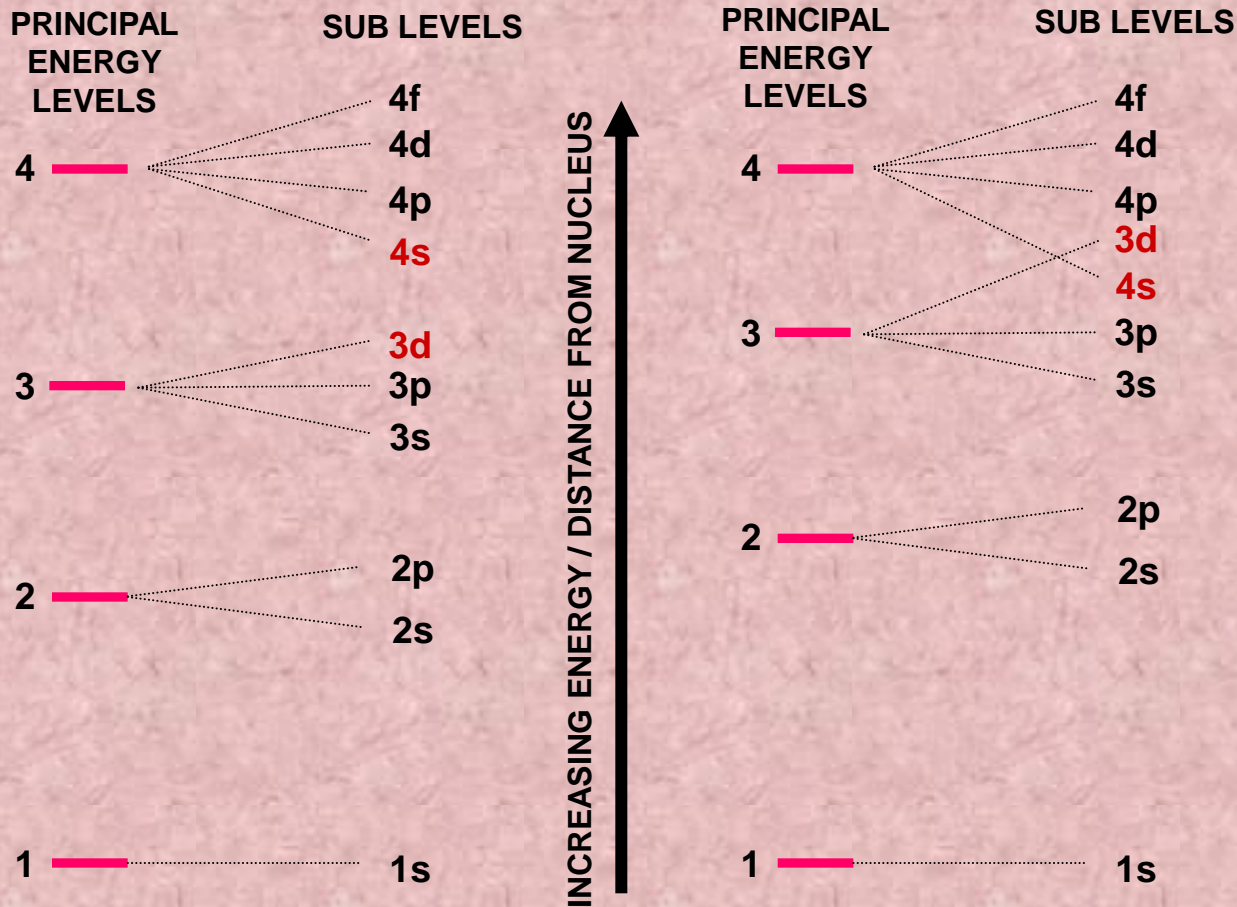
Orbitals are not filled in numerical order because the principal energy levels get closer together as you get further from the nucleus. This results in overlap of sub levels. The first example occurs when the **4s orbital is filled before the 3d orbitals**.

ORDER OF FILLING ORBITALS



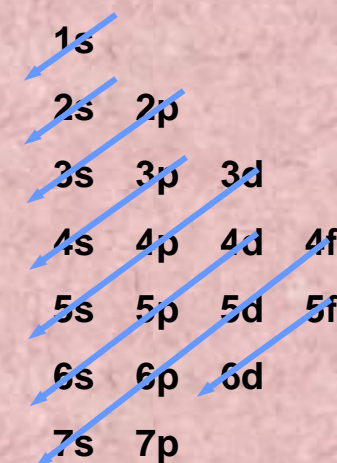
Orbitals are not filled in numerical order because the principal energy levels get closer together as you get further from the nucleus. This results in overlap of sub levels. The first example occurs when the **4s orbital is filled before the 3d orbitals**.

ORDER OF FILLING ORBITALS



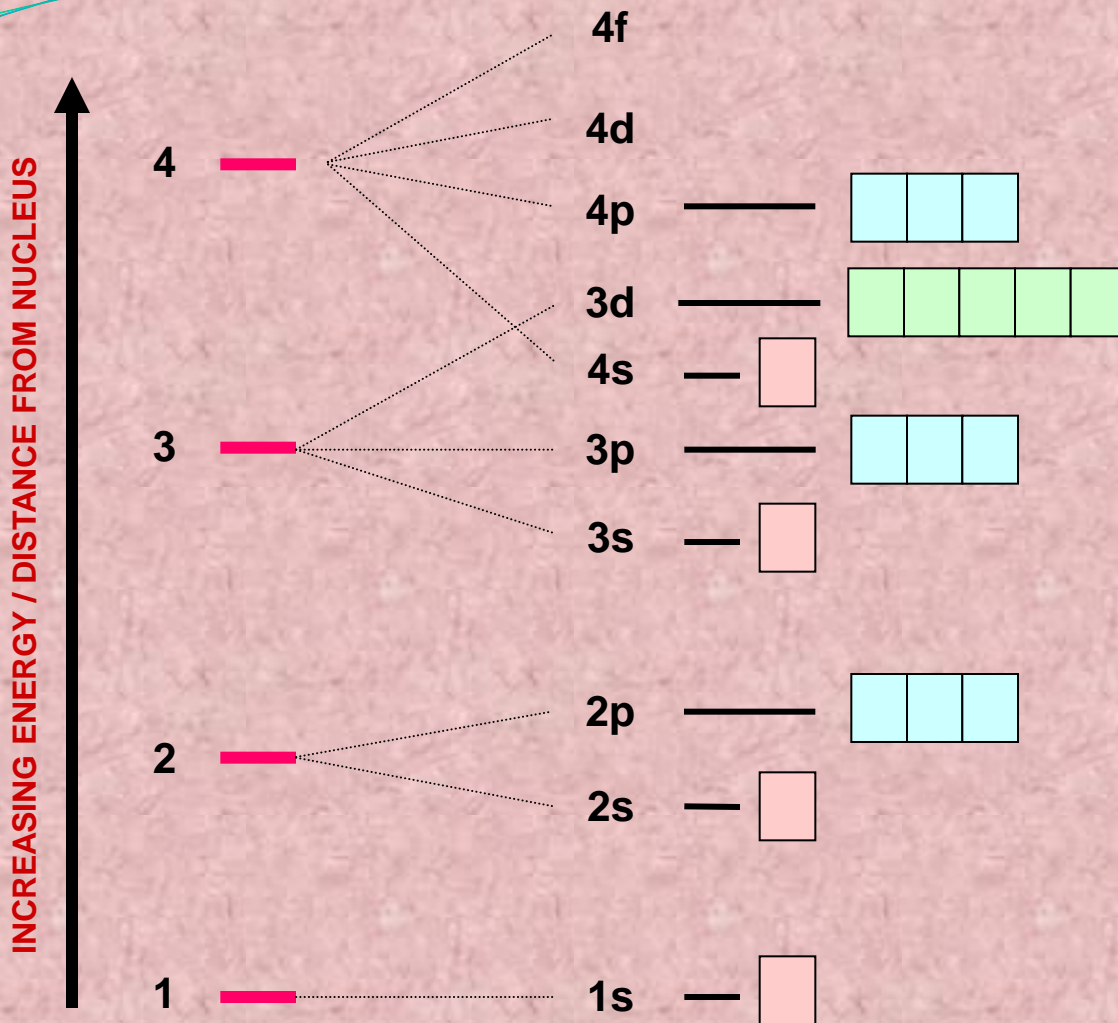
HOW TO REMEMBER ...

THE FILLING ORDER



Orbitals are not filled in numerical order because the principal energy levels get closer together as you get further from the nucleus. This results in overlap of sub levels. The first example occurs when the **4s orbital is filled before the 3d orbitals**.

THE 'AUFBAU' PRINCIPAL



This states that...

**“ELECTRONS ENTER THE
LOWEST AVAILABLE
ENERGY LEVEL”**

The following sequence will show the ‘building up’ of the electronic structures of the first 36 elements in the periodic table.

Electrons are shown as half headed arrows and can spin in one of two directions

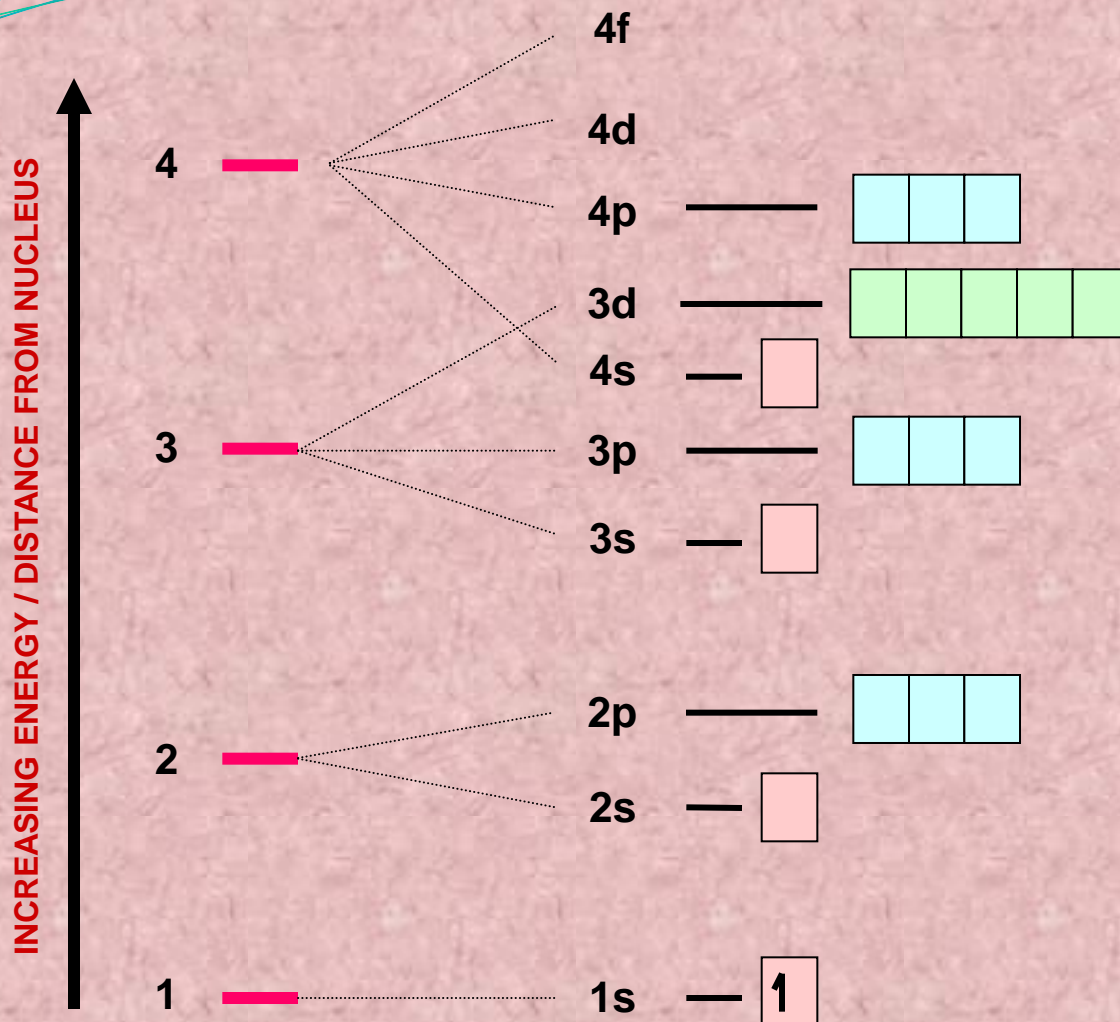
↑ or ↓

s orbitals

p orbitals

d orbitals

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



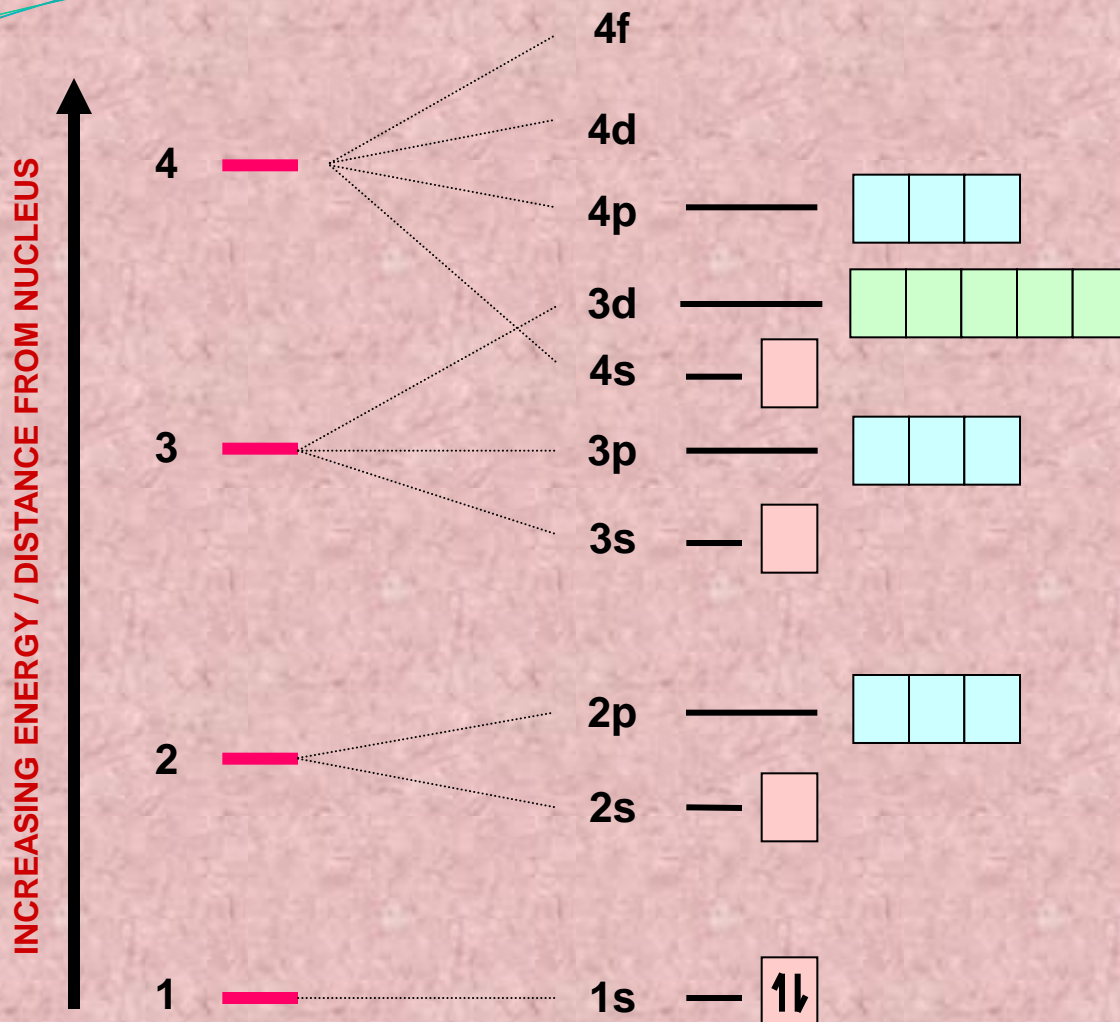
HYDROGEN



Hydrogen atoms have one electron. This goes into a vacant orbital in the lowest available energy level.

**'Aufbau'
Principle**

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



HELIUM



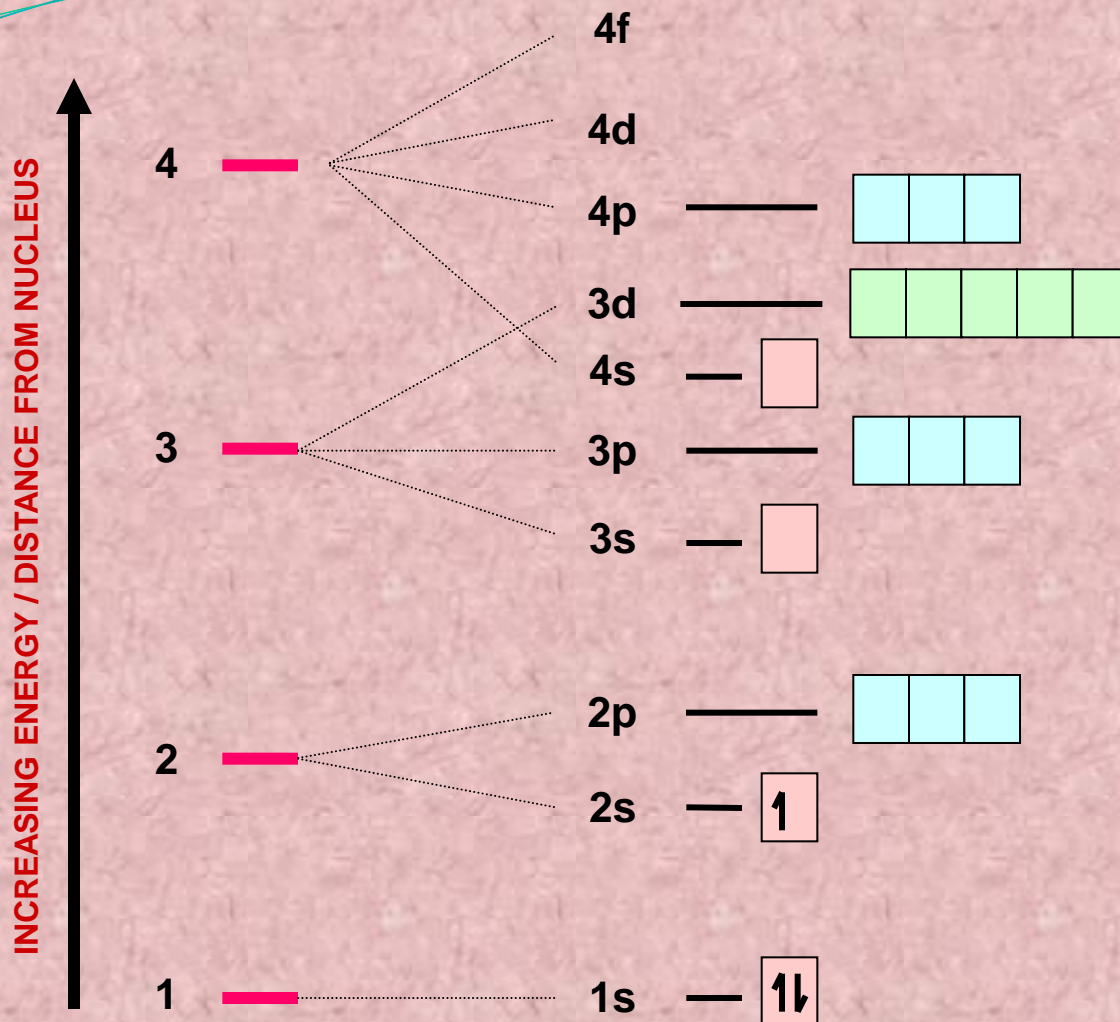
Every orbital can contain 2 electrons, provided the electrons are spinning in opposite directions. This is based on...

PAULI'S EXCLUSION PRINCIPLE

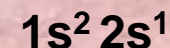
The two electrons in a helium atom can both go in the 1s orbital.

'Aufbau' Principle

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



LITHIUM

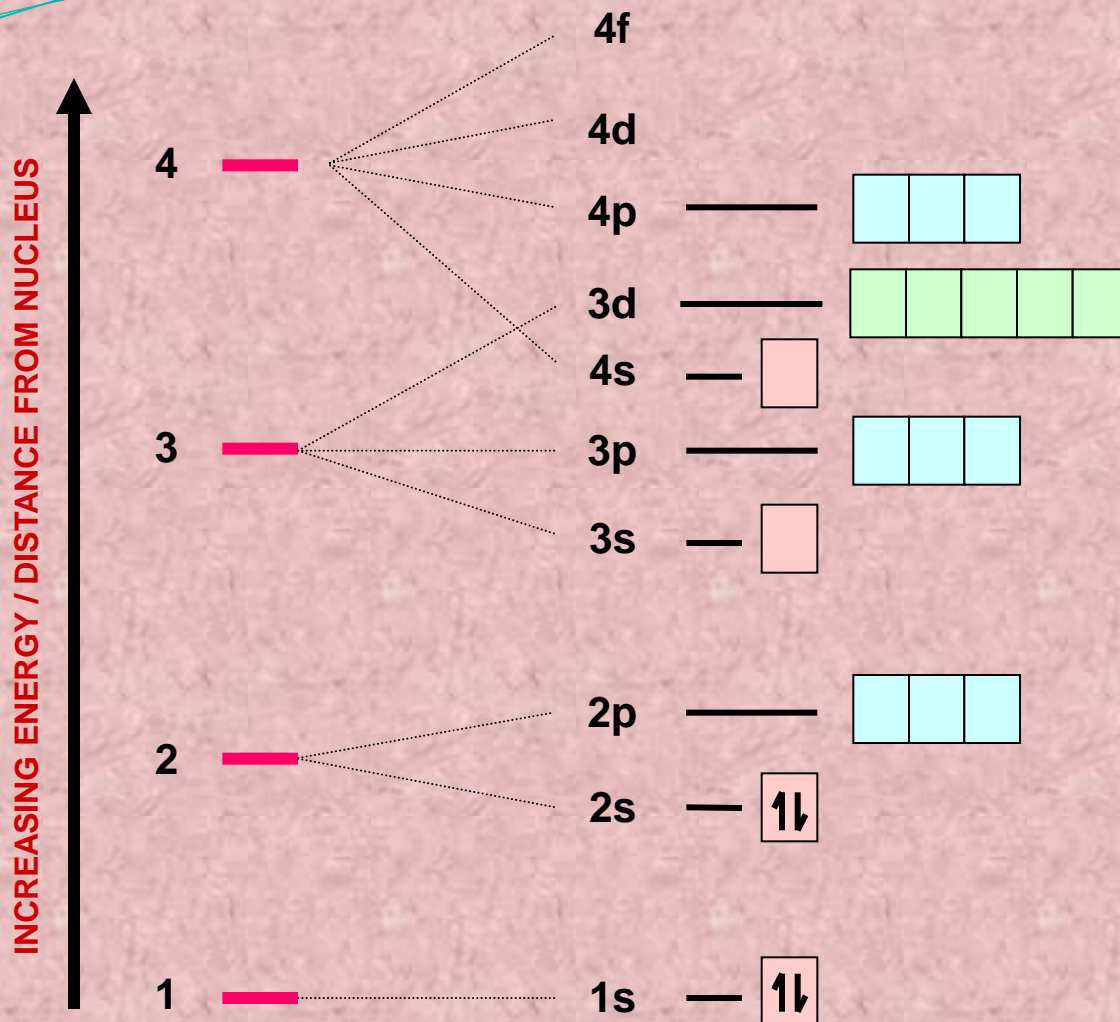


1s orbitals can hold a maximum of two electrons so the third electron in a lithium atom must go into the next available orbital of higher energy. This will be further from the nucleus in the second principal energy level.

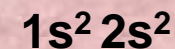
The second principal level has two types of orbital (s and p). An s orbital is lower in energy than a p.

**'Aufbau'
Principle**

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



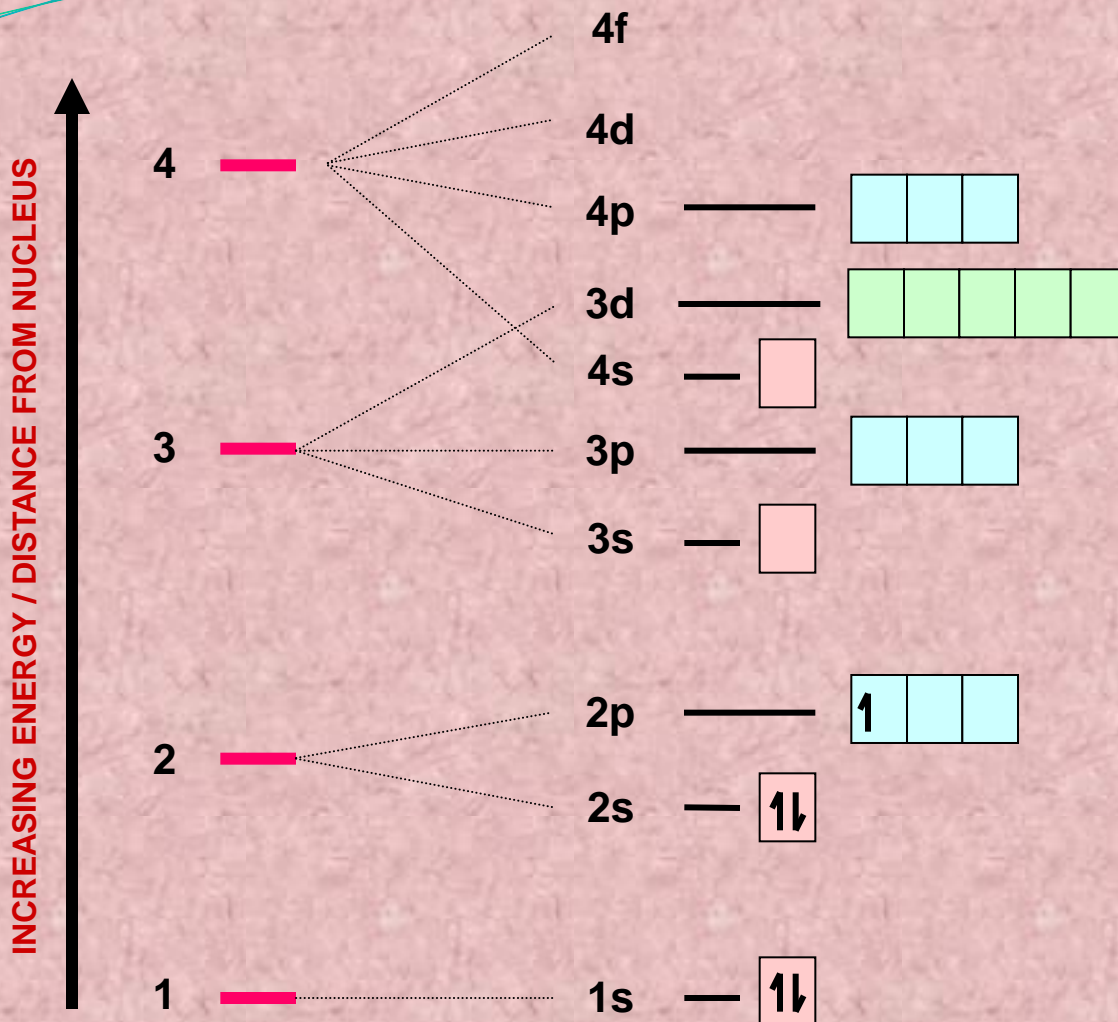
BERYLLIUM



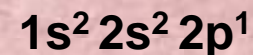
Beryllium atoms have four electrons so the fourth electron pairs up in the 2s orbital. The 2s sub level is now full.

'Aufbau'
Principle

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



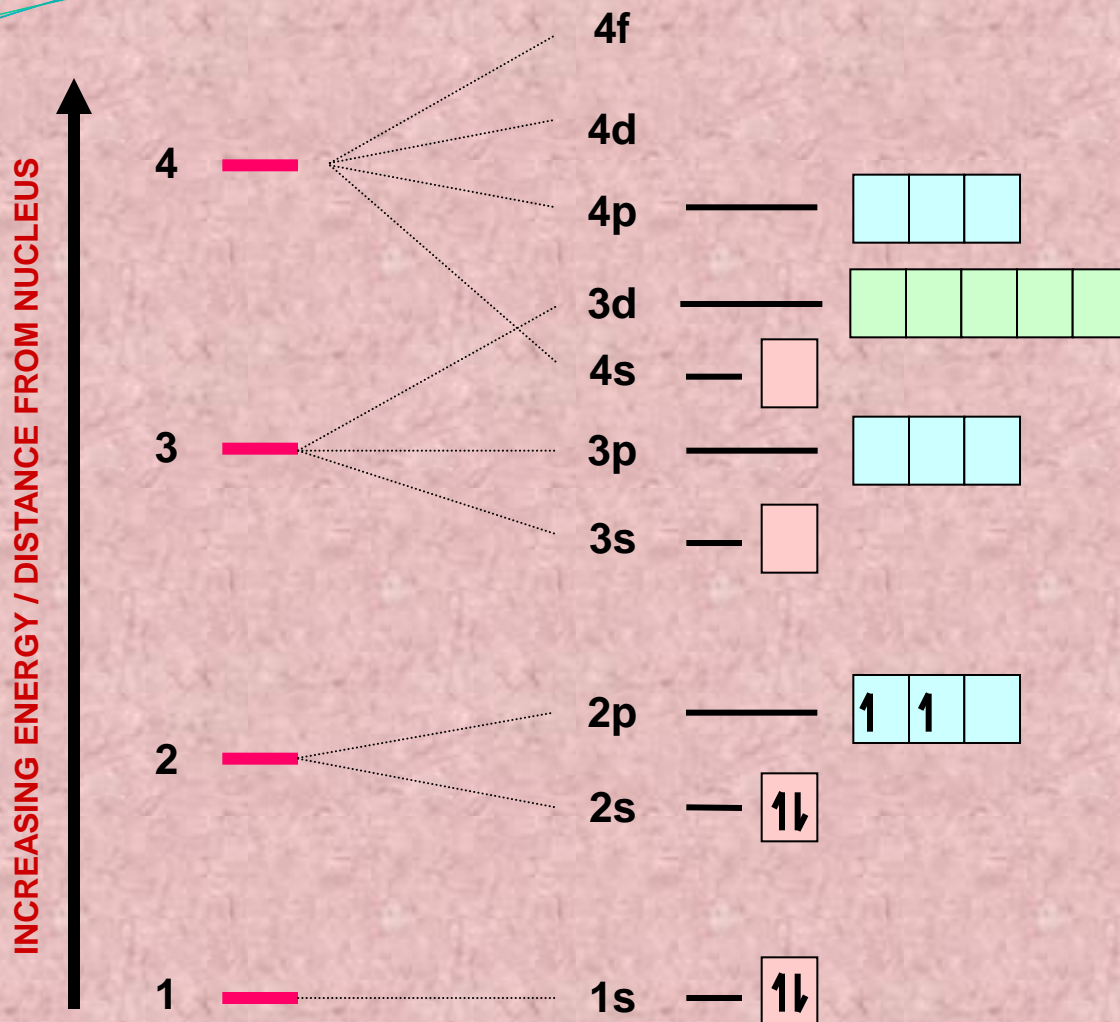
BORON



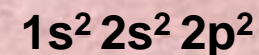
As the 2s sub level is now full, the fifth electron goes into one of the three p orbitals in the 2p sub level. The 2p orbitals are slightly higher in energy than the 2s orbital.

**'Aufbau'
Principle**

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



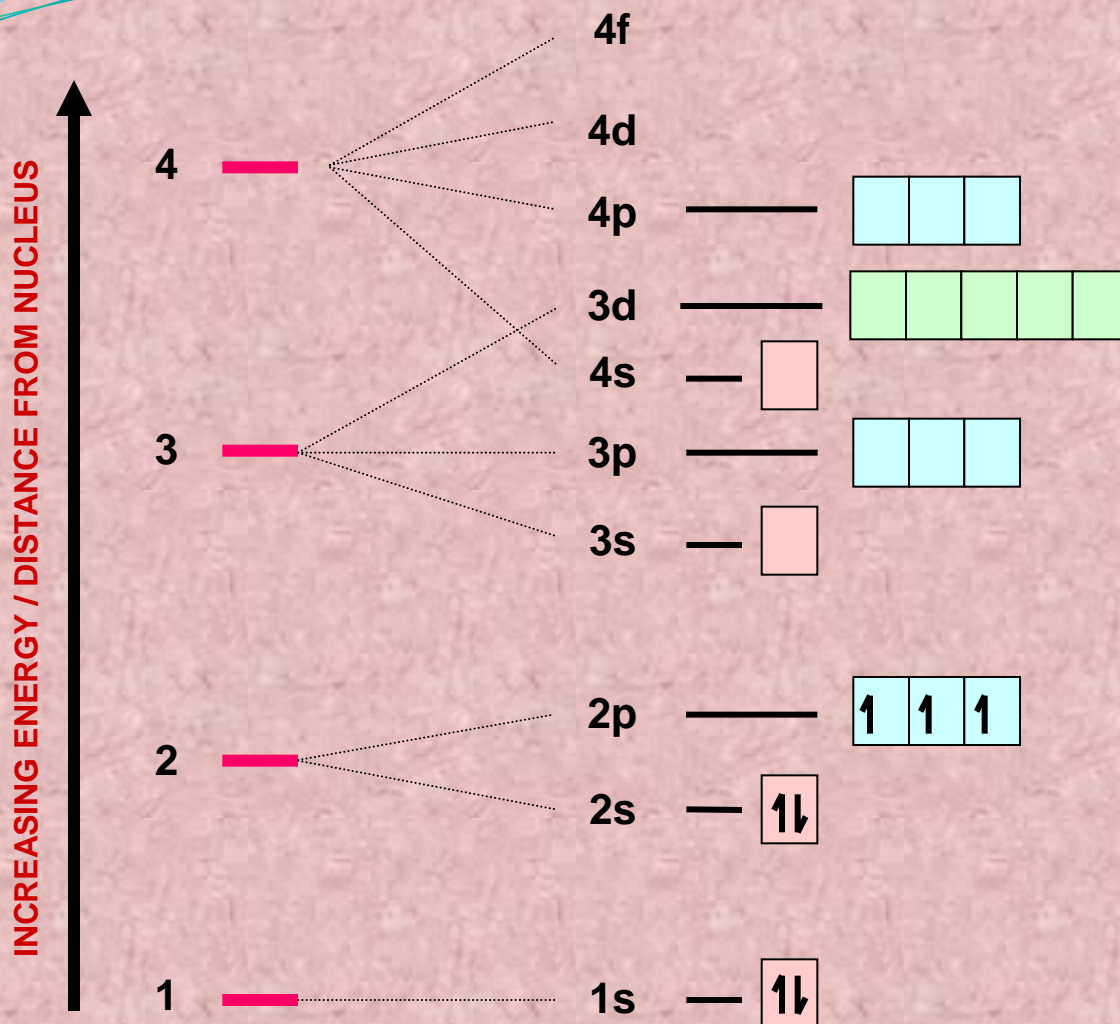
CARBON



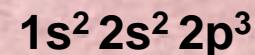
The next electron in doesn't pair up with the one already there. This would give rise to **repulsion between the similarly charged species**. Instead, it goes into another p orbital which means **less repulsion, lower energy and more stability**.

**HUND'S RULE
OF
MAXIMUM MULTIPLICITY**

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



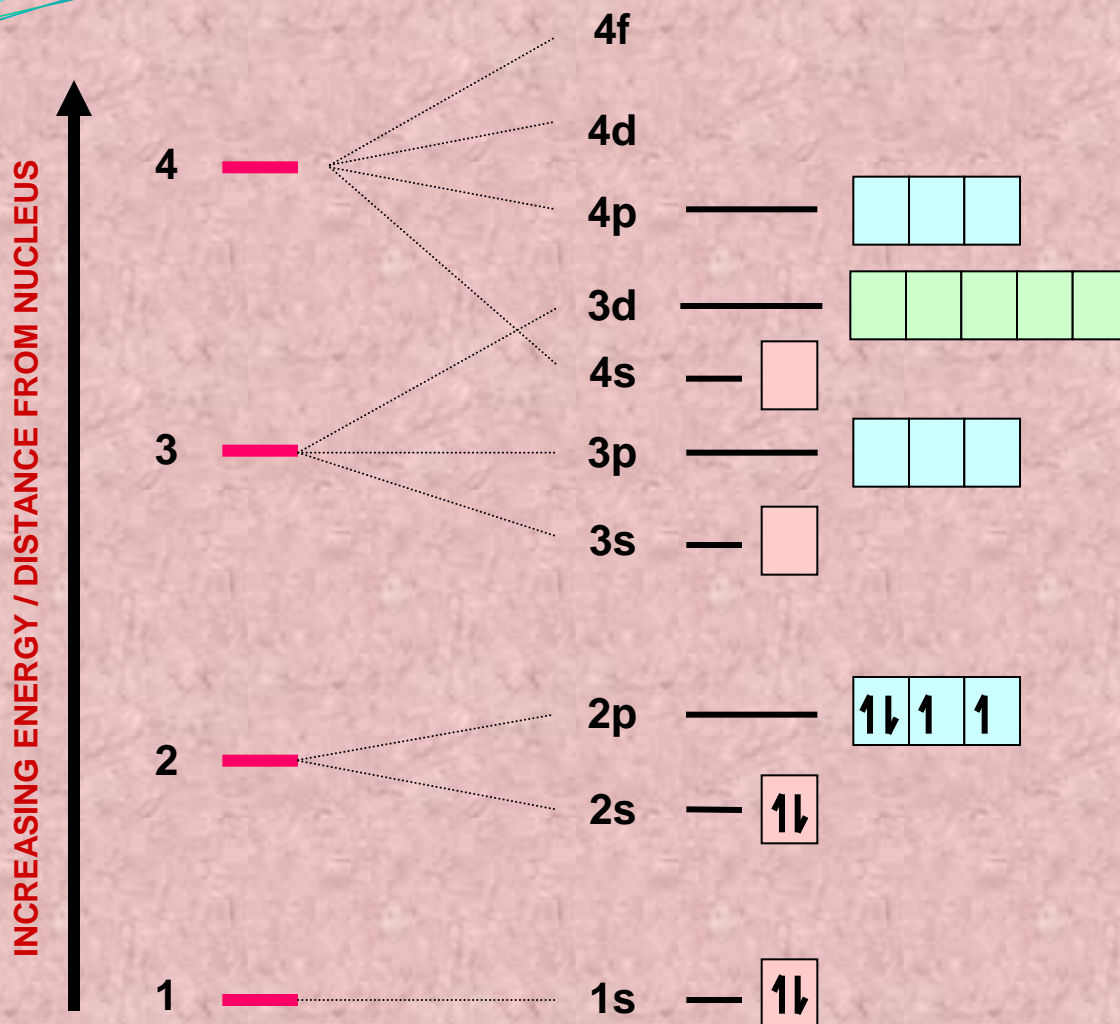
NITROGEN



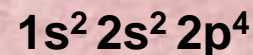
Following Hund's Rule, the next electron will not pair up so goes into a vacant p orbital. All three electrons are now unpaired. This gives **less repulsion, lower energy and therefore more stability.**

HUND'S RULE
OF
MAXIMUM MULTIPLICITY

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



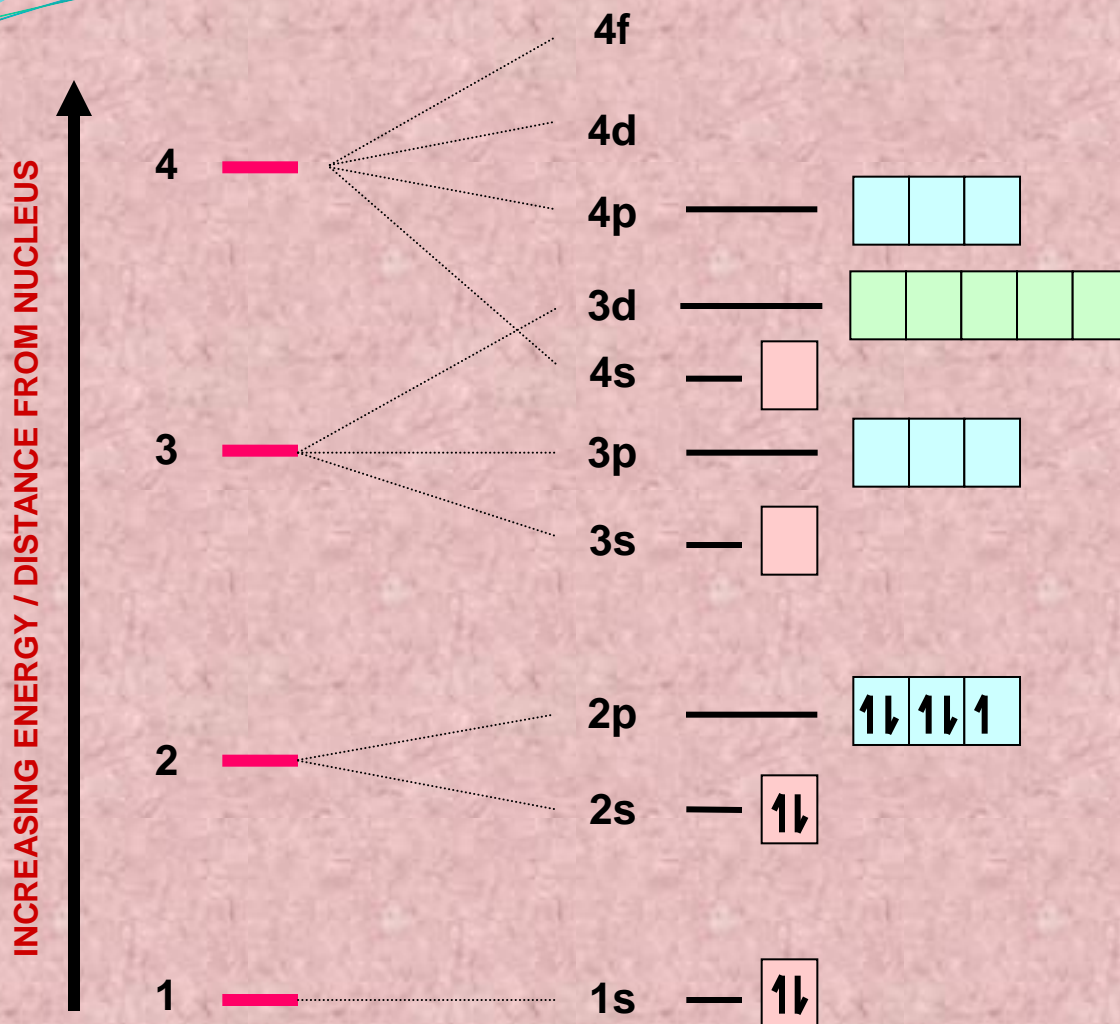
OXYGEN



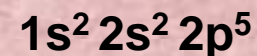
With all three orbitals half-filled, the eighth electron in an oxygen atom must now pair up with one of the electrons already there.

'Aufbau' Principle

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS

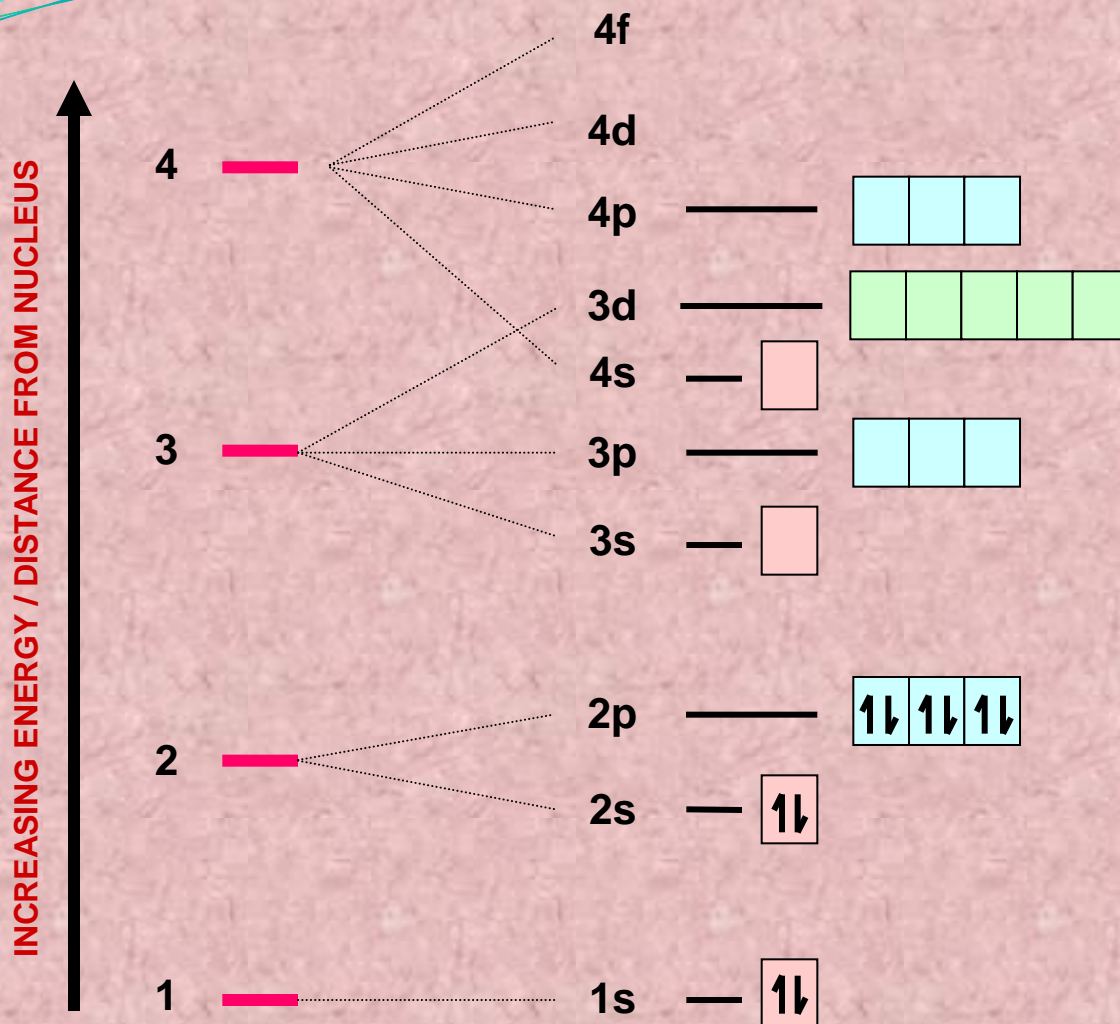


FLUORINE

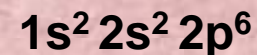


The electrons continue to pair up with those in the half-filled orbitals.

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



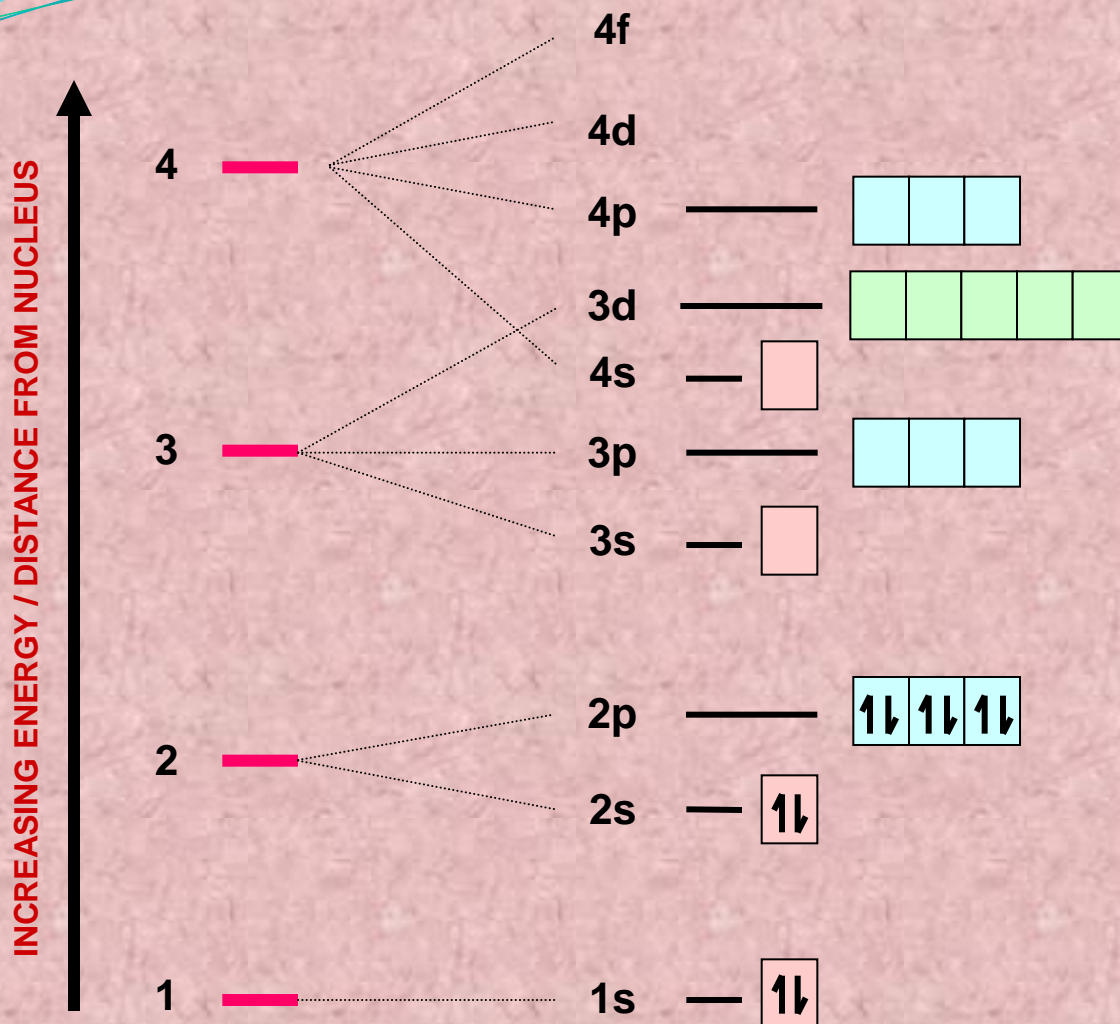
NEON



The electrons continue to pair up with those in the half-filled orbitals. The 2p orbitals are now completely filled and so is the second principal energy level.

In the older system of describing electronic configurations, this would have been written as 2,8.

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



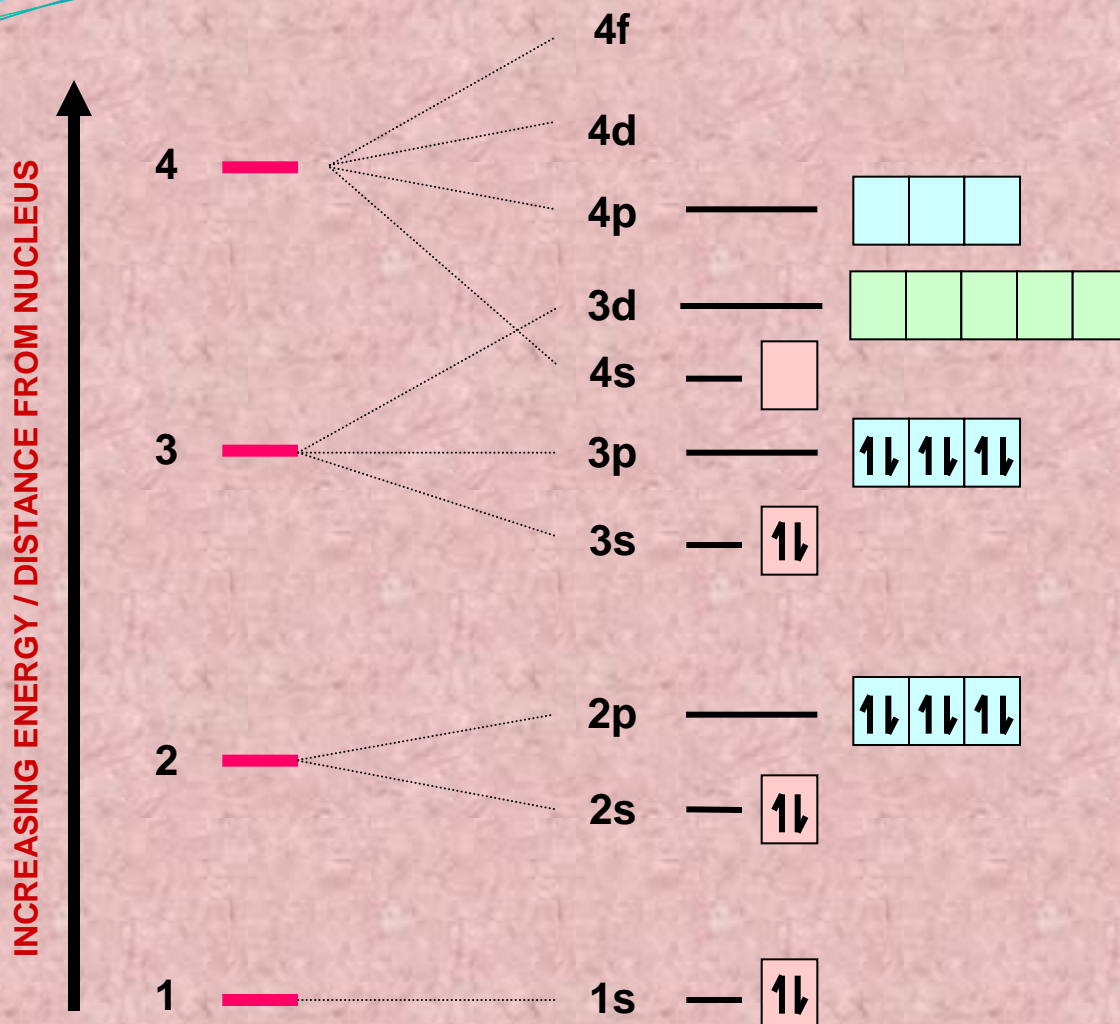
SODIUM - ARGON

With the second principal energy level full, the next electrons must go into the next highest level. The third principal energy level contains three types of orbital; s, p and d.

The 3s and 3p orbitals are filled in exactly the same way as those in the 2s and 2p sub levels.

**'Aufbau'
Principle**

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS

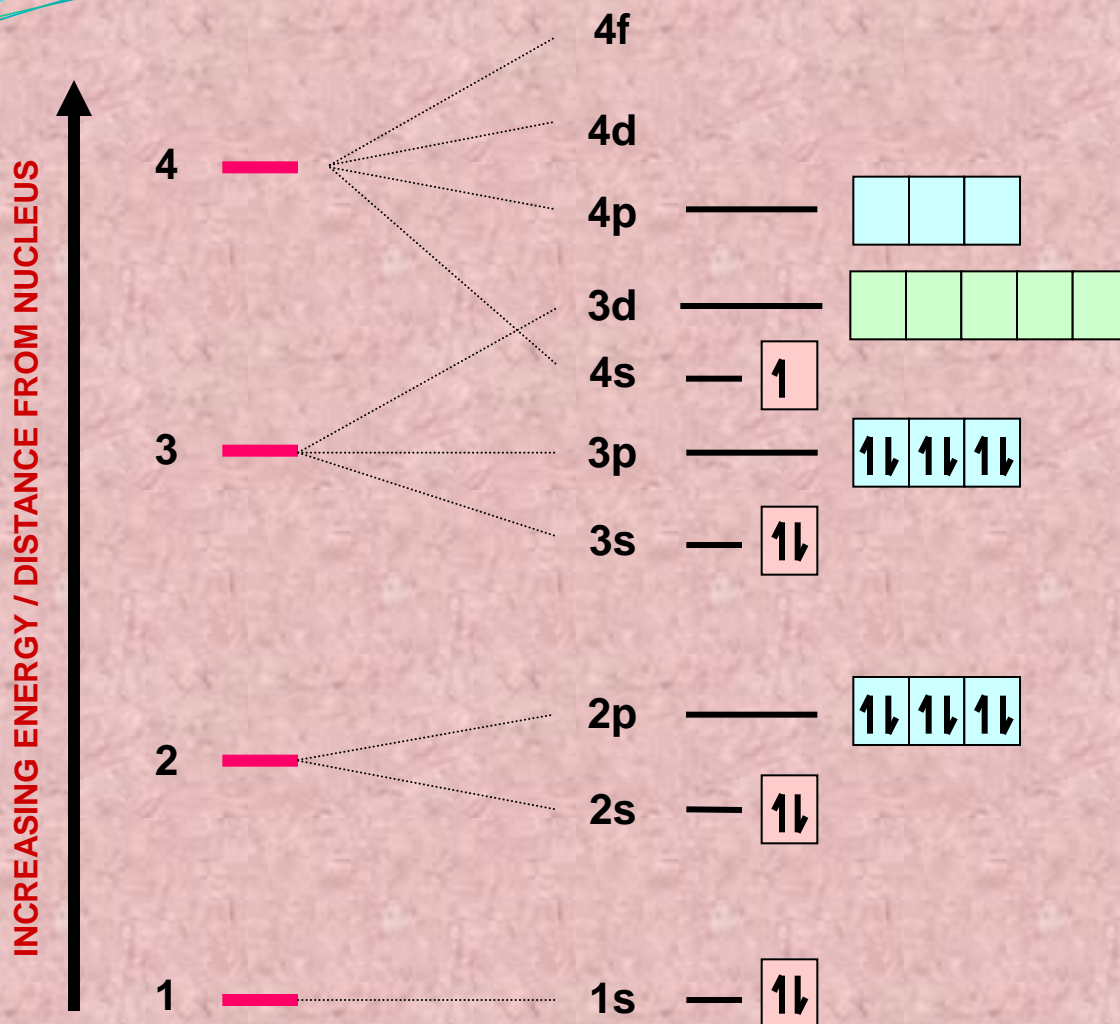


SODIUM - ARGON

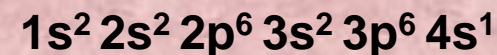
Na	$1s^2 2s^2 2p^6 3s^1$
Mg	$1s^2 2s^2 2p^6 3s^2$
Al	$1s^2 2s^2 2p^6 3s^2 3p^1$
Si	$1s^2 2s^2 2p^6 3s^2 3p^2$
P	$1s^2 2s^2 2p^6 3s^2 3p^3$
S	$1s^2 2s^2 2p^6 3s^2 3p^4$
Cl	$1s^2 2s^2 2p^6 3s^2 3p^5$
Ar	$1s^2 2s^2 2p^6 3s^2 3p^6$

Remember that the 3p configurations follow Hund's Rule with the electrons remaining unpaired to give more stability.

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



POTASSIUM

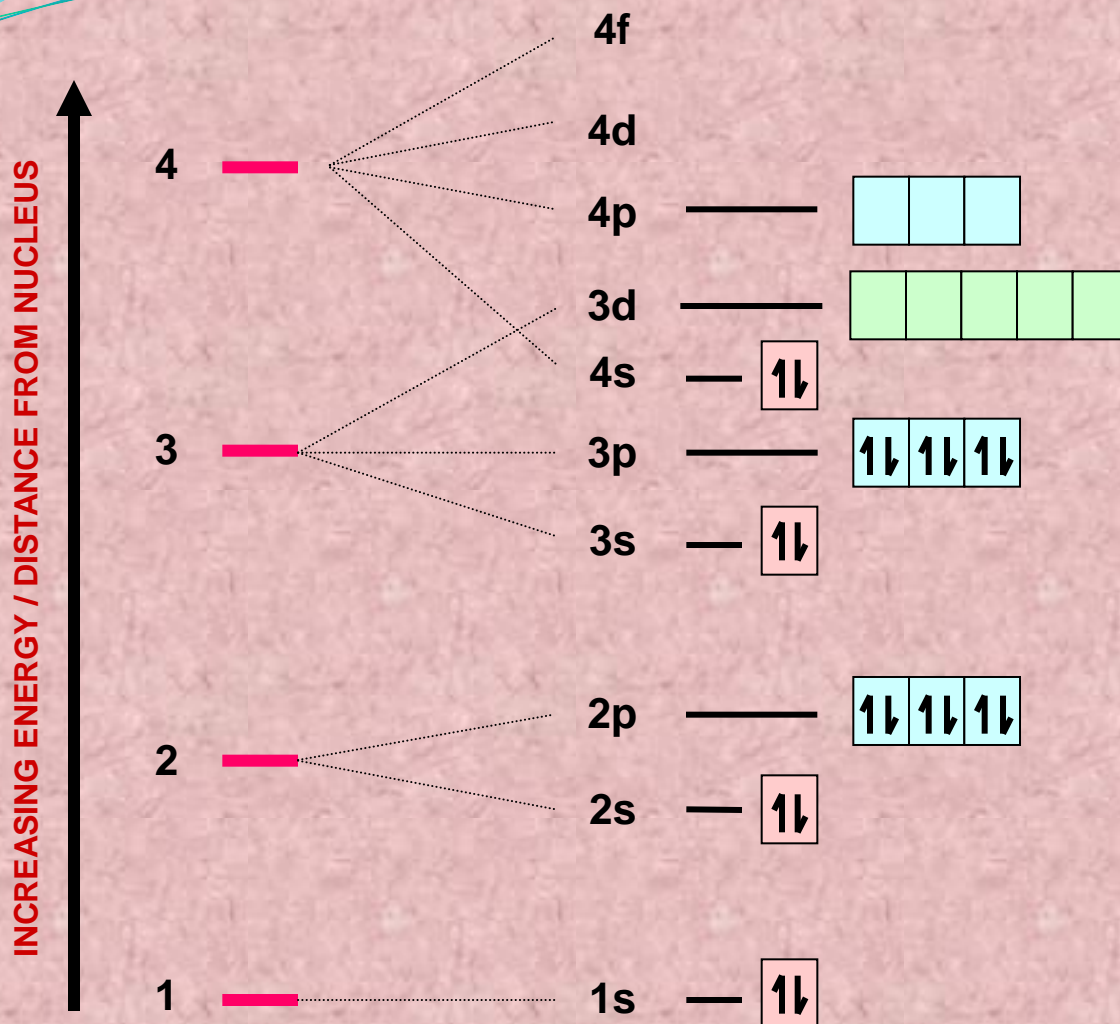


In numerical terms one would expect the 3d orbitals to be filled next.

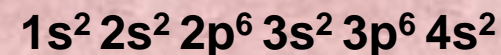
However, because the **principal energy levels get closer together as you go further from the nucleus** coupled with the **splitting into sub energy levels**, the **4s orbital is of a LOWER ENERGY than the 3d orbitals** so gets filled first.

**'Aufbau'
Principle**

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



CALCIUM

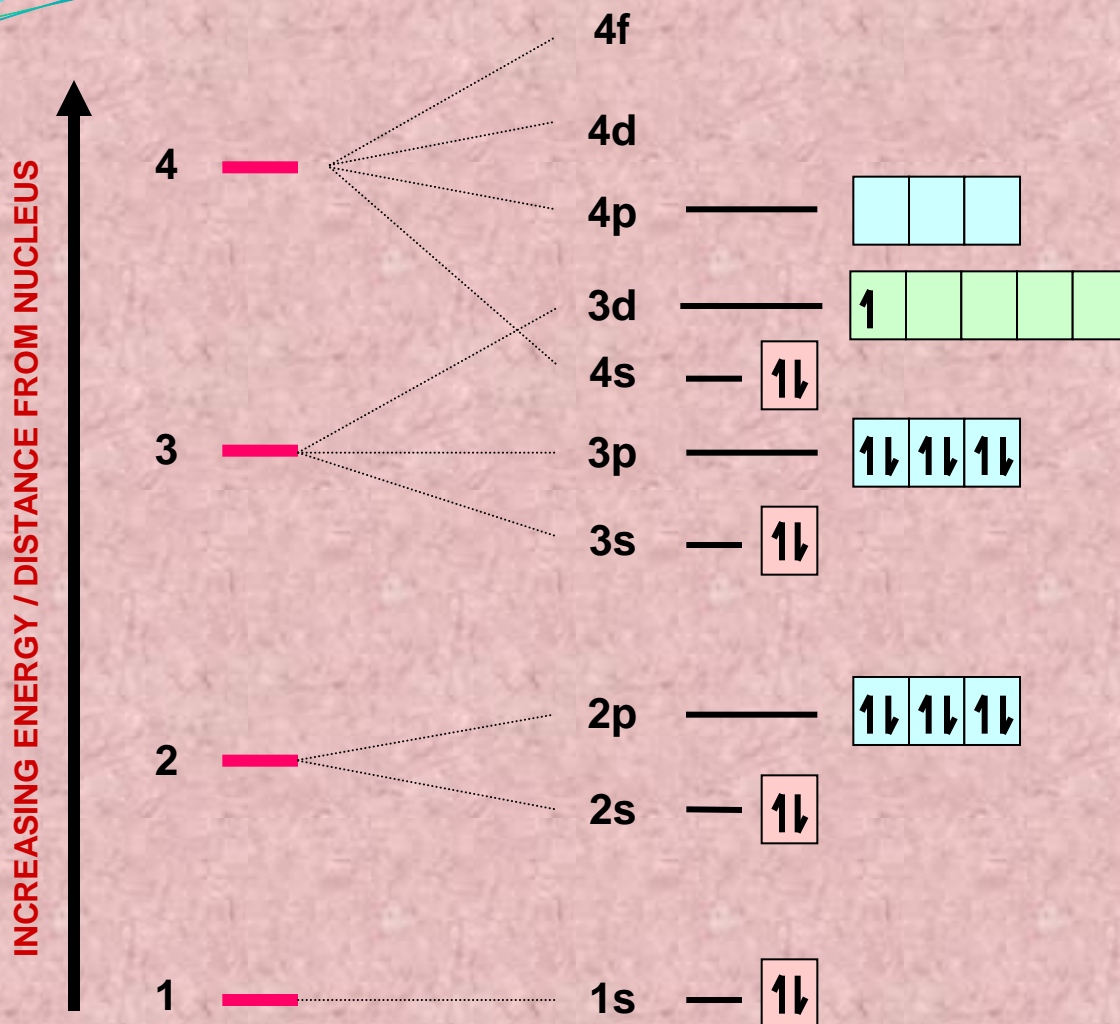


As expected, the next electron pairs up to complete a filled 4s orbital.

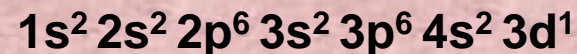
This explanation, using sub levels fits in with the position of potassium and calcium in the Periodic Table. All elements with an $-s^1$ electronic configuration are in Group I and all with an $-s^2$ configuration are in Group II.

**'Aufbau'
Principle**

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



SCANDIUM

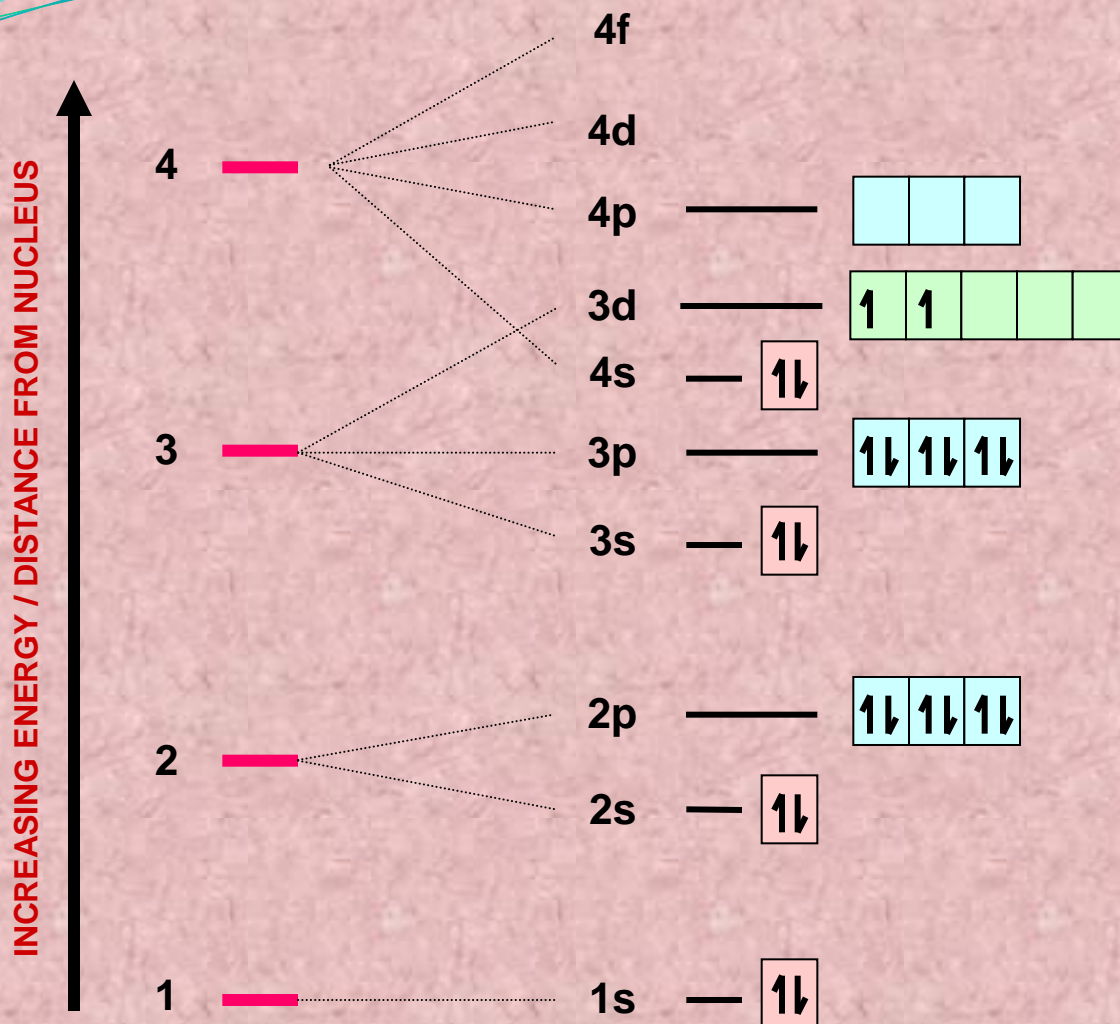


With the lower energy 4s orbital filled, the next electrons can now fill the 3d orbitals. There are five d orbitals. They are filled according to Hund's Rule -

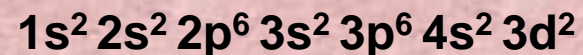
BUT WATCH OUT FOR TWO SPECIAL CASES.

**HUND'S RULE
OF
MAXIMUM MULTIPLICITY**

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



TITANIUM

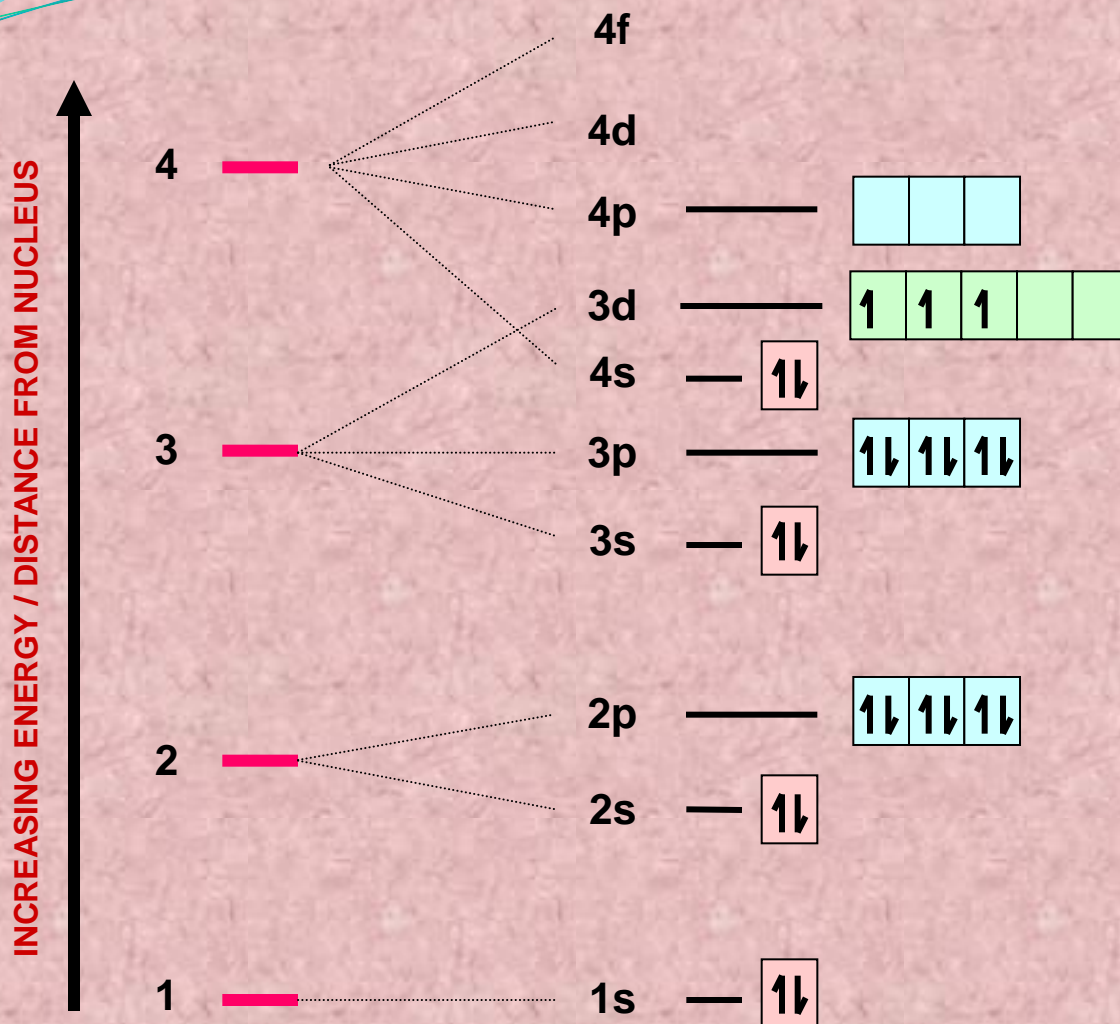


The 3d orbitals are filled according to Hund's rule so the next electron doesn't pair up but goes into an empty orbital in the same sub level.

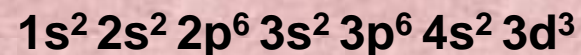
HUND'S RULE

OF
MAXIMUM MULTIPLICITY

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



VANADIUM



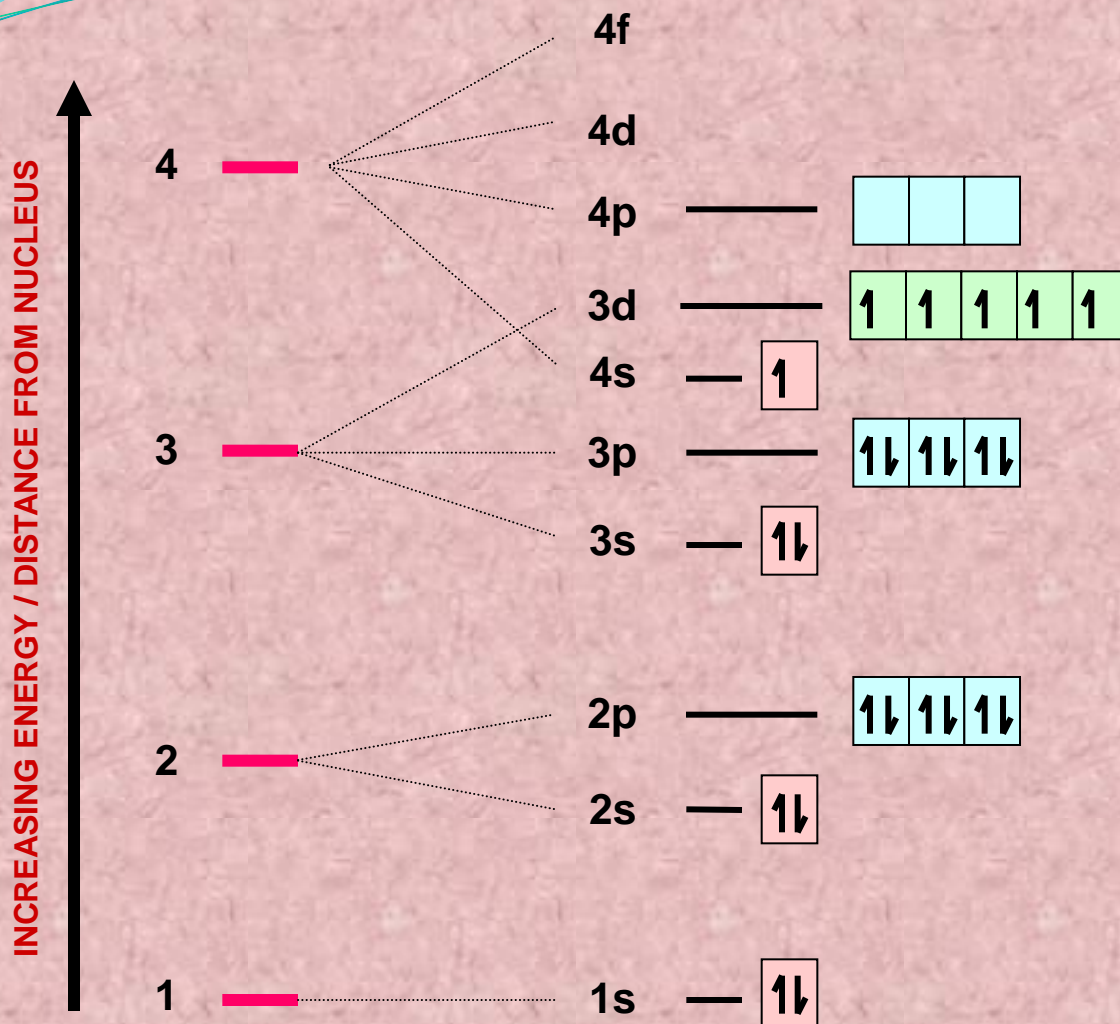
The 3d orbitals are filled according to Hund's rule so the next electron doesn't pair up but goes into an empty orbital in the same sub level.

HUND'S RULE

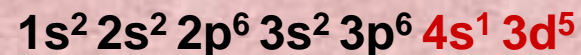
OF

MAXIMUM MULTIPLICITY

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



CHROMIUM



One would expect the configuration of chromium atoms to end in $4s^2 3d^4$.

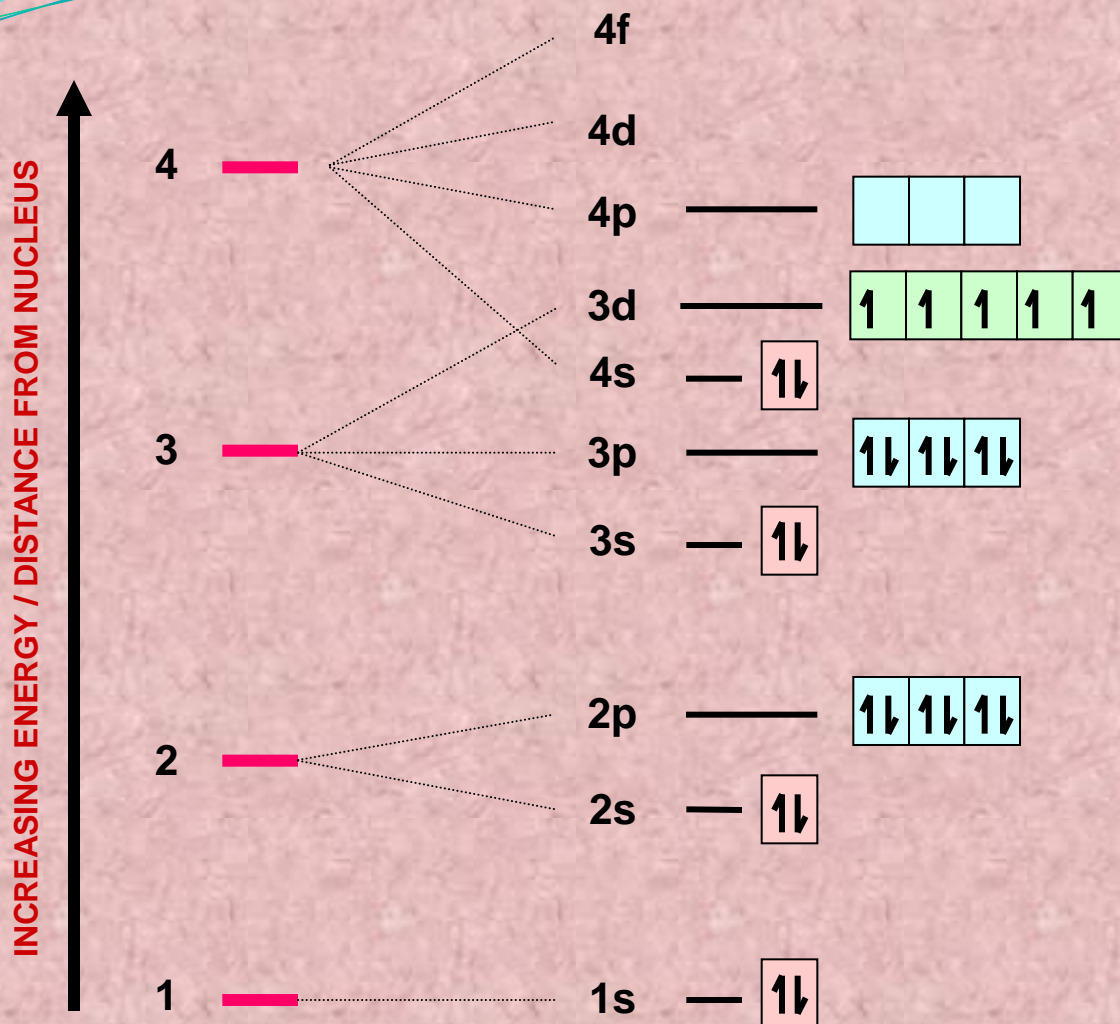
To achieve a **more stable arrangement of lower energy**, one of the 4s electrons is promoted into the 3d to give **six unpaired electrons with lower repulsion**.

HUND'S RULE

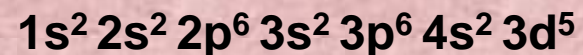
OF

MAXIMUM MULTIPLICITY

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



MANGANESE



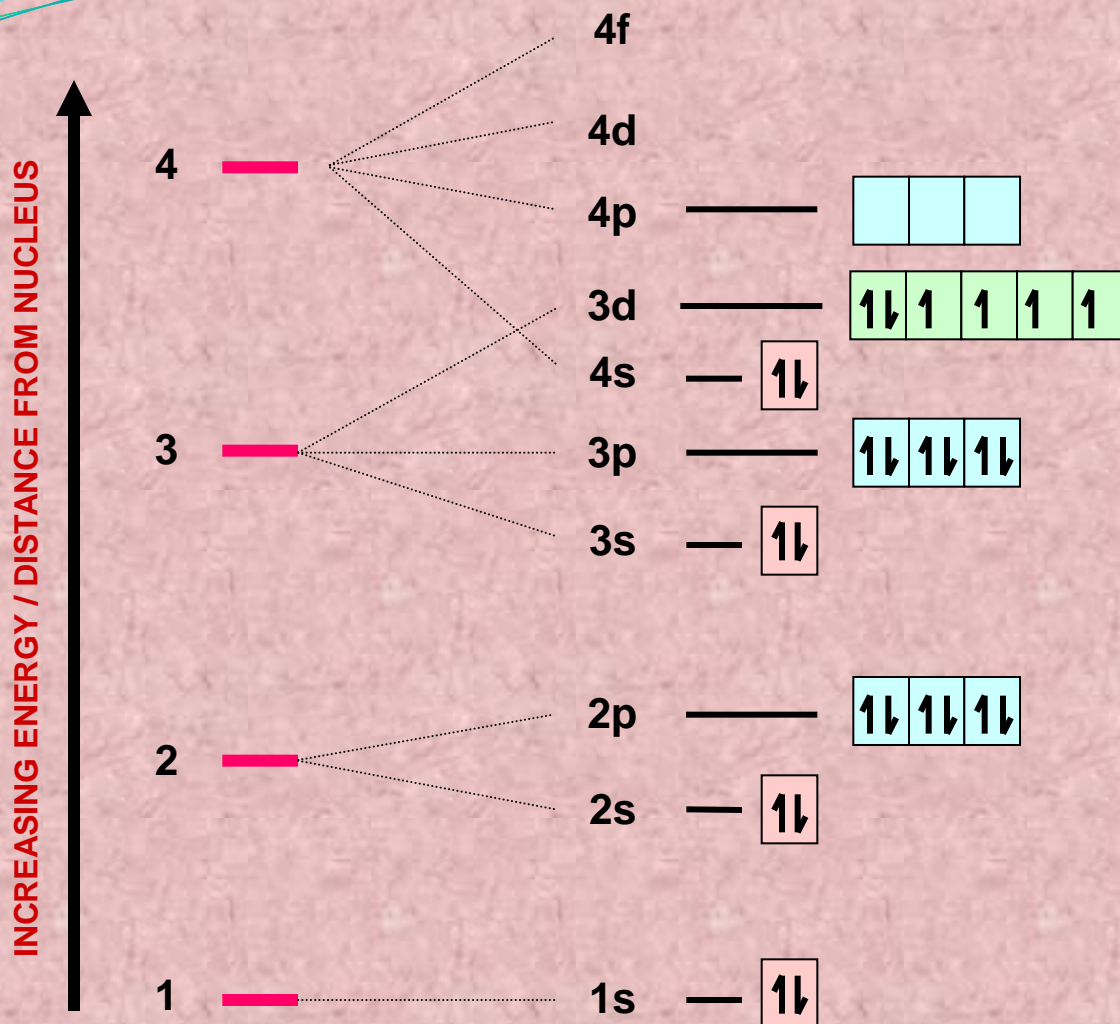
The new electron goes into the 4s to restore its filled state.

HUND'S RULE

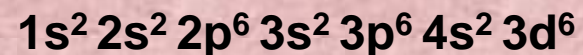
OF

MAXIMUM MULTIPLICITY

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



IRON

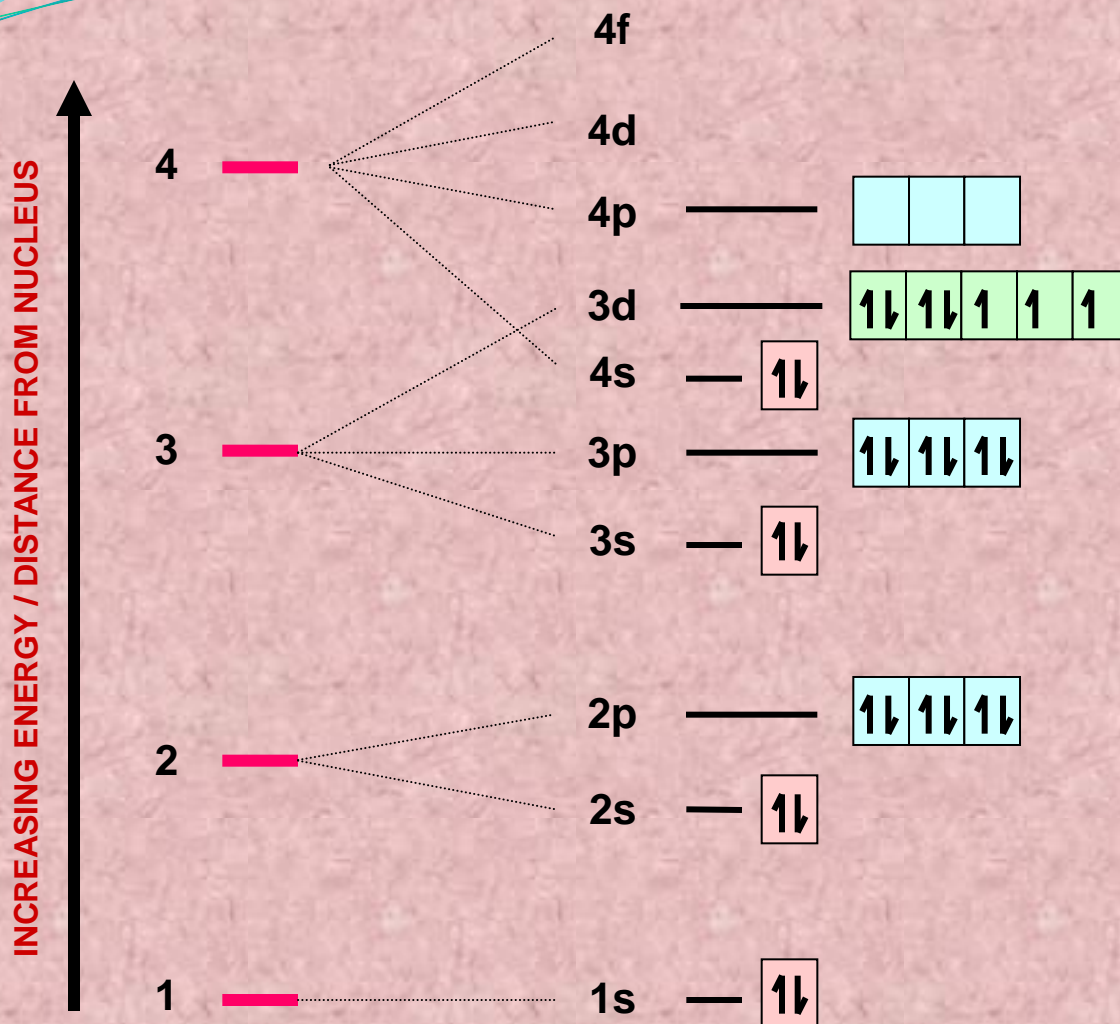


Orbitals are filled according to Hund's Rule. They continue to pair up.

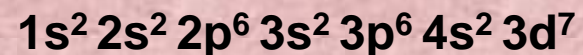
HUND'S RULE

OF
MAXIMUM MULTIPLICITY

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



COBALT

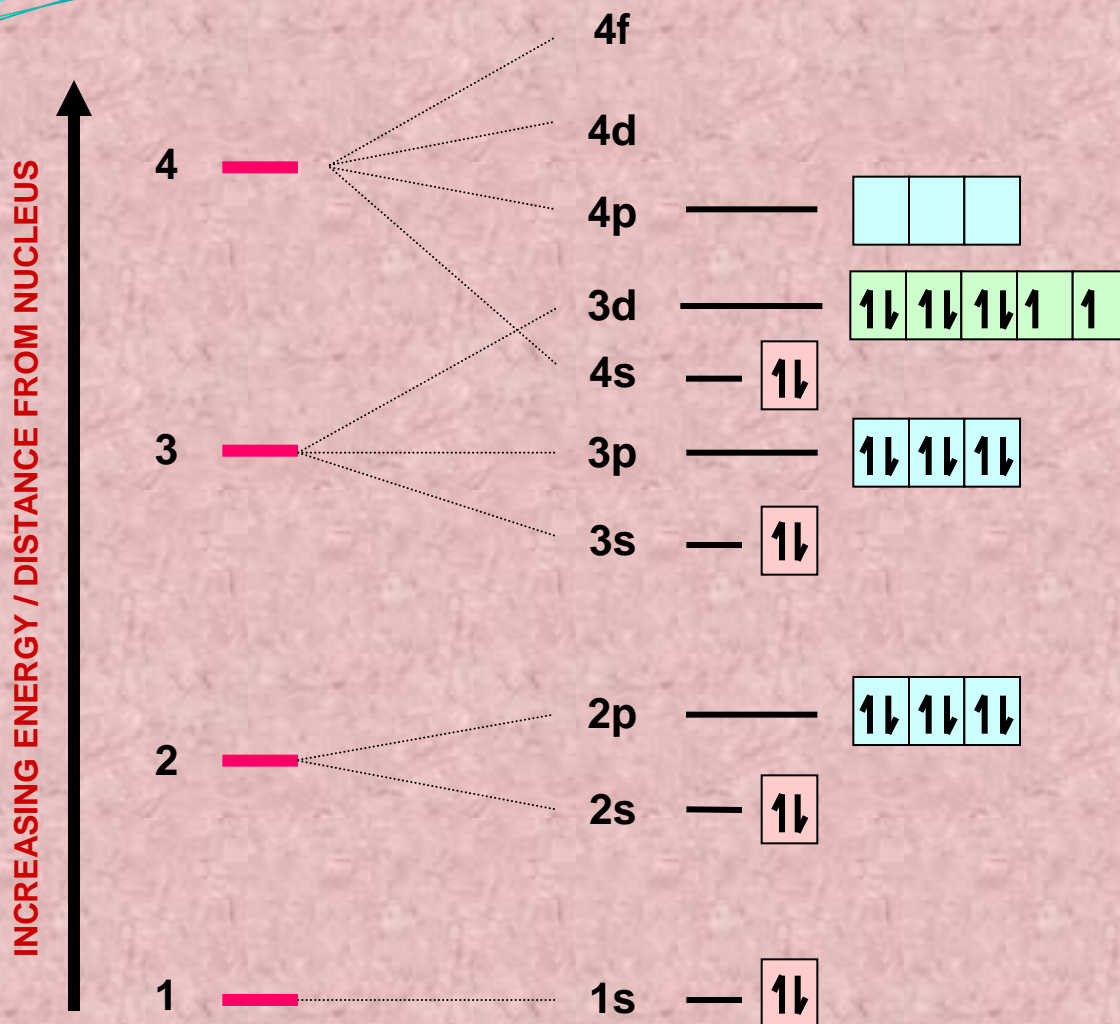


Orbitals are filled according to Hund's Rule. They continue to pair up.

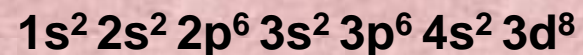
HUND'S RULE

OF
MAXIMUM MULTIPLICITY

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



NICKEL

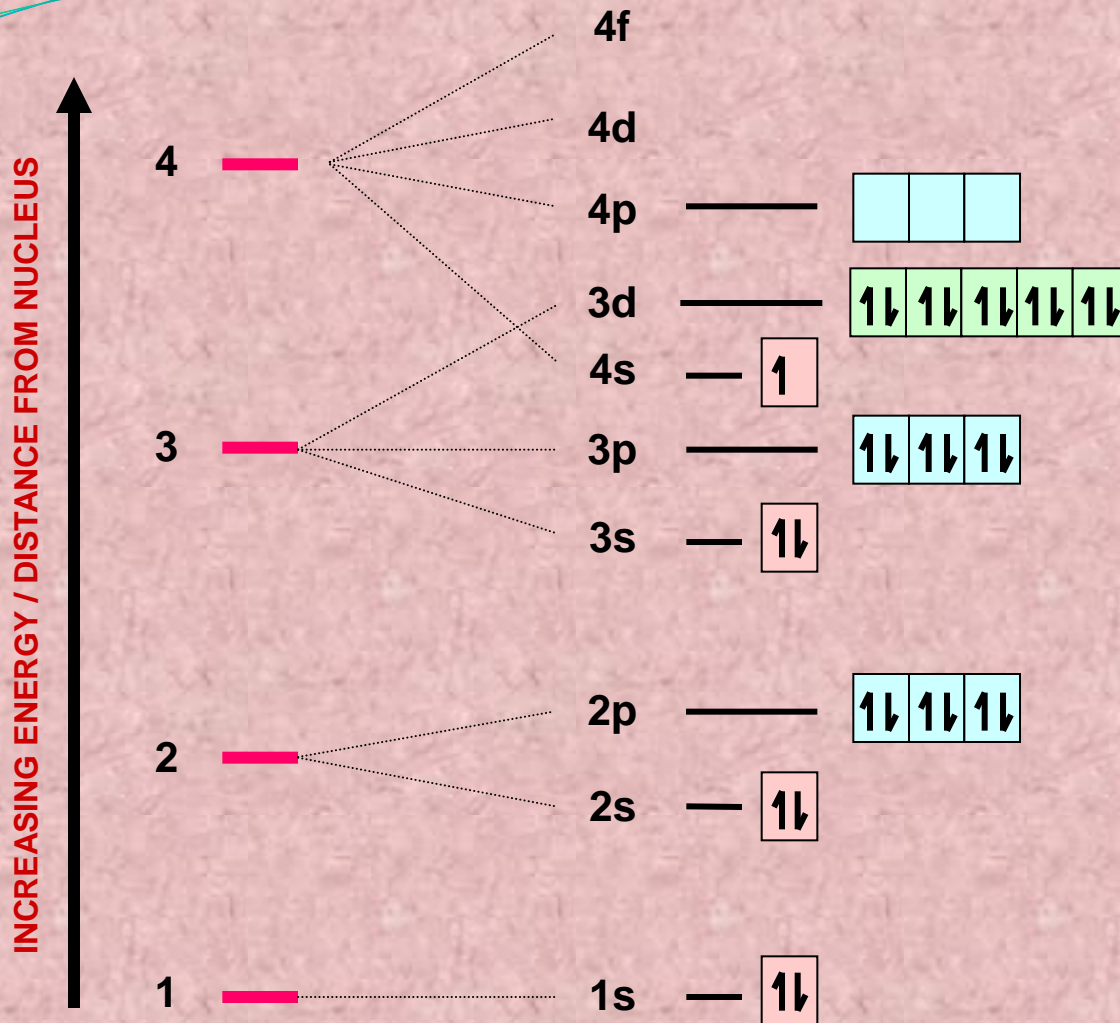


Orbitals are filled according to Hund's Rule. They continue to pair up.

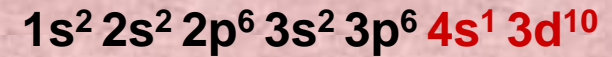
HUND'S RULE

OF
MAXIMUM MULTIPLICITY

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



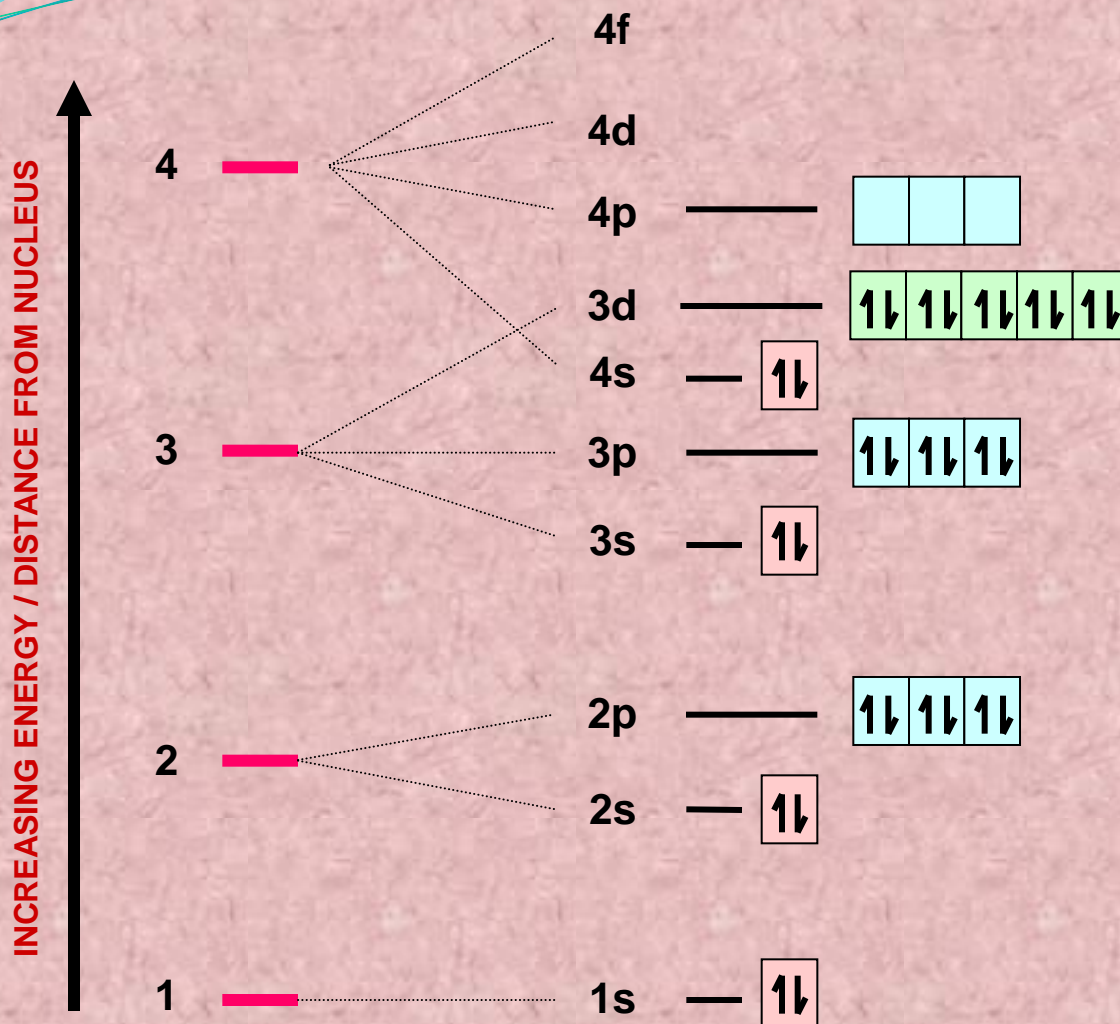
COPPER



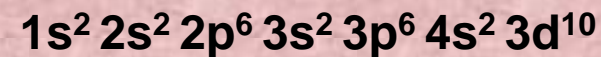
One would expect the configuration of chromium atoms to end in $4s^2 3d^9$.

To achieve a **more stable arrangement of lower energy**, one of the 4s electrons is promoted into the 3d.

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS

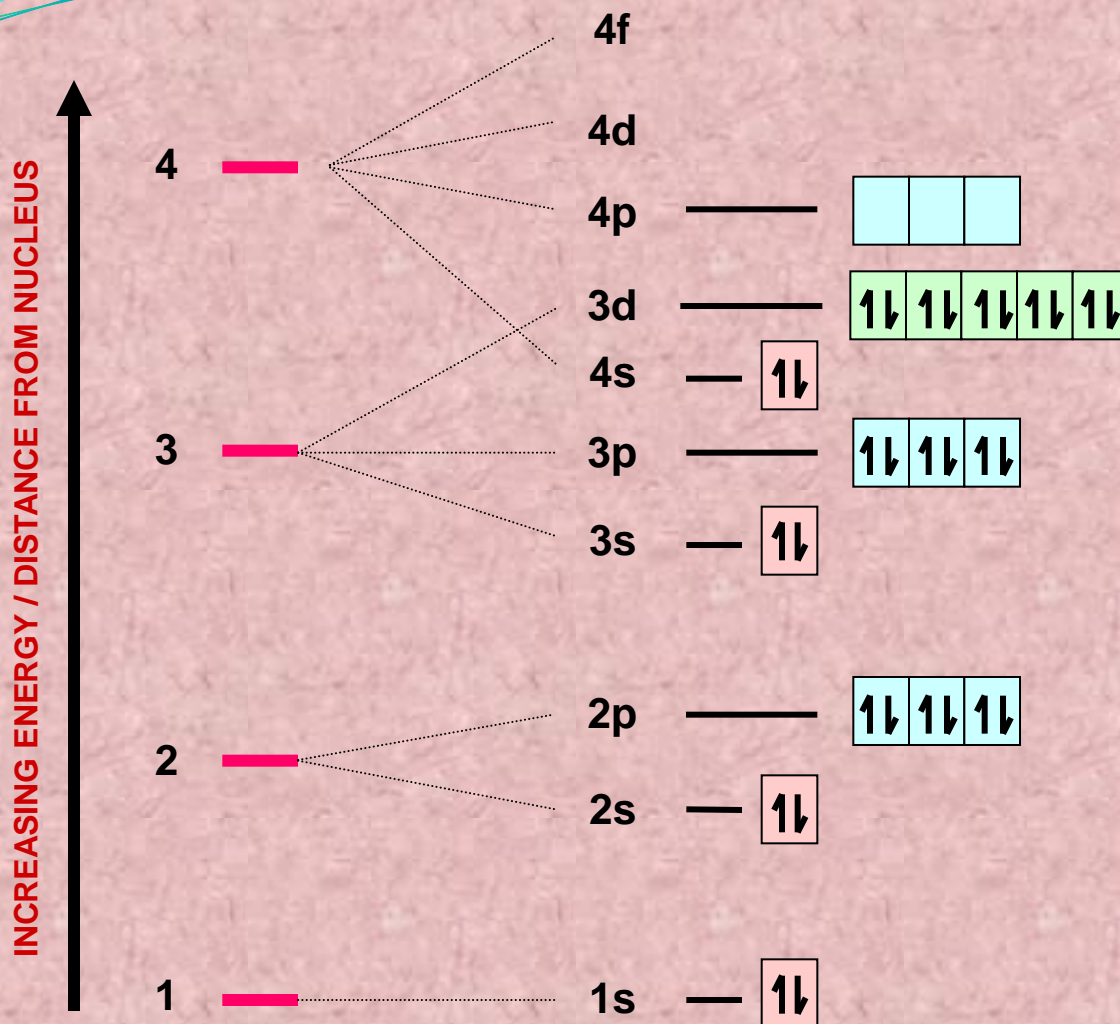


ZINC



The electron goes into the 4s to restore its filled state and complete the 3d and 4s orbital filling.

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS

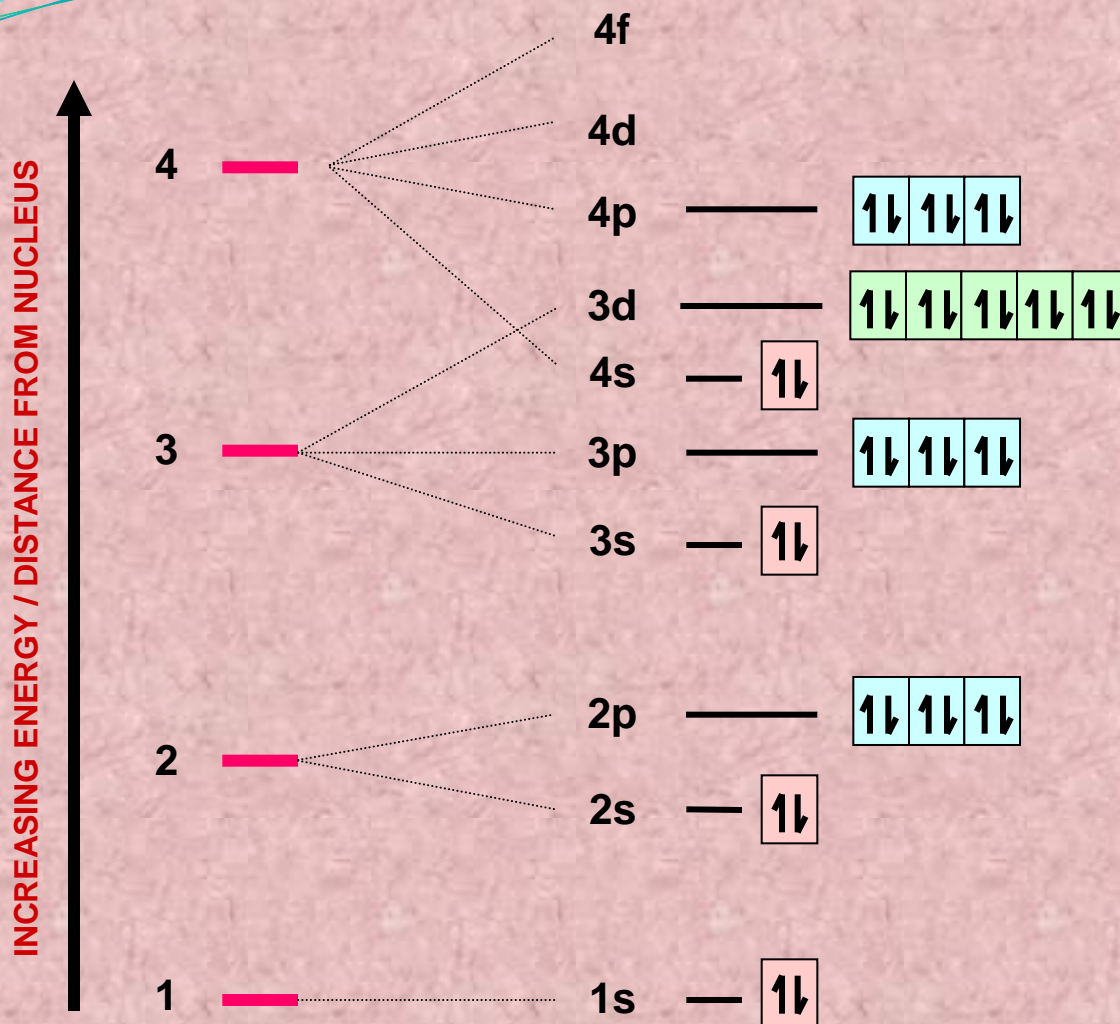


GALLIUM - KRYPTON

The 4p orbitals are filled in exactly the same way as those in the 2p and 3p sub levels.

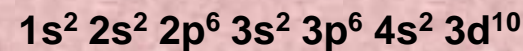
HUND'S RULE
OF
MAXIMUM MULTIPLICITY

THE ELECTRONIC CONFIGURATIONS OF THE FIRST 36 ELEMENTS



GALLIUM - KRYPTON

Prefix with...



Ga - $4p^1$

Ge - $4p^2$

As - $4p^3$

Se - $4p^4$

Br - $4p^5$

Kr - $4p^6$

Remember that the 4p configurations follow Hund's Rule with the electrons remaining unpaired to give more stability.

ELECTRONIC CONFIGURATIONS OF ELEMENTS 1-30

H	↑												$1s^1$
He	↑↓												$1s^2$
Li	↑↓	↑											$1s^2 2s^1$
Be	↑↓	↑↓											$1s^2 2s^2$
B	↑↓	↑↓	↑										$1s^2 2s^2 2p^1$
C	↑↓	↑↓	↑	↑									$1s^2 2s^2 2p^2$
N	↑↓	↑↓	↑	↑	↑								$1s^2 2s^2 2p^3$
O	↑↓	↑↓	↑↓	↑	↑								$1s^2 2s^2 2p^4$
F	↑↓	↑↓	↑↓	↑↓	↑								$1s^2 2s^2 2p^5$
Ne	↑↓	↑↓	↑↓	↑↓	↑↓								$1s^2 2s^2 2p^6$
Na	↑↓	↑↓	↑↓	↑↓	↑↓	↑							$1s^2 2s^2 2p^6 3s^1$
Mg	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓							$1s^2 2s^2 2p^6 3s^2$
Al	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑						$1s^2 2s^2 2p^6 3s^2 3p^1$
Si	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	↑					$1s^2 2s^2 2p^6 3s^2 3p^2$
P	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	↑	↑				$1s^2 2s^2 2p^6 3s^2 3p^3$
S	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	↑	↑			$1s^2 2s^2 2p^6 3s^2 3p^4$
Cl	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑				$1s^2 2s^2 2p^6 3s^2 3p^5$
Ar	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓				$1s^2 2s^2 2p^6 3s^2 3p^6$
K	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑			$1s^2 2s^2 2p^6 3s^2 3p^6 4s^1$
Ca	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓			$1s^2 2s^2 2p^6 3s^2 3p^6 4s^2$
Sc	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑		$1s^2 2s^2 2p^6 3s^2 3p^6 4s^2 3d^1$
Ti	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	↑	$1s^2 2s^2 2p^6 3s^2 3p^6 4s^2 3d^2$
V	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	↑	$1s^2 2s^2 2p^6 3s^2 3p^6 4s^2 3d^3$
Cr	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	↑	$1s^2 2s^2 2p^6 3s^2 3p^6 4s^1 3d^5$
Mn	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	↑	$1s^2 2s^2 2p^6 3s^2 3p^6 4s^2 3d^5$
Fe	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	$1s^2 2s^2 2p^6 3s^2 3p^6 4s^2 3d^6$
Co	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	$1s^2 2s^2 2p^6 3s^2 3p^6 4s^2 3d^7$
Ni	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	$1s^2 2s^2 2p^6 3s^2 3p^6 4s^2 3d^8$
Cu	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑	$1s^2 2s^2 2p^6 3s^2 3p^6 4s^1 3d^{10}$
Zn	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	↑↓	$1s^2 2s^2 2p^6 3s^2 3p^6 4s^2 3d^{10}$

ELECTRONIC CONFIGURATIONS

THE END

