

Paramagnetism and Diamagnetism

- Atoms with unpaired \uparrow electrons are called ***paramagnetic***.
 - Paramagnetic atoms are attracted to a magnet.
- Atoms with paired $\uparrow\downarrow$ electrons are called ***diamagnetic***.
 - Diamagnetic atoms are repelled by a magnet.

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- The number of orbitals per n level is given by n^2 .
 - The maximum number of electrons per n level is $2n^2$.

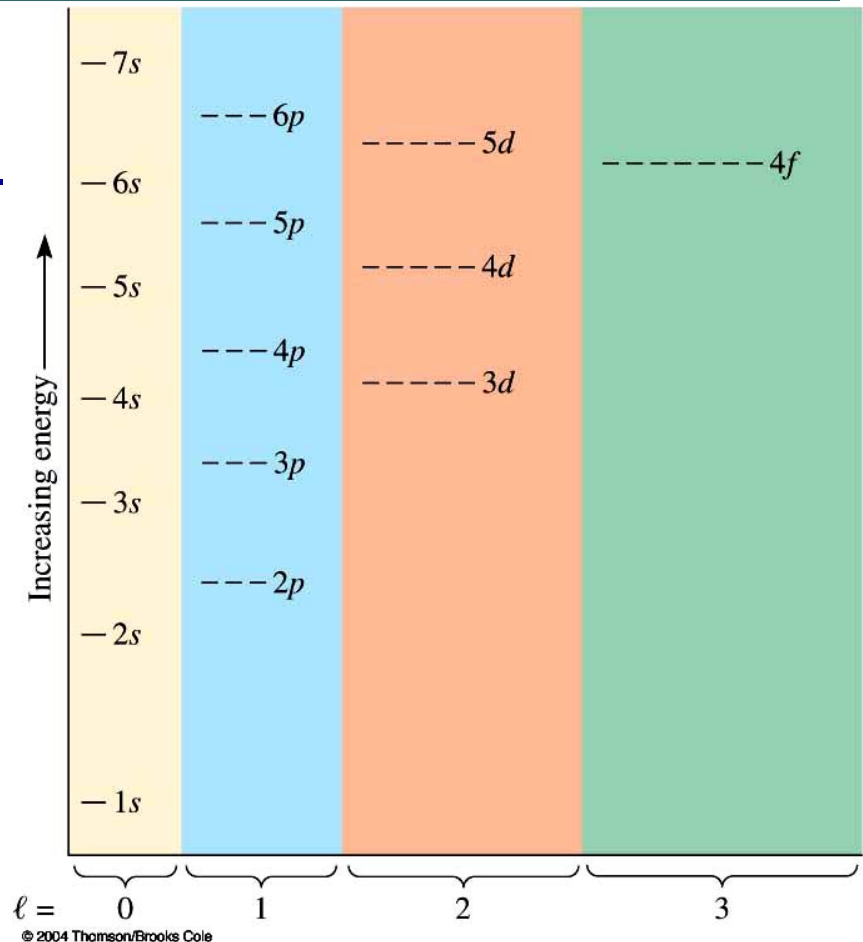
<u>Energy Level</u>	<u># of Orbitals</u>	<u>Max. # of e⁻</u>
n	n²	2n²
1	1	2
2	4	8
3	9	18
4	16	32

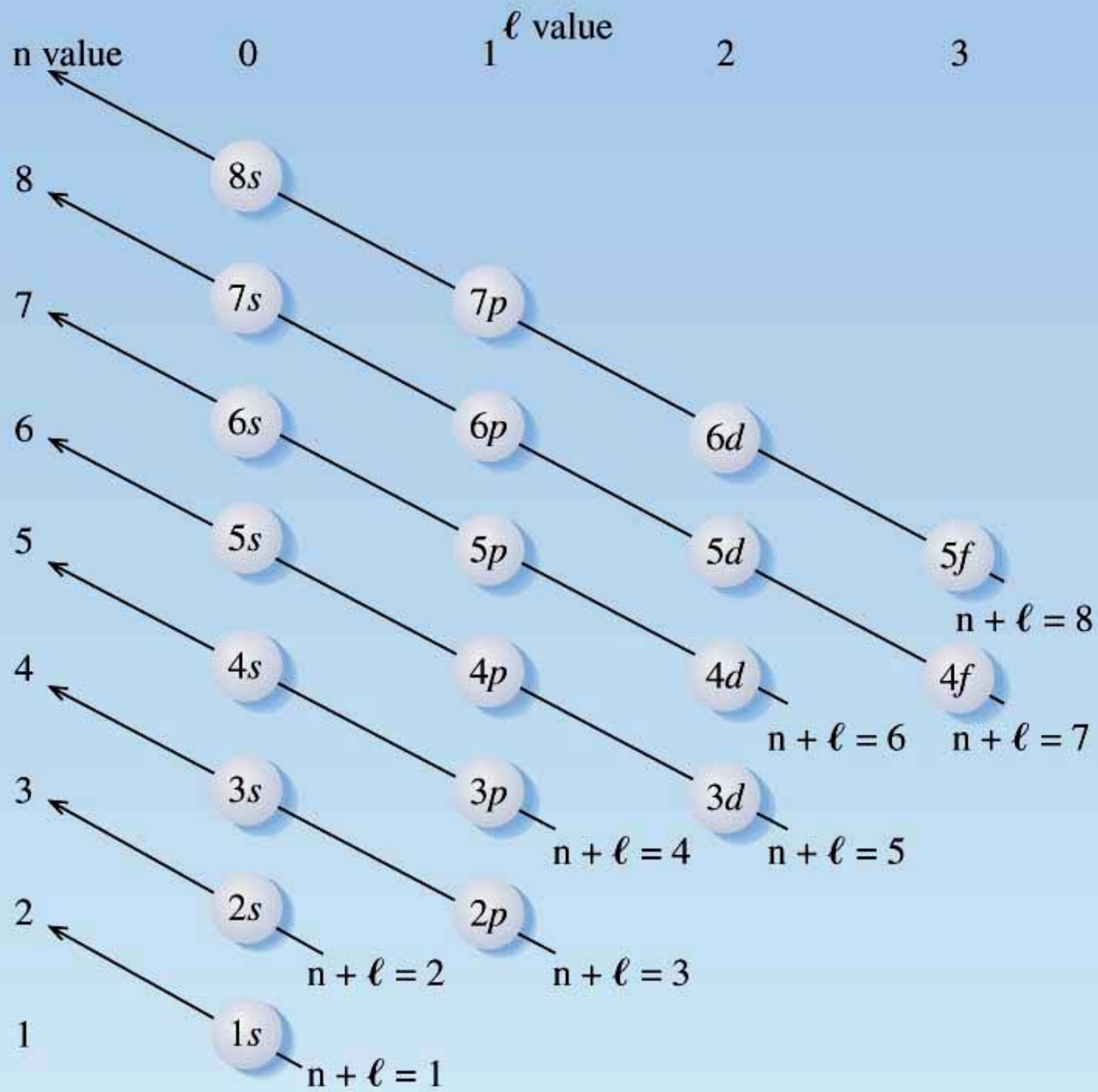
The Periodic Table and Electron Configurations

- The principle that describes how the periodic chart is a function of electronic configurations is the **Aufbau Principle**.
- The electron that distinguishes an element from the previous element enters the lowest energy atomic orbital available.

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- The Aufbau Principle describes the electron filling order in atoms.





Hund's rule

- Hund's rule tells us that the electrons will fill the p orbitals by placing electrons in each orbital singly and with same spin until half-filled. Then the electrons will pair to finish the p orbitals.

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- 1st row elements

	<u>1s</u>	<u>Configuration</u>
${}_1\text{H}$	<u>↑</u>	$1s^1$

	<u>1s</u>	<u>Configuration</u>
${}_1\text{H}$	<u>↑</u>	$1s^1$
${}_2\text{He}$	<u>↑↓</u>	$1s^2$

- 2nd row elements

	<u>1s</u>	<u>2s</u>	<u>2p</u>	<u>Configuration</u>
₃ Li	<u>↑↓</u>	<u>↑</u>	— — —	1s ² 2s ¹
₄ Be	<u>↑↓</u>	—	— — —	1s ² 2s ²
₅ B	<u>↑↓</u>	<u>↑↓</u>	<u>↑</u> — —	1s ² 2s ² 2p ¹
₆ C	<u>↑↓</u>	<u>↑↓</u>	<u>↑</u> <u>↑</u> —	1s ² 2s ² 2p ²
₇ N	<u>↑↓</u>	<u>↑↓</u>	— — —	1s ² 2s ² 2p ³
₈ O	<u>↑↓</u>	<u>↑↓</u>	<u>↑↓</u> <u>↑</u> <u>↑</u>	1s ² 2s ² 2p ⁴
₉ F	<u>↑↓</u>	<u>↑↓</u>	<u>↑↓</u> <u>↑↓</u> <u>↑</u>	1s ² 2s ² 2p ⁵
₁₀ Ne	<u>↑↓</u>	<u>↑↓</u>	<u>↑↓</u> <u>↑↓</u> <u>↑↓</u>	1s ² 2s ² 2p ⁶

- 3rd row elements

		<u>3s</u>	<u>3p</u>	<u>Configuration</u>
₁₁ Na	[Ne]	<u>↑</u>	<u>— — —</u>	[Ne]3s ¹
₁₂ Mg	[Ne]	<u>—</u>	<u>— — —</u>	[Ne]3s ²
₁₃ Al	[Ne]	<u>↑↓</u>	<u>↑ — —</u>	[Ne]3s ² 3p ¹
₁₄ Si	[Ne]	<u>↑↓</u>	<u>↑ ↑ —</u>	[Ne]3s ² 3p ²
₁₅ P	[Ne]	<u>↑↓</u>	<u>↑ ↑ ↑</u>	[Ne]3s ² 3p ³
₁₆ S	[Ne]	<u>↑↓</u>	<u>— — —</u>	[Ne]3s ² 3p ⁴
₁₇ Cl	[Ne]	<u>↑↓</u>	<u>↑↓ ↑↓ ↑</u>	[Ne]3s ² 3p ⁵
₁₈ Ar	[Ne]	<u>↑↓</u>	<u>↑↓ ↑↓ ↑↓</u>	[Ne]3s ² 3p ⁶

There is an extra measure of stability associated with half-filled or completely filled orbitals.

	<u>3d</u>	<u>4s</u>	<u>4p</u>	<u>Configuration</u>
$_{19}\text{K}$ [Ar]	— — — — —	↑	— — — — —	[Ar]4s ¹
$_{20}\text{Ca}$ [Ar]	— — — — —	↑↓	— — — — —	[Ar]4s ²
$_{21}\text{Sc}$ [Ar]	↑ — — — —	↑↓	— — — — —	[Ar]4s ² 3d ¹
$_{22}\text{Ti}$ [Ar]	↑ ↑ — — —	↑↓	— — — — —	[Ar]4s ² 3d ²
$_{23}\text{V}$ [Ar]	↑ ↑ ↑ — —	↑↓	— — — — —	[Ar]4s ² 3d ³
$_{24}\text{Cr}$ [Ar]	↑ ↑ ↑ ↑ ↑	↑	— — — — —	[Ar]4s ¹ 3d ⁵

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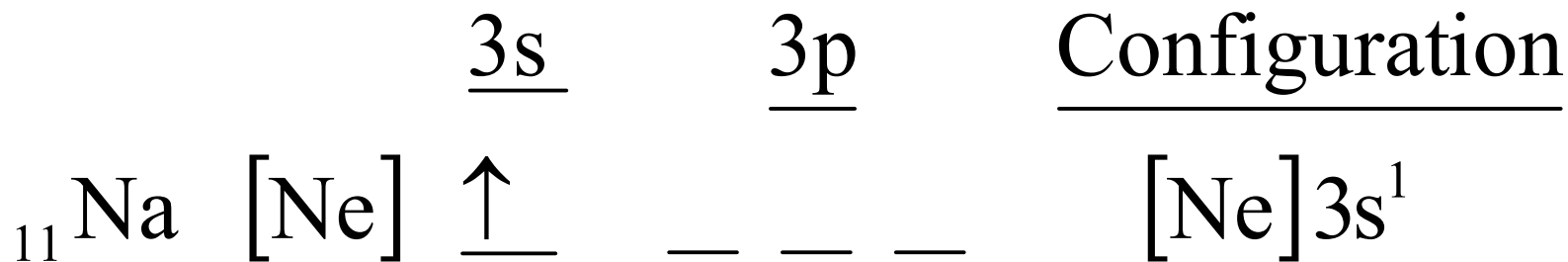
		<u>3d</u>	<u>4s</u>	<u>4p</u>	<u>Configuration</u>
₂₅ Mn	[Ar]	↑ ↑ ↑ ↑ ↑	↑↓	— — —	[Ar]4s ² 3d ⁵
₂₆ Fe	[Ar]	↑↓ ↑ ↑ ↑ ↑	↑↓	— — —	[Ar]4s ² 3d ⁶
₂₇ Co	[Ar]	↑↓ ↑↓ ↑ ↑ ↑	↑↓	— — —	[Ar]4s ² 3d ⁷
₂₈ Ni	[Ar]	↑↓ ↑↓ ↑↓ ↑ ↑	↑↓	— — —	[Ar]4s ² 3d ⁸
₂₉ Cu	[Ar]	↑↓ ↑↓ ↑↓ ↑↓ ↑↓	↑	— — —	[Ar]4s ¹ 3d ¹⁰
₃₀ Zn	[Ar]	↑↓ ↑↓ ↑↓ ↑↓ ↑↓	↑↓	— — —	[Ar]4s ² 3d ¹⁰

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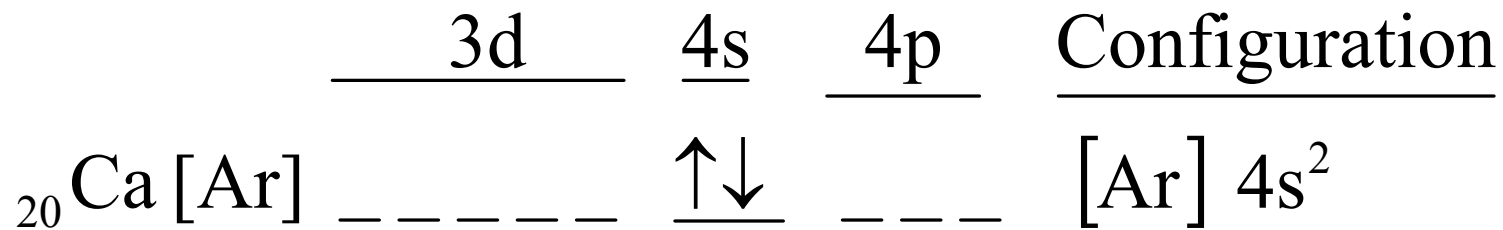
		<u>3d</u>	<u>4s</u>	<u>4p</u>	<u>Configuration</u>
₃₁ Ga	[Ar]	<u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u>	<u>↑↓</u>	<u>↑</u> <u>—</u> <u>—</u>	[Ar] 4s ² 3d ¹⁰ 4p ¹
₃₂ Ge	[Ar]	<u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u>	<u>↑↓</u>	<u>↑</u> <u>↑</u> <u>—</u>	[Ar] 4s ² 3d ¹⁰ 4p ²
₃₃ As	[Ar]	<u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u>	<u>↑↓</u>	<u>↑</u> <u>↑</u> <u>↑</u>	[Ar] 4s ² 3d ¹⁰ 4p ³
₃₄ Se	[Ar]	<u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u>	<u>↑↓</u>	<u>↑↓</u> <u>↑</u> <u>↑</u>	[Ar] 4s ² 3d ¹⁰ 4p ⁴
₃₅ Br	[Ar]	<u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u>	<u>↑↓</u>	<u>↑↓</u> <u>↑↓</u> <u>↑</u>	[Ar] 4s ² 3d ¹⁰ 4p ⁵
₃₆ Kr	[Ar]	<u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u> <u>↑↓</u>	<u>↑↓</u>	<u>↑↓</u> <u>↑↓</u> <u>↑↓</u>	[Ar] 4s ² 3d ¹⁰ 4p ⁶

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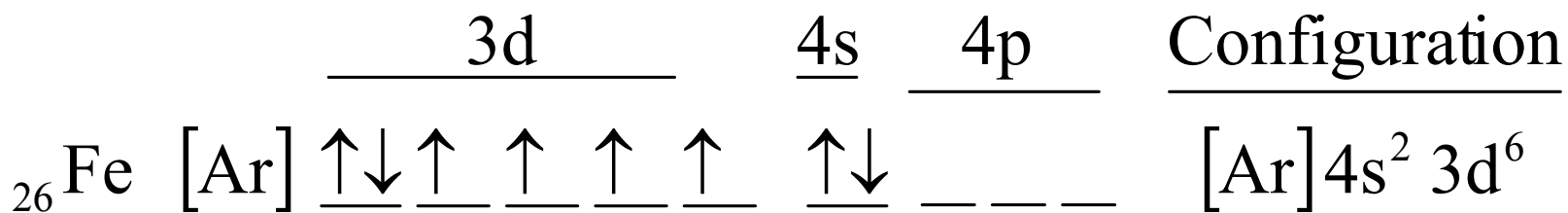
- Now we can write a complete set of quantum numbers for all of the electrons in these three elements as examples.
 - Na
 - Ca
 - Fe



	<u>n</u>	<u>l</u>	<u>m_l</u>	<u>m_s</u>	
<u>1st e⁻</u>	1	0	0	+ 1/2	}
<u>2nd e⁻</u>	1	0	0	- 1/2	
<u>3rd e⁻</u>	2	0	0	+ 1/2	}
<u>4th e⁻</u>	2	0	0	- 1/2	
<u>5th e⁻</u>	2	1	-1	+ 1/2	}
<u>6th e⁻</u>	2	1	0	+ 1/2	
<u>7th e⁻</u>	2	1	+ 1	+ 1/2	
<u>8th e⁻</u>	2	1	- 1	- 1/2	
<u>9th e⁻</u>	2	1	0	- 1/2	
<u>10th e⁻</u>	2	1	+ 1	- 1/2	
<u>11th e⁻</u>	3	0	0	+ 1/2	}
					3 s electron



	<u>n</u>	<u>l</u>	<u>m_l</u>	<u>m_s</u>	
$[\text{Ar}]$ <u>$19^{\text{th}} e^-$</u>	4	0	0	+1/2	} 4 s electrons
<u>$20^{\text{th}} e^-$</u>	4	0	0	-1/2	



	<u>n</u>	<u>l</u>	<u>m_l</u>	<u>m_s</u>	
$[\text{Ar}] \underline{19^{\text{th}} e^-}$	4	0	0	+1/2	}
$\underline{20^{\text{th}} e^-}$	4	0	0	-1/2	
$\underline{21^{\text{st}} e^-}$	3	2	-2	+1/2	
$\underline{22^{\text{nd}} e^-}$	3	2	-1	+1/2	
$\underline{23^{\text{rd}} e^-}$	3	2	0	+1/2	
$\underline{24^{\text{th}} e^-}$	3	2	+1	+1/2	
$\underline{25^{\text{th}} e^-}$	3	2	+2	+1/2	
$\underline{26^{\text{th}} e^-}$	3	2	-2	-1/2	

Chemistry is fun!