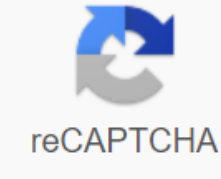




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## Wave mechanical atomic model

The second quantum number is often called the azimuthal quantum number ( $l$ ). The  $l$  value describes the shape of the space area occupied by the electron. The allowable values of  $l$  depend on the value  $n$  and can range from 0 to  $n-1$  ( $l=0,1,2,\dots,n-1$ ) for example, if it is  $n=1$ ,  $l$  can be only 0; For a given atom, all wave functions that have the same values as both  $n$  and  $l$  form a subshell. Areas of space occupied by electrons in a subshell usually have the same shape, but are oriented differently in space. The third quantum number is the magnetic quantum number ( $m_l$ ). The value ( $m_l$ ) describes the orientation of the area in the space occupied by an electron according to a functional magnetic field. Allowed values ( $m_l$ ) depend on the value  $l$ :  $m_l$  can range from  $-l$  to  $l$  in the integral stages:  $m_l = -l, -(l-1), \dots, 0, \dots, (l-1), l$  for example, if ( $l=0$ ), ( $m_l$ ) can only be 0; If  $l=1$ ,  $m_l$  can be  $-1, 0, \text{ or } +1$ ; Each wave function with a permissible combination of values  $n, l$ , and  $m_l$  describes an atomic orbital, a specific spatial distribution for an electron. For a given set of quantum numbers, each main shell has a fixed number of subshells, and each subshell has a fixed number of orbitals. Example:  $n=4$  Shell Structure How many subshells and orbitals are contained within the principal shell with  $n=4$ ? According to:  $N$  value asked for: Count number of sub-shells and orbitals in the main shell strategy: According to  $n=4$ , calculate the allowable values  $l$ . From these allowed values, count the number of subshells. Calculate the allowable amounts of  $m_l$  for each allowable value  $l$ . The total number of orbitals per subshell is the number of orbitals in the main shell. Solution: A We know that  $l$  can have all integral values from 0 to  $n-1$ . If  $n=4$ , then  $l$  can be 0, 1, 2 or 3. Because the shell has four  $l$  values, it has four sub-shells, each of which will contain a different number of orbitals, depending on the allowed values of  $m_l$ . B For  $l=0$ , the  $m_l$  can be only 0, and thus the subshell  $l=0$  has only one orbital. For  $l=1$ ,  $m_l$  can be 0 or  $\pm 1$ ; So  $l=1$  subshell has three orbitals. For  $l=2$ ,  $m_l$  can be 0,  $\pm 1$ , or  $\pm 2$ , so there are five orbitals in subshell  $l=2$ . The last allowable value is  $l=3$ , for which  $m_l$  can be 0,  $\pm 1$ ,  $\pm 2$ , or  $\pm 3$ , resulting in seven orbitals in subshell  $l=3$ . The total number of orbitals in the shell  $n=4$  is the main total number of orbitals per subshell and is equal to  $n^2=16$  Exercise:  $n=3$  multi-shell structure shell and orbital in the main shell with  $n=3$ ? Reply to three sub-shells; Not orbital instead of specifying all meanings  $n$  and  $l$  every time we talk about a subshell or orbital, chemists use a shortened system in lowercase to note the value  $l$  for a specific subshell or orbital:  $l=0,1,2,3$  Designation  $s, p, d, f$  the original quantum number is first named, followed by the letter  $s, p, d, \text{ or } f$  as appropriate. (These orbital determinations derive from historical terms for the corresponding spectroscopy characteristics: sharp, original, dumbbical, and fundamental.) A  $1s$  orbital has  $n=1$  and  $l=0$ ; a  $2p$  subshell has  $n=2$  and  $l=1$  (and has three  $2p$  orbitals, corresponding to  $m_l = -1, 0, \text{ and } +1$ ); a  $3d$  subshell has  $n=3$  and  $l=2$  (and has five  $3d$  orbitals, corresponding to  $m_l = -2, -1, 0, +1, \text{ and } +2$ ); and so is it. We can summarize the relationships between quantum numbers and the number of sub-shells and orbitals as follows (Table 6.5.1): Each main shell has  $n$  subshells. For  $n=1$ , only one unit under is possible ( $1s$ ); for  $n=2$ , there are two subshells ( $2s, 2p$ ); for  $n=3$ , there are three subshells ( $3s, 3p, \text{ and } 3d$ ); and so is it. Each shell has a  $ns$  subshell, each shell with  $n \geq 2$  also has a  $np$  subshell, and each shell with  $n \geq 3$  also has a  $nd$  subshell. Since a  $2d$  subshell requires both  $n=2$  and  $l=2$ , which is not the allowed value of  $l$  for  $n=2$ , there is no  $2d$  subshell. Each subshell has  $2l+1$  orbitals. This means that all  $ns$  subshells are included in a single orbital  $s$ , all  $np$  subshells contain three orbitals  $p$ , all  $nd$  subshells contain five orbitals  $d$  subshells, and all  $nf$  subshells contain seven orbitals  $f$ . Each main shell has  $n$  subshells, and each subshell has  $2l+1$  orbitals. Table: Values of  $n, l$ , and  $m_l$  through  $n=4$  Subshell Designation ( $m_l$ ) Number of Orbitals in Subshell Number of Orbitals in Shell  $1, 0, 1, 1, 1, 2, 0, 2s, 0, 1, 4, 1, 2p, -1, 0, 1, 3, 0, 3s, 0, 1, 9, 1, 3p, -1, 0, 1, 3, 2, 3d, -2, -1, 0, 1, 2, 5, 4, 0, 1, 1, 1, 1, 4p, -1, 0, 1, 3, 2, 4d, -2, -1, 0, 1, 2, 5, 3, 4f, -3, -2, -1, 0, 1, 2, 3, 7$  The newsflash interrupts your favorite TV program. At the First National Bank one is kept. The suspect fled in a car and is believed to be somewhere in the downtown area. The robber can be located only in a specific area – the police do not have an exact location, just a general idea as to whereabouts of the thief. In 1926, Austrian physicist Erwin Schrödinger (1887–1961) used the duality of electron wave-particles to develop and solve a complex mathematical equation that accurately described the behavior of electrons in a hydrogen atom. The quantum mechanical model of the atom comes from the solution of the Schrödinger equation. Quantization of electron energy is a need to solve the equation. This is in contrast to the Bohr model in which quantization was simply assumed to be without a mathematical basis. Remember that in the Bohr model, the exact electron was confined to very well-defined circular orbits around the nucleus. The quantum mechanical model is a radical departure from it. The solutions of the Schrödinger wave equation, called wave functions, are only likely to find electrons at a given point around the nucleus. Electrons do not travel around the nucleus in simple circular circuits. Figure 1. An electron cloud: The darker area nearer the nucleus indicates a high probability of finding electrons, while the lighter area is greater than the nucleus indicating a lower probability of finding electrons. The location of electrons in the quantum mechanical model of the atom is often referred to as the electron cloud. The electron cloud can be imagined as follows: imagine putting a square piece of paper on the ground where the point in the circle represents the nucleus. Now visit a marker and drop it on paper over and over again, making small marks anywhere marker. If you drop many markers, the overall pattern of the point will be almost circular. If you aim towards the center well, there will be more dots near the core and gradually less point as you move away from it. Each point represents the place where the electron can be at any moment. Because of the principle of uncertainty, there is no way to know exactly where the electron is. An electron cloud has variable densities: the high density in which the electron will most likely be, and the low density in which the electron is the least likely (Fig. 1). For a specific definition of cloud shape, it is customary to refer to an area of space within which there is a 90% chance of finding an electron. This is called an orbital, three-dimensional space zone that indicates where there is a high probability of finding an electron. Schrödinger wave equation summary replaced Bohr ideas about electron location with a factor of uncertainty. The location of the electron can only be given as a possibility that the electron is somewhere in a specific region. To answer the following questions, use the link below What was a problem with the Bohr Atom model? What did Heisenberg show about electrons? What did Schrödinger derive? What does it take to revisit the perspective of quantum mechanics of atoms? What's the wave function? What does a high-density super-electron suggest? Super-electron: The location of electrons in the quantum mechanical model is atom. Orbital: The 3D space area that indicates where the electron is likely to be found. Quantum Mechanics Model: A model of atom derived from the Schrödinger wave equation and deals with probabilities. Wave function: Only the probability of finding an electron at a given point around I'm a chemistry student and I'm doing a project about Erwin Schrödinger that came with the atomic model of the mechanical wave. Let's get started! Erwin Schrödinger was born on August 12, 1887, and died on January 4, 1961. He was an Austrian physicist who made many important theories about quantum theory. He won the 1933 Nobel Prize in Physics and the Max Planck Medal in 1937. Irwin's father came from a Bavarian family who had settled in Vienna generations earlier. In 1926, he combined the Bohr model Erwin Schrödinger with the hypothesis de Broglie. He suggested that the electron is a 3D waveform that bypasses the nucleus at a full number of wavelengths, allowing the waveform to replicate itself as a stable standing wave that represents the energy levels of the Bohr model. A standing wave is a wave that does not convey energy or move but undergoes resonance. This means it can absorb energy from a nearby source that is fluctuating at a suitable frequency. A standing wave must also have wavelengths so that a full number of parts of the wave fit within the setting. If the number of wave sections is a complete number, then the wave collapses. Schrödinger suggested that de Broglie was correct about matter waves and that electrons were located in atomic space according to the frequencies of standing waves. Therefore, the energy needed to change from one standing wave to another must be quantized to maintain the total number of wavelengths and avoid collapse. In support of his hypothesis, Schrödinger developed a mathematical equation to describe electron-like wave behavior. The Schrödinger wave equation not only gave the correct levels of energy to the hydrogen atom, but was somewhat useful in atoms with more than one electron. His theory was that mathematical equations could be used to find the probability of an electron's location. This atomic theory predicts the chances of electron location. This model is known as quantum mechanics model. The quantum mechanics model does not define the exact path of an electron, but predicts the chances of the location of the electron. This model can be portrayed as a nucleus surrounded by a super-electron. Where the cloud is densest, electrons are more likely to find the largest, and on the contrary, electrons are less likely to be placed in a less dense area than the cloud. In this way, this model introduced the concept of sub-energy levels. The two-notch experiment is a show where matter and energy can display the characteristics of both waves and particles, revealing the probable nature of electrons. The experiment belongs to a general class of double track experiments in which a wave is divided into two separate waves that are later re-combined. Changes along the path of both waves lead to The change causes the pattern of interference. The Schrödinger equation is a partial DDL equation that describes how the quantum state of some physical systems changes with time. Formulated in late 1925, it was published in 1926 by Austrian physicist Erwin Schrödinger. schrodinger .