



Short Note

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On Filling Energy Levels and Sub-Levels

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As it is known, the filling of energy levels and sub-levels is produced in accordance with V. M. Klechkowski I and II rules [1-3], which are based on the two solutions of the E. Schroedinger equation [4-5] – the principal n and the orbital l quantum numbers, but nothing is said about the third solution – the magnetic m_l quantum number. It is true that the order obtained by increasing the energy corresponds to the experimental data, but in our opinion it is better to use a single unified rule in which to the sum of the quantum numbers (n and l) will be added the third term. The third solution of the Schroedinger

equation m_l determines the number of orbitals on the sub-levels $k_{m_l} [k_{m_l}(s) = 1, k_{m_l}(p) = 3, k_{m_l}(d) = 5, k_{m_l}(f) = 7]$. Our input is to add the inverse values of these numbers (1, 3, 5 and 7) $\frac{1}{k_{m_l}} - 1, 0.33, 0.2$ and 0.14 to this sum. With this in mind, the sum of three numbers $\left(\Sigma = n + l + \frac{1}{k_{m_l}}\right)$ for all sub-levels of the 7 periods of the periodic table of the elements is given in Table 1.

Table 1: The sum of three numbers for all sub-levels of the 7 periods of the periodic table of the elements is given in the table.

| sub-level | n | l | k_{m_l} | $\frac{1}{k_{m_l}}$ | Σ |
|-----------|---|---|-----------|---------------------|----------|
| 1s | 1 | 0 | 1 | 1 | 2 |
| 2s | 2 | 0 | 1 | 1 | 3 |
| 2p | 2 | 1 | 3 | 0.33 | 3.33 |
| 3s | 3 | 0 | 1 | 1 | 4 |
| 3p | 3 | 1 | 3 | 0.33 | 4.33 |
| 3d | 3 | 2 | 5 | 0.2 | 5.2 |
| 4s | 4 | 0 | 1 | 1 | 5 |
| 4p | 4 | 1 | 3 | 0.33 | 5.33 |
| 4d | 4 | 2 | 5 | 0.2 | 6.2 |
| 4f | 4 | 3 | 7 | 0.14 | 7.14 |
| 5s | 5 | 0 | 1 | 1 | 6 |
| 5p | 5 | 1 | 3 | 0.33 | 6.33 |
| 5d | 5 | 2 | 5 | 0.2 | 7.2 |
| 5f | 5 | 3 | 7 | 0.14 | 8.14 |
| 6s | 6 | 0 | 1 | 1 | 7 |
| 6p | 6 | 1 | 3 | 0.33 | 7.33 |
| 6d | 6 | 2 | 5 | 0.2 | 8.2 |
| 6f | 6 | 3 | 7 | 0.14 | 9.14 |
| 7s | 7 | 0 | 1 | 1 | 8 |
| 7p | 7 | 1 | 3 | 0.33 | 8.33 |
| 7d | 7 | 2 | 5 | 0.2 | 9.2 |
| 7f | 7 | 3 | 7 | 0.14 | 10.14 |

Table 2: The layout of the sub-levels according to the sum increase is given in the table.

| sub-level | n | l | k_{m_l} | $\frac{1}{k_{m_l}}$ | Σ |
|-----------|---|---|-----------|---------------------|----------|
| 1s | 1 | 0 | 1 | 1 | 2 |
| 2s | 2 | 0 | 1 | 1 | 3 |
| 2p | 2 | 1 | 3 | 0.33 | 3.33 |
| 3s | 3 | 0 | 1 | 1 | 4 |
| 3p | 3 | 1 | 3 | 0.33 | 4.33 |
| 4s | 4 | 0 | 1 | 1 | 5 |
| 3d | 3 | 2 | 5 | 0.2 | 5.2 |
| 4p | 4 | 1 | 3 | 0.33 | 5.33 |
| 5s | 5 | 0 | 1 | 1 | 6 |
| 4d | 4 | 2 | 5 | 0.2 | 6.2 |
| 5p | 5 | 1 | 3 | 0.33 | 6.33 |
| 6s | 6 | 0 | 1 | 1 | 7 |
| 4f | 4 | 3 | 7 | 0.14 | 7.14 |
| 5d | 5 | 2 | 5 | 0.2 | 7.2 |
| 6p | 6 | 1 | 3 | 0.33 | 7.33 |

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|----|---|---|---|------|-------|
| 7s | 7 | 0 | 1 | 1 | 8 |
| 5f | 5 | 3 | 7 | 0.14 | 8.14 |
| 6d | 6 | 2 | 5 | 0.2 | 8.2 |
| 7p | 7 | 1 | 3 | 0.33 | 8.33 |
| 6f | 6 | 3 | 7 | 0.14 | 9.14 |
| 7d | 7 | 2 | 5 | 0.2 | 9.2 |
| 7f | 7 | 3 | 7 | 0.14 | 10.14 |

The layout of the sub-levels according to the sum increase is given in Table 2.

The new filling rule will be formulated as follows:

The electron sub-levels are filled in accordance with the sum of the principal n , the orbital l quantum numbers and the inverse values of the number of orbitals on the sub-level - $\Sigma = n + l + \frac{1}{k_{m_l}}$.

Based on this, the principle of least energy is formed:

The electronic sub-levels are arranged in the following order according to the energy increase:

$1s < 2s < 2p < 3s < 3p < 4s < 3d < 4p < 5s < 4d < 5p < 6s < 4f < 5d < 6p < 7s < 5f < 6d < 7p < 6f < 7d < 7f$.

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