Volume Charge Density Functional Theory (VDFT)

[OR: Volume Density Functional Theory (VDFT)]

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"I may be wrong and you may be right, and by an effort, we may get nearer to the truth."

----------Sir Karl Raymond Popper

The Basic Points of View:

Volume density functional theory (VDFT), is about the electronic structure (Law of charge distribution of extra-nuclear) latest theory, is a research on concentric spherical layers model [1], multi electron atoms (Z >1) system based on the electronic structure (structure and charge distribution of extra-nuclear) quantization mouthed, the main goal is to replace the *Electron density* and the *Wave function* for basic quantity research by *Volume charge density*, and build a final accord with reality, decisive, causality, locality and completeness, multiple causes multiple effects of quantum theory.

The total charge Q = Ze; *Volume* Charge density functional ρ_i =F（E，H，Q),
Frequency-Field Function v_i=f（ρi)=f[F（E，H，Q）],
Levels represents a function Ei=hv_i=h.f（ρi)=h.{f[F（E，H，Q）]} .

Sphere center symmetry, isotropic, $\rho_i=q_i/V_{icl}$； Distribution in the radius ρ;
Divergence on the radial gradient; Fine Structure ------ gradient divergence;
Hyperfine structure ------second gradient divergence; between the various qi cited, balance and nuclear repulsion between.

Ei-level relationship with the natural frequency of the v_i and charge density ρ_i of:
Ei=hv_i=h.f（ρ_i)=h.{f[F（E，H，Q）]}.

Reference

[1] Concentric spherical layer-type model of an atom
http://vixra.org/abs/1402.0104

< The Causes and Mechanism of Atomic Energy Levels quantization >