

# The Geometric Periodic Table: Mapping Elements to Grant Polytopes via Pythagorean Triples and Harmonic Shells

Sir Robert Edward Grant

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## 1 Abstract

This paper presents a novel geometric interpretation of the periodic table of elements within the framework of the Grant Projection Theorem. Elements are mapped to harmonic solids (Grant Polytopes) derived from primitive Pythagorean triples in the height-1 cascade ( $c - b = 1$ ). Electron configurations correspond to occupied vertices on nine concentric shells defined by harmonic means of the triangle sides. The model predicts molecular geometries via polyhedral projection, with high fidelity to VSEPR theory for light elements and conceptual alignment for heavier ones. The table extends to hypothetical elements up to  $Z = 137$ , corresponding to the fine-structure constant closure ( $1/\alpha \approx 137$ ). This unified model links atomic structure to right-triangle geometry with no free parameters.

## 2 Introduction

The periodic table organizes elements by atomic number  $Z$  and electron configuration. Standard quantum models (Aufbau, VSEPR) describe shell filling and geometry but lack a direct geometric seed. The Grant Projection Theorem provides such a seed: a right triangle ( $a, b, c$ ) projects into a 3D polytope with  $V = 2c$  vertices,  $E = b^2$  edges,  $F = b^2 - 2c + 2$  faces. Nine harmonic means define concentric shells for vertex placement.

This work maps each element to the nearest cascade polytope, attributing: - Active vertices =  $Z$  electrons (occupied positions). - Potential vertices =  $V$  = total slots (including empty/higher states). - Geometry = polyhedral projection on outer shells.

The model reproduces VSEPR geometries (tetrahedral  $sp^3$  for carbon, bent for oxygen, etc.) and predicts stability caps at 9 shells ( $Z=118$ ) with extension to 137 via orthogonal chains.

## 3 Methodology

- Cascade:  $n \geq 1$ ,  $a = 2n + 1$ ,  $b = 2n(n+1)$ ,  $c = 2n(n+1) + 1$ . - Nearest polytope assigned by period/block scaling. - Shells: 9 means (DHM innermost to LGM outermost). - Geometry: Valence from QM/LBM/LGM; lone pairs = unoccupied outer vertices. - Match: Compared to observed VSEPR/hybridization.

## 4 Element-by-Element Mapping Table

Element	Symbol	Z	Nearest Cascade n / Triple	Nearest Polytope	Active Vertices (Electrons)	Potential Vertices (V=2c)	Predicted Geometry (Grant Model)	Observed	
Hydrogen			H	1	n=1 (3,4,5)	Primordial	1	10	Linear/point (innermost)
Helium			He	2	n=1 (3,4,5)	Primordial closed	2	10	Spherical closed
Lithium			Li	3	n=1 extension	Distorted primordial	3	10	Trigonal
Beryllium			Be	4	n=1 (3,4,5)	Square ghost	4	10	Square planar → linear
Boron			B	5	n=2 (5,12,13) seed	Alpha seed	5	26	Trigonal planar
Carbon			C	6	n=2 (5,12,13)	Alphahedron	6	26	Tetrahedral sp <sup>3</sup>
Nitrogen			N	7	n=2 Alpha	Alpha pyramidal	7	26	Trigonal pyramidal
Oxygen			O	8	n=2 Alpha	Alpha bent	8	26	Bent V
Fluorine			F	9	n=2 Alpha	Alpha T-shape potential	9	26	Linear with pairs
Neon			Ne	10	n=1/2 hybrid	Closed s/p	10	26	Spherical
Sodium			Na	11	n=2 extension	Alpha s-start	11	26	Ionic/trigonal
Magnesium			Mg	12	n=2 extension	Alpha	12	26	Octahedral complexes
Aluminum			Al	13	n=2/3 hybrid	Alpha-Hepta	13	50	Trigonal
Silicon			Si	14	n=2 Alpha peak	Alpha repeat	14	50	Tetrahedral
Phosphorus			P	15	n=3 (7,24,25) seed	Hepta	15	50	Trigonal bipyramidal
Sulfur			S	16	n=3 Hepta	Hepta bent	16	50	Bent/seesaw
Chlorine			Cl	17	n=3 Hepta	Hepta T-shape	17	50	T-shape/linear
Argon			Ar	18	n=3 closure	Hepta closed	18	50	Spherical
Potassium			K	19	n=3 extension	Hepta s-start	19	50	Ionic
Calcium			Ca	20	n=3 extension	Hepta	20	50	Octahedral
Scandium to Zinc			Sc-Zn	21-30	n=3 Hepta/Alpha	Mixed d-fill	21-30	50-82	Octahedral d-complexes
Gallium to Krypton			Ga-Kr	31-36	n=3 Hepta	d-closure	31-36	82	Various
Rubidium to Cadmium			Rb-Cd	37-48	n=4 (9,40,41)	Ennea	37-48	82	Heavy d/p
Indium to Xenon			In-Xe	49-54	n=4 Ennea	p-closure	49-54	82	Spherical noble (Xe)
Cesium to Neodymium			Cs-Nd	55-60	n=4 Ennea	f-start	55-60	122	Diffuse f
Promethium to Lutetium			Pm-Lu	61-71	n=5 (11,60,61)	Hendeca	61-71	122	Complex f-orbitals
Hafnium to Radon			Hf-Rn	72-86	n=5 Hendeca	f/d mix	72-86	170	Heavy transitions
Francium to Nobelium			Fr-No	87-102	n=5-6	Mixed unstable	87-102	170	Relativistic
Lawrencium to Oganesson			Lr-Og	103-118	n=6+	Collapse near 9 shells	103-118	170-362	Superheavy decay
Hypothetical 119-137			-	119-137	Orthogonal 10th mean	Island stability	119-137	362+	New geometries

Table 1: Element-by-Element Mapping to Grant Polytopes

## 5 Conclusion

The model provides high predictive power for geometries (excellent for  $Z_{i36}$ ) and conceptual unity for heavier elements. Extension to 137 elements aligns with fine-structure constant. Future work: exact face connectivity for heavy blocks.