

Prediction of Neutrino Masses from Information in the Electrovacuum

A Numerical Method Obtaining Particle Quantities from Classical Equations

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Abstract: It is demonstrated how to obtain quantities of free particles and masses of nuclei from known equations of the electrovacuum around the particle. A sampling method is used based on a higher number of numerical tests. The single computation ends at a geometric limit, which involves eigenvalues. These eigenvalues, significantly correlating with known particle values (mass, spin, electric charge, magnetic moment) and also masses of nuclei, are detected. This allows predicting particle quantities being unknown up to date. In particular, possible masses of neutrinos are predicted. Obtained values are 0.068 eV, 0.095 eV, 0.155 eV, 0.25 eV, 0.31 eV, 0.39 eV, 0.56 eV, 1.63 eV, 2.88 eV, 5.7 eV. The algorithm is explained to some detail.

Keywords: Theory of Relativity · Rainich Theory · Algorithm · Numerical Simulations · Fundamental Quantities

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1 Introduction

The standard theory being used for describing structure of matter on nuclear and subatomic level is based on considerations of symmetry and was successful in constructing a classification scheme of sub-atomic particles. This theory, although considered as being “the best we have”, has a number of shortcomings. It cannot be unified with general relativity and there is no way to compute masses of elementary particles by a method based on first-principles. Instead, masses of sub-atomic and elementary particles (as well as other properties) have to be introduced as adaptable parameters. Although there were attempts in the past to overcome this problem, these approaches have not been considered in mainstream physics up to date. One of these approaches is the unification of electromagnetism with general relativity by the EINSTEIN-MAXWELL theory which the work of this paper is based on. The equations and the resulting geometry were found by RAINICH [2, 3] already in 1924 and are explained in the next section. In section three, details of the computational method for obtaining quantities of elementary particles are discussed. The results for neutrinos are presented, and confronted with results for nuclei and the electron in section four. While the masses of neutrinos could not be determined exactly up to now, we present a prediction based on our calculations which lies within the range safely known by experiment. The method works well on the basis of inherent information in the electrovacuum around the particle.

2 The Equations

The theory is based on the relativistic tensor equations [5]

$$R_{ik} = \kappa \left(\frac{1}{4} g_{ik} F_{ab} F^{ab} - F_{ia} F_k{}^a \right) \quad , \quad (1)$$

$$F_{ij,k} + F_{jk,i} + F_{ki,j} = 0 \quad , \quad (2)$$

$$F^{ia}{}_{;a} = 0 \quad , \quad (3)$$

in which g_{ik} are the components of metrics, R_{ik} those of the RICCI tensor and F_{ik} those of the electromagnetic field tensor. κ is EINSTEIN’s gravitation constant. If we express the field tensor by a vector potential with

$$F_{ik} = A_{i,k} - A_{k,i} \quad , \quad (4)$$

equation (2) is identically fulfilled. Thus, we can base our calculations on quantities having the character of potentials that are metrics and the electromagnetic vector potential.

These equations are known as EINSTEIN-MAXWELL equations. The energy-momentum tensor of electrodynamics is equated to the energy-momentum tensor of EINSTEIN's theory [1]. In detail, here are used the *homogeneous* MAXWELL equations for force equilibrium and conservation of energy and momentum, mathematically expressed by the BIANCHI identities. These equations describe the electrovacuum around a particle and involve geometry described by the EINSTEIN part (equation (1)) of the equations. These equations and the involved geometry were found by RAINICH already in the year 1924 [2, 3]. The sources of related inhomogeneous equations are replaced by integration constants [4]. Mass, spin, electric charge, and magnetic moment are the first integration constants.

The geometric equations do *not* define causality, because causality is not a geometric category. The geometric equations yield only 10 independent equations for 14 components g_{ik}, A_i . We will consider only the diagonal elements of g_{ik} plus two off-diagonal elements for practical calculation, what reduces the number of equations.

3 Numerical Simulations

Analytic solutions (different from zero) based on integration constants lead commonly to singularities. There are two types of singularities. The first type is singularities assumed for point masses and charges in order to simplify the equations so that analytical solutions are feasible. This is often considered as a deficit when comparing a calculation with the situation in reality. However, in our calculations, these formal singularities are placed into the inner of the particle (according to observer's coordinates) which is not subject of calculation. With spherical coordinates, the formal singularity is at the centre.

The basic idea of calculation is as follows. The equations (1,3,4) are evaluated on a radial grid from outer to inner and so one approaches the unknown inner area successively. At a certain radius, the calculation starts to diverge because the central singularity becomes predominant. It is important to notice that this radius of divergence is clearly separated from the central singularity so a second type of "singularity" here appears. A kind of chaos

takes effect as well, what would be matter of research.

These numerical simulations according to the EINSTEIN-MAXWELL equations show that this singularity is irrelevant, for the following reason: Numerical simulations using iterative, non-integrating methods lead always to a boundary at the conjectural particle radius. As a result, the second singularity is always within a geometric limit. The area within this geometric limit *according to observer's coordinates* is not accessible to further investigation but this is even not required. The geometric limit is the mathematical reason for the existence of discrete solutions. This has also to do with chaos [4]. These discrete solutions involve discrete values of the integration constants, i.e. eigenvalues. We shall see that the electrovacuum is able to produce such eigenvalues.

In order to gain eigenvalues, one has to do lots of tests, because the particle quantities are integration constants and have to be inserted into the initial conditions (for more details see [4]), which are defined for the electrovacuum around the particle.

As already mentioned, the basis for computations are equations (1,3,4). For the sake of simplicity, we restrict equations (1,3,4) to time independence and rotation symmetry. That results, with spherical coordinates

$$x^1 = r , x^2 = \vartheta , x^3 = \varphi , x^4 = jct ,$$

in 6 independent equations for 8 components with potential character $A_3, A_4, g_{11}, g_{12}, g_{22}, g_{33}, g_{34}, g_{44}$, the rest vanishes. In order to override the indeterminacy by the two missing equations, we define

$$g_{12} = 0 \quad (\text{and, consequently, } g^{12} = 0) \quad (5)$$

and

$$g = \det|g_{ik}| = r^4 \sin^2 \vartheta \quad . \quad (6)$$

These conditions are arbitrary, in which the second is taken from the free-field Minkowski metric. They are in combination leading to reasonable results.

The integration constants from equations (1,3,4) result from a series expansion. The first coefficients of expansion are the input for the simulations and are inserted into the initial conditions [4]. The output is the number of grid points along the radius until divergence occurs, which is a measure for the stability of the solution.

The first coefficients (integration constants) are

$$c_1 = - \frac{\kappa m}{4\pi} \implies \frac{\kappa m}{4\pi} \quad (7)$$

(mass),

$$c_2 = j \frac{\kappa s}{4\pi c} \implies \frac{\kappa s}{4\pi c} \quad (8)$$

(spin),

$$c_3 = -j \frac{\mu_o^{\frac{1}{2}} Q}{4\pi} \implies \frac{\kappa^{\frac{1}{2}} \mu_o^{\frac{1}{2}} Q}{4\pi} \quad (9)$$

(charge), and

$$c_4 = - \frac{\varepsilon_o^{\frac{1}{2}} M}{4\pi} \implies \frac{\kappa^{\frac{1}{2}} \varepsilon_o^{\frac{1}{2}} M}{4\pi} \quad (10)$$

(magnetic moment).

As explained, these follow from a comparison of series expansion from the EINSTEIN-MAXWELL equations (homogeneous MAXWELL equations) with the solutions of equivalent inhomogeneous equations, see [4]. The dimensionless terms after the arrow are taken for computation, and have positive values. The imaginary unit is eliminated. The radius unit ($r = 1$) corresponds to 10^{-15} m. By this, the initial conditions become, using $T = \frac{\pi}{2} - \vartheta$,

$$g_{11} = 1 + \frac{c_1}{r} - \frac{1}{2} \left(\frac{c_3}{r}\right)^2 + \frac{\left(\frac{c_4}{r^2}\right)^2 (1 + \cos^2 T)}{10}, \quad (11)$$

$$g_{22} = r^2 \left\{ 1 + \left(\frac{c_4}{r^2}\right)^2 \left(\frac{1}{3} \cos^2 T - \frac{3}{10}\right) \right\}, \quad (12)$$

$$g_{33} = r^2 \cos^2 T \left\{ 1 + \left(\frac{c_4}{r^2}\right)^2 \left(\frac{\cos^2 T}{15} - \frac{3}{10}\right) \right\}, \quad (13)$$

$$g_{44} = 1 - \frac{c_1}{r} + \frac{1}{2} \left\{ \left(\frac{c_3}{r}\right)^2 + \left(\frac{c_4}{r^2}\right)^2 \sin^2 T \right\}, \quad (14)$$

$$g_{34} = r \cos^2 T \left(\frac{c_2}{r^2} - \frac{1}{2} \frac{c_3 c_4}{r^3} \right), \quad (15)$$

$$A_3 = r \cos^2 T \frac{c_4}{r^2}, \quad (16)$$

$$A_4 = \frac{c_3}{r}. \quad (17)$$

The actual computation is done with quantities performed from “physical components”, i.e. the unity summands in the equations are eliminated. This

is because the physical components of metrics have a magnitude of ca 10^{-40} and thus divergence is detectable at all. It is defined for example

$$g_{11} = 1 + g_{(11)} , \quad (18)$$

$$g_{22} = r^2 (1 + g_{(22)}) , \quad (19)$$

$$g_{33} = r^2 \sin^2 \vartheta (1 + g_{(33)}) , \quad (20)$$

$$g_{44} = 1 + g_{(44)} , \quad (21)$$

correspondingly with the CHRISTOFFEL symbols. The indices in parenthesis mean “physical” parts of equations (11-14). So the unities vanish in the equations before one evaluates them.

We have to insert the values of the integration constants into the modified initial conditions (with physical components), see program in the supplementary data package (available at author’s current website³). The conversion of physical into normalized (dimensionless) values and vice versa is described in detail in [4, 7]. Table 1 shows some values with radius unit of 10^{-15}m . These examples allow for convenient conversion.

Table 1: Physical and normalized values for conversion

	physical value	norm. value
proton mass	$1.672 \times 10^{-24}\text{g}$	2.48×10^{-39}
electr. mass	$0.911 \times 10^{-27}\text{g}$	1.35×10^{-42}
\hbar	$1.054 \times 10^{-27}\text{cm}^2\text{g/s}$	5.20×10^{-40}
elem. charge	$1.602 \times 10^{-19}\text{As}$	1.95×10^{-21}
μ_B	$1.165 \times 10^{-27}\text{Vs cm}$	3.70×10^{-19}

Higher moments are missing in the equations because of lack of knowledge, their influence is estimated to be rather small. In the results section we will insert known values and values deviating from them, and compare the results.

³http://www.bruchholz-acoustics.de/physics/neutrino_data.tar.gz

The algorithm for evaluating the equations requires numerical differentiation. We do this by separating the quantity with highest radius index at the left-hand side, and all previously evaluated quantities are at the right-hand side. These quantities come from equations (1) and (3) using (4). For example when we calculate spherical shells from outside to inside, the new quantity is $f_{m+2,n}$. In the following difference equations f stands for *any* potential-like quantity:

$$\left. \frac{\partial f}{\partial r} \right|_{r_m, T_n} = \frac{f_{m-1,n} - f_{m+1,n}}{2 \, dr} , \quad (22)$$

$$\left. \frac{\partial^2 f}{\partial r^2} \right|_{r_m, T_n} = \frac{f_{m+2,n} - 2f_{m,n} + f_{m-2,n}}{(2 \, dr)^2} , \quad (23)$$

$$\left. \frac{\partial f}{\partial T} \right|_{r_m, T_n} = \frac{f_{m,n+1} - f_{m,n-1}}{2 \, dT} , \quad (24)$$

$$\left. \frac{\partial^2 f}{\partial T^2} \right|_{r_m, T_n} = \frac{f_{m,n+1} - 2f_{m,n} + f_{m,n-1}}{dT^2} . \quad (25)$$

This equation set is linearly resolvable to the new quantities in equation (23), and the whole adherent tensor equations are non-linear. Detailed formulae are available in the Pascal code. (The Pascal code is in the supplementary data.) This method is made possible by the fact that 2nd derivatives in the tensor equations appear always linearly. Therefore the doubled difference in equation (23).

When the program runs, the values of the several components are successively quantified in one spherical shell after the other. The computation is done for all components along the inclination (ϑ values) at a given radius, and along the radius (with all inclination values) from outside to inside step by step until geometric limits are reached. After starting the procedure, we get the values as expected from the initial conditions. Suddenly, the values grow over all limits. At this point geometric limits are reached and the calculation is stopped.

The step count (number of iterations) up to the first geometric limit of a metrical component (where the absolute value of the “physical” component becomes unity) depends on the inserted values of the integration constants. A relatively coarse grid reflects strong dependencies, however, the referring values of the integration constants are imprecise. Computations with finer grid lead to smaller contrast of the step counts, but the values are more precise.

The resulting eigenvalues of the integration constants are obtained where the step count until divergence is at maximum. Round-off errors have to be respected because these can be in the order of step count differences for the formulae.

In order to see the eigenvalues, lots of tests were done with parameters more and less deviating from reference values. The output parameter (used for the plots discussed in the results section) is the mentioned step count. In order to make visible the differences, the step count above a “threshold” is depicted in resulting figures by a more or less fat “point”.

Though neutrinos are uncharged, one has to use always the full EINSTEIN-MAXWELL equations to account for the inherent non-linearity. Because the information is in the entire field outside the geometric boundary, one has to do so even if charge and magnetic moment are zero.

4 Computational Results

4.1 Spins, Electric Charges, Magnetic Moments

Tests including parameters different from mass had to be run with an initial radius close to the conjectural particle radius. Here, the influences of the four relevant parameters to metrics (about 10^{-40}) are comparable.

The best result has been achieved with the free electron, see [4, 7]. The magnetic moment of the electron arises specially sharply, due to the dominant influence.

Unfortunately, the mass gets lost in the “noise” from rounding errors. Only cases with charge and mass together can be made visible in exceptional cases, see for example [6].

4.2 Masses

4.2.1 *Masses of Nuclei*

The influence of mass to metrics prevails in a certain distance from the conjectural particle or nucleus radius, respectively. It proves being possible to set the remaining parameters to zero. Figs. 1 and 2 show related tests, with possible assignment of maxima in the figures to nuclei [6].

It was necessary in the tests according to Figs. 1, 2 to “pile up” the data. For this purpose, several test series with *slightly* different parameters (mostly initial radius) have been run, and the related step counts (the output) have

been added. So the “noise” from rounding errors is successively suppressed. With 80 bit floating point registers, the rounding error is in the 20th decimal. As well, the relative deviation of difference quotients from related differential quotients *in the first step* is roughly 10^{-20} – that is the limit, where the onset of chaotic behaviour can be seen. Consequently, simulations with only 64 bit (double) lead to no meaningful results.

One can see certain patterns in the figures, which could arise from errors by neglecting other parameters.

4.2.2 *Masses of Leptons*

It is principally possible to deduce the masses of all free particles, if they are stable to some extent. Since the electron mass is relatively small, one needs an initial radius of about 4×10^{-13} m in order to be able to neglect the influence of spin, charge, magnetic moment to some extent, see Fig. 3 [7]. One step count maximum (piled) appears fairly correctly at the experimental value, flanked by adjoining maxima, possibly caused by the neglected parameters.

The success in detecting known masses gives good reasons for trying a prediction of neutrino masses. That implies that neutrinos are stationary particles, i.e. have rest mass at all. Then they can never reach light speed.

The Particle Data Group [8] commented in the year 2002:

There is now compelling evidence that neutrinos have nonzero mass from the observation of neutrino flavor change, both from the study of atmospheric neutrino fluxes by SuperKamiokande, and from the combined study of solar neutrino cross sections by SNO (charged and neutral currents) and SuperKamiokande (elastic scattering).

The neutrino has the advantage of being electromagnetically neutral. As well, the spin does not perceptibly influence other components of metrics than those for the spin itself. So we can unscrupulously neglect the spin, and search for quite tiny masses.

Quoting the Particle Data Group (in 2002) again [8]:

Mass⁴ $m < 3$ eV.

Interpretation of tritium beta decay experiments is complicated by anomalies near the endpoint, and the limits are not without ambiguity.

Newer experiments re-verify this ambiguity, just providing multiple mass bounds.

⁴of electron neutrino

Ten plausible maxima have been found in the area for the electron neutrino, see Figs. 4, 5, 6, 7, 8, and the supplementary data. Obtained values are 0.068 eV, 0.095 eV, 0.155 eV, 0.25 eV, 0.31 eV, 0.39 eV, 0.56 eV, 1.63 eV, 2.88 eV, 5.7 eV. Smaller values (Fig. 4) are less convincing.

The mentioned ambiguity gets along with the fact that multiple mass values have been detected. It could be possible that the set of values is reduced by computation with spin. The precision with 80 bit registers is not sufficient for such calculations. However, it could well be possible interpreting some values as composites from smaller values. Here we could have comparable circumstances like in nuclei so that there is no reason for the assumption that only one value can exist. This conclusion is supported by multiple experimental mass bound values.

Many mass values are integer multiples of ~ 0.08 eV, within the tolerances of the method. At the place of this value there is a hole in the figure, flanked by maxima at 0.068 eV and 0.095 eV. This could be:

- 1) a methodical error, or
- 2) both values are a kind of basic values, where the other values are composite from.

Other interpretations cannot be precluded.

5 Conclusion

It has been shown in this paper that neutrino masses can be predicted by numerical calculations based on EINSTEIN-MAXWELL theory. The resulting masses for electron neutrinos come out to lie in the range being known by experiments. This is probably the first time that neutrino masses are predicted by a theory based on first principles.

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References

- [1] A. EINSTEIN, *Grundzüge der Relativitätstheorie*. A back-translation from the Four Lectures on Theory of Relativity. Akademie-Verlag Berlin, Pergamon Press Oxford, Friedrich Vieweg & Sohn Braunschweig (1969).
- [2] G. Y. RAINICH, Electrodynamics in the General Relativity Theory. *Proc. N.A.S.*, **10**, 124 (1924).
- [3] G. Y. RAINICH, Second Note Electrodynamics in the General Relativity Theory. *Proc. N.A.S.*, **10**, 294 (1924).
- [4] U. E. BRUCHHOLZ, Key Notes on a Geometric Theory of Fields, *Progress in Physics*, **5** (2), 107 (2009).
- [5] U. E. BRUCHHOLZ, Geometry of Space-Time, *Progress in Physics*, **5** (4), 65 (2009).
- [6] U. E. BRUCHHOLZ, Masses of Nuclei Constituted from a Geometric Theory of Fields, *Adv. Studies Theor. Phys.*, **7** no.19, 901 (2013).
- [7] U. E. BRUCHHOLZ, How to Deduce Masses and Further Parameters of Particles and Nuclei, *International Journal of Knowledge Based Computer System*, **1** issue 2 (2013)
- [8] K. HAGIWARA et al. (Particle Data Group), *Phys. Rev. D* **66**, 010001 (2002) (URL: <http://pdg.lbl.gov>)

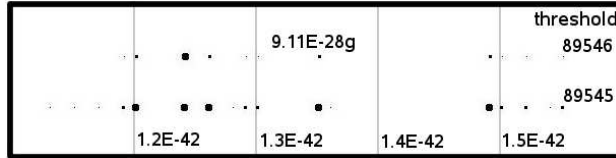


Figure 3: Tests for the free electron. Initial radius 400, 51 values, 9 times piled (459 tests)

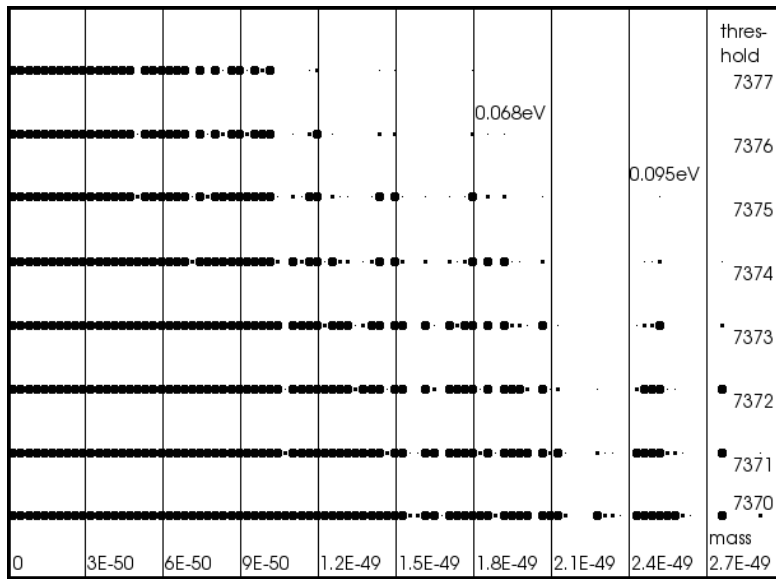


Figure 4: Tests for the electron neutrino, masses $< 0.11eV$. Initial radius 5, 100 values, 9 times piled (900 tests)

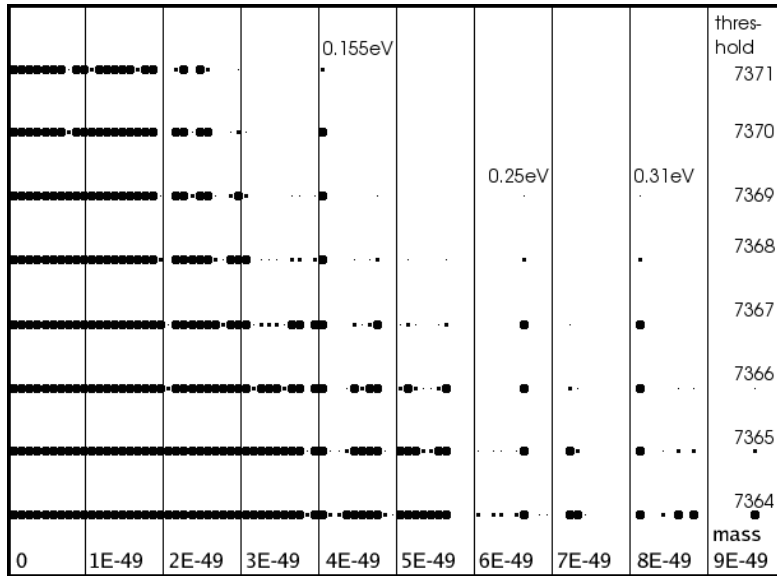


Figure 5: Tests for the electron neutrino, masses $< 0.4\text{eV}$. Initial radius 5, 100 values, 9 times piled (900 tests)

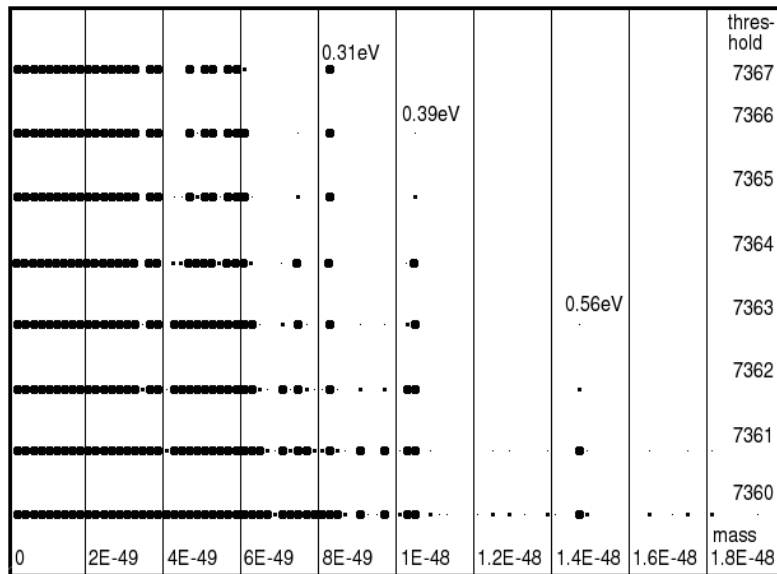


Figure 6: Tests for the electron neutrino, masses $< 1\text{eV}$. Initial radius 5, 99 values, 9 times piled (891 tests)

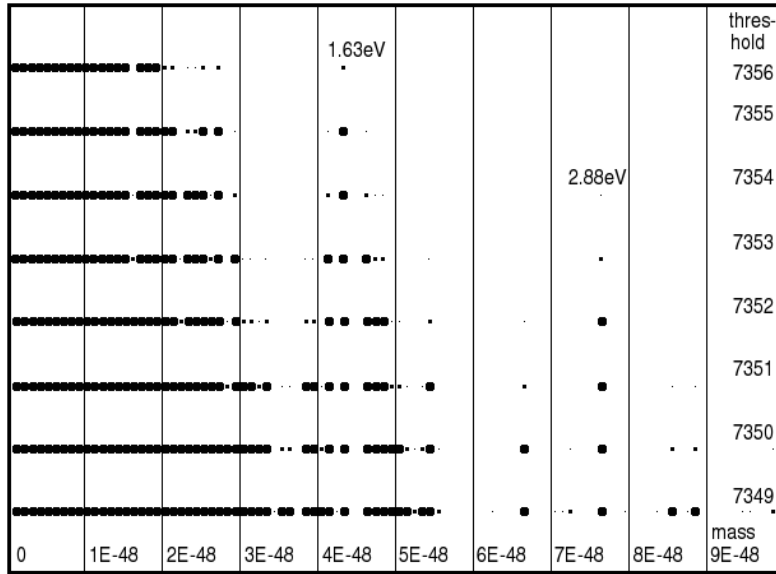


Figure 7: Tests for the electron neutrino, masses $< 4\text{eV}$. Initial radius 5, 99 values, 9 times piled (891 tests)

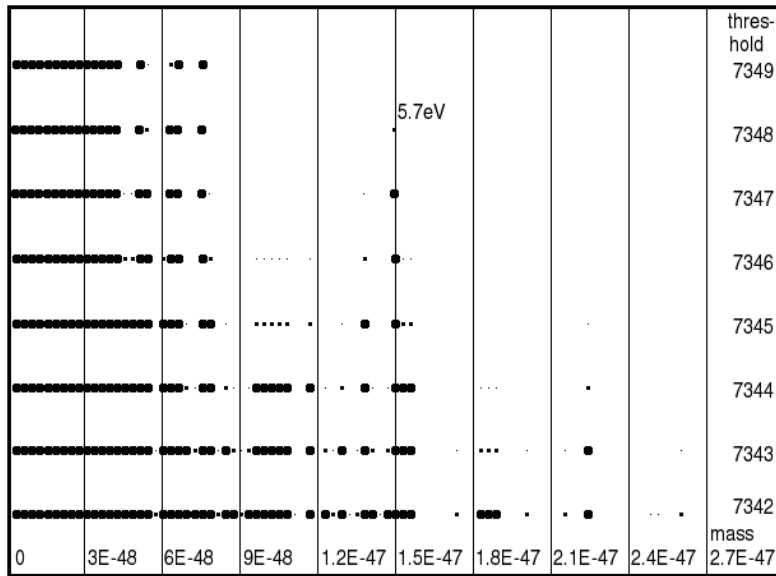


Figure 8: Tests for the electron neutrino, masses $< 11\text{eV}$. Initial radius 5, 100 values, 9 times piled (900 tests)