

FYS3500 MIDTERM HOME EXAM

Candidate number: 15522

In this assignment I mainly collaborated with candidate 155003. In addition I had some minor discussions with 15521, 15506, 15517 and I shared with them diagrams and figures.

1 Some shorter questions

a

This result can be explained by means of the nuclear shell model. According to this model each sub-level of the nucleus can be identified by 4 quantum numbers $\{n, j, l, m_l\}$. For fixed $\{n, j, l\}$ one has that the $2l + 1$ sub-levels given by $m_l = -l, -l + 1, \dots, l - 1, l$ are degenerate and each of these sub-levels can contain two nucleons with different spin z-projection ($m_s = \pm 1/2$).

At first one can notice that if one sub-shell is filled with two nucleons (of the same type) then the spin contribution of that sub-shell is 0. Having an even number of protons and neutrons means that they combine in couples in the sub-shell in a way such that the contribution to the spin of each sub-shell is 0. For what concerns the parity, each particle in the nucleus gives a contribute of $(-1)^l$ due to the orbital motion (the intrinsic parity of the nucleons is +1) and the total parity is the product of single parities times the orbital motion contributions. The product of the parities of two protons (neutrons) in the same sub-shell is always +1 since they, in particular, share the same quantum number l . Hence it is legitimate to expect an even-even nucleus to be in the state 0^+ .

Answer: spin 0, parity +

b

The process is governed by the strong interaction, hence quark flavour numbers and baryon numbers must be conserved. By observing the final state it is immediate to see that the flavour number is 0, and the same holds for the baryon number, hence the hadron must be a meson.

By looking at the [table](#) of the allowed mesons one can conclude that the isospin I must be either 0 or 1. On the other side in the final state $I_3 = 0$ and because of isospin components conservation law one concludes that $I_3 = 0$ also for the decaying pion. In conclusion the admitted couples (I, I_3) of isospins are $(0, 0)$, $(1, 0)$.

c

The nuclear potential is charge independent, in the sense that protons and neutrons feel the same nuclear force. It is repulsive for very short distances and attractive for longer distances (but the force range is short). In addition a proton feels the effect of the Coulomb interaction with the nucleus, a factor that does not affect the neutron since it is electric neutral. Figure 1 reports a sketch plot of the potentials.

d

The only two other possibilities are the combinations uuu and ddd . Since we are dealing with fermions, the *total* wavefunction must be antisymmetric, which means that it must be antisymmetric when taking account of all the degrees of freedom. This means that the spatial and spin parts of the wavefunction can be both symmetric or antisymmetric (hence symmetric in total), provided that the other contributions give an antisymmetric result. Since for $l = 0$ the spatial contribution is symmetric, the spin contribution must be itself symmetric. But this is possible only if the spins are aligned for a total of $J = 3/2$ which contradicts the hypothesis.

e

The energy to steal a neutron can be calculated by computing the difference between the binding energy $B(A, Z)$ before and after the removal (see exercise 2b for more details).

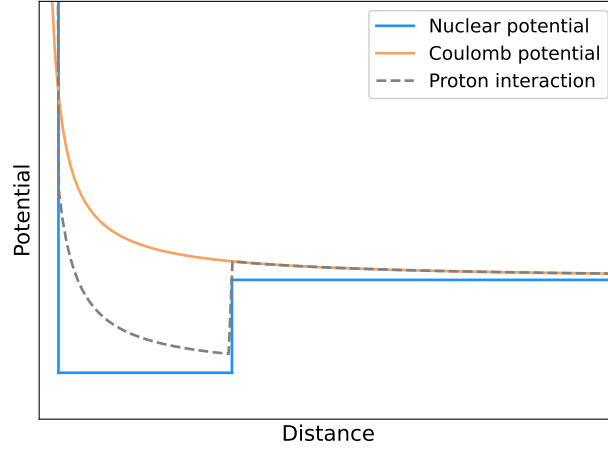


Figure 1: Nucleons interaction. The plot reports a simplified model of the nuclear potential between two nucleons (blue) and the Coulomb repulsion (orange) which is present only for the proton. The grey dotted line represents an approximated potential for the proton, which, in addition, feels also the Coulomb repulsion. The plot is purely qualitative and not in scale.

1. $^{118}\text{Sn} \rightarrow B(118, 50) - B(117, 50) \approx 9.03 \text{ MeV}/c^2$
2. $^{16}\text{O} \rightarrow B(16, 8) - B(15, 8) \approx 15.80 \text{ MeV}/c^2$
3. $^{119}\text{Sn} \rightarrow B(119, 50) - B(118, 50) \approx 6.54 \text{ MeV}/c^2$

hence it is harder to steal a neutron from ^{16}O .

f

The number of events N , the luminosity \mathcal{L} and the cross section σ are connected through the relation

$$N = \sigma L$$

and the correct number of events, that is the value that takes account of the efficiency ϵ of the detector and of the background events number N_{back} , is

$$N = \frac{N_{obs} - N_{back}}{\epsilon}$$

so that

$$\sigma = \frac{(N_{obs} - N_{back})}{\epsilon \mathcal{L}} = \frac{984}{0.25 \cdot 5} \text{ pb} \approx 0.787 \text{ pb}$$

g

Each vertex in the diagram introduces a multiplicative factor of $\alpha = 1/137$ in the probability of the decay to occur, where α is the fine structure constant. Hence one can expect the probability of a second decay to occur to be approximately $1/137$ of the first one, and the probability of a third decay to be $1/(137)^2$ of the first one.

The cross section is a direct measure of the probability of the reaction to occur, hence ratios are expected to be approximately

$$1 : \frac{1}{137} : \frac{1}{137^2}$$

2 Nuclear binding energy

a

The mass of the atom is given by

$$M(^{48}_{20}\text{Ca}^{28}) = Z(m_p + m_e) + (A - Z)m_n - \frac{B(A, Z)}{c^2}$$

where B indicates the binding energy, which in turns can be calculated via the semi-empirical formula (atomic units)

$$B(A, Z) = a_v A - a_s A^{2/3} - a_c \frac{Z(Z-1)}{A^{1/3}} - a_{sym} \frac{(A-2Z)^2}{A} + \delta(A, Z) \quad (1)$$

where

$$\delta(A, Z) = \begin{cases} \frac{a_p}{A^{1/2}} & \text{if } Z \text{ and } N \text{ are even} \\ 0 & \text{if } A \text{ is odd} \\ -\frac{a_p}{A^{1/2}} & \text{if } Z \text{ and } N \text{ are odd} \end{cases}$$

One can simply pop in the values of A and Z into the formula using the fitted coefficients (source Wikipedia) in units of MeV/c^2

$$a_v = 15.8 \quad a_s = 18.3 \quad a_c = 0.714 \quad a_{sym} = 23.2 \quad a_p = 12$$

and obtains $B(A, Z) = 412.84 \text{ MeV}/c^2$. Hence, inserting the result into 2 one obtains

$$M(^{48}\text{Ca}) \simeq 44670.64 \text{ MeV}/c^2 \simeq 47.96 \text{ u}$$

The result is close to the experimental value (47.952). One could justify eventual differences by observing that the SEMF is not a complete theoretical model, but rather a fit to experimental data of a model with some theoretical foundations. As such, it is reasonable that the model can give an overall good description of the data but cannot perfectly describe each of them: this is especially the case of lighter atoms for which the quantum shell structures of the nucleus should be taken into account.

b

One can calculate the energy to remove a neutron as the difference in energy between the atom's state with the neutron and without the neutron. The energy is the sum of two terms, one that take accounts of the mass of the particles and it is simply the sum of all the masses. The second term, instead, is the binding energy. In the initial and final state the mass energy contributions are the same (we are not making any particle disappear, we are just moving one away), hence the difference is only due to the binding energy contribution. One can calculate this contribution as

$$B_{neutron} = B(A, Z) - B(A-1, Z)$$

and using expression 1 with $A = 44$ and $Z = 20$ to calculate the two quantities one ends up with an estimation for the energy required to extract the neutron

$$\Delta E_{neutron} \approx 11.14 \text{ MeV}/c^2$$

c

One can search for the most stable atom with fixed A searching for the minimum of the mass formula

$$M_A(Z) = Z(m_p + m_e) + (A - Z)m_n - B_A(Z)$$

Since one does not know if Z is even or odd in advance, one can set the pairing term in the binding energy to 0: this will lead to a non-integer Z and the true minimum value will be either the closest greater or smaller integer (one has to check which of the two has minimum energy). By direct calculation

$$0 = \frac{dM_A}{dZ} = m_p + m_e - m_n - \frac{a_c}{A^{1/3}} (A - 2Z) + 4 \frac{a_{sym}}{A} (A - 2Z)$$

and rearranging terms

$$Z = \frac{m_n - (m_p + m_e) + 4a_{sym}A + a_c A^{2/3}}{2a_c A^{2/3} + 8a_{sym}} \approx 56.59 \quad \longrightarrow \quad Z = 56$$

Figures 2 and 3 report a plot of the nucleus masses as a function of the atomic number Z for fixed A using the mass semi-empirical formula. The odd A plot contains only one parabola, while the even A plot contains two parabolas: this is because of the pairing term in the MSEF, which is either positive if Z is even or negative if Z is odd. In both cases, atoms which do not lie in the minimum tend to decay via β -decay towards the most stable configuration (the minimum of the parabola) by changing neutrons into protons and vice-versa.

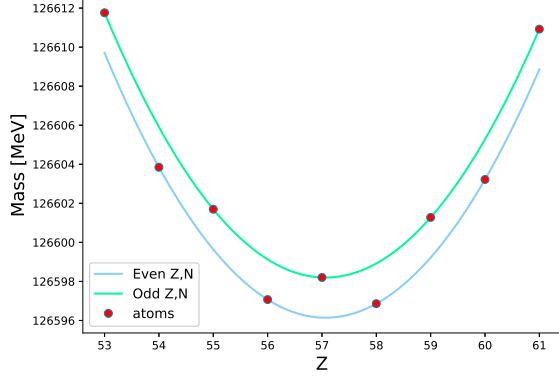


Figure 2: Even A

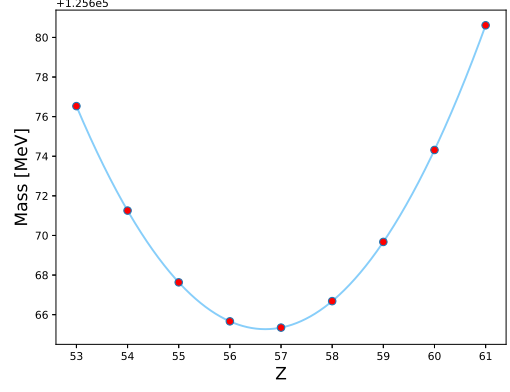


Figure 3: Odd A

3 Higgs boson decay

a

In the resting frame of the Higgs boson the invariant mass is simply the Higgs boson mass m_H . The invariant mass is a Lorentz invariant, hence this quantity is the same in the laboratory frame. More specifically, one can calculate W at the end of the process using the electrons data (laboratory frame)

$$W = \sqrt{(\sum_i E_i)^2 - |\sum \mathbf{p}_i|^2}$$

where the summation index i runs over all the electrons. By equating this expression to m_H and inserting the given values together with the electrons' masses, one obtains

$$m_H \simeq 125.20 \text{ GeV}/c^2$$

b

In principle each pair of electron-positron can be generated by the Z boson, in the sense that no rules would be violated. In other words there are 4 different ways to combine the couples of electron-positron to the bosons

1. $Z \rightarrow e_1^+ + e_1^-$, $Z^* \rightarrow e_2^+ + e_2^-$
2. $Z \rightarrow e_1^+ + e_2^-$, $Z^* \rightarrow e_2^+ + e_1^-$
3. $Z \rightarrow e_2^+ + e_1^-$, $Z^* \rightarrow e_1^+ + e_2^-$
4. $Z \rightarrow e_2^+ + e_2^-$, $Z^* \rightarrow e_1^+ + e_1^-$

One can now impose the conservation of the 4-momentum in the decay and, in particular for the Z boson, this means

$$E_Z = \sqrt{m_Z^2 + p_Z^2} = E_{e^+} + E_{e^-}$$

and by rearranging terms

$$p_Z^2 = (E_{e^+} + E_{e^-})^2 - m_Z^2$$

where

$$E_{e^+} + E_{e^-} = \sqrt{m_{e^+}^2 + p_{e^+}^2} + \sqrt{m_{e^-}^2 + p_{e^-}^2}$$

Every physical particle satisfies these relation with their physical mass but virtual particles do not (that is why they are called "off-shell"). Hence one can assume the true mass of Z and check if the corresponding momentum satisfies this relation (or if it is "on-shell") or, equivalently, one can assume the momentum by the conservation law and calculate the mass through the last relation which should result greater or equal than the true mass. I chose the first approach. By explicit calculation for the 4 cases, inserting the experimental value of the **Z boson mass**, one obtains

1. $p_Z^2 \simeq -2414 \text{ GeV}/c$
2. $p_Z^2 \simeq -7350 \text{ GeV}/c$
3. $p_Z^2 \simeq 744 \text{ GeV}/c$
4. $p_Z^2 \simeq -5872 \text{ GeV}/c$

hence only in the third case (that is the one presented in the text of the exercise) the particle can be regarded as physical and not virtual.

c

The three equations needed to solve the problem are the invariance of the invariant mass W and the shell relation $E^2 = p^2 + m^2$. More specifically

$$\begin{cases} m_H = E_Z + E_{Z^*} \\ E_Z^2 = m_Z^2 + p_Z^2 \\ E_{Z^*}^2 = m_{Z^*}^2 + p_Z^2 \end{cases}$$

this system can be solved for E_Z, E_{Z^*}, p_Z and the two energies are

$$E_{Z^*} = \frac{m_H^2 + m_{Z^*}^2 - m_Z^2}{2m_H}$$

$$E_Z = m_H - \frac{m_H^2 + m_{Z^*}^2 - m_Z^2}{2m_H}$$

d

Every quantity can in principle be computed by the conservation of 4-momentum. This means that all the four components must be conserved in each step of the process and in the center of mass frame this means that

$$(m_H, \mathbf{0}) = (E_Z + E_{Z^*}, \mathbf{p}_Z + \mathbf{p}_{Z^*}) = \left(\sum_i E_i, \sum_i \mathbf{p}_i \right)$$

where the last summation index runs over the electrons. In addition, for each physical particle, one can use the shell relation $E = \sqrt{p^2 + m^2}$. All these equations can be put together and the system of equations can be solved for the energies of the electrons.

4 Lepton universality

a

Let us consider the two decays

$$\tau^- \rightarrow \mu^- \bar{\nu}_\mu \nu_\tau$$

$$\mu^- \rightarrow e^- \bar{\nu}_e \nu_\mu$$

represented by the two Feynman diagrams

The probability for a particle to get scattered by a certain potential is proportional to the squared of the

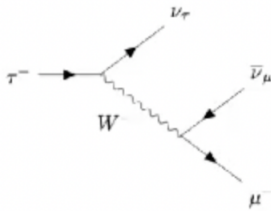


Figure 4: Feynman diagram for the reaction $\tau^- \rightarrow \mu^- \bar{\nu}_\mu \nu_\tau$

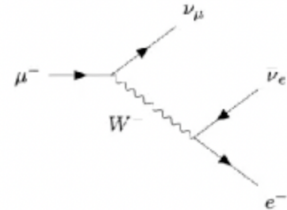


Figure 5: Feynman diagram for the reaction $\mu^- \rightarrow e^- \bar{\nu}_e \nu_\mu$

so called *amplitude* \mathcal{M} . Assuming that the coupling constant g^2 of the Yukawa potential (atomic units)

$$V(r) = -\frac{g^2}{4\pi} \frac{e^{-r/R}}{r}$$

is small compared to 4π (this means that the potential can be considered as a perturbation to the free particle solution), then the amplitude \mathcal{M} of the proces can be computed by means of the Born approximation

$$\mathcal{M}(\mathbf{p}) = \int V(\mathbf{r}) \exp(i\mathbf{r} \cdot \mathbf{r}) d^3\mathbf{r}$$

where $\mathbf{p} = \mathbf{p}_i - \mathbf{p}_f$ denotes the momentum difference between the initial and final state. By direct computation one can prove that

$$\mathcal{M}(\mathbf{p}) = \mathcal{M}(p^2) = -\frac{g^2}{p^2 + m_x^2} \quad (2)$$

where m_x denotes the mass of the propagator. A multiplicative factor of $\sqrt{2}$ must be added if the spin interaction is taken into account.

In addition, the force carrier that governs the interaction is the W boson, the mass of which is many times bigger than all the other involved masses. Hence a more suitable form of 2 is

$$\mathcal{M}(\mathbf{p}) = \mathcal{M}(p^2) = -\sqrt{2} \frac{g^2}{m_x^2} \approx 1.166 \cdot 10^{-5} (GeV)^{-2} \equiv G_F$$

and G_F is the Fermi coupling constant. This results is known as *leptons universality* and states that the amplitude of a lepton's decay process is a constant that does not depend on the chosen lepton.

At this point one can introduce the decay rate Γ which is proportional to the probability of the process to occur

$$\Gamma \approx K M^2 A$$

where A is a constant whose units are those of an energy to the fifth power, because $[M^2] = (GeV)^{-4}$. Since we expect the decay rate to depend on the generating particle, a reasonable choice for A is $A = m_i^5$ where indeed m_i denotes the mass of the generating particle. Hence

$$\Gamma \approx -\sqrt{2} K \frac{g^2}{m_x^2} m_i^5$$

Applying it to the particular case of the two given processes one obtains

$$\frac{\Gamma(\tau^- \rightarrow \mu^- \bar{\nu}_\mu \nu_\tau)}{\Gamma(\mu^- \rightarrow e^- \bar{\nu}_e \nu_\mu)} \approx \left(\frac{m_\tau}{m_\mu}\right)^5 \approx 1.34 \cdot 10^6 \quad (3)$$

b

The two relevant experimental quantities are the lifetime of the particles τ and the branching ratios of the reactions. Let us define $\Gamma_\tau \equiv \Gamma(\tau \rightarrow X \nu_\tau)$ and $\Gamma_\mu \equiv \Gamma(\tau \rightarrow Y \nu_\mu)$. Equation 3 can be rewritten as

$$\frac{\Gamma(\tau^- \rightarrow \mu^- \bar{\nu}_\mu \nu_\tau)}{\Gamma(\mu^- \rightarrow e^- \bar{\nu}_e \nu_\mu)} = \frac{\Gamma(\tau^- \rightarrow \mu^- \bar{\nu}_\mu \nu_\tau)}{\Gamma_\tau} \frac{\Gamma_\mu}{\Gamma(\mu^- \rightarrow e^- \bar{\nu}_e \nu_\mu)} \frac{\Gamma_\tau}{\Gamma_\mu} = \frac{B(\tau^- \rightarrow \mu^- \bar{\nu}_\mu \nu_\tau)}{B(\mu^- \rightarrow e^- \bar{\nu}_e \nu_\mu)} \frac{\tau_\mu}{\tau_\tau}$$

which is a more relevant form for the experimental quantities.

5 Quantum numbers

a

Parity

Let us consider a wavefunction $\psi(\mathbf{r})$. One can define the parity operator \hat{P} as an operator such that

$$\hat{P} \psi(\mathbf{r}) = \psi(-\mathbf{r})$$

Since

$$\hat{P}^2 \psi(\mathbf{r}) = \hat{P} \psi(-\mathbf{r}) = \psi(\mathbf{r}) \quad (4)$$

the eigenvalues of the parity operator are ± 1 and the corresponding eigenfunctions are the odd (eigenvalue -1) and even (eigenvalue $+1$) wavefunctions. If the wavefunction describes the state of particle, and the state is an eigenstate of the parity operator, the corresponding eigenvalue is also said to be the (intrinsic) parity of the particle. Formally

$$\hat{P}\psi(\mathbf{r}) = \pi\psi(\mathbf{r})$$

and π is the (intrinsic) parity of the particle.

Let us consider a system of two particles $A + B$ described by a wavefunction $\psi_{AB}(\mathbf{r}_A, \mathbf{r}_B)$. It can be proven that the parity of the system is given by

$$\hat{P}\psi_{AB} = \pi_A\pi_B(-1)^l\psi_{AB}$$

where l denotes the orbital angular momentum quantum number of the relative motion. Hence in a reaction of the type $A + B \rightarrow C + D$ described by a hamiltonian \hat{H} that commutes with parity, parity is conserved or, in other words

$$\pi_A\pi_B(-1)^{l_{AB}} = \pi_C\pi_D(-1)^{l_{CD}}$$

This, for example, is not the case of the weak interaction where, in general, the hamiltonian operator does not commute with the parity operator.

Charge conjugation

The charge conjugation operator is an operator that changes the sign of all the forces' quantum charges, specifically electric charge, baryon number, lepton number, flavor charges, isospin, ... In particular, if applied to a particle, the C-parity operator transforms the particle into its antiparticle. If a particle is in a state $|\psi\rangle$ then the action of the conjugation operator reads

$$\hat{C}|\psi\rangle = |\bar{\psi}\rangle$$

where $|\bar{\psi}\rangle$ represents the antiparticle's state.

The eigenstates of the C-parity operator are the systems neutral to all force charges like the photon or particle-antiparticle bound states. In this case

$$\hat{C}|\psi\rangle = C_\psi|\psi\rangle$$

and C_ψ is said to be the C-parity of the particle. As for the the parity operator, the eigenvalues can be only ± 1 .

Both P-parity and C-parity are conserved in the processes governed by all the fundamental forces **but the weak force**.

b

The notation 3S_1 stands for $L = 1$ $S = 1$ and $J = 1$. The J/Ψ Meson is a bound state of a charm quark and an anti-charm quark. By applying the C-parity operator on the state $|J/\Psi\rangle$ and noting that each quark simply swap their roles, one obtains that the C-parity of the J/ψ meson is -1 (the only contribution is due to the angular momentum).

$$\hat{C}|J/\Psi\rangle = (-1)^J|J/\Psi\rangle = -|J/\Psi\rangle$$

c

^{15}N atom has 7 protons and 8 neutrons. Since the neutrons (in the ground state) are all coupled (each sub-shell admits two nucleons) they do not contribute to the spin of the nucleus and the parity contribution is $+1$. Instead the protons configuration of the ^{15}N atom in the ground state is $(1s)^2(1p_{3/2})^4(1p_{1/2})^1$: only the last proton contributes to the interested quantities. Since the orbital p represents the quantum number $l = 1$ the parity of the the nucleus is the product of the neutrons an protons contribution, that is $(+1) \cdot (-1)^1 = -1$. The spin is then $1/2$.

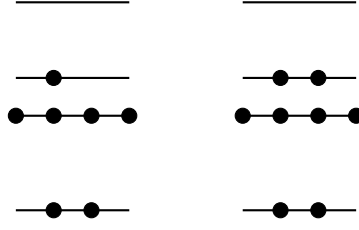


Figure 6: Ground state ^{15}N nucleus configuration

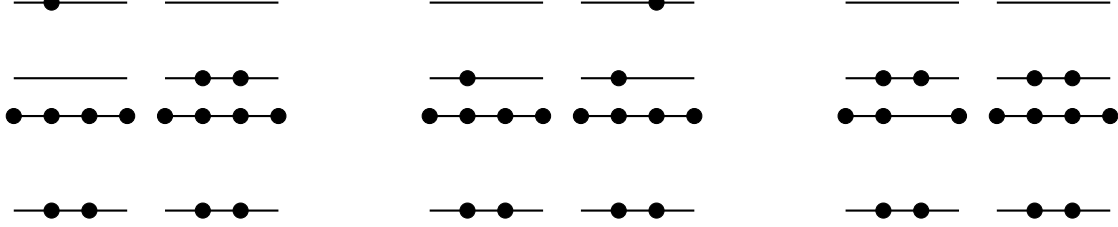


Figure 7: First three excited states of a ^{15}N nucleus

There are three possible alternatives for the first excited state, all of which consist in moving a proton or a neutron to another level starting from the ground state configuration.

1. Move the $1p_{1/2}$ proton to the $1d_{5/2}$ level (figure 7 on the left)
 \rightarrow the new proton configuration is $(1s)^2(1p_{3/2})^4(1d_{5/2})^1$ and the spin-parity is $5/2^+$.
2. Move the $1p_{1/2}$ neutron to the $1d_{5/2}$ level (figure 7 in the middle)
 \rightarrow the new neutron configuration is $(1s)^2(1p_{3/2})^4(1p_{1/2})^1(1d_{5/2})^1$. The parity here is non-trivial because of the angular momemnta addition rules.
3. Move the $1p_{3/2}$ proton to the $1p_{1/2}$ level (figure 7 on the right)
 \rightarrow the new proton configuration is $(1s)^2(1p_{3/2})^3(1p_{1/2})^2$ and the spin-parity is $3/2^-$.

By looking at the suggested table, one can notice that the first excited state is described by configuration 1, the second excited state by configuration 2 and the third excited state is described by configuration 3.

d

The third excited state is the one that has $J^p = \frac{3}{2}^-$. From this state the nucleus can decay into the second excited state, the first excited state and the ground state. Let us first consider the decay to the ground state. When the excited proton returns to the $1p_{3/2}$ level there is no change in parity and one has that $\Delta J = J_{exc} - J_{ground} = 1$. This means that one must have odd-L magnetic fields and even-L electric fields.

Because of the general selection rules in the gamma decays one has that, said L the photon's angular momentum

$$|J_{exc} - J_{ground}| \leq L \leq J_{exc} + J_{ground} \quad \rightarrow \quad 1 \leq L \leq 2$$

The allowed states are

$$M_1, E_2$$

For $A = 15$ one has that (reference chapter 10.3 Krane)

$$\frac{\lambda(E_2)}{\lambda(M_1)} \approx 10^{-3} \quad (5)$$

hence we expect the M_1 decay to be the most probable one, but with a non-neglectable contribution of E_2 .

In an analogous way one can find that for the decays to the second and first excited states the relations reported in table 1. For the transition $2 \rightarrow 1$ one has that the dominant field is E_2 ($2 \leq L \leq 3$ and no parity change) while for the decay $2 \rightarrow 0$ the dominant field is E_1 ($0 \leq L \leq 1$ and parity change). Hence from the second excited state the nucleus can decay to the ground state via E_1 radiation, or it can decay to the first excited state via E_2 radiation. By comparing probabilities (reference chapter 10.3 Krane)

$$\frac{P_{2 \rightarrow 0}}{P_{2 \rightarrow 1}} \approx \frac{\lambda(E_1)}{\lambda(E_2)} = \frac{1}{7.3} \cdot 10^7 \cdot A^{-2/3} > 1$$

Transition	ΔE	$\Delta\pi$	L	Fields	Dominants
$3 \rightarrow 0$	$\approx 6.3 \text{ MeV}$	no	$1 \leq L \leq 2$	M_1, E_2	M_1, E_2
$3 \rightarrow 1$	$\approx 1 \text{ MeV}$	yes	$1 \leq L \leq 4$	E_1, M_2, E_3, M_4	E_1
$3 \rightarrow 2$	$\approx 1 \text{ MeV}$	yes	$1 \leq L \leq 2$	E_1, M_2	E_1

Table 1: γ transitions from the 3rd excited state

6 Allowed, suppressed and forbidden processes

I use the definition of decay stated [here](#).

I do not follow the order of the questions as given, but everything is reported here.

Process	Type	Interaction	Allowed	Suppressed	Reason not allowed
1	decay	electroweak	yes	no	charge
2			no		
3	decay	electroweak	yes	yes	
4	decay	weak	yes	no	lepton number
5	decay	strong	yes	no	
6	reaction	electroweak + strong	yes	no	
7	decay	weak	yes	no	
8			no		

Process	Interaction	Typical Interaction time	Lifetime
1	electromagnetic	$10^{-14} \sim 10^{-20}$	Not found
3	weak	$10^{-8} \sim 10^{-13}$	10^{-12}
4	weak	$10^{-8} \sim 10^{-13}$	10^{-10}
5	strong	$< 10^{-22}$	10^{-24}

Process 3 is not the most common decay for the bottom quark, hence extremely rare and suppressed. Processes 7 and 8 differ by the presence of 2 antineutrinos. The process 8, namely the neutrinoless double beta decay, is not allowed according to the Standard model because it violates the lepton number conservation. The observation of such decay would indeed confirm the Majorana nature of the neutrinos and the lepton number conservation rule would be violated.

When the neutron decays into 2 protons and 2 electrons without the neutrino, the energy is equally splitted to the generated particles according to their masses. On the opposite, when the antineutrinos are produced, part of the energy would be transferred to them, but the portion is not fixed, hence it would produce a continuous energy spectrum.

7 Radioactive decay

a

Said $N(t)$ the number of atoms present at time t , the radioactive decay law states that

$$\mathcal{A} = -\frac{dN}{dt} = \lambda N \quad (6)$$

where λ is called the decay constant.

By integrating it in time one obtains

$$N(t) = N_0 e^{-\lambda t} \quad (7)$$

Now if we consider a chain decay of the type $A \rightarrow B \rightarrow C$ with initial conditions

$$N_A(t=0) = N_0 \quad N_B(t=0) = N_C(t=0) = 0$$

one can immediately obtain the number of A atoms as a function of time by applying [8](#)

$$N_A(t) = N_0 e^{-\lambda_A t}$$

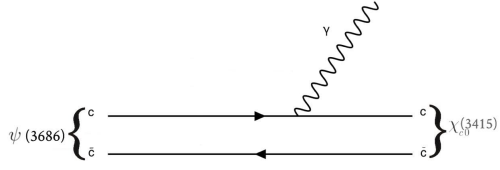


Figure 8: $\psi(3686) \rightarrow \chi_0(3145) + \gamma$

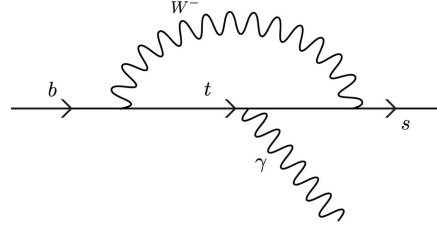


Figure 9: $b \rightarrow s\gamma$

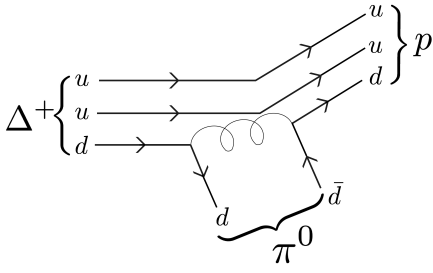


Figure 10: $\Delta^+ \rightarrow p\pi^0$

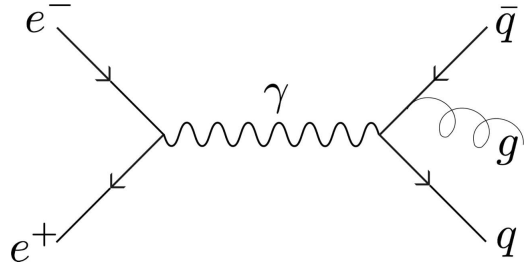


Figure 11: $e^+e^- \rightarrow q\bar{q}g$

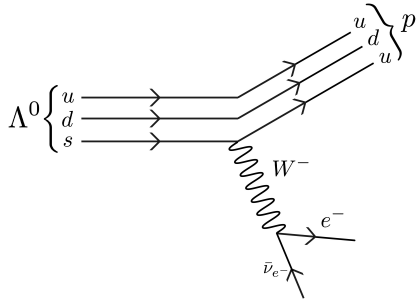


Figure 12: $\Lambda^0 \rightarrow p + e^- + \bar{\nu}_e$

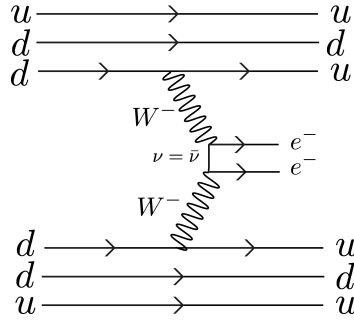


Figure 13: $2n \rightarrow 2p + 2e^-$

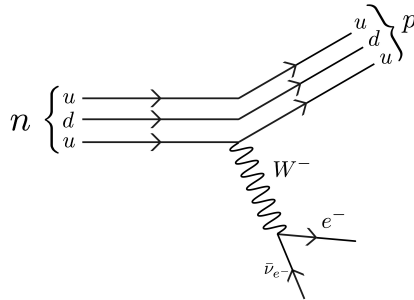


Figure 14: $2n \rightarrow 2p + 2e^- + 2\bar{\nu}_e$

To study the number of atoms of type B one has to take account of two factors: the number of "created" atoms by means of A 's decay, and the number of atoms decayed into C . This means that the radioactive law reads

$$\frac{dN_B(t)}{dt} = -\lambda_B N_B(t) + \lambda_A N_{A \rightarrow B}(t)$$

where the first term on the right side member represents the B 's decay and the the last term takes account of the number of B 's obtained by A 's decay.

One can now substitute 7 obtaining

$$\frac{dN_B(t)}{dt} + \lambda_B N_B(t) = \lambda_A N_0 e^{-\lambda_A t}$$

The general solution of this differential equation is

$$N_B(t) = e^{-A(t)} \left(N_B(0) + \int_0^t \lambda_A N_0 e^{-\lambda_A s} e^{A(s)} ds \right)$$

where $A(s) = \lambda_B \int dt = \lambda_B t$. Hence

$$\begin{aligned} N_B(t) &= \lambda_A e^{-\lambda_B t} \int_0^t N_0 e^{-(\lambda_A - \lambda_B)s} ds = \frac{N_0 \lambda_A e^{-\lambda_B t}}{\lambda_A - \lambda_B} \left(1 - e^{-(\lambda_A - \lambda_B)t} \right) = \\ &= N_0 \lambda_A \frac{e^{-\lambda_A t} - e^{-\lambda_B t}}{\lambda_B - \lambda_A} \end{aligned}$$

b

Let us consider the single-decay law

$$N(t) = N_0 e^{-\lambda t}$$

One can notice here that λ 's physical units are inverse of time. The time $t_{1/2}$ is defined as the time necessary to halve the number of atoms from N_0 . This can be found by imposing $N(t) = N_0/2$ and solving for t one obtains

$$t_{1/2} = \frac{1}{\lambda} \log 2 \equiv \tau \log 2$$

where I introduced the mean lifetime $\tau \equiv 1/\lambda$. This last term, which has units of time, has a particular meaning. To understand this, one can first think of

$$p(t) = \frac{N(t)}{\int N(t) dt}$$

as a probability distribution function related to the probability of a particle to decay. One can then calculate the expected lifetime of such particle

$$\langle t \rangle = \frac{\int t N(t) dt}{\int N(t) dt} = \frac{\int_0^\infty t e^{-\lambda t} dt}{\int_0^\infty e^{-\lambda t} dt} = \frac{1}{\lambda}$$

and it is exactly τ .

Another interesting quantity is the natural decay width Γ . One can make plot the probability of a decay to occur as function of the decaying particle's energy: this function is peaked around a most probable value E_0 with a distribution described by the Breit Wigner formula. The half width at half maximum of this distribution is the factor Γ which has units of energy and is related to the average lifetime of the particle through the relation energy-time uncertainty relation

$$\Delta E \cdot \Delta t \approx \frac{\hbar}{2}$$

which in this particular case reads

$$\frac{\Gamma}{2} \cdot \tau \approx \frac{\hbar}{2}$$

or

$$\tau \approx \frac{\hbar}{\Gamma}$$

c

$1mCi = 37MBq$. Taking the solution

$$N_1(t) = N_0 e^{-\lambda t}$$

and imposing $\mathcal{A}(t=0) = -\frac{dN_1(t)}{dt}|_{t=0} = 37MBq$ one obtains

$$N_0 = \frac{37MBq}{\lambda} \simeq 29689798324$$

In this exercise I found some ambiguity with the definition of the activity in a decay chain. In fact, for a single decay, it is always true that

$$\frac{dN}{dt} = -\lambda N \quad (8)$$

hence one can calculate the activity both as $-\frac{dN}{dt}$ or $-\lambda N$. But when it comes to decay chains, it is not anymore true that for each species a relation as 8 sussists, hence there are two possible definitions of the activity. Marting and Shaw (chapter 1.6.4 third edition) defines tha activity as $\mathcal{A} = -\frac{dN}{dt}$, but other sources, such as [this](#), use $\mathcal{A} = -\lambda N$ as definition. I chose to attach to the latter.

d

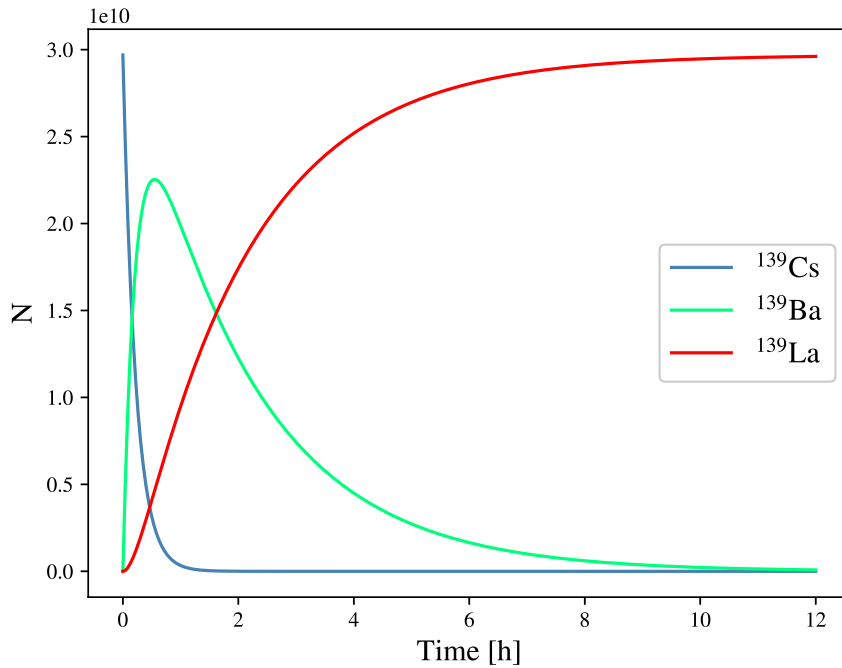


Figure 15: Number of atoms as a function of time in the decay chain

Figures 15 and 16 report respectively the number of atoms and the activities as functions of time. Of course at time $t = 0$ the only present nuclei are ^{139}Cs , and as the time passes the number of ^{139}Cs nuclei decreases, while the two other elements start forming in chain. The activity of the first reaction is proportional to the the number of atoms of ^{139}Cs , hence it is logical that the maximum activity of the first reaction is at the beginning of the simulation. For what concerns ^{139}Ba , the activity grows as more ^{139}Ba atoms are produced but a certain point the rate of production of ^{139}La overcomes the rate of decay of ^{139}Cs , and the ^{139}Ba activity starts decreasing. Finally, the activity of ^{139}La is always 0 because no decays are observed

e

The maximum activity of ^{139}Ba is registered after 33 minutes, with a maximum value of $3.1391 MBq = 0.0848 mCi$.

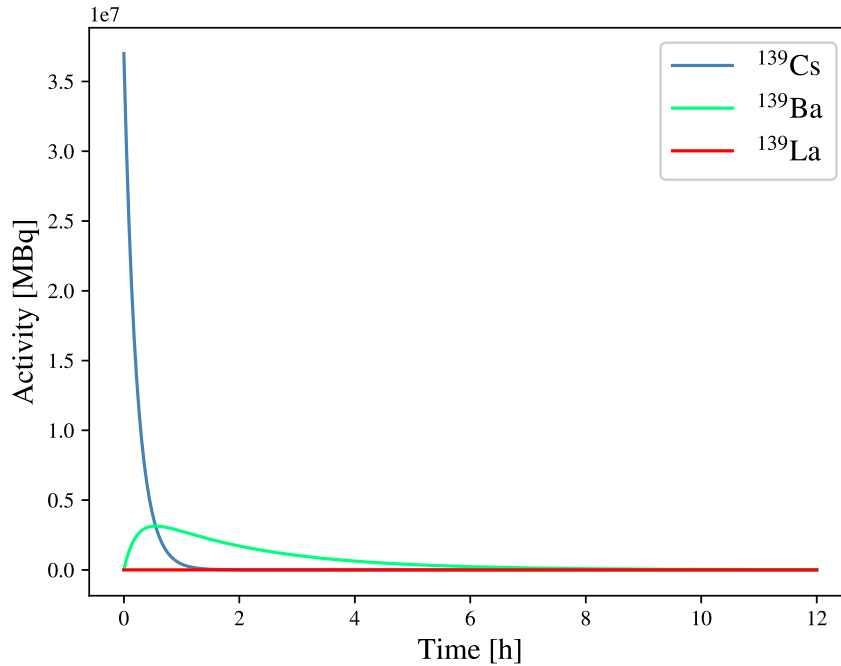


Figure 16: Activities as a function of time in the decay chain

f

The activities of ^{139}Cs and ^{139}Ba become equal after 0.33 minutes (see script below), that is in the peak of the activity of ^{139}Ba .

8 Code for exercise 1

```
from matplotlib import pyplot as plt
import numpy as np
```

```
# LSQ Coefficients (source Wikipedia)
```

```
a_v = 15.8
```

```
a_s = 18.3
```

```
a_c = 0.714
```

```
a_symm = 23.2
```

```
a_p = 12
```

```
# Other constants
```

```
mp = 938.272 # Proton mass in MeV/c^2
```

```
me = 0.511 # Electron mass in MeV/c^2
```

```
mn = 939.565 # Neutron mass in MeV/c^2
```

```
MeV_to_u = 931.5 # conv factor from MeV/c^2 to atomic units (divide)
```

```
def B(A, Z):
```

```
    delta = 0
```

```
    if A%2 == 0 and Z%2 == 0:
```

```
        delta = a_p/np.sqrt(A)
```

```
    elif A%2 == 0 and Z%2 == 1:
```

```
        delta = -a_p/np.sqrt(A)
```

```
    return + a_v*A - a_s*A**(2/3) - a_c*Z*(Z-1)/A**(1/3) - a_symm*(A-2*Z)**2/A + delta
```

```
def Coulomb(r, a=1):
```

```
    return a/r
```

```

def protonPotential(r, a=1):
    if r<0.5:
        return 100
    elif r<5:
        return Coulomb(r, a) - 10
    else:
        return Coulomb(r, a)

# Repulsion vertical line
repulsion = (1000, -10)
xrepulsion = (0.5, 0.5)
# Well horizontal line
well = (-10, -10)
xwell = (0.5, 5)
# Attraction vertical line
attraction = (-10, 0)
xattraction = (5, 5)
# Horizontal line
horiz = (0, 0)
xhoriz = (5, 100)
# Coulomb curve
xline = np.linspace(0.01, 100, 1000)
yline = Coulomb(xline, a=10)
# Proton potential
zline = [protonPotential(xx, a=10) for xx in xline]

plt.plot(xrepulsion, repulsion, color='dodgerblue',
         label="Nuclear potential", linewidth=2.0)
plt.plot(xwell, well, color='dodgerblue', linewidth=2.0)
plt.plot(xattraction, attraction, color='dodgerblue', linewidth=2.0)
plt.plot(xhoriz, horiz, color='dodgerblue', linewidth=2.0)
plt.plot(xline, yline, color='sandybrown',
         label='Coulomb potential', linewidth=2.0)
plt.plot(xline, zline, '—', color='gray',
         label='Proton interaction', linewidth=2.0)
plt.legend(fontsize=14)
plt.xlim(0, 15)
plt.ylim(-15, 30)
plt.xlabel('Distance', fontsize=16)
plt.ylabel('Potential', fontsize=16)
plt.xticks([], [])
plt.yticks([], [])
plt.show()

# 1e) neutron energy
print("118 Sn:", B(118, 50) - B(117, 50))
print("16 O:", B(16, 8) - B(15, 8))
print("119 Sn:", B(119, 50) - B(118, 50))

```

9 Code for exercise 2

```

import numpy as np
from matplotlib import pyplot as plt

# LSQ Coefficients (source Wikipedia)
a_v = 15.8
a_s = 18.3
a_c = 0.714
a_symm = 23.2
a_p = 12

```

```

# Other constants
mp = 938.272 # Proton mass in MeV/c^2
me = 0.511 # Electron mass in MeV/c^2
mn = 939.565 # Neutron mass in MeV/c^2
MeV_to_u = 931.5 # conv factor from MeV/c^2 to atomic units (divide)

# Mass semi-empirical formula
def M(A, Z, force_even_Z=False, force_odd_Z=False):
    bind = B(A, Z, force_even_Z=force_even_Z, force_odd_Z=force_odd_Z)
    return Z*(mp+me) + (A-Z)*mn - bind

# Binding energy
def B(A, Z, force_even_Z=False, force_odd_Z=False):
    mass_term = a_v*A - a_s*A**(2/3) - a_c*Z*(Z-1)/A**(1/3)
    delta = 0
    if (A%2 == 0 and Z%2 == 0 and not force_odd_Z) or force_even_Z:
        delta = a_p/np.sqrt(A)
    elif (A%2 == 0 and Z%2 == 1) or force_odd_Z:
        delta = -a_p/np.sqrt(A)
    return mass_term - a_symm*(A-2*Z)**2/A + delta

# Find most stable atom for fixed A
def most_stable(A):
    Zbest_rounded = int((mn - mp - me + 4*A*a_symm + a_c*A**(2/3)) /
        (2*a_c*A**(2/3) + 8*a_symm))
    if M(A, Zbest_rounded) > M(A, Zbest_rounded + 1):
        return Zbest_rounded + 1
    else:
        return Zbest_rounded

print("Mass of 48Ca:", M(48, 20)/MeV_to_u)
print("Energy to extract a neutron from 44Ca:", (B(44, 20) - B(43, 20)))
print("Most stable Z for fixed A=136:", most_stable(136))

# Z values
Zline = np.linspace(53, 61, 1000)
Zvals = np.arange(53, 62, 1)

# Calculate masses line
Mvals_oddA = M(135, Zline)
Mvals_evenA_evenZ = [M(136, zz, force_even_Z=True) for zz in Zline]
Mvals_evenA_oddZ = [M(136, zz, force_odd_Z=True) for zz in Zline]
# Discrete masses value
M_oddA = [M(135, zz) for zz in Zvals]
M_evenA = [M(136, zz) for zz in Zvals]

# Odd A
plt.plot(Zline, Mvals_oddA, linewidth=1.8, color='lightskyblue')
plt.plot(Zvals, M_oddA, '.', markersize=14,
    markerfacecolor='red', linewidth=1.8)
plt.xlabel('Z', fontsize=16)
plt.ylabel('Mass [MeV]', fontsize=16)
plt.show()

# Even A
plt.plot(Zline, Mvals_evenA_evenZ, label="Even Z,N",
    linewidth=1.8, color='lightskyblue')
plt.plot(Zline, Mvals_evenA_oddZ, label="Odd Z,N",
    linewidth=1.8, color='mediumspringgreen')
plt.plot(Zvals, M_evenA, '.', markersize=14, markerfacecolor='red',

```

```

        linewidth=1.8, label='atoms')
plt.xlabel('Z', fontsize=16)
plt.ylabel('Mass [MeV]', fontsize=16)
plt.legend(fontsize=12)
plt.show()

```

10 Code for exercise 3

```

import numpy as np

p1 = [0.814, -13.810, -8.409]
p2 = [-27.649, -1.511, -20.665]
p3 = [38.632, 13.361, 44.775]
p4 = [-13.443, 2.514, -5.853]

m_Z = 91.188 #GeV
m_e = 5.1e-4 #Gev

def norm3(x):
    return np.sqrt(x[0]**2 + x[1]**2 + x[2]**2)

def Energy(p, m):
    return np.sqrt(m**2 + norm3(p)**2)

en_Z = Energy(p2, m_e) + Energy(p4, m_e)
print("Squared momentum of Z:", en_Z**2 - m_Z**2)

En_sum = Energy(p1, m_e) + Energy(p2, m_e) + Energy(p3, m_e) + Energy(p4, m_e)
Etot = np.sqrt((En_sum)**2 - norm3(p1+p2+p3+p4)**2)

print("Higgs boson mass:", Etot)

```

11 Code for exercise 7

```

from matplotlib import pyplot as plt
import numpy as np
import matplotlib

matplotlib.rcParams['mathtext.fontset'] = 'stix'
matplotlib.rcParams['font.family'] = 'STIXGeneral'

T_half_Cs = 9.27/60 # Half life time of 139Cs in hours
T_half_Ba = 82.93/60 # Half life time of 139Ba in hours
lam_Cs = np.log(2)/T_half_Cs # Decay width of 139Cs
lam_Ba = np.log(2)/T_half_Ba # Decay width of 139Ba

N0 = 37e6/(lam_Cs/60/60)

print("N0:", int(N0))

# First decay
def N1(N0, t, lam1):
    return N0*np.exp(-lam1*t)

# Second decay
def N2(N0, t, lam1, lam2):
    return N0 * lam1 * (np.exp(-lam1*t) - np.exp(-lam2*t)) / (lam2 - lam1)

# Third decay
def N3(N0, t, lam1, lam2):

```



```

    return N0*(lam1*(1-np.exp(-lam2*t))-lam2*(1-np.exp(-lam1*t)))/(lam1-lam2)

# Activity of the first reaction
def A1(N0, t, lam1):
    return lam1/3600*N1(N0, t, lam1)

def A2(N0, t, lam1, lam2):
    return lam2/3600*N2(N0, t, lam1, lam2)

def A3(N0, t, lam1, lam2):
    return [0 for _ in range(len(t))]

# Time values to evaluate the numbers of species
time_ticks = np.linspace(0, 12, 100000)

# N(t)
N_Cs = N1(N0, time_ticks, lam_Cs)
N_Ba = N2(N0, time_ticks, lam_Cs, lam_Ba)
N_La = N3(N0, time_ticks, lam_Cs, lam_Ba)

# Activities
A_Cs = A1(N0, time_ticks, lam_Cs)
A_Ba = A2(N0, time_ticks, lam_Cs, lam_Ba)
A_La = A3(N0, time_ticks, lam_Cs, lam_Ba)

# Finding requested quantities
max_A_Ba = print("Maximum activity: %.3f at time %.3f hours" %
(max(A_Ba), time_ticks[np.where(A_Ba == max(A_Ba))[0][0]]))
diff = np.abs(A_Ba - A_Cs)
print("Activities of Cs and Ba are equal at time %.3f hours" %
(time_ticks[np.where(diff == min(diff))]))

# Plotting N(t)
plt.plot(time_ticks, N_Cs, label='$^{139}$Cs', color='steelblue')
plt.plot(time_ticks, N_Ba, label='$^{139}$Ba', color='springgreen')
plt.plot(time_ticks, N_La, label='$^{139}$La', color='red')
plt.xlabel('Time [h]', fontsize=14)
plt.ylabel('N', fontsize=14)
plt.legend(fontsize=14)
plt.show()

# Plotting activities
plt.plot(time_ticks, A_Cs, label='$^{139}$Cs', color='steelblue')
plt.plot(time_ticks, A_Ba, label='$^{139}$Ba', color='springgreen')
plt.plot(time_ticks, A_La, label='$^{139}$La', color='red')
plt.xlabel('Time [h]', fontsize=14)
plt.ylabel('Activity [MBq]', fontsize=14)
plt.legend(fontsize=14)
plt.show()

```