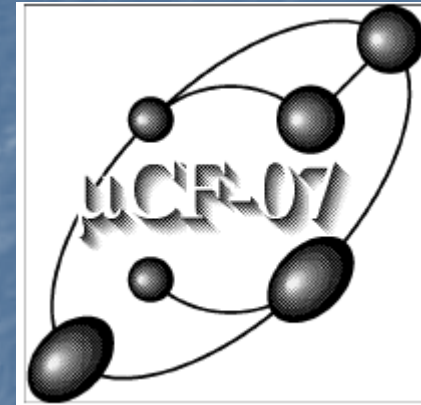


Calculation of Muonic Atom cascade Dynamics in D-T Mixtures

JINR, Dubna, Mucf07
19 June 2007

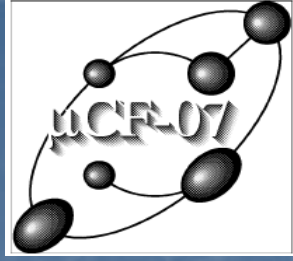
S. Z. Kalantari and M. Raisi
Department of Physics
Isfahan University of Technology
Isfahan, IRAN



- **Introduction**
- **Cascade processes of muonic atoms**
- **Kinetics of Cascade processes of muonic atoms**
- **Cascade Processes of Kaonic atoms**
- **Forthcoming Research**



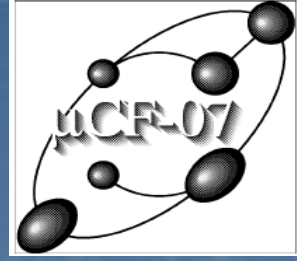
Introduction



- Why the calculations of kinetics of cascade processes of muonic atoms are important.
 1. Cascade processes of muonic atoms affect the muon catalyzed fusion.
 2. It can help to realize the kinetics of cascade processes of the other Exotic atoms.



Introduction



$$\pi^- \longrightarrow \mu^-(4.3 \text{ MeV}) + \bar{\nu}$$

I) Ionization $\mu + (X)_{gs} \longrightarrow \mu + e + X^+$

II) Inelastic collision $\mu + (X)_{gs} \longrightarrow \mu + (X)_{nl}$

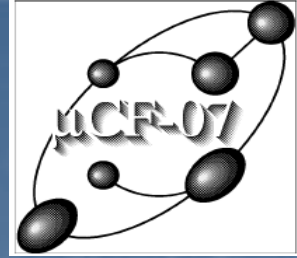
III) Captureing $\mu + (X)_{gs} \longrightarrow (\mu X)_{nl} (1\text{eV} \sim 10\text{eV}) + e \quad \text{if } V_\mu \sim V_e$

$$x = D, D_1$$

In high energy with the Born approximation, the predicted time for stopping muon is in the order of : 10^{-12} s to 10^{-11} s



Introduction



$$r_n = \frac{n^2 \hbar^2}{\bar{m}_\mu z e^2} \quad E_n = -\frac{z e^2}{2 r_n} = -\frac{\bar{m}_\mu z^2 e^4}{2 n^2 \hbar^2} \quad V_n = \frac{z e^2}{n \hbar}$$

$$\bar{m}_\mu, z$$

are Muon reduced mass and charge of nuclues, Resp.

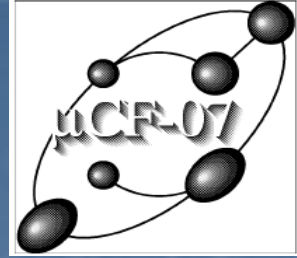
They have a peak around the following values. (J.S. Cohen, Phys. Rev. ,27(1983)167.

$$n_{peak} = n_\mu^{\frac{1}{2}} \approx 12$$

$$l_{peak} = (2 m_\mu) \left(1 - \frac{m_\mu}{2 m}\right)^{\frac{1}{2}}$$



Cascade processes of muonic atoms

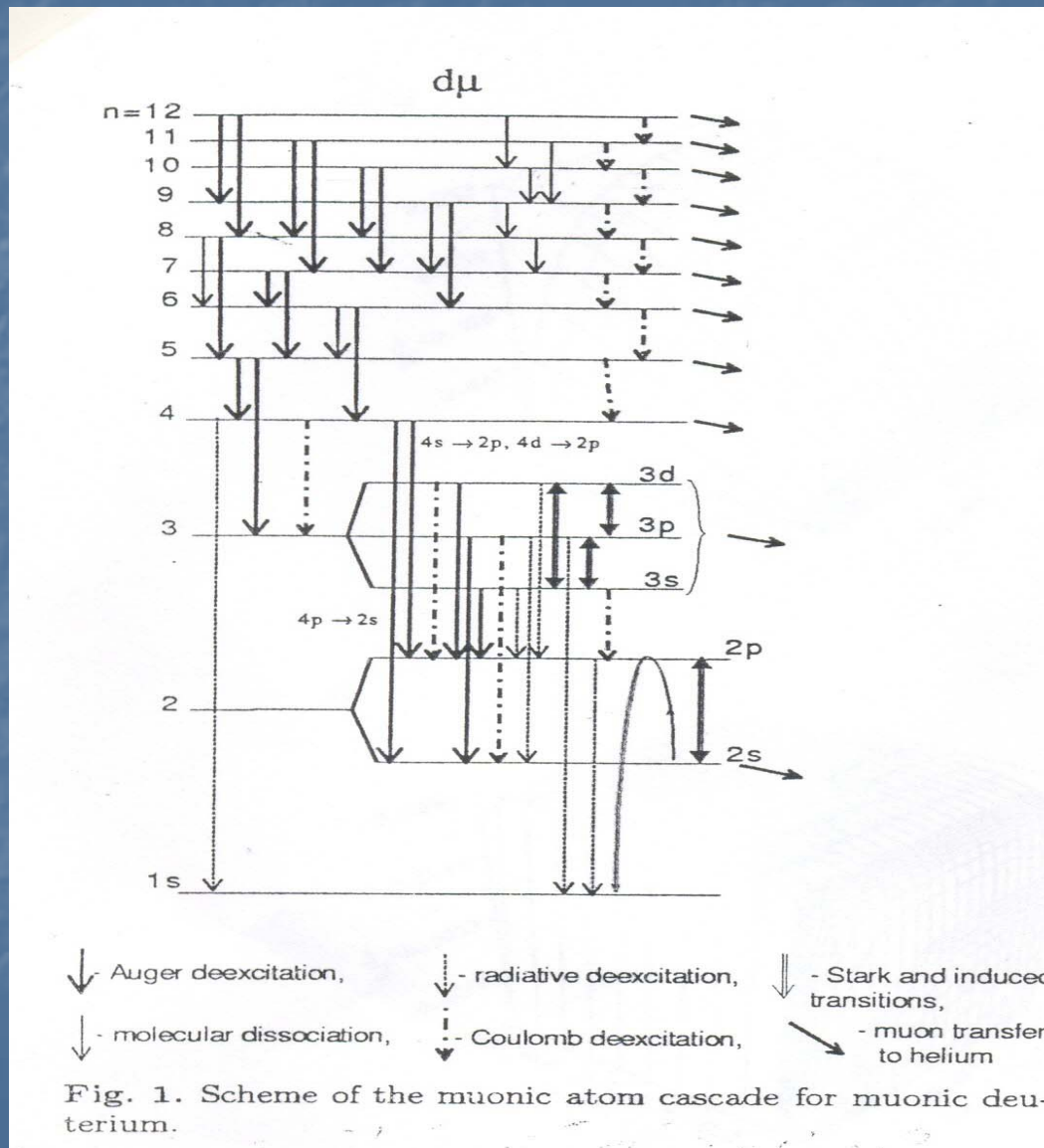
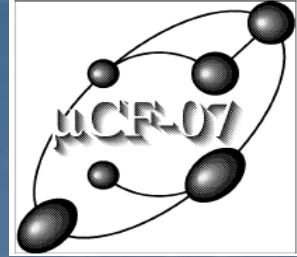


■ Cascade processes

Mechanism	Process	Energy Dependence	Density Dependence
Chemical dissociation	$(x\mu)_{n_i} + x'_2 \rightarrow x'^+ + x^- + (x\mu)_{n_f}, n_f < n_i, l_f = l_i$	Weak	Linear
Radiative	$(x\mu)_{n_i} \rightarrow (x\mu)_{n_f} + \gamma$	None	None
Auger	$(x\mu)_{n_i} + x'_2 \rightarrow x'^+ + e + (x\mu)_{n_f}$	Weak	Linear
Stark mixing	$(x\mu)_{n_l} + X \rightarrow (x\mu)_{n_l'} + X$	Moderate	Linear
Scattering	$(x\mu)_n + X \rightarrow (x\mu)_n + X'$	Strong	Linear
Coulomb	$(x\mu)_{n_i} + x' \rightarrow (x\mu)_{n_f} + x'$	Strong	Linear
Transfer	$(x\mu)_n + x' \rightarrow (x'\mu)_n + x$	Strong	Linear

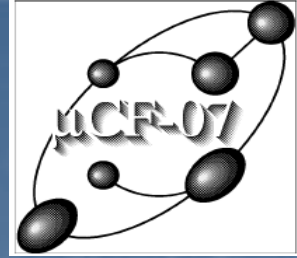


Cascade processes of muonic atoms

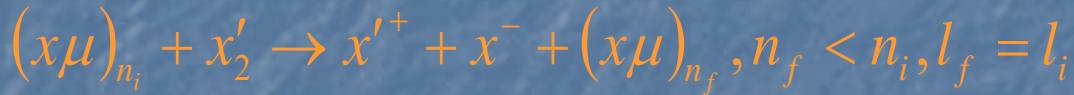




Cascade processes of muonic atoms



I) Chemical Dissociation



$$\Gamma_{nn'} = \frac{N}{2} v \pi a_n^2 (s^{-1}), \Delta E_{nn'} = E_n - E_{n'}$$

II) Radiative Transition



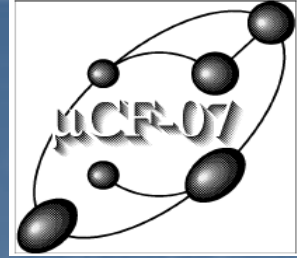
$$\Gamma_{nn'} = \frac{4}{3} (\Delta E)^3 \left| R_n^{n'} \right|^2 \overline{M}^{-2} \times 1.60 \times 10^{-10} (s^{-1})$$

Average over l-value of initial state and summed over the final state.

Bethe H.A. and Salpeter E.E.1975, Quantum Mechanics of one and two electron atoms



Cascade processes of muonic atoms



III) External Auger effect

$$\lambda_{if} = N\nu_i\sigma_{if} = 4.3 \times 10^{15} (\text{sec}^{-1}) (R^{n'l'}_{nl})^2 \overline{M}^{-2} (2\Delta E + 1.39)^{-1/2}$$

$$R_{nl} = \left(\frac{z}{a_0}\right)^{\frac{2}{3}} e^{zr/na_0} f_{nl}\left(\frac{zr}{a_0}\right) \quad a_0 = \frac{1}{M}$$

$$\Delta E = \Delta E_{x\mu} - \delta_H$$

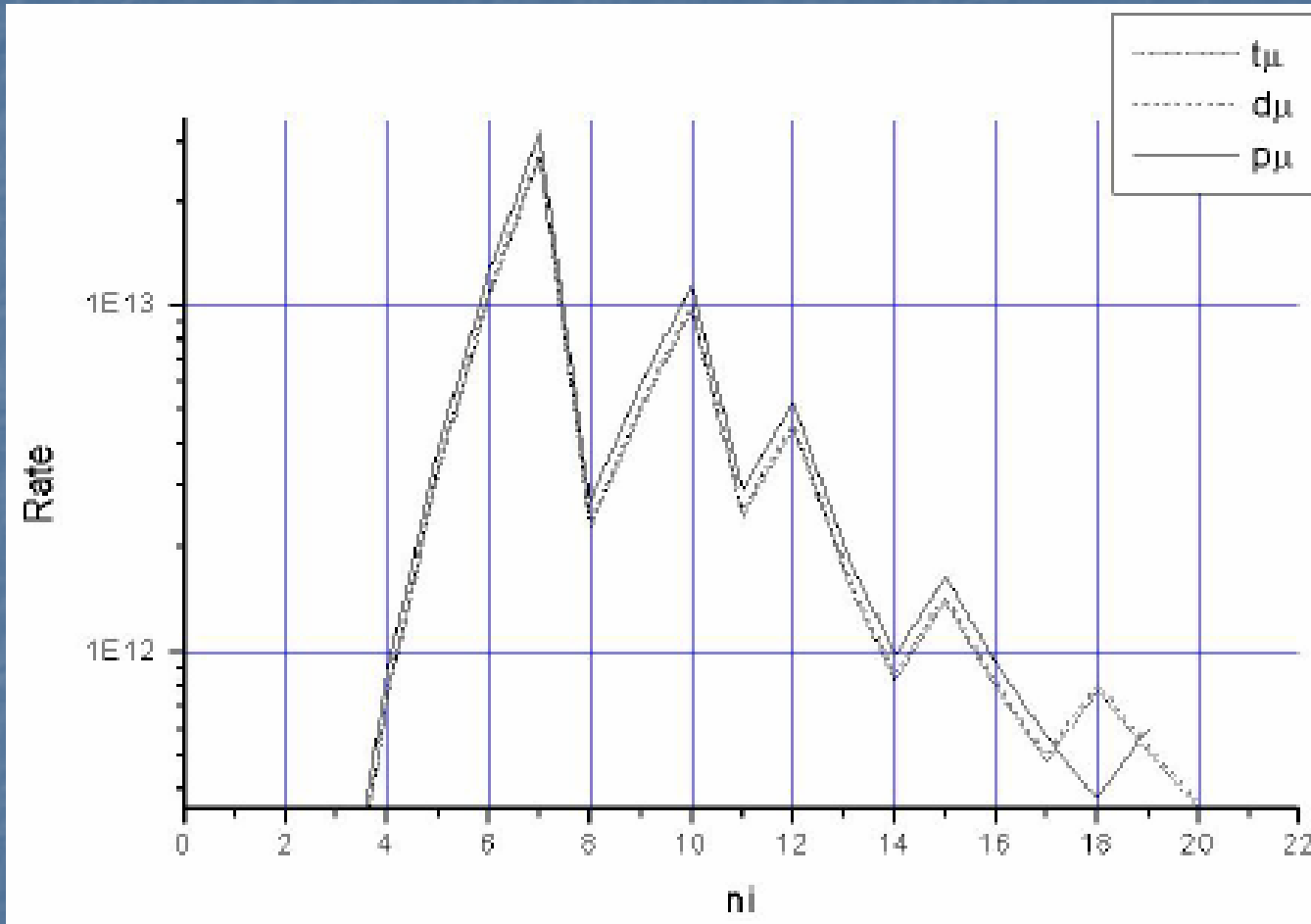
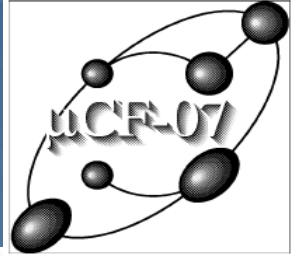
$\Delta E_{x\mu}$ **Is the excitation energy of $x\mu$ atom**

$\delta_H = 15.6 \text{ eV}$ Is the Ionization energy of hydrogen atom

Leon M. and Bethe H., Phys. Rev., 127(1962)636



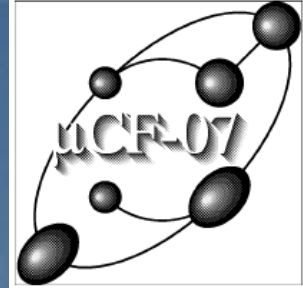
Cascade processes of muonic atoms



External Auger effect rates as a function of initial state.

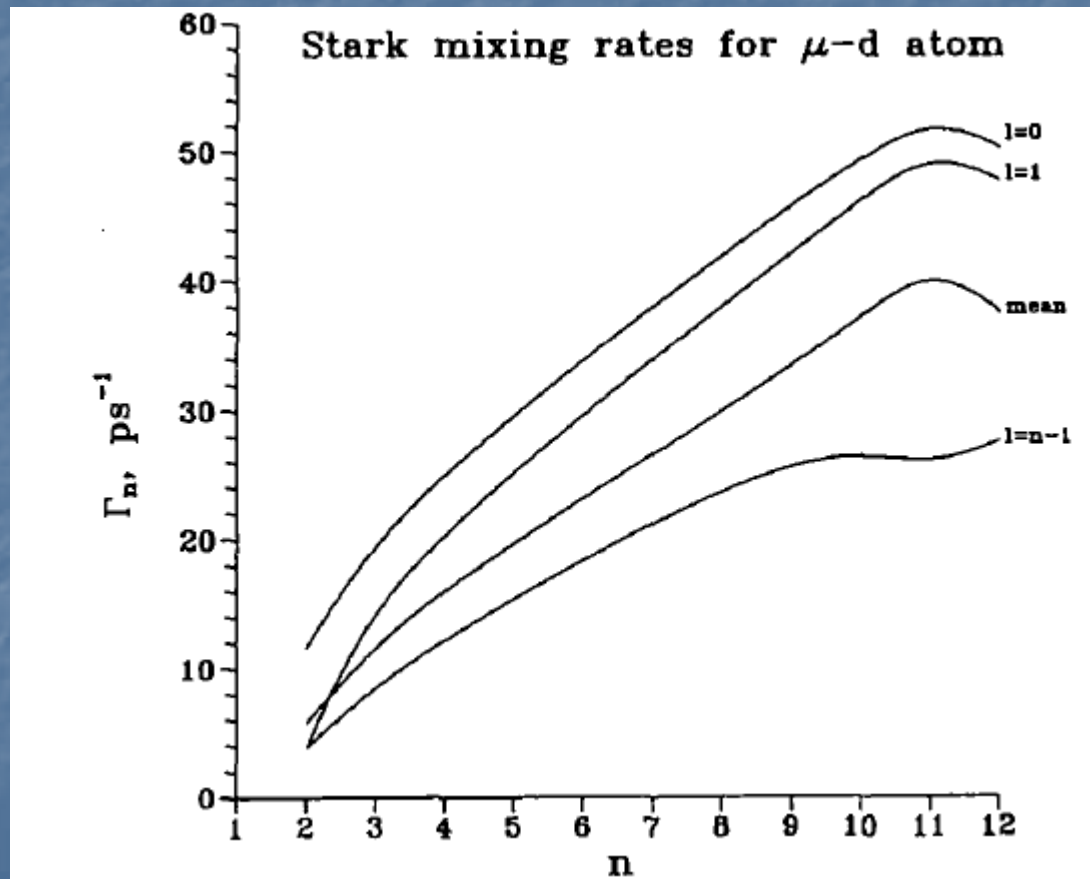


Cascade processes of muonic atoms



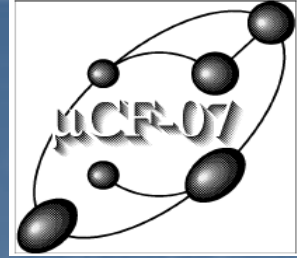
IV) Stark Mixing

$$(\mu x)_{nl} + x' \longrightarrow (\mu x)_{nl} + x'$$

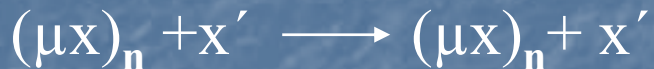




Cascade processes of muonic atoms



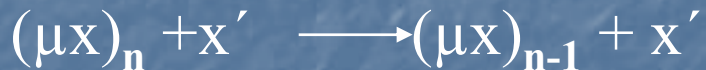
V) Deceleration (elastic scattering)



$$\chi_n^{\text{dec}} = N_0 v \frac{2M_\mu M_x}{(M_\mu + M_x)^2} \sigma_n^{\text{el}}(E) \quad \sigma_n^{\text{el}} = \frac{\pi^2 \hbar^2 (n^2 - 1)}{4M_\mu E} (\text{cm}^2)$$

L.I. Menshikov and L.I. Ponomarev, Z. Phys. D 2(1986)1

VI) Coulomb deexcitation



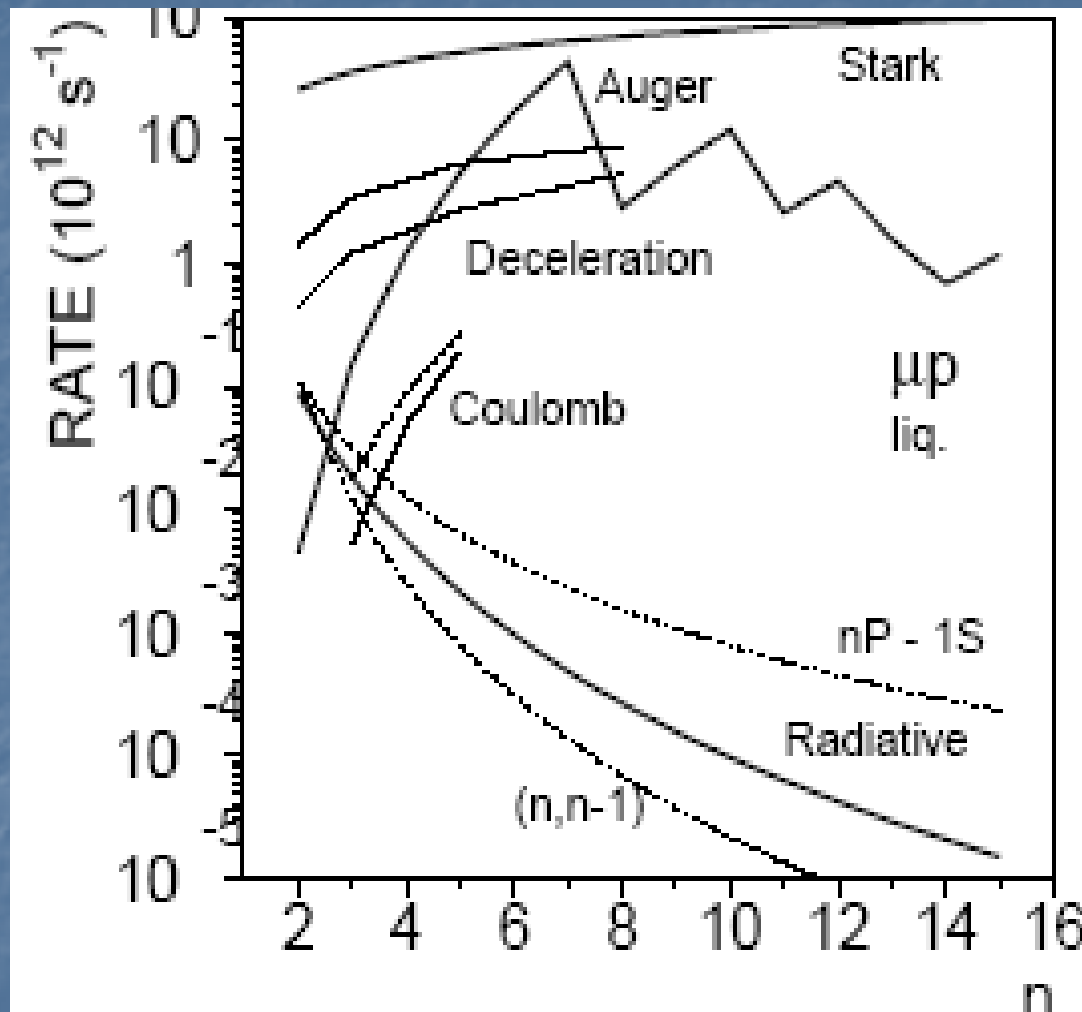
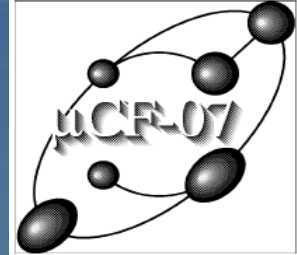
The released energy in this process is about $5\text{keV}/n(n-1)^2$.

$$\Gamma^{\text{coul}}_{n \rightarrow n'} = \frac{Nv}{20} \pi a_n^2 \left(1 + \frac{m_{\mu d}}{En^2}\right), n' = n - 1$$

Barcci L.I., et al, Nuovo Cimento A, 43(1978)9.



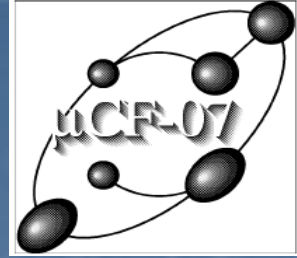
Cascade processes of muonic atoms



Comparison of the rates of competitive Cascade processes of muonic atoms.



Kinetics of cascade processes of muonic atoms

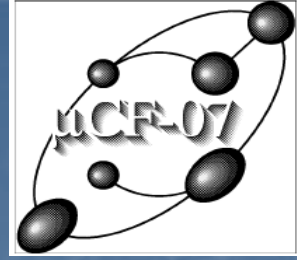


We have suggested two method for Kinetics calculattions

1. Find the **kinetics equations** and solve them by neumerical methods
2. Simulate the kinetics of cascade processes by **Monte-Carlo** method



Kinetics of cascade processes of muonic atoms



1. Kinetics equations

***Calculation of the kinetic energy of muonic atoms after collision**

$$\vec{v}'_a = \vec{v}_{cm} + \gamma_b \vec{v}'_{rel}$$

$$\vec{v}'_b = \vec{v}_{cm} - \gamma_a \vec{v}'_{rel}$$

a and b denotes the muonic and the kind of target isotopes, respectively

$$\gamma_b = \frac{m_b}{m_a + m_b}$$

$$\gamma_a = \frac{m_a}{m_a + m_b}$$

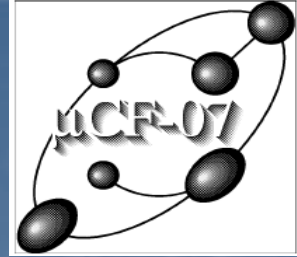
$$E_a + E_b = E'_a + E'_b + \Delta u$$

$$|\vec{v}'_{rel}| = |\vec{v}_{rel}| \sqrt{1 + \frac{\Delta u}{E}}$$

Δu and E are the released energy in reaction and relative kinetic energy in collision, respectively.

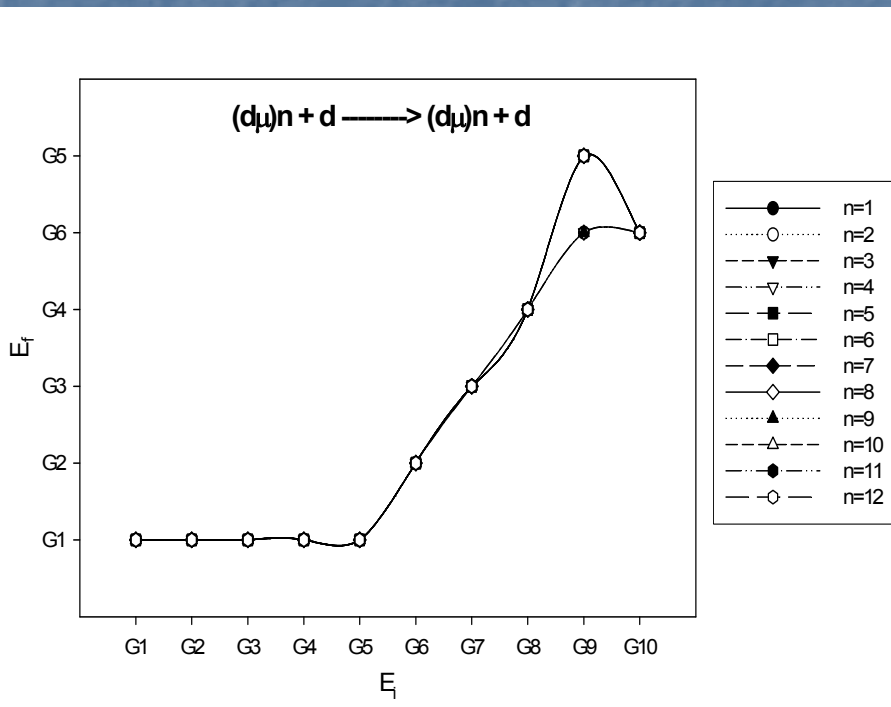


Kinetics of cascade processes of muonic atoms



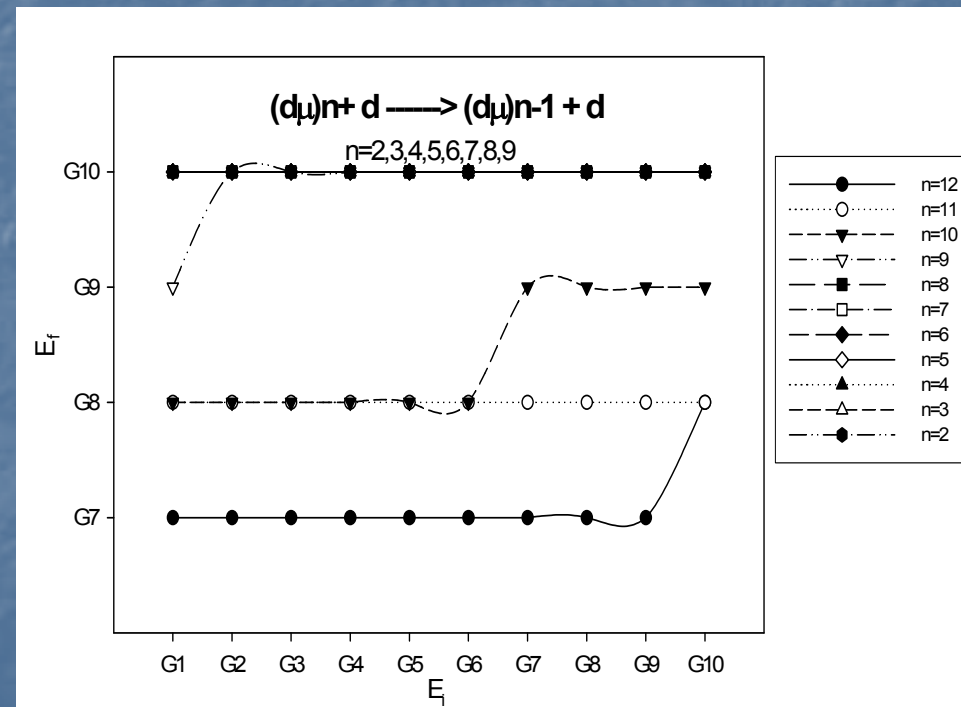
1. Kinetics equations

*For consideration of the energy dependency of the cross sections we have used a multi group method.



Elastic Collision

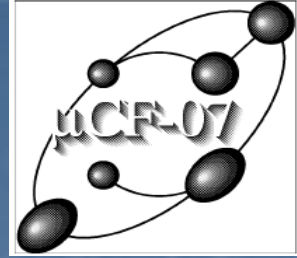
Determination of the number of energy group after collision



Coulomb de-excitation



Kinetics of cascade processes of muonic atoms



1. Kinetics equations

$$\begin{aligned}
 \frac{dP_{(n,m)_{d\mu}}}{dt} = & \sum_{i(i>n)} \lambda_{ra}^{\left\{ \begin{smallmatrix} i \rightarrow n \\ m \rightarrow m \end{smallmatrix} \right\}} P_{(i,m)_{d\mu}} - \sum_{i(i<n)} \lambda_{ra}^{\left\{ \begin{smallmatrix} n \rightarrow i \\ m \rightarrow m \end{smallmatrix} \right\}} P_{(n,m)_{d\mu}} + \sum_{i(i>n)} \phi \lambda_{au}^{\left\{ \begin{smallmatrix} i \rightarrow n \\ m \rightarrow m \end{smallmatrix} \right\}} P_{(i,m)_{d\mu}} - \sum_{i(i<n)} \phi \lambda_{au}^{\left\{ \begin{smallmatrix} n \rightarrow i \\ m \rightarrow m \end{smallmatrix} \right\}} P_{(n,m)_{d\mu}} \\
 & + \sum_l \phi(c_d) \lambda_{coul}^{\left\{ \begin{smallmatrix} n+1 \rightarrow n \\ l \rightarrow m \end{smallmatrix} \right\}} P_{(n+1,l')_{d\mu}} - \sum_l \phi(c_d) \lambda_{coul}^{\left\{ \begin{smallmatrix} n \rightarrow n-1 \\ m \rightarrow l \end{smallmatrix} \right\}} P_{(n,m)_{d\mu}} + \sum_l \phi(c_t) \lambda_{coul}^{\left\{ \begin{smallmatrix} n+1 \rightarrow n \\ l \rightarrow m \end{smallmatrix} \right\}} P_{(n+1,l)_{d\mu}} - \sum_l \phi(c_t) \lambda_{coul}^{\left\{ \begin{smallmatrix} n \rightarrow n-1 \\ m \rightarrow l \end{smallmatrix} \right\}} P_{(n,m)_{d\mu}} \\
 & + \sum_l \phi(c_d) \lambda_{tr}^{\left\{ \begin{smallmatrix} n \rightarrow n \\ l \rightarrow m \end{smallmatrix} \right\}} P_{(n,l)_{t\mu}} - \sum_l \phi(c_t) \lambda_{tr}^{\left\{ \begin{smallmatrix} n \rightarrow n \\ m \rightarrow l \end{smallmatrix} \right\}} P_{(n,m)_{d\mu}} + \sum_l \phi(c_d) \lambda_{tr}^{\left\{ \begin{smallmatrix} n+1 \rightarrow n \\ l \rightarrow m \end{smallmatrix} \right\}} P_{(n,l)_{t\mu}} - \sum_l \phi(c_t) \lambda_{tr}^{\left\{ \begin{smallmatrix} n \rightarrow n-1 \\ m \rightarrow l \end{smallmatrix} \right\}} P_{(n,m)_{d\mu}} \\
 & + \sum_{l(l>m)} \phi(c_d) \lambda_{dec}^{\left\{ \begin{smallmatrix} n \rightarrow n \\ l \rightarrow m \end{smallmatrix} \right\}} P_{(n,l)_{d\mu}} - \sum_{l(l<m)} \phi(c_d) \lambda_{dec}^{\left\{ \begin{smallmatrix} n \rightarrow n \\ m \rightarrow l \end{smallmatrix} \right\}} P_{(n,m)_{d\mu}} + \sum_{l(l>m)} \phi(c_t) \lambda_{dec}^{\left\{ \begin{smallmatrix} n \rightarrow n \\ l \rightarrow m \end{smallmatrix} \right\}} P_{(n,l)_{d\mu}} - \sum_{l(l<m)} \phi(c_t) \lambda_{dec}^{\left\{ \begin{smallmatrix} n \rightarrow n \\ m \rightarrow l \end{smallmatrix} \right\}} P_{(n,m)_{d\mu}}
 \end{aligned}$$

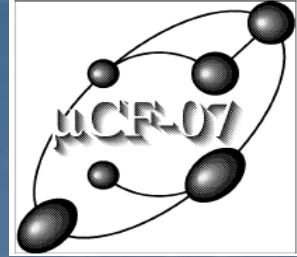
Principal quantum number: $n=12$

Number of energy group: $m=10$

These 240 equations are coupled linear differential equations. To solve them we use the Runge-Kutta method in the fourth order.

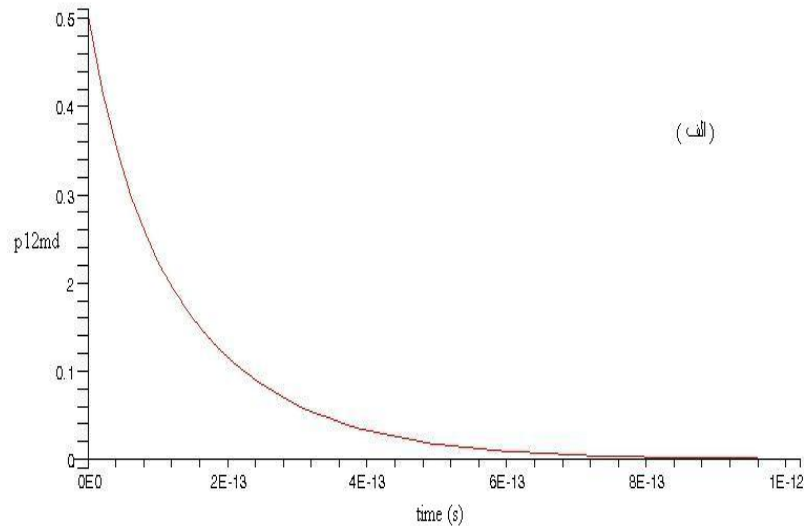


Kinetics of cascade processes of muonic atoms

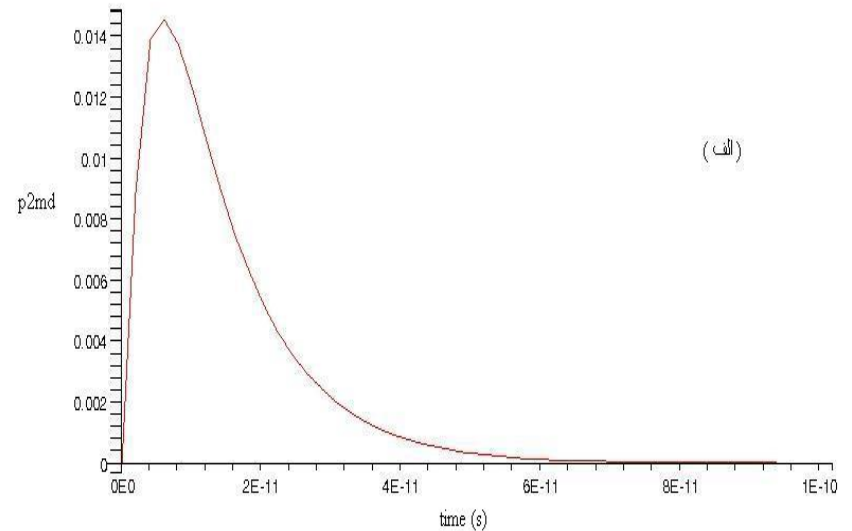


1. Kinetics equations: Some of our results

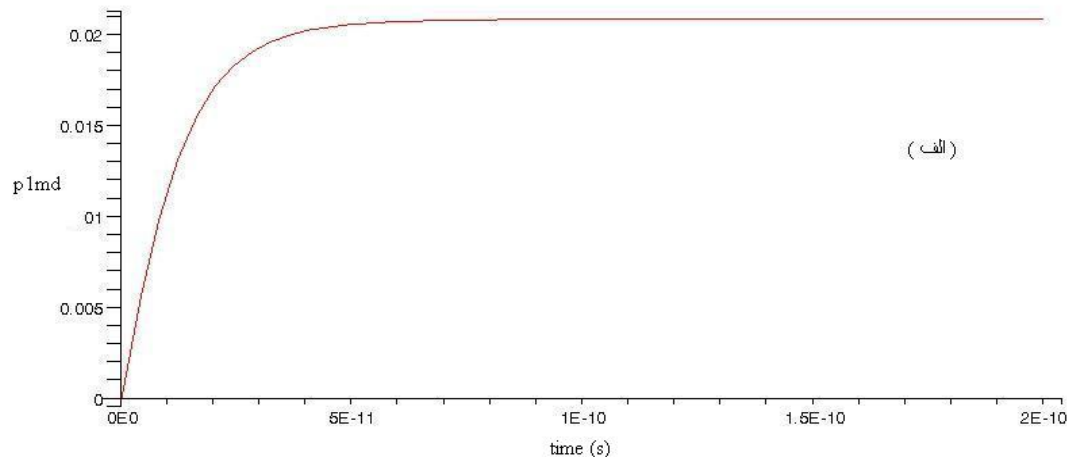
dm n=12 G=1-10 p12md f=1 cd=0.5



dm n=2 G=1-10 p2md f=1 cd=0.5



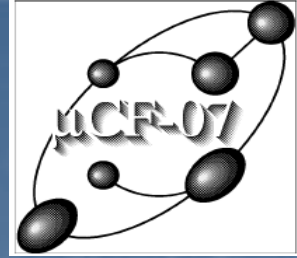
dm n=1 G=1-10 p1md f=1 cd=0.5



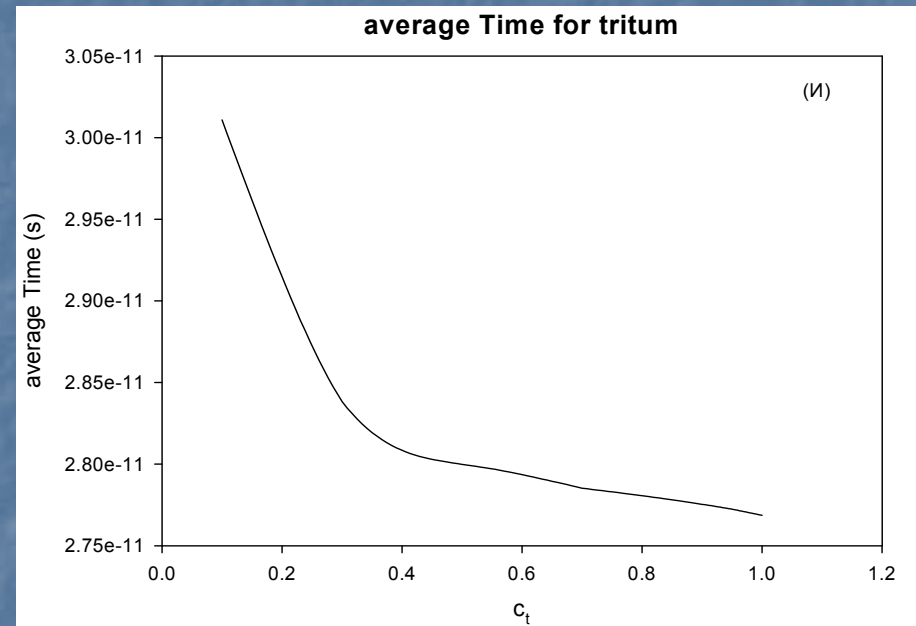
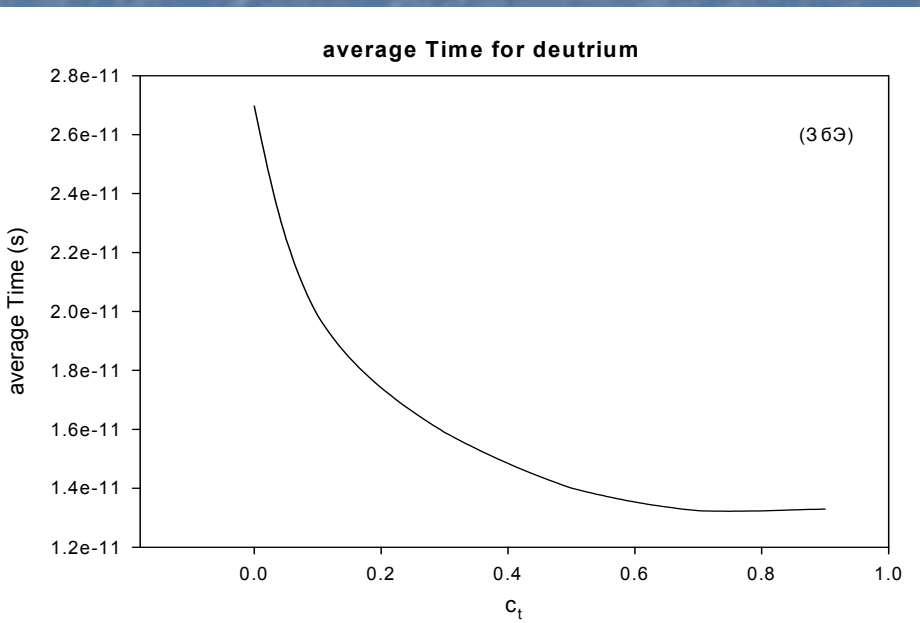
Populations of muonic atoms



Kinetics of cascade processes of muonic atoms



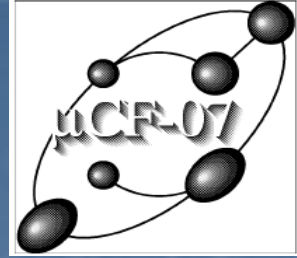
1. Kinetics equations: Some of our results



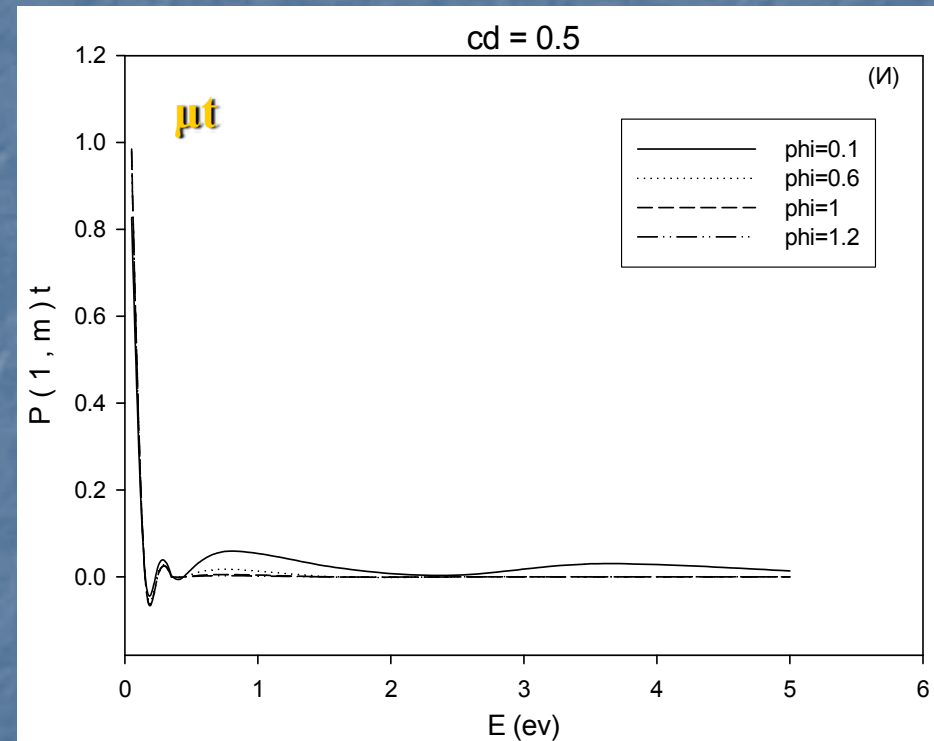
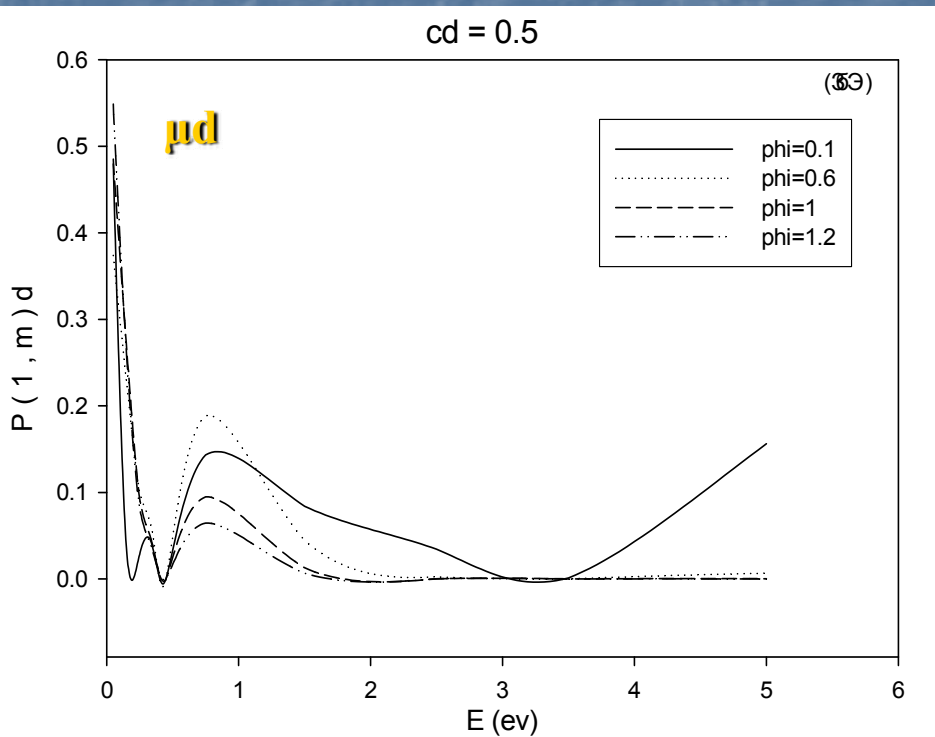
Average time of cascade processes



Kinetics of cascade processes of muonic atoms



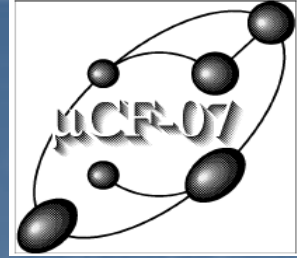
1. Kinetics equations: Some of our results



**Kinetic Energy distribution of muonic atoms in the 1s state.
It is not Maxwellian.**



Kinetics of cascade processes of muonic atoms



1. Kinetics equations

Advantages of this method:

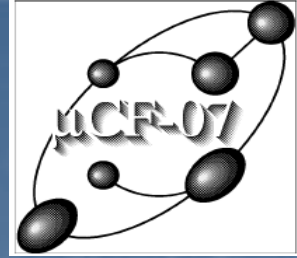
1. We can consider the energy dependency of the cross sections.
2. We can solve the problem as a function of time.
3. We can solve the problem in different physical conditions.

Disadvantages:

We should write many kinetics equations especially if the number of groups is chosen larg.



Kinetics of cascade processes of muonic atoms

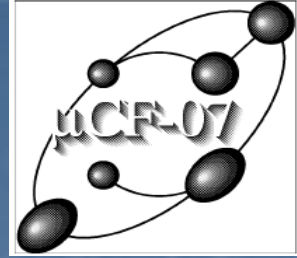


We have suggested two method for Kinetics calculations

1. Find the **kinetics equations** and solve them by numerical methods
2. Simulate the kinetics of cascade processes by **Monte-Carlo** method



Kinetics of cascade processes of muonic atoms



2. Monte-Carlo method

* Total transition rate from excited state n is:

$$\lambda_{total}^n(E) = \sum_{n'=1}^{n-1} \lambda_{columb}^{n \rightarrow n'}(E) + \sum_{n'=1}^{n-1} \lambda_{radiation}^{n \rightarrow n'} + \lambda_{scattering}^{n \rightarrow n}(E) + \sum_{n'=n-1}^n \lambda_{transfer}^{n(x\mu) \rightarrow n'(x'\mu)}(E) \\ + \lambda_{stark}^{(n,l \rightarrow n,l')_{Ave}}(E) + \sum_{n'=1}^{n-1} \lambda_{auger}^{n \rightarrow n'}(E)$$

* Transition probability from $n \rightarrow n'$ ($n' < n$) due to i th mechanism in a kinetic energy E is :

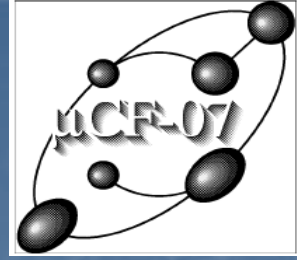
$$P_i^{n \rightarrow n'}(E) = \lambda_i^{n \rightarrow n'}(E) / \lambda_{total}^n(E)$$

* Cumulative distribution function:

$$PDF(n, j, E) = \sum_{i=1}^j \sum_{n'=1}^{n-1} P_i^{n \rightarrow n'}(E)$$



Kinetics of cascade processes of muonic atoms



2. Monte-Carlo method

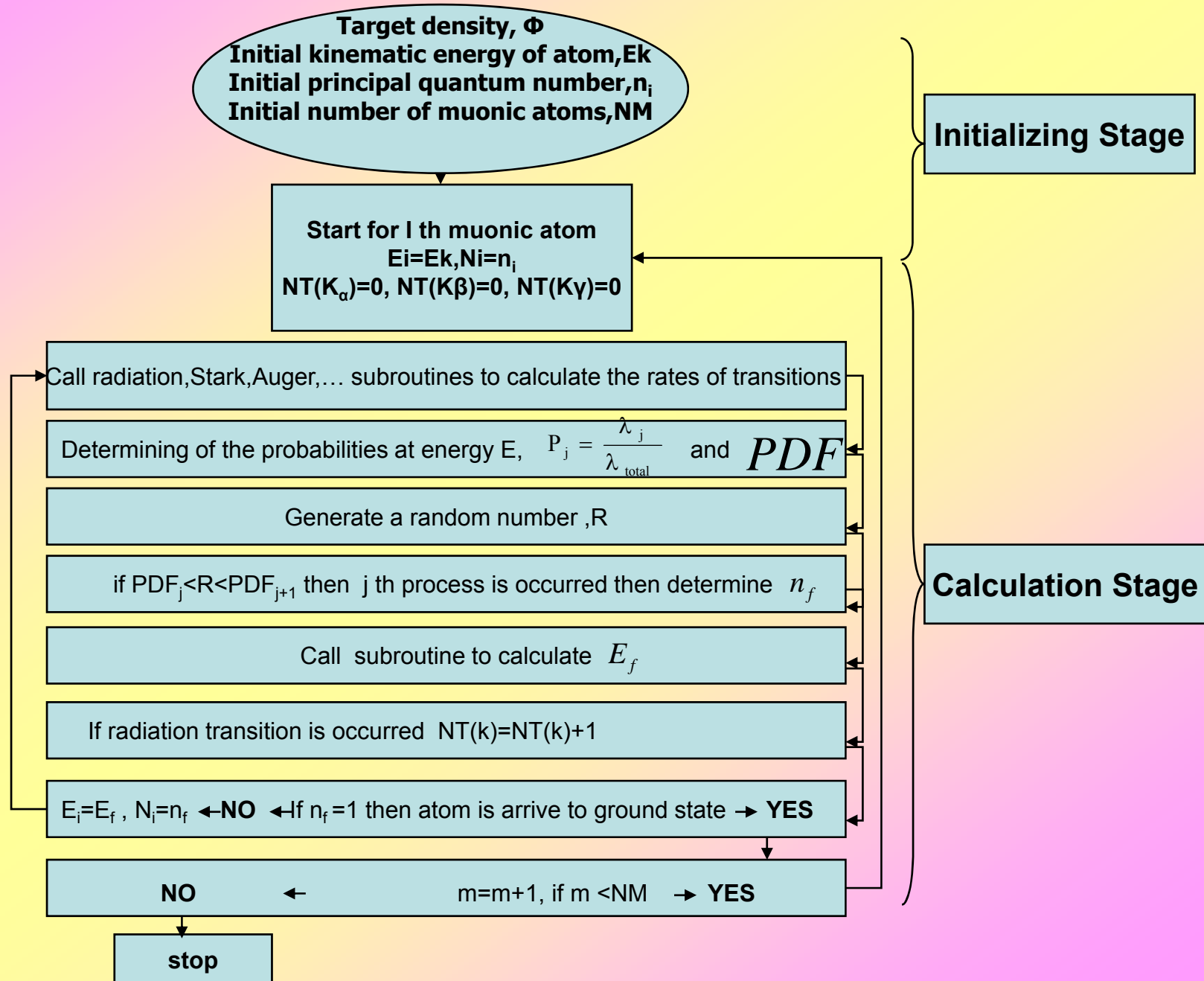
I) Calculate the cumulative distribution function at specific energy PDF(n, j, E).

II) Generate a random number R in the Interval $(0, 1)$.

III) If $\text{PDF}(n, j, E) < R < \text{PDF}(n, j+1, E)$ then j th mechanism transition is accepted. Then by the same proceed we can determine n'

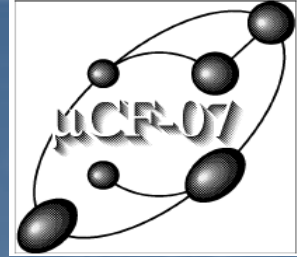
If the collisional process is every one of the scattering, Coulomb deexcitation or transfer, We should determine the kinetic energy of muonic atom after collision. Otherwise, continue from step (III) for the new n state. We should also determine the time of transition.

Monte-Carlo Simulation Flow Chart

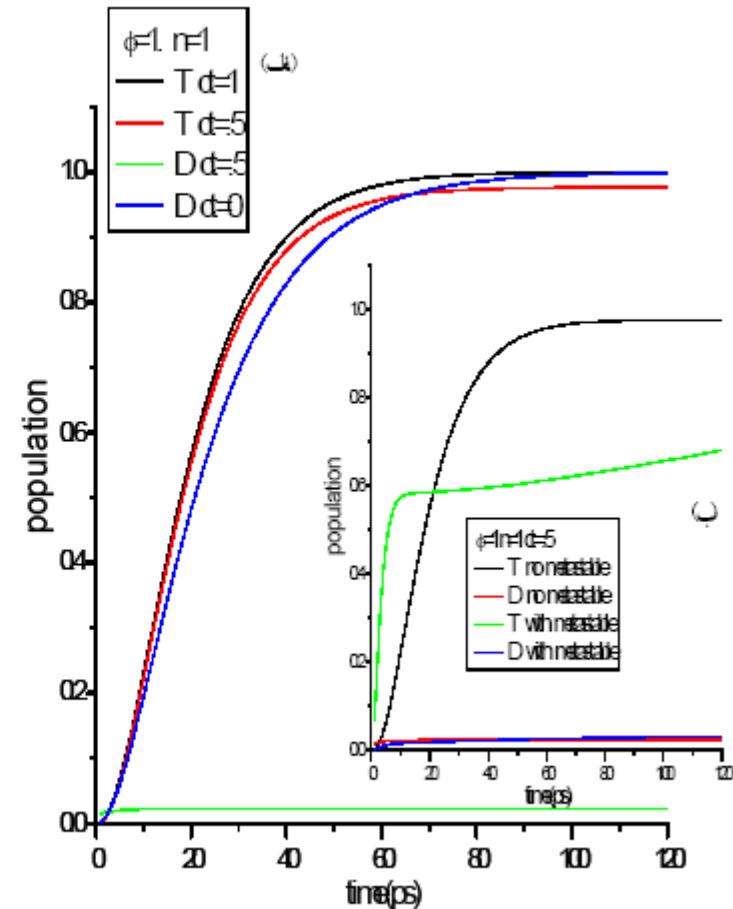
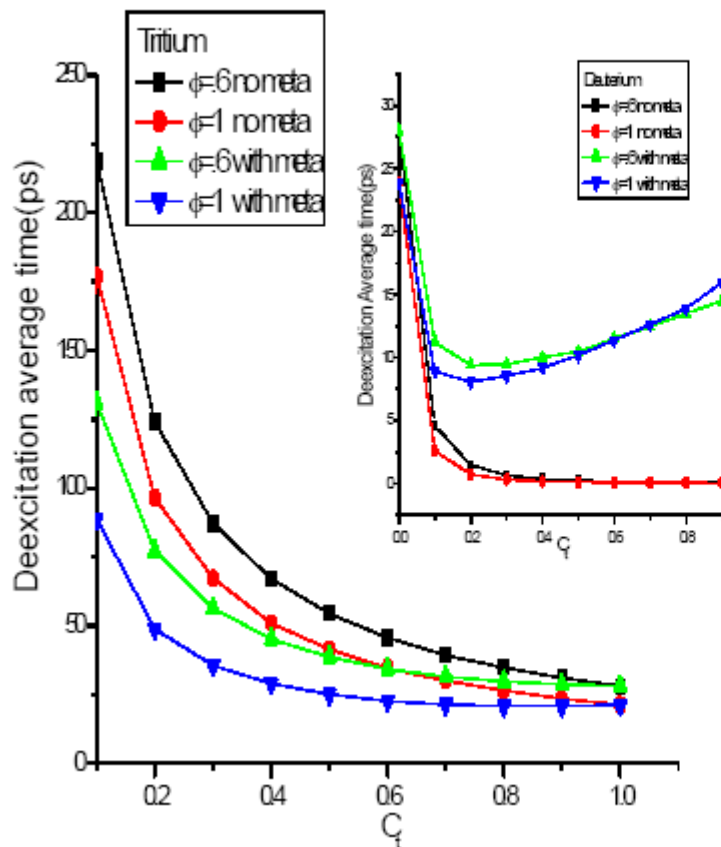




Kinetics of cascade processes of muonic atoms

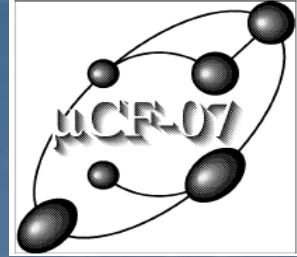


2. Monte-Carlo method: Some of our results

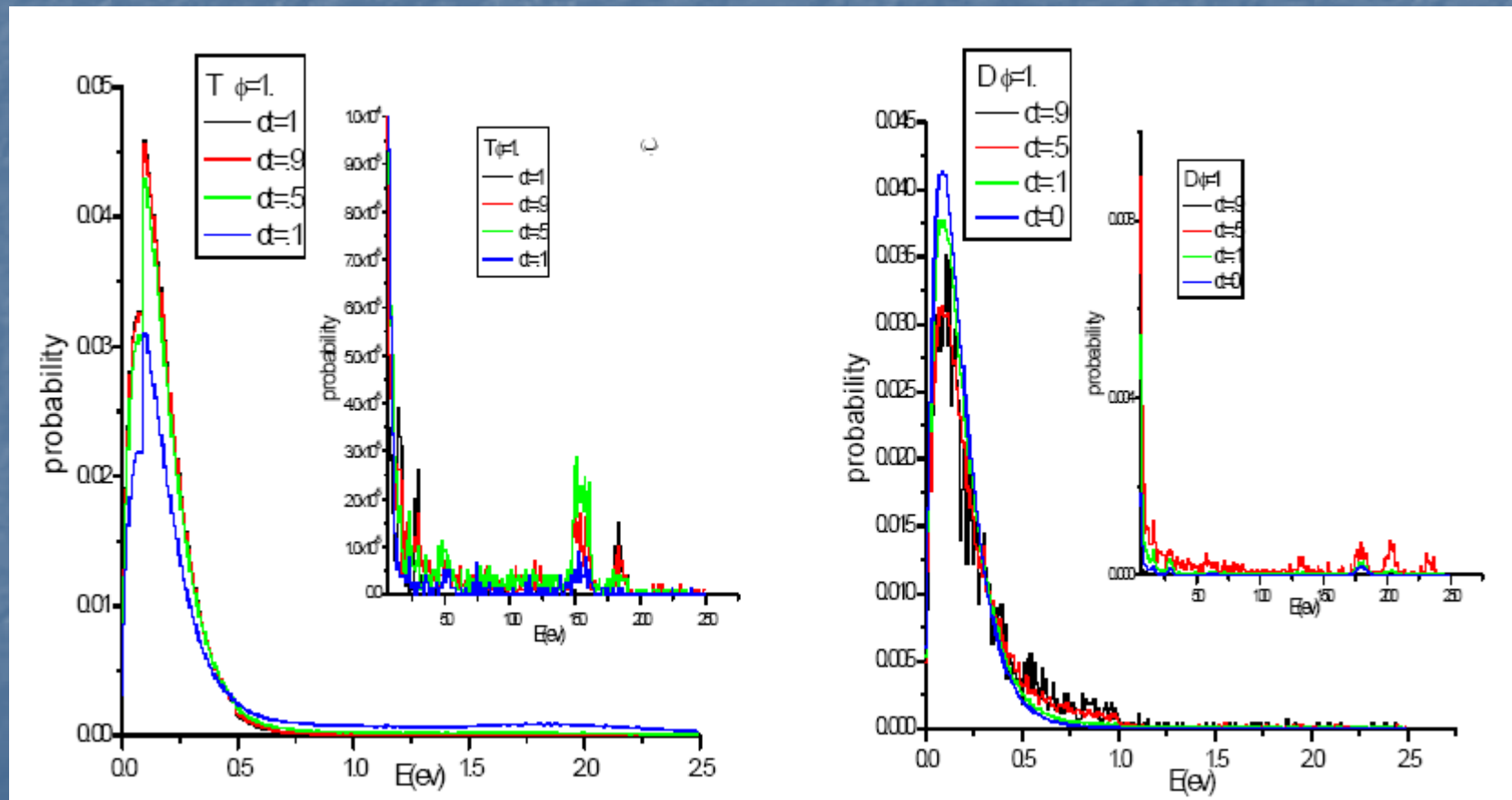




Kinetics of cascade processes of muonic atoms



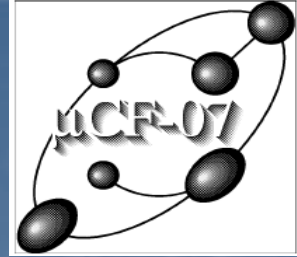
2. Monte-Carlo method: Some of our results



Kinetic energy distribution of muonic atoms

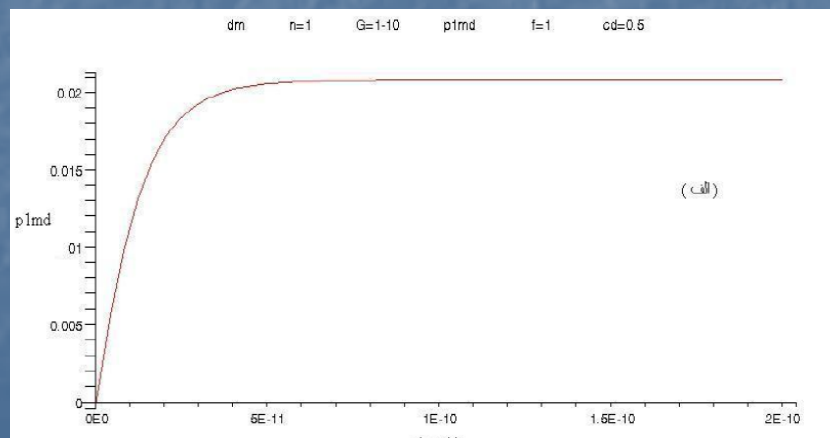


Kinetics of cascade processes of muonic atoms

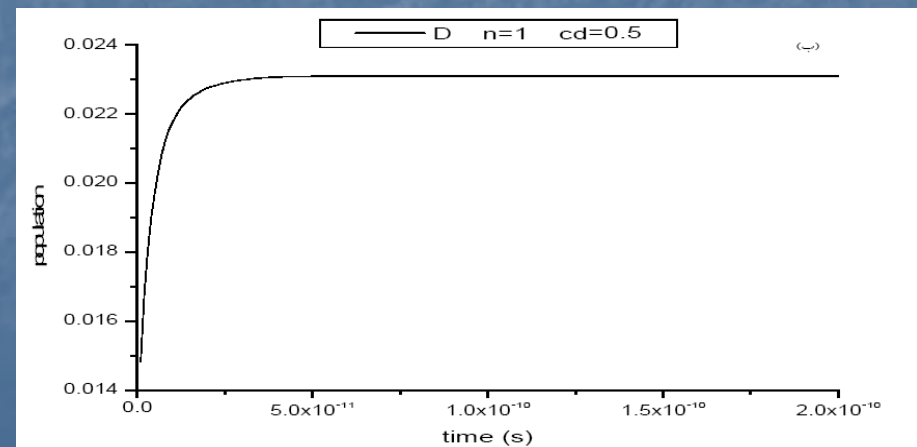


Comparison of the results from two method

Average deexcitation time	μd	μt
Solve of kinetics equations	$2.7 \times 10^{-11} s$	$2.8 \times 10^{-11} s$
Monte-Carlo simulation	$2.4 \times 10^{-11} s$	$2.1 \times 10^{-11} s$



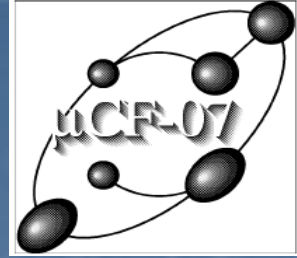
Kinetics method



Mont Carlo method

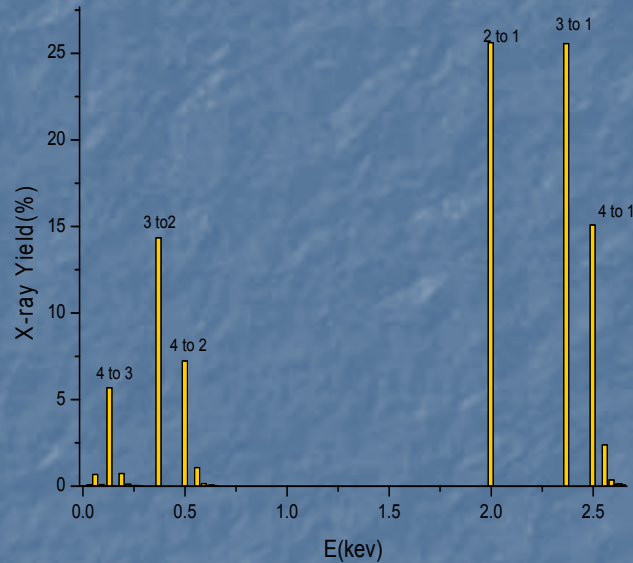


Kinetics of cascade processes of muonic atoms

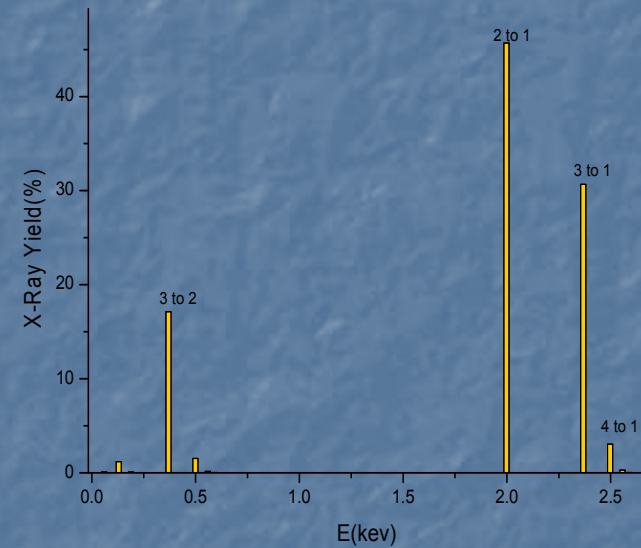


X-ray yields from muonic atoms by Monte-Carlo method

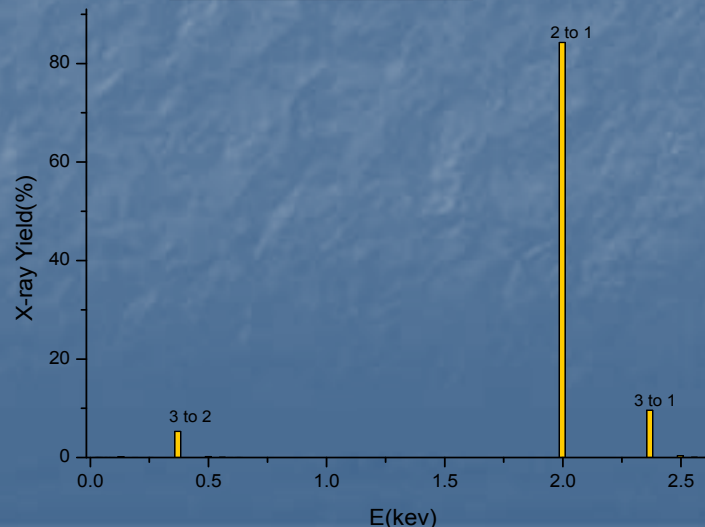
Muonic Deuterium $\phi=.01$



Muonic Deuterium $\phi=.1$

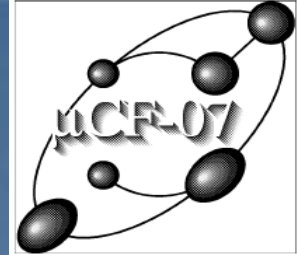


Muonic Deuterium $\phi=1.$

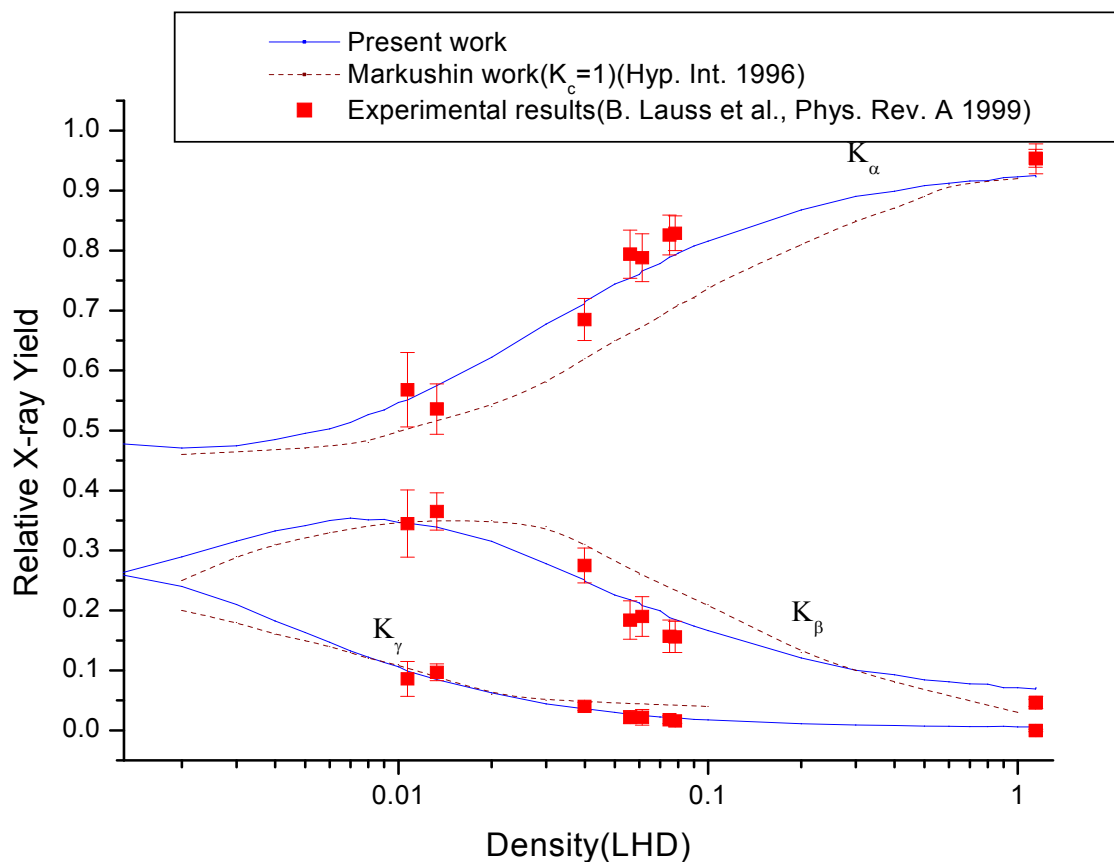


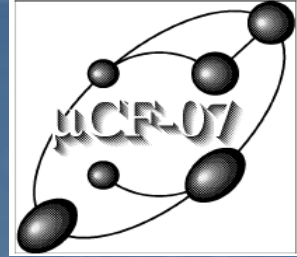


Kinetics of cascade processes of muonic atoms



Comparison our results with the other results





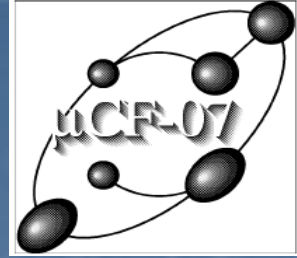
Cascade processes of Kaonic atoms

■ Cascade processes

Mechanism	Process	Energy Dependence	Density Dependence
Radiative	$(xK)_{n_i} \rightarrow (xK)_{n_f} + \gamma$	None	None
Auger	$(xK)_{n_i} + x'_2 \rightarrow x'^{+}_2 + e + (xK)_{n_f}$	Weak	Linear
Stark mixing	$(xK)_{n_l} + X \rightarrow (xK)_{n_l'} + X$	Moderate	Linear
Scattering	$(xK)_n + X \rightarrow (xK)_n + X'$	Strong	Linear
Coulomb	$(xK)_{n_i} + x' \rightarrow (xK)_{n_f} + x'$	Strong	Linear
Transfer	$(xK)_n + x' \rightarrow (x'K)_n + x$	Strong	Linear
Nuclear absorption	$(xK)_n + X \rightarrow \pi^0 + \Lambda + X$?	?
Nuclear reaction	$(xK)_n \rightarrow \pi^0 + \Lambda$	None	None
Weak decay	$K^- \rightarrow \mu^- + \bar{\nu}_\mu$	None	None



Simulation of cascade processes of K \bar{p} atoms



X-ray yields from cascade of kaonic atoms is important to test of strong interaction and the theory of QCD in low energies.

As the kaon interacts strongly with the nucleus the 1s energy level is shifted and broadened

$$\varepsilon_1 = (E_{K_\alpha})_{\text{exp}} - (E_{K_\alpha})_{\text{em}}$$

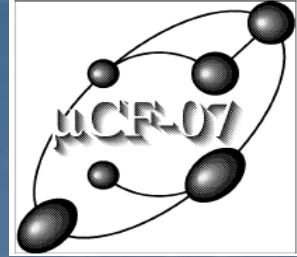
Density of the target can affect the X-ray yields.

Because the X-ray yields are rare therefore we should find an optimum condition of the target for the experiment.

Shift of energy in 1s state and width of 1s and 2p and ... may change the X-ray yields. We can use them in our calculation as a free parameter.



Simulation of cascade processes of K-p atoms

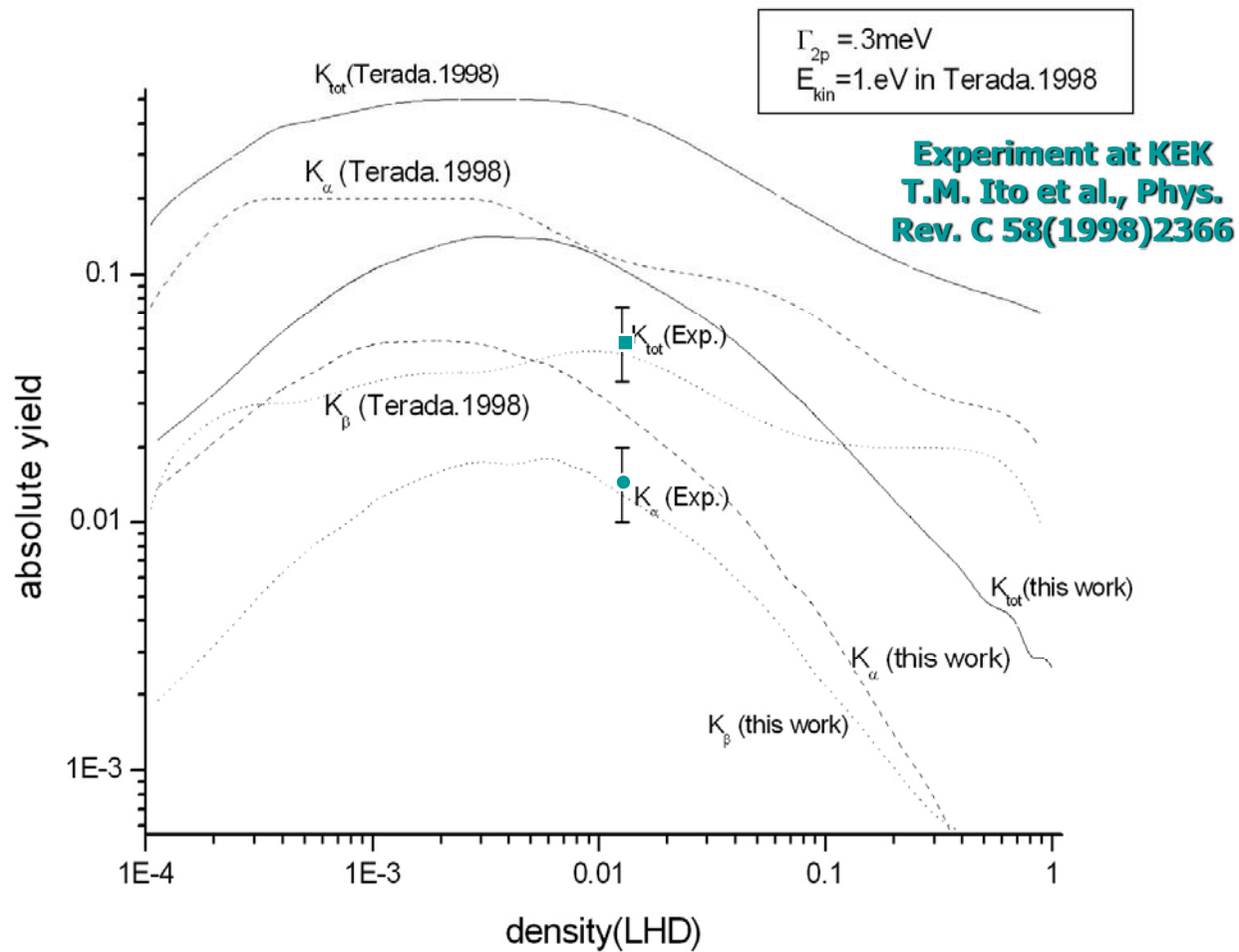
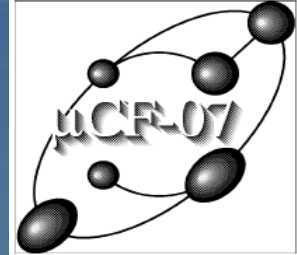


Our plan to do the simulation

- I. Determining of the cross sections or rates of the Stark transitions, external Auger effect, Coulomb deexcitation, elastic scattering and absorption as a function of kinetic energy of kaonic atoms and the rates of radiative transition and nuclear reactions of kaonic atoms in any excited states.
- II. Preparing a computer code based on the Monte-Carlo method to simulate the cascade processes in kaonic atoms. The program begins by a highly excited kaonic atom and calculates its life history as they cascade down through the nl states. The program determines the kinetic energy of the kaonic atom after any collisional processes such as Coulomb deexcitation and elastic scattering in each step of time. The program repeat for 10^6 kaonic atoms and take an average of any quantity.
- IV. If during the simulation any radiative transition is take place, the energy and time of emitted X-ray is recorded. Finally the X-ray spectrum can be calculated in any time. These calculations can be done for any physical conditions such as density target.

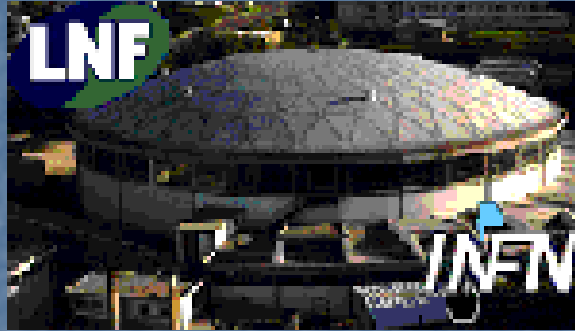
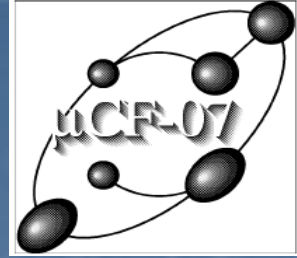


Simulation of cascade processes of K-p atoms





Simulation of cascade processes of K-d atoms in SIDDHARTA



We hope to determine an optimum condition for the SIDDHARTA experiment (For K-d atoms at LNF) to achieve the highest X-ray yields.

Then In order to study the strong Interactionour results will be compared with the experiment.

