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ACK DECAY OF MUONIC MOLECULAR
RESONANCES AND THE MEASURED
VALUE OF $d_{\mu}d$ -FORMATION RATE
IN MUON-CATALYZED FUSION
IN DEUTERIUM

A. Gula*, A. Adamczak** and M. Bubak***

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BACK DECAY OF MUONIC MOLECULAR RESONANCES AND THE
MEASURED VALUE OF $d\mu d$ - FORMATION RATE IN
MUON-CATALYZED FUSION IN DEUTERIUM

ROZPAD REZONANSÓW MOLEKULARNYCH W KANAŁ WEJŚCIOWY
A MIERZONA WARTOŚĆ SZYBKOŚCI TWORZENIA $d\mu d$
W MIUNOWEJ KATALIZIE SYNTEZY JĄDROWEJ

РАСПАД МО-МОЛЕКУЛЯРНЫХ РЕЗОНАНСОВ ПО ЭХОДНОМУ КАНАЛУ
А ЭКСПЕРИМЕНТАЛЬНАЯ ВЕЛИЧИНА СКОРОСТИ ОБРАЗОВАНИЯ $d\mu d$
В МИУОННОМ КАТАЛИЗЕ РЕАКЦИИ ДЕФУТОГО СИНТЕЗА

by

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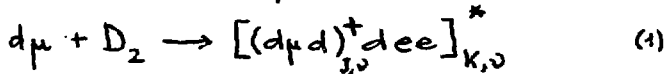
Abstract

It is shown that the experimental values of $d_{\mu d}$ formation rate, obtained without taking into account the decay of the μ -molecular resonance compound $[(d_{\mu d})^+ dee]^*$ back to the formation channel $d_{\mu} + D_2$, are underestimated. The correction depends on the rate of this resonance back decay and the rates of processes leading to fusion in $d_{\mu d}$.

For their current estimates the correction significantly exceeds the experimental error of the uncorrected $d_{\mu d}$ formation rate $\lambda_m^{obs} = 2.76 \pm 0.08 \mu s^{-1}$ reported recently. It is argued that back decay may lead to variation of λ_m^{obs} with target density which may provide useful information on the parameters of muon-catalyzed fusion.

When a negative muon enters a deuterium target it forms a muonic deuterium atom ($d\mu$) and subsequently a tightly bound ($d\mu d$)⁺ system (r.m.s. $\sim 10^{-11}$ cm) in which the deuterons rapidly undergo nuclear fusion [1,2]. The muon is then most likely freed ($f \approx 87\%$) and becomes available to catalyze another fusion reaction.

It is now well established that ($d\mu d$)⁺ formation proceeds mainly via creation of a resonance μ -molecular compound [3-5]:



Such process is possible due to existence of a weakly bound ($\epsilon_b = 1.91$ eV) rovibrational state of $d\mu d$: $J = 1, \nu = 1$ [2]. The energies are tuned in (1) by the vibrational ($\epsilon_\nu \sim 10^2$ meV) and rotational ($\epsilon_K \sim 1$ meV) excitations of the compound and by the relative collision energy ($\epsilon_T \sim 10$ meV) [2,6]. It has been commonly assumed so far, that - unless the muon decays - (1) is followed inevitably by fusion in $d\mu d$. However, as pointed out in [7], the decay of compound (1) back to the formation channel may seriously compete with transitions leading to fusion. In this note we investigate an important consequence of this observation.

Fig. 1 presents the main processes forming one muon catalyzed fusion (MCF) cycle in deuterium with inclusion of the resonance back decay (RBD) shown as a dashed line. Other important decay modes are [7]: (i) direct fusion in the ($J = 1, \nu = 1; K, \nu$) state, (ii) de-excitation of $d\mu d$ to a lower $J = 1, \nu = 0$ state, (iii) vibrational de-excitation of compound (1) in collisions with D_2 molecules. In (ii) the resonance object is destroyed as $\Delta \epsilon_b \approx 2 \cdot 10^2$ eV exceeds the ionization energy and in (iii) already a $\nu \rightarrow \nu - 1$ transition brings the compound below the elastic threshold ($\Delta \epsilon_\nu \gg \epsilon_T$). Although the analogous collisional K -transitions are expected to be very fast [8] we assume tentatively, as in [7], that they do not compete with RBD. In fact, the $d\mu d$ -fusion data can be described by a single narrow resonance at $\epsilon_T = \epsilon_0 \approx 50$ meV taking into account only vibrational excitation [3]. Adding the rotational excitations on top leaves the formation channel open in the ensuing

transitions, thus, the general structure of the kinetic graph in Fig. 1 is not affected.

To see how RBD influences the kinetics of MCF in deuterium let us consider the Laplace transform, $F_1(s)$, of the time distribution of fusion events ending the first cycle, $F_1(t) = dN_{11}/dt$. Without RBD it reads [9-11]:

$$F_1(s) = \frac{s \phi(s)}{\Lambda_\mu + s} \frac{\lambda_a}{\Lambda_{d\mu} + s} \frac{\lambda_m}{\Lambda_{\mu mr} + s} H(\lambda_d, \lambda_e, \lambda_f, \lambda'_f; s) \quad (2)$$

where λ_i are transition rates defined in Fig. 1, $s \cdot \phi(s)$ represents the incoming muon flux, Λ_i are sums of the transition rates over all lines leaving an i -th vertex (note: $\Lambda_{(1,1;K,\nu)} = \Lambda_{\mu mr}$) and $H(s)$ describes the structure of the subgraph linking the μ and F nodes (channels leading to fusion in $d\mu d$) [11].

One can easily verify, either by directly solving the first-cycle kinetic equations [10] or by applying the Mason topological formula [11], that the $d\mu \rightleftharpoons \mu mr$ loop created in the kinetic graph by RBD couples $\Lambda_{d\mu}$ and $\Lambda_{\mu mr}$ in eq. (2) so that

$$\begin{aligned} (\Lambda_{d\mu} + s)(\Lambda_{\mu mr} + s) &\longrightarrow (R_{d\mu} + s)(R_{\mu mr} + s) = \\ &= (\Lambda_{d\mu} + s)(\Lambda'_{\mu mr} + s) - \lambda_b \lambda_m \end{aligned} \quad (3)$$

The other Λ_i - hence also $H(s)$ - remain unchanged and, with inclusion of RBD, $\Lambda_{\mu mr} \rightarrow \Lambda'_{\mu mr} = \Lambda_{\mu mr} + \lambda_b$. Since the time distribution of fusion events ending an arbitrary k -th cycle is given by $F_k(s) = s^{k-1} [F_1(s)]^k$, inclusion of RBD amounts to replacing in all kinetic formulae: $\Lambda_{d\mu}, \Lambda_{\mu mr} \rightarrow R_{d\mu}, R_{\mu mr}$. In particular, the large time behaviour of $F_k(t)$ is determined by the position of the rightmost pole in (2) [10]. This is now $(-R_{d\mu})$ since

$$R_{d\mu} = \lambda_0 + \frac{1}{2} [\lambda_m + \lambda_f + \lambda_b - \sqrt{(\lambda_f + \lambda_b - \lambda_m)^2 + 4\lambda_b \lambda_m}] < \Lambda_{d\mu} \quad (4)$$

and $\Lambda_{d\mu} < \Lambda_{i \neq d\mu}$ [2]. We put above: $\lambda_F = \lambda_e + \phi \lambda_d$.

$\lambda_e = \lambda_f + \lambda_t$, $\lambda_m = \varphi \lambda_m^0$. λ_0 is muon decay rate, φ is target density expressed in units of liquid hydrogen density (LHD) and index 0 denotes the rates normalized to LHD. For $\lambda_b \gg \lambda_m$ eq. (4) becomes

$$\lambda_m^{\text{obs}} = (R_{d\mu} - \lambda_0) \varphi^{-1} = \frac{\lambda_f}{\lambda_f + \lambda_b} \lambda_m^0 \quad (4a)$$

where λ_m^{obs} is the quantity identified so far with λ_m^0 .

An immediate consequence of eq. (4) is that neglecting RBD in the determination of λ_m^0 from the large time slope of $F_1(t)$ [4] leads to underestimated values of this important parameter. To estimate this effect let us remark that upon the thermal averaging [7]:

$$\lambda_b = \text{const } e^{\frac{\epsilon_0}{kT}} (kT)^{\frac{3}{2}} \lambda_m^0(T) = \alpha(T) \lambda_m^0(T). \quad (5)$$

In contrast to all previous analyses where $R_{d\mu} \rightarrow \Lambda_{d\mu} = \lambda_0 + \lambda_m$ (e.g., [3,4]), λ_m^0 depends now on λ_b and α , the latter being determined by the resonance energy ϵ_0 and the rotational quantum numbers K_{D2} and K , which count the number of states contributing to relation (5). Neglecting the K -structure and using $\epsilon_0 = 53 \text{ meV}$ [12] one obtains [7] $\alpha(293\text{K}) \approx 300$. This value together with $\lambda_f = 870 \mu\text{s}^{-1}$ [13], $\lambda_t = 190 \mu\text{s}^{-1}$ [14] and $\lambda_d^0 = 10 \mu\text{s}^{-1}$ [7] applies to the result of [4] $\lambda_m^{\text{obs}} = 2.76 \pm 0.08 \mu\text{s}^{-1}$ ($\varphi = 0.1$) gives $\lambda_m^0(293\text{K}) = 12.6 \pm 1.7 \mu\text{s}^{-1}$. Muon transitions to impurity atoms and nonresonant $d\mu$ production which have been neglected in writing down eqs. (4,5) may change this number by $\approx 5\%$ [15,16]. If, more realistically, the $K_{D2}=0 \rightarrow K=1$ transitions are assumed to dominate ($\alpha \rightarrow \alpha/3 = 100$) then

$$\lambda_m^0 = 3.7 \pm 0.14 \mu\text{s}^{-1}.$$

To illustrate the importance of the correction for back decay let us notice that for λ_m^0 to coincide with $\lambda_m^{\text{obs}} = 2.76 \pm 0.08 \mu\text{s}^{-1}$

within two standard deviations, one should have, according to (4a): $\lambda_{\text{F}}/\alpha > 50 \mu\text{s}^{-1}$, while with λ_{F} of [7,13,14] and $\alpha = 100$ this ratio is only $\sim 10 \mu\text{s}^{-1}$. Thus, if λ_{F} is not grossly underestimated, $\lambda_{\text{m}}^{\circ}$ is actually much higher than the existing theoretical prediction [12]. However, the results of [4,5] already challenge this estimate and indications appear that the μ -molecular formation rates may significantly differ from those reported previously [17].

More accurate data on $\bar{r}_{\text{x}}(t)$ at small t which are sensitive to λ_{F} [10,16] would greatly help to resolve these ambiguities. Another interesting and relatively simple possibility is to measure $\lambda_{\text{m}}^{\text{obs}}$ at various target densities. Fig. 2 shows $\lambda_{\text{m}}^{\text{obs}}$ as a function of φ for several values of α , λ_{e} and $\lambda_{\text{d}}^{\circ}$ with $\lambda_{\text{m}}^{\text{obs}}$ fixed at the most accurately measured value ($2.76 \pm 0.08 \mu\text{s}^{-1}$ at $\varphi = 0.103$) of [4]. A wide range of $\lambda_{\text{d}}^{\circ}$ is covered to include the possibility that the rotational de-excitations also compete with RBD as conjectured in [8]. (In fact, $\lambda_{\text{d}}^{\circ}$ for vibrational de-excitation depends strongly on the assumptions used in its determination and a wide spectrum of values can also be expected [18].)

Curves for $\lambda_{\text{d}}^{\circ} < 100 \mu\text{s}^{-1}$ deviate very weakly from $\lambda_{\text{m}}^{\text{obs}} = \text{const}$ and are not shown. The recent data of [4] published for other densities (2.4 ± 0.6 , 2.86 ± 0.11 , 2.74 ± 0.14 , $2.84 \pm 0.10 \mu\text{s}^{-1}$ at $\varphi = 0.06$, 0.08 , 0.09 , 0.105 , respectively) and the estimate of [5] ($\sim 3 \mu\text{s}^{-1}$ at $\varphi \approx 0.3 - 0.7$) are not conclusive with the present errors. They agree with $\lambda_{\text{d}}^{\circ} \lesssim 100 \mu\text{s}^{-1}$, although higher $\lambda_{\text{d}}^{\circ}$ are not excluded if α is appropriately lowered or λ_{e} increased. On the other hand, the result of [3], $0.76 \pm 0.11 \mu\text{s}^{-1}$, cannot be reconciled with these data with any choice of parameters. As pointed out in [7] $\lambda_{\text{d}}^{\circ}$ should grow by 3-4 orders of magnitude between 400 and 2000 K and quite pronounced effects may, thus, appear at somewhat higher temperatures. Hence, back decay should be taken into account also in consideration of other factors leading to non-trivial density dependence of MCF parameters. Therefore, the measurements of $\lambda_{\text{m}}^{\text{obs}}$ in pure deuterium over a wide range of temperatures and densities may provide very useful information.

After this work been completed we learned that L.I. Menshikov estimated the influence of RBD on the μ - molecular resonance formation rates, using a probability argument which led him to eq. (4a).

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References

1. F.C. Frank: *Nature* 160 (1947) 525,
A.D. Sakharov: Report of the Phys. Inst. Acad. Sci. USSR,
Moscow, 1948,
Ya.B. Zeldovich: *Dokl. Akad. Nauk SSSR* 95 (1954) 493,
2. For introduction and a recent review see:
L. Bracci, G. Fiorentini: *Phys. Rep.* 86 (1982) 170,
L.L. Ponomarev: *Atomkernenergie/Kerntechnik* 43 (1983) 175,
3. V.M. Bystritsky, V.P. Dzhelepov et al.: *Zh. Eksp. Teor. Fiz.* 76
(1979) 460 *Sov. Phys. JETP* 49 (1979) 232 .
4. D.V. Balin, A.A. Vorobev et al.: *Phys. Lett.* 141 B (1984) 173;
D.V. Balin, A.A. Vorobev et al.: *Zh. Eksp. Teor. Fiz. - Pisma* 40
(1984) 318,
5. S.E. Jones: Talk at Ninth International Conference on Atomic
Physics, Seattle, July, 1984,
6. M. Leon: *Phys. Rev. Lett.* 52 (1984) 605,
7. M. Lane: *Phys. Lett.* 98A (1983) 337,
8. V.N. Ostrovski, V.I. Ustimov: *Zh. Eksp. Teor. Fiz.* 79 (1980) 1228,
9. V.M. Bystritsky, A. Guła, J. Woźniak: *Atomkernenergie/Kerntechnik*
45 (1984) 197,
10. M. Bubak, V.M. Bystritsky, A. Guła: Preprint JINR E1-84-164,
Dubna, 1984 (to be published in *Acta Physica Polonica*),
11. A. Guła: Preprint JINR E1-84-301, Dubna, 1984 (to be published
in *Acta Physica Polonica*),
12. S.I. Vinitsky, L.L. Ponomarev et al. : *Zh. Eksp. Teor. Fiz.* 74
(1978) 849 [*Sov. Phys. JETP* 47 (1978) 444] .
13. L.M. Bogdanova et al.: *Phys. Lett.* 115B (1982) 171.
14. Value obtained by M.P. Faifman using the algorithm of L.L. Pono-
marev, M.P. Faifman: *Zh. Eksp. Teor. Fiz.* 71 (1976) 1689 [*Sov.*
Phys. JETP 44 (1976) 886] ,

15. D.V. Balin et al. : Preprint LNPI-964, Leningrad, 1984,
16. A. Gula, M.Bubak, J. Niwicki: Report JINR E1-84-560, Dubna, 1984; preprint INT 189/P, Kraków, 1985 (to be published in Atomkernenergie/Kerntechnik),
17. J.S. Cohen, R.L. Martin : Phys. Rev. Lett. 53 (1984) 738 and M. Leon. *ibid.*, p. 739,
18. We are grateful to L.L. Menshikov for bringing this point to our attention,

Figure captions

Fig. 1. Kinetic graph presenting the processes forming one muon-catalyzed fusion cycle in deuterium.

The nodes represent : (μ) - free muon, ($d\mu$) - deuterium μ -atom, ($1,1; K, \nu$) $\equiv (\mu m r)$ - μ -molecular resonance compound in a $J = 1, \nu = 1, K, \nu$ state, ($1,1; K, \nu - 1$) - the same compound after vibrational de-excitation, ($1,0$) - $d\mu d$ in the $J = 1, \nu = 0$ state. (F) is a subsidiary vertex representing fusion in $d\mu d$.

Fig. 2. $\lambda_m^{obs} = \varphi^{-1}(R_{d\mu} - \lambda_0)$ as a function of target density (in units of LHD) for different α , λ_0 and λ_d^0 listed in the upper-left corner. All curves correspond to $\lambda_m^{obs}(\varphi = 1.035) = 2.76 (\pm 0.08) \mu s^{-1}$ of [4]. The inset (b) (bottom-right) shows in more detail the region $\varphi < 0.1$ (bottom and lower-right scales). The numbers in brackets in the inset are the "true" values of λ_m^0 .

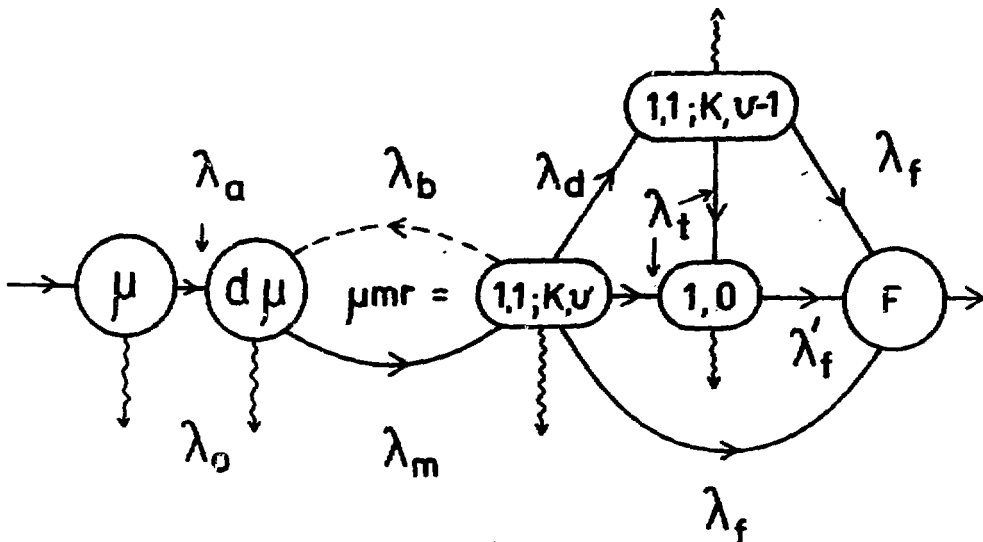


Fig. 1

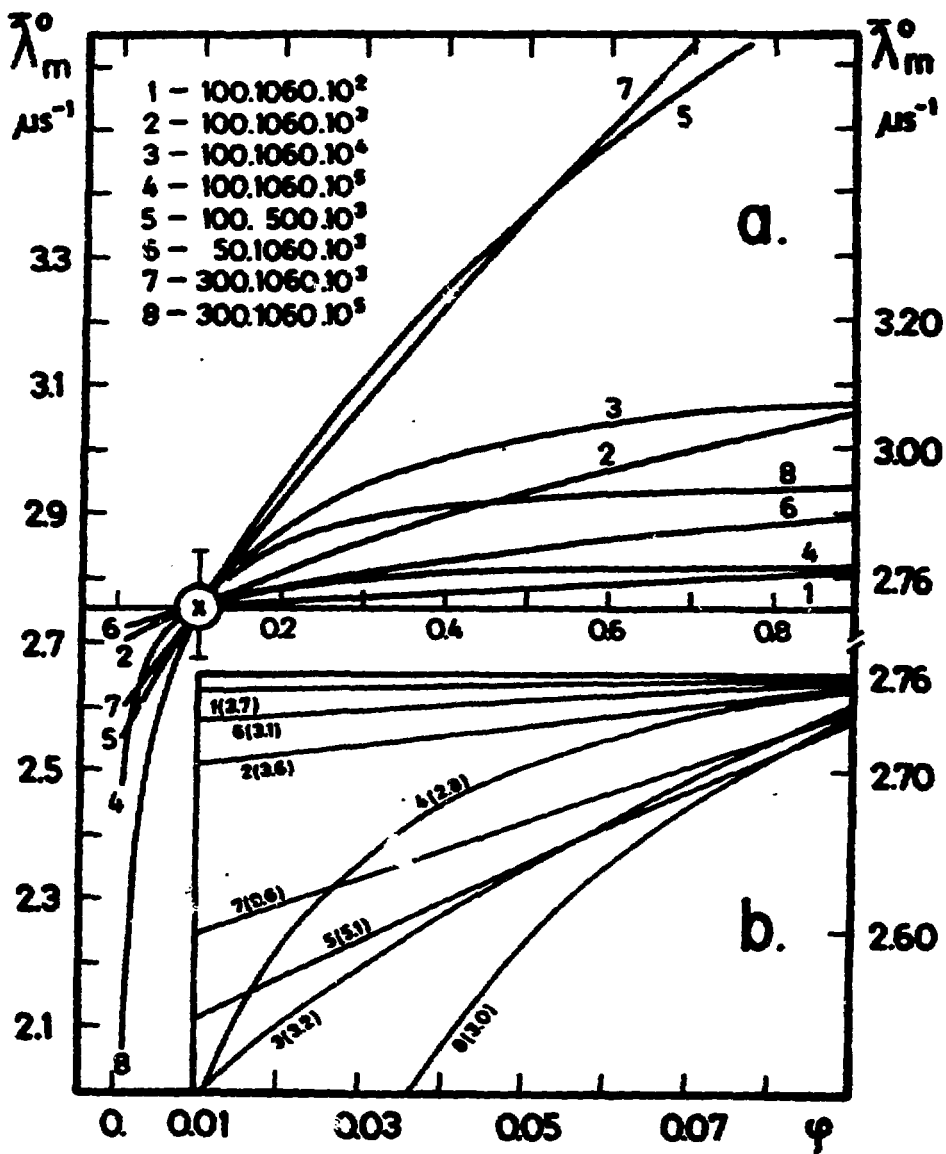


Fig. 2

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