

Reply to “Comment on ‘Effect of rotation on the elastic moduli of solid ^4He ’”

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We respond to the Comment by Pushkarov. We find that, in solid ^4He , ^3He impurity atoms that pin dislocation segments are not impuritons, but randomly walking particles. The interpretation for the rotation effect, which was based on the impuriton model, does have to be totally reconsidered, taking into account the motion of randomly walking ^3He in the strong elastic potential near the core of edge dislocations.

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The elastic anomaly of solid ^4He discovered by Day and Beamish has been elucidated well as an effect of ^3He impurity atoms to the dynamics of dislocation network [1]. We have found that the elastic anomaly, an increase in shear modulus below 0.2 K, is suppressed when the solid ^4He sample is rotated unidirectionally up to 4 rad/s [2]. Since the dynamics of the dislocation network seems to be hardly influenced by sample rotation with several rad/s, we have concluded that the pinning probability of ^3He impurities are suppressed several orders of magnitude by rotation [2]. Past studies suggested that ^3He impurity atoms behave as so-called impuritons, a Bloch-like wave in the periodic potential of a solid ^4He matrix. We have presented a possible mechanism based on the impuriton picture that the circular motion of ^3He impuritons by Coriolis force can decrease the probability of pinning of ^3He on dislocation segments [2]. Pushkarov criticizes our interpretation concerning the Coriolis force acting on ^3He impuritons [3]. The claims are summarized as follows:

(1) We proposed that the Coriolis force exerted to ^3He impuritons can be enhanced by effective mass, but no enhancement occurs with gravitational and centrifugal forces. This is in contradiction with equality of inertial and gravitational masses. (2) The quantity responsible for inertial effects of ^3He impuritons in solid ^4He is not the effective mass, but the bare mass of ^3He . Therefore, the enhancement of Coriolis force proposed by us does not occur. (3) Pushkarov proposed in previous theoretical work that diffusion of vacancies in solid ^4He should show strong spatial anisotropy when the solid ^4He sample is rotated. We mentioned this theoretical result as a possible example of rotation effect on defect dynamics in solid ^4He . Pushkarov has applied the theory to the case of ^3He impurity and concluded that the rotation effect does not exist in our current experimental condition.

We agree with the claims: The equivalence principle should hold even in the effective band mass of impuritons. As to claim (2), we point out a counterexperiment that showed that the macroscopic motion of a neutron in a Si crystal is determined by the band effective mass of the neutron [4]. However, this experiment has also confirmed that the equivalence principle holds between gravitational and inertial

effective masses. Since the effective mass of ^3He impuritons can reach 10^6 times the bare mass, the enhancement of ^3He mass would lead to an unrealistic spatial distribution in solid ^4He samples. Within the constraint of equivalence principle, it is difficult to conclude that the Coriolis force surpasses other forces. We therefore abandon the idea that the motion of ^3He is governed by the impuriton effective mass. In the last claim (3), Pushkarov develops his theory to the case of ^3He impuritons, and concludes that the rotation effect is absent in our experimental condition. We also agree with this conclusion, although we did not apply this theory to our experimental result in our publication [2].

Although the elasticity of solid ^4He is dominated by interaction between ^3He impurity and dislocation segment, the interaction was not theoretically studied at all. We find that, from a simple discussion shown below, the impuriton picture is inappropriate in discussing pinning of ^3He to dislocations because of the strong elastic interaction between dislocation and ^3He impurity.

Under external potential $U(r)$, the state of ^3He impurity is determined by the magnitude relation between impuriton bandwidth Δ and the change in potential when impurity moves from a site to a neighbor site, $a|\nabla U(r)|$, where a is a lattice constant [5]. The impuriton picture is valid at position r_0 only if the condition $\Delta > a|\nabla U(r)|_{r=r_0}$ is satisfied [5]. On the other hand, if $\Delta < a|\nabla U(r)|_{r=r_0}$, i.e., external force is large or bandwidth is small, the impuriton picture is not justified. Landesman, Winter, and Yamashita showed that in this case ^3He can move from site to site by random walk [6]. Since ^3He atoms are attracted to the dislocation core by elastic interaction, the impuriton picture breaks down near a dislocation segment. We schematically explain the situation.

Suppose a straight edge dislocation in the z direction as shown in Figs. 1(a) and 1(b). The origin and xy axes are taken in the figure. The ^3He dislocation elastic interaction, which is caused by the volume difference of a ^3He atom in the solid ^4He , is given by the following formula [7]:

$$W(x, y) = \frac{4}{3} \left(\frac{1+v}{1-v} \right) \frac{\mu b \epsilon r_0^3 y}{x^2 + y^2 + r_0^2}, \quad (1)$$

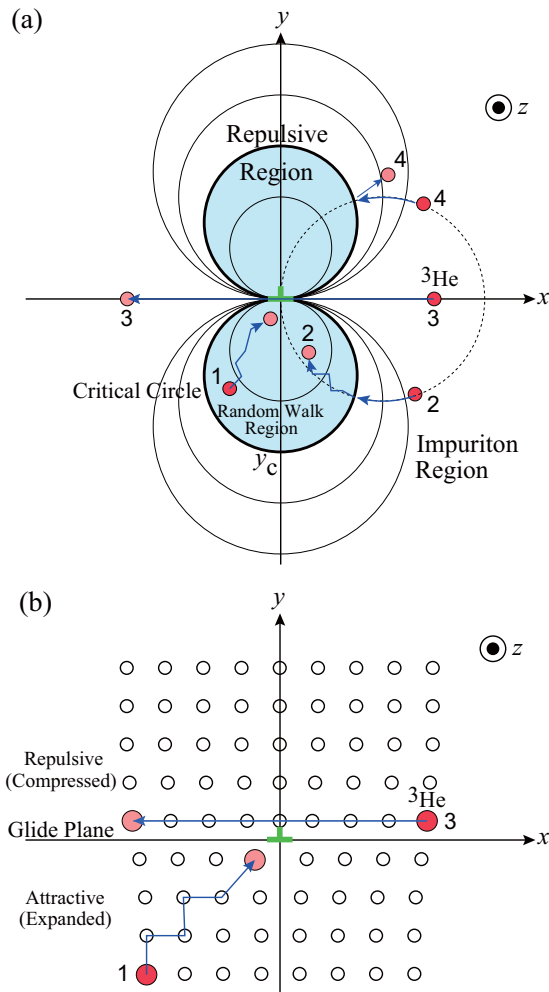


FIG. 1. (a) Solid circles show equipotential lines of the elastic interaction Eq. (1) between a ^3He atom and a straight edge dislocation located at the origin (shown as a mark \perp) [7]. Bold solid circles show the “critical circles” at which $\Delta = a|\nabla W(x, y)|$ [5]. The impuriton picture is valid only outside the circles, while inside them (light-blue region) ^3He atoms can move only by random walk [6]. Red and orange circles schematically indicate ^3He atoms, with numbers 1–4. Atom 1 can approach the dislocation core by random walk. Atom 2 propagates on a potential gradient shown by dashed circle, and is converted to randomly walking atom. Atom 3 propagates along the glide plane as an impuriton, and passes through the dislocation core. Atom 4 propagates as an impuriton in the repulsive region, and is reflected at the critical circle. (b) The dislocation core region on atomic scale. The numbering of ^3He atoms has the same meaning as (a). Atom 1 is trapped at the most expanded lattice site after random walk. Atom 3 runs on the glide plane and passes through the core.

where ν is Poisson’s ratio (~ 0.33), μ is the shear modulus (~ 15 MPa), and b is the Burgers vector. When the radii of

^4He and ^3He atoms are given by r_4 and r_3 respectively, ϵ is defined as $\epsilon = (r_3 - r_4)/r_4$. $r'_0 \sim 2b/3$ is a correction for violation of Hooke’s law at the dislocation core. Iwasa and Suzuki [8] estimated the depth of the potential at the core of an edge dislocation to be about 0.6 K, which is very close to the value of the pinning energy of ^3He , 0.7 K, which was determined from recent elastic measurements [1]. Therefore, Eq. (1) is valid for solid ^4He .

Equipotential lines of this interaction are shown in Fig. 1(a). ^3He is attracted by dislocation in the lower half region, while in the upper half one the interaction is repulsive. The equipotential lines are circles bunched at the origin (dislocation core). Taking $x = 0$, we can estimate the “critical diameter” of the circle y_c , at which the relation $\Delta = a|\nabla W(x = 0, y)|_{y=y_c}$ holds. For our two solid samples with pressures $P = 3.6$ and 5.4 MPa, the diameter y_c is estimated to be about 3×10^{-7} and 1×10^{-6} m, respectively. These values are 1/10–1/30 of the average length of dislocation segments, i.e., separation of segments, $L_N \sim 1 \times 10^{-5}$ m.

The impuriton picture is valid only outside the “critical circles” of diameter y_c , which are shown by light blue circles (the circle in the upper half plane can also be the critical circle). As shown in Fig. 1(b), ^3He atoms are trapped at the sites where the lattice is most expanded. We easily see that ^3He cannot approach the dislocation core as an impuriton, because $a|\nabla W(x, y)|$ exceeds Δ inside the critical circle. ^3He can approach the trapping site only as a randomly walking particle, which is shown by an illustration number 1 in Figs. 1(a) and 1(b). The impuriton propagating outside the critical circle (shown by the number 2) can be converted to the randomly walking particle. When ^3He moves along the gliding plane as shown by an atom number 3 in Figs. 1(a) and 1(b), it can behave as an impuriton. However, since there is no potential gradient in the direction of the glide plane, the impuriton cannot be trapped to the dislocation but passes through the dislocation core region. Eventually, there is no possibility that impuritons are directly trapped by dislocations. Note that ^3He impuritons in the repulsive region are reflected at the critical circle, as shown by an atom number 4 in Fig. 1(a).

In real solid ^4He , the critical circles will exist around the edge and mixed dislocation segments in a three-dimensional dislocation network, but not around screw dislocations, which do not trap ^3He because of no volume expansion. The random walk of ^3He will be assisted by phonons and potential-energy fluctuations caused by movement of dislocations.

To conclude, ^3He atoms that pin dislocation segments are not impuritons, but randomly walking particles. Therefore, the interpretation for the rotation effect found by us [2] has to be totally reconsidered taking into account the motion of randomly walking ^3He in the strong elastic potential near-edge dislocation core.

[1] J. Day and J. Beamish, *Nature (London)* **450**, 853 (2007); I. Iwasa, *J. Low Temp. Phys.* **171**, 30 (2013); S. Balibar *et al.*, *C. R. Phys.* **17**, 264 (2016).

[2] T. Tsuiki, D. Takahashi, S. Murakawa, Y. Okuda, K. Kono, and K. Shirahama, *Phys. Rev. B* **97**, 054516 (2018).

- [3] D. I. Pushkarov, preceding comment, *Phys. Rev. B* **99**, 066501 (2019).
- [4] K. Raum, M. Koellner, A. Zeilinger, M. Arif, and R. Gähler, *Phys. Rev. Lett.* **74**, 2859 (1995); K. Raum, M. Weber, R. Gähler, and A. Zeilinger, *J. Phys. Soc. Jpn. Suppl. A* **65**, 277 (1996); K. Raum, J. Felber, M. A. Horne, P. Geltenbort, and A. Zeilinger (unpublished), available at <http://www.univie.ac.at/qfp/publications3/pdf/1998-18.pdf> (1998).
- [5] A. F. Andreev and I. M. Lifshitz, *Sov. Phys. JETP* **29**, 1108 (1969); A. F. Andreev, *Sov. Phys. Usp.* **19**, 137 (1976).
- [6] A. Landesman and J. M. Winter, *Proceedings of LT13*, edited by K. D. Timmerhaus (Plenum Press, New York, 1973), p. 73; A. Landesman, *Phys. Lett. A* **54**, 137 (1975); These articles did not explicitly describe that ^3He impurities take random walks in large elastic potential gradients. The random-walk picture was clearly stated in the following article: Y. Yamashita, *Bussei Kenkyu (Condensed Matter Research)* **28**, F26 (1977) (in Japanese). Article available at <http://hdl.handle.net/2433/89402>.
- [7] H. Saka, *Classical Theory of Crystal Dislocations: From Iron to Gallium Nitride* (World Scientific, Singapore, 2017), p. 93; A. H. Cottrell, *Dislocations and Plastic Flow in Crystals* (Oxford University Press, New York, 1953), p. 57.
- [8] I. Iwasa and H. Suzuki, *J. Phys. Soc. Jpn.* **49**, 1722 (1980).