

FROM THE WARDEN
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Dear Robin,

Many thanks for your paper for submission to Chemical Physics Letters. I have read it carefully and shall send it off at once with a recommendation that it should be published.

It is a pretty piece of work and I congratulate you on having got it out so nicely. I have talked about this possibility several times with my colleagues here, but we have never had the initiative to make it work, though you may have seen a short paper we wrote to Molecular Physics some years ago about the relative relaxation times of the components of an AB quartet. I think it is a beautiful method. Do you think there is any hope of reducing the computing time, which seems to me to be prohibitive for general use at the moment?

Yours sincerely,



R. E. Richards

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