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Journal Name:	Physical Science International Journal
Manuscript Number:	**Ms_PSIJ_27742**
Title of the Manuscript:	**The Dielectric behavior of Acetone and Dimethylformamide in Electric Field**
Type of the Article	**Original Research Article**

General guideline for Peer Review process:

This journal's peer review policy states that **NO** manuscript should be rejected only on the basis of '**lack of Novelty'**, provided the manuscript is scientifically robust and technically sound.
To know the complete guideline for Peer Review process, reviewers are requested to visit this link:

(http://www.sciencedomain.org/page.php?id=sdi-general-editorial-policy#Peer-Review-Guideline)

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PART 1: Review Comments

	Reviewer's comment	**Author's comment** *(if agreed with reviewer, correct the manuscript and highlight that part in the manuscript. It is mandatory that authors should write his/her feedback here)*
Compulsory REVISION comments	**Abstract** Result indicates that at frequency 0.1GHz and temperature 10oC, the calculated value for acetone was 22.4637 and that of DMF was 40.3380. This shows that an increase in dielectric constant caused a corresponding decrease in the loss factor, which means an inverse relation between the frequency and temperature. These results compared well with experimental values for both acetone and DMF at the same frequency temperature of 22.21 ± 0.04 and 40.34 ± 0.06 respectively. *-It is unclear about what parameter values are given in the abstract and there is no specific result or values have been discussed in the result or conclusion.* *-The author also has been stated about simulation butit is not clear about the simulation result and the experimental results in the presentation.* **Introduction** produced by the applied **voltage (Craig, 1995).** *-Follow one type of reference writing format. In this case I suggest that the author to restructure the writing format and provide the reference accordingly.* *-Re-label the figures in order.*	

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	Total Polarization P_e is the orientation polarization P_i is ionic polarization P_o is the orientation polarization P_s is the space charge polarization *-rectify the mistake in explanation* *Abstract and Conclusion needs attention.*	
Minor REVISION comments		
Optional/General comments		

Reviewer Details:

Name:	**Subramaniam Jahanadan**
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