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SDI Review Form 1.6

Journal Name:	Asian Journal of Physical Sciences
Manuscript Number:	**Ms_AJOPS_30695**
Title of the Manuscript:	**THE COHESIVE ENERGY CALCULATIONS OF SOME BCC (Li, Cr, Fe, Mo) LATTICES USING DENSITY FUNCTIONAL THEORY**
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)

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		
Minor REVISION comments	(i) The reason for only one alkali and three transition metals, is not clear. (ii) Line 39 is not clear. (iii) Input and output parameters in a tabular form will do good. (iv) The reasons for overestimation are not mentioned, in particular, for Cr and Fe. A concluding remark about it should be added.	
Optional/General comments		

Reviewer Details:

Name:	**Amitava Ghorai**
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