



SDI Review Form 1.6

Journal Name:	International Research Journal of Pure and Applied Chemistry
Manuscript Number:	Ms_IRJPAC_37695
Title of the Manuscript:	3D-QSAR Modeling and Molecular Docking Studies on a series of 1,2,4 triazole containing diarylpyrazolyl carboxamide as CB1 cannabinoid receptor ligand
Type of the 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	The manuscript is of great interest for computational chemistry readers. It is well written with minor revisions necessary in language.	Done
Optional/General comments	Only <i>in silico</i> studies are involved in this paper whereas the structure activity relationship must involve the <i>in vivo</i> or <i>in vitro studies</i> also. So the author may include those studies or may add the reference for it so that visible relation between the two can be observed	A reference [21] is added as in vivo study for CB1 receptor binding affinity