# SCIENCEDOMAIN international

www.sciencedomain.org



# **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 <u>NO</u> manuscript should be rejected only on the basis of '<u>lack of Novelty'</u>, 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	In the Introduction section, the authors must not showed previous conclusions on the present manuscript (page 2, line 44-46).	
	In the Introduction section, the authors need to update the information, once must of the references are very old (page 14 - 1987, 1991, 1992, 1993)	
	Please, provide the empirical scoring function used in the docking calculations. In addition, do the authors used the re-docking calculations to validate the scoring function?	
	I recommend that the authors check the English language, because there are some mistakes, as example page 3 and 4 lines 75 and 89, respectively.	
Minor REVISION comments		
Optional/General comments		

# **Reviewer Details:**

Name:	Otávio Augusto Chaves
Department, University & Country	Department Of Chemistry, Universidade Federal Rural Do Rio De Janeiro, Brazil

Created by: EA Checked by: ME Approved by: CEO Version: 1.6 (07-06-2013)