

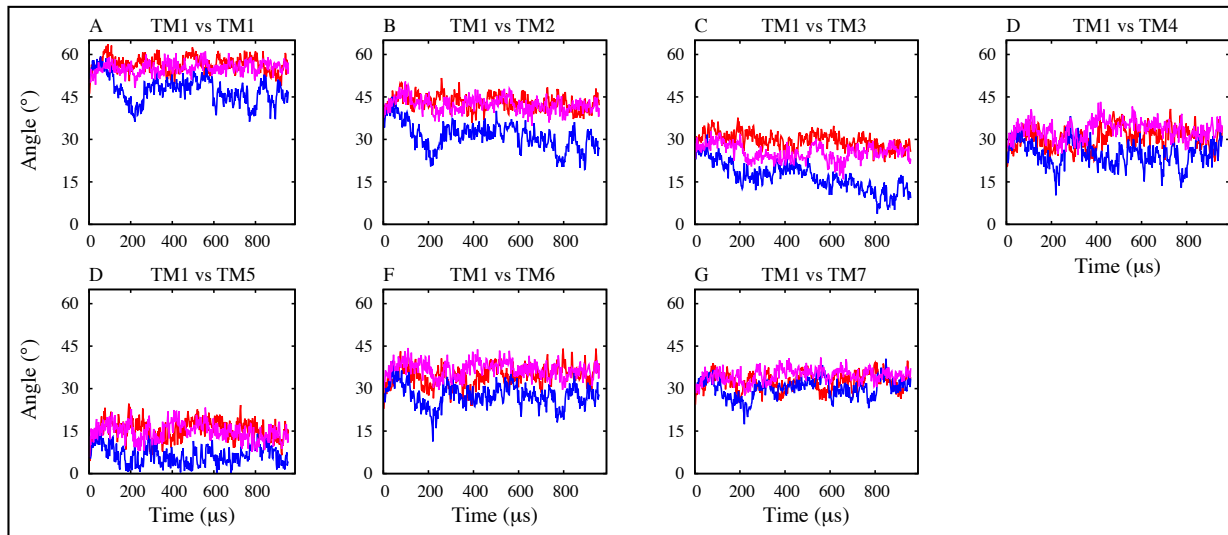
**Supporting Information: Cholesterol Dependence of the Conformational  
Changes in Metabotropic Glutamate Receptor 1**

Ugochi H. Isu, Shadi A. Badiiee, Adithya Polasa, Seyed H. Tabari, Mortaza

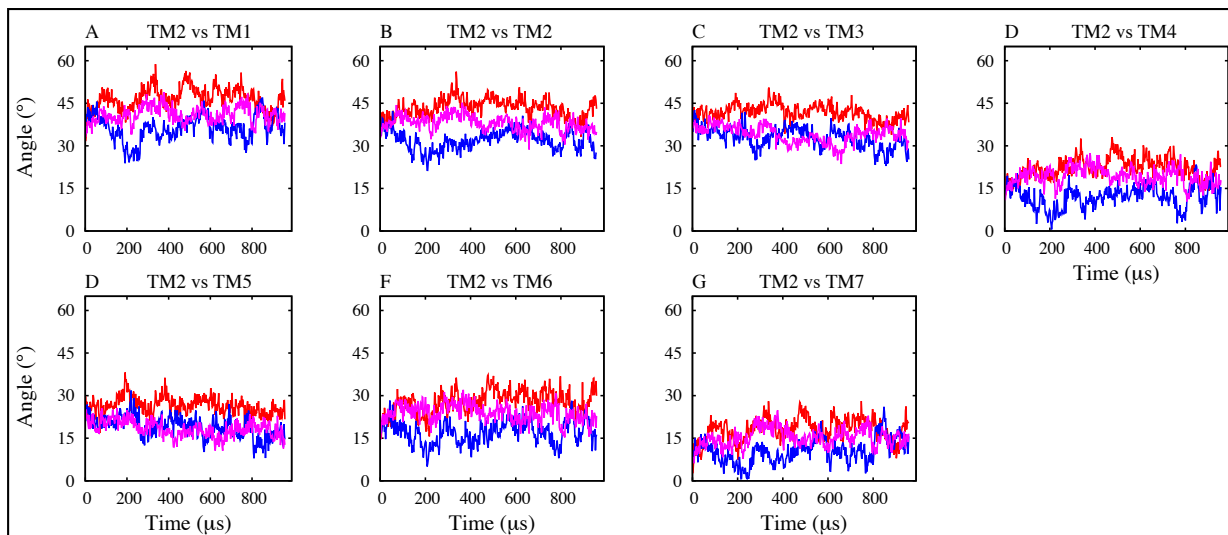
Derakhshani-Molayousefi, and Mahmoud Moradi \*

Department of Chemistry and Biochemistry, University of Arkansas, Fayetteville, AR 72701

E-mail: moradi@uark.edu



**Fig. S1.** Time series plot showing the interhelical angle throughout the simulations for simulation set 1. The interhelical angle between TM1 and TM2-TM7 has been calculated for each frame of the simulation.



**Fig. S2.** Time series plot showing the interhelical angle throughout the simulations for simulation set 1. The interhelical angle between TM2 and TM1, TM3-TM7 has been calculated for each frame of the simulation.

## Tables

**Table S1: Occupancy (%) of Hydrogen Bond Interactions of All Protomers from Simulation Set 1**

System	0 % CHOL		10 % CHOL		25 % CHOL	
	A	B	A	B	A	B
<b>Y672-T794</b>	3	12	13	8	<b>98</b>	6
<b>T768-Y792</b>	0	0	0	0	<b>51</b>	0
<b>S627-K678</b>	<b>19</b>	59	<b>35</b>	61	68	73
<b>M791-T768</b>	94	95	96	95	<b>34</b>	95
<b>Y770-N680</b>	95	96	96	<b>51</b>	98	<b>33</b>
<b>R661-E728</b>	3	46	<b>84</b>	13	14	8
<b>E783-K834</b>	74	98	97	97	94	73