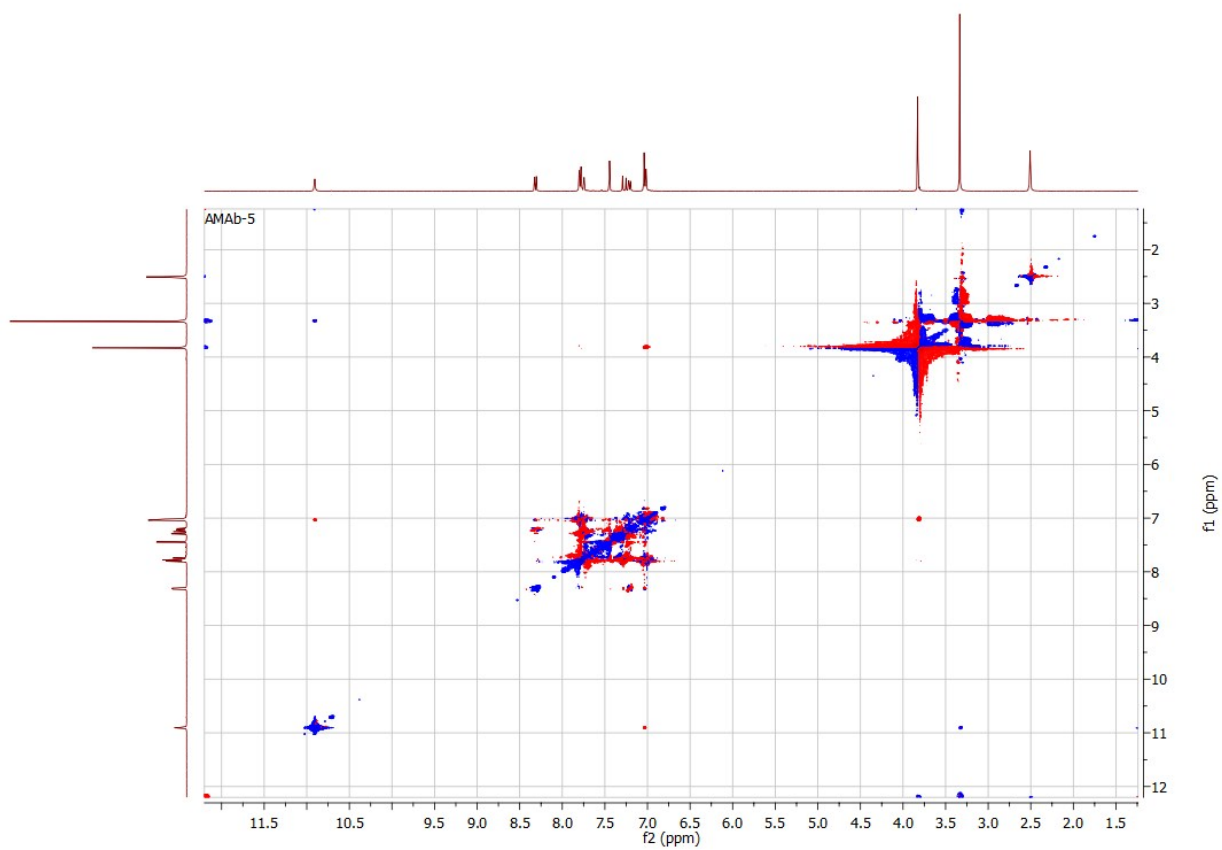
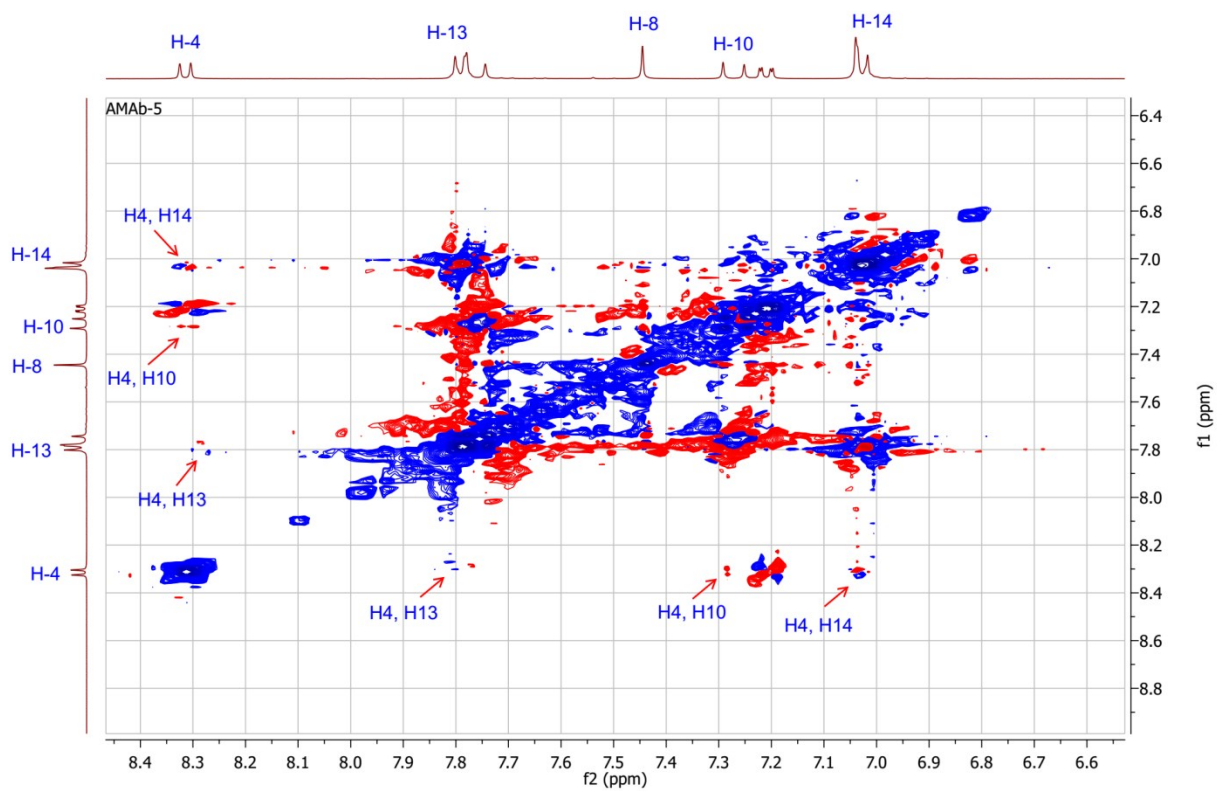


Supplementary material

Saccharomonosporine A Inspiration; Synthesis of Potent Analogues as Potential PIM Kinases Inhibitors

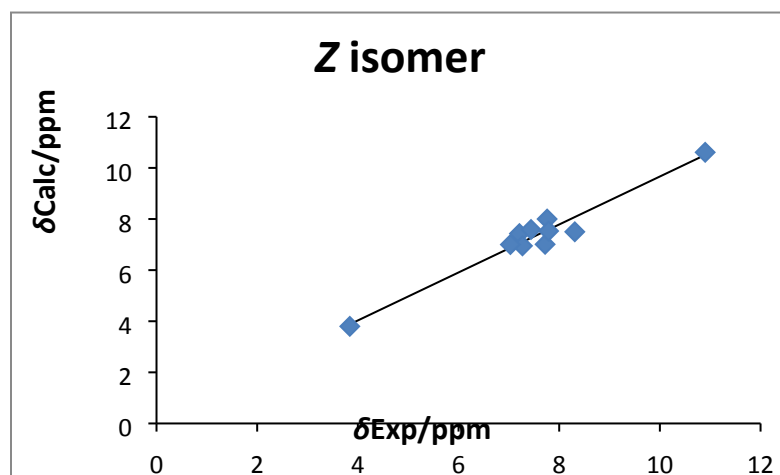


SI. ^1H - ^1H NOESY chart for compound 5.

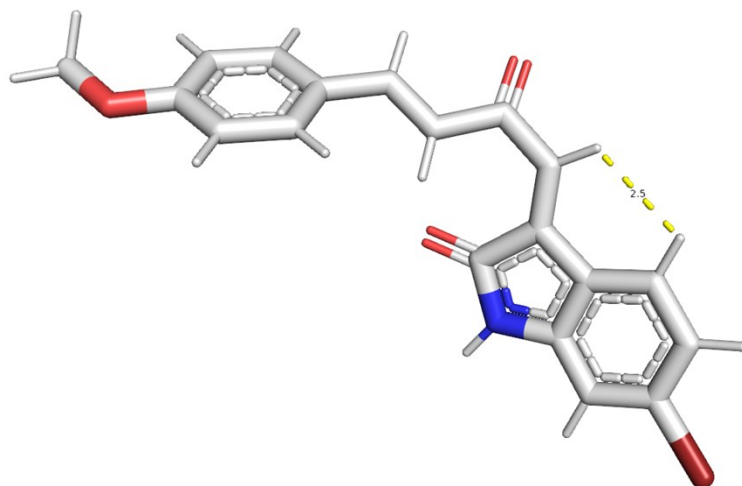


S2. ^1H - ^1H NOESY chart for compound **5**.

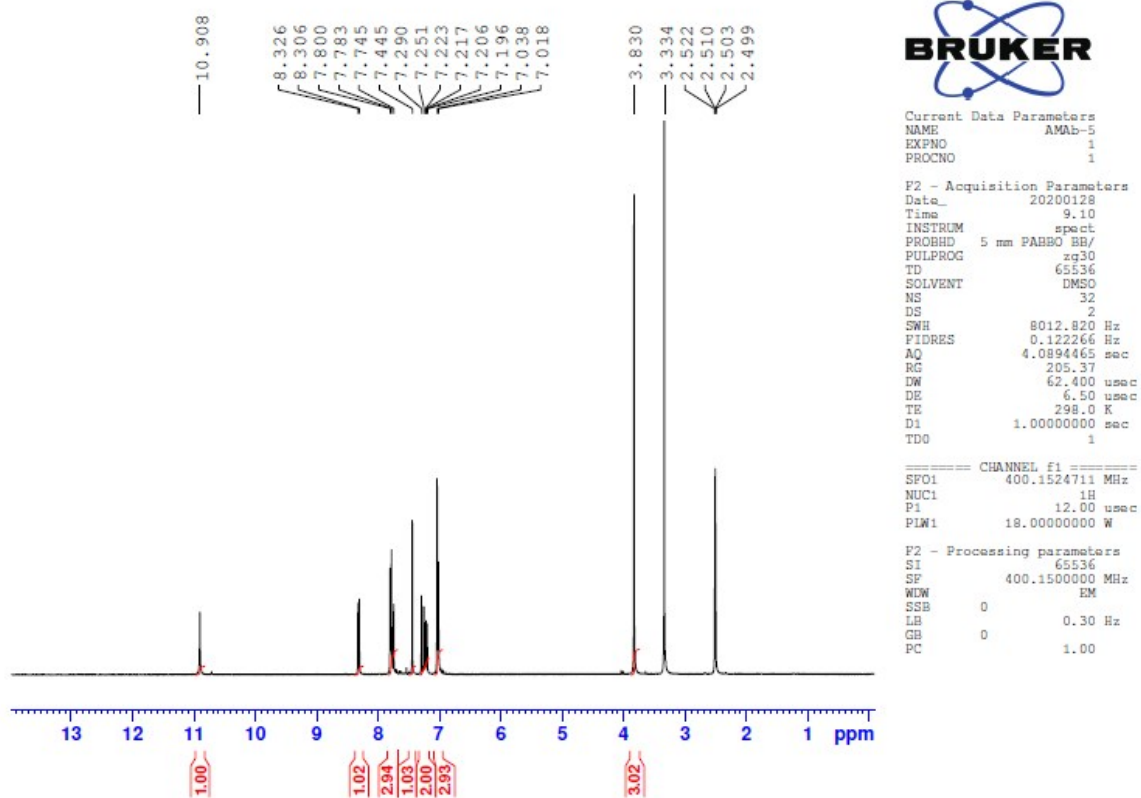
No.	$\delta_{\text{Exp/ppm}}$	$\delta_{\text{Calc/ppm}}$ (<i>E</i> isomer)	$\delta_{\text{Calc/ppm}}$ (<i>Z</i> isomer)
1	10.9	10.61	10.61
2	-	-	-
3	-	-	-
3a	-	-	-
4	8.31	7.4	7.5
5	7.21	7.4	7.43
6	-	-	-
7	7.44	7.58	7.59
7a	-	-	-
8	7.72	7.66	7.01
9	-	-	-
10	7.27	7.06	6.95
11	7.76	8.06	8
12	-	-	-
13	7.79	7.53	7.53
14	7.03	7.01	7
15	-	-	-
16	3.84	3.81	3.8
Calculated Relative Energy		MMFF Energy= 40.1537 kcal/mol	MMFF Energy= 45.2583 kcal/mol



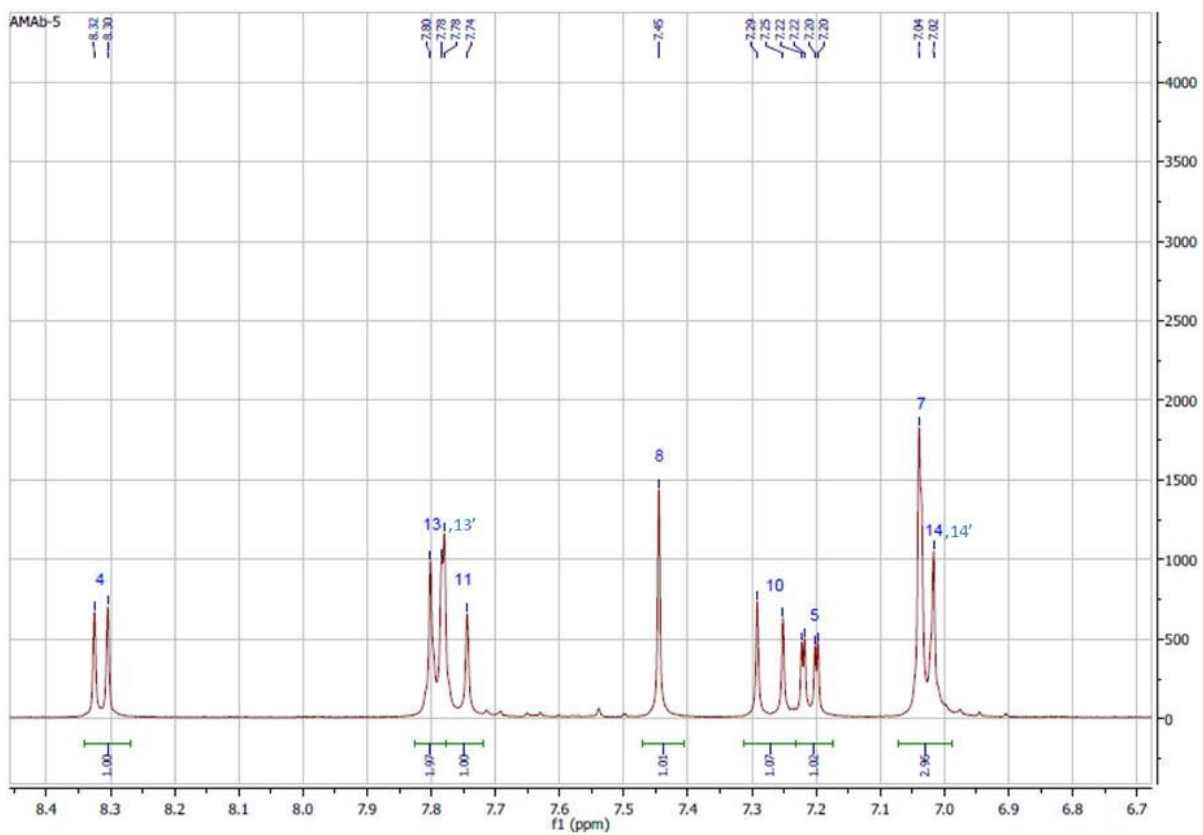
S3. Theoretical calculations of ^1H NMR chemical shifts of compound 5 in *E* isomer.



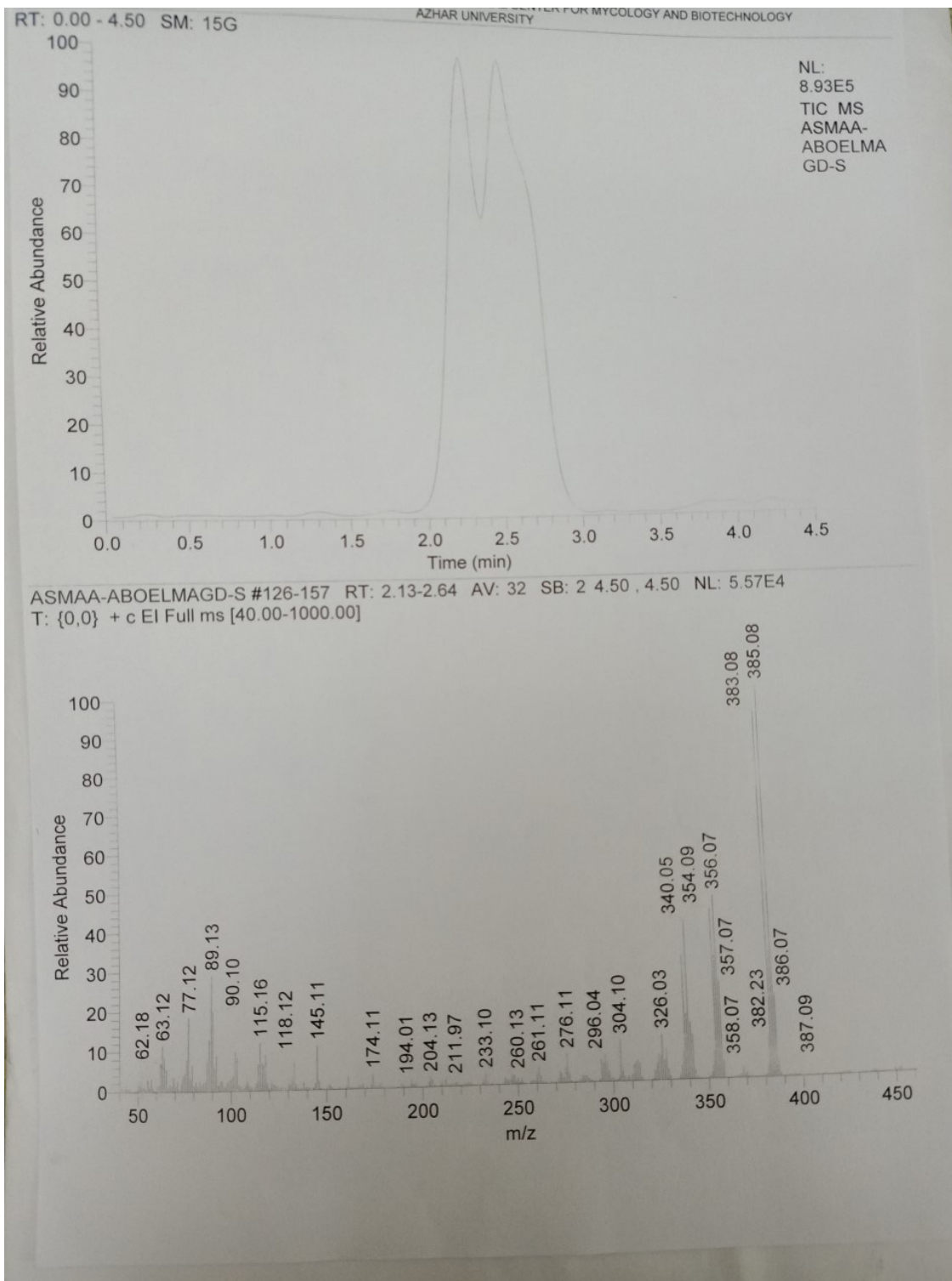
S4: ^1H - ^1H correlations of cis isomer of compound 5.



S5: ^1H NMR of compound 5.



S6: ^1H NMR of compound **5**.



S7: Mass spectrum of compound 5.