

Synthesis of Amantadine clubbed *N*-Aryl amino thiazoles as potent Urease, α -Amylase & α -Glucosidase inhibitors and their kinetic and molecular docking studies.

Synthesis of Amantadine clubbed *N*-Aryl amino thiazoles as potent Urease, α -Amylase & α -Glucosidase inhibitors, kinetic and molecular docking studies.

Fatima Tuz Zahra^a, Aamer Saeed^{*a}, Atteeque Ahmed^a, Hammad Ismail^b, Muhammad Umar Ijaz^c, Fernando Albericio,^{*d,e}

^a *Department of Chemistry Quaid-i-Azam University-45320, Islamabad, Pakistan*

^b *Department of Biochemistry and Biotechnology, University of Gujrat, Gujrat, 50700, Pakistan*

^c *Department of Zoology, Wildlife and Fisheries, University of Agriculture, Faisalabad 38040, Pakistan*

^d *Peptides Science Laboratory, School of Chemistry and Physics, University of KwaZulu-Natal, Westville, Durban 4000, South Africa.*

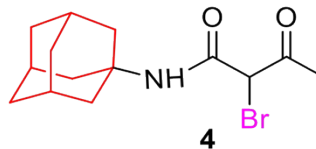
^e *CIBER-BBN, Networking Centre on Bioengineering, Biomaterials and Nanomedicine, and Department of Organic Chemistry, University of Barcelona, 08028 Barcelona, Spain.*

Correspondence

Prof. Dr. Aamer Saeed

Email: asaheed@qau.edu.pk ; ORCID: <http://orcid.org/0000-0002-7112-9296>

Contact number: +92-51-9064-2128, Fax: +92-51-9064-2241



Chemical Formula: $C_{14}H_{20}BrNO_2$

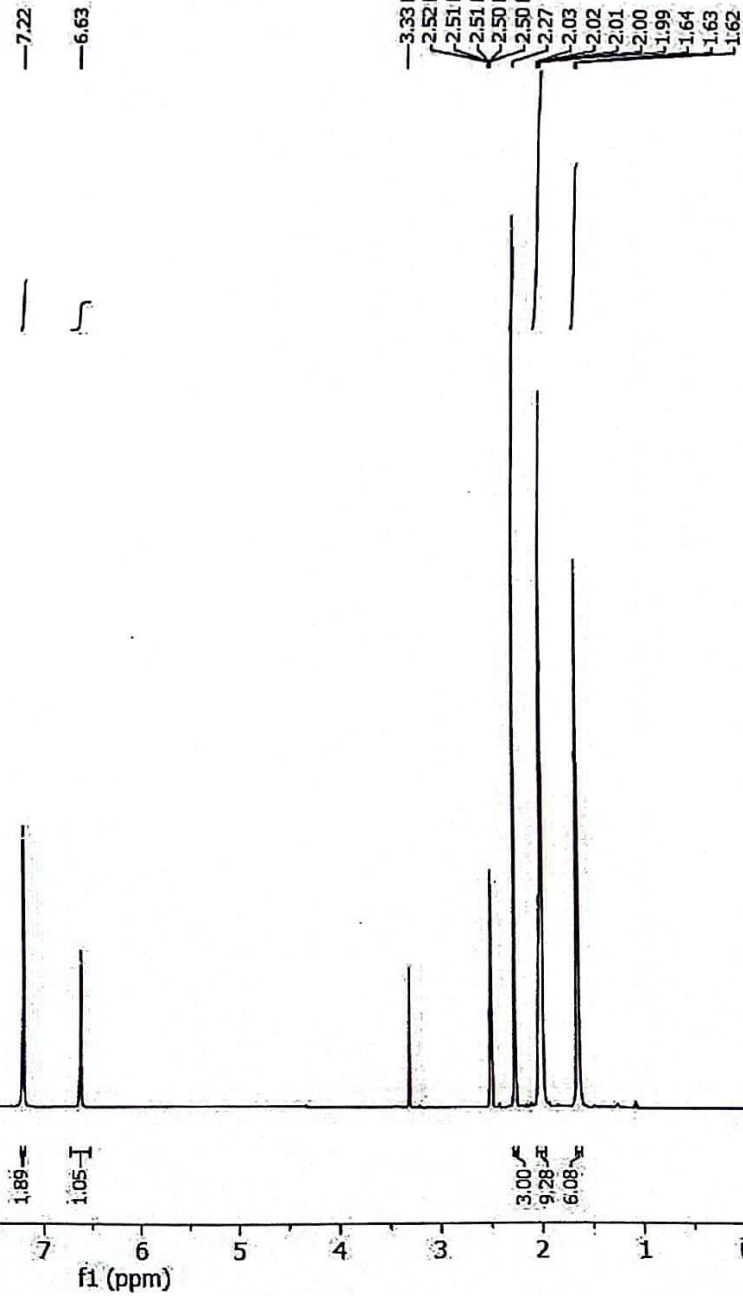
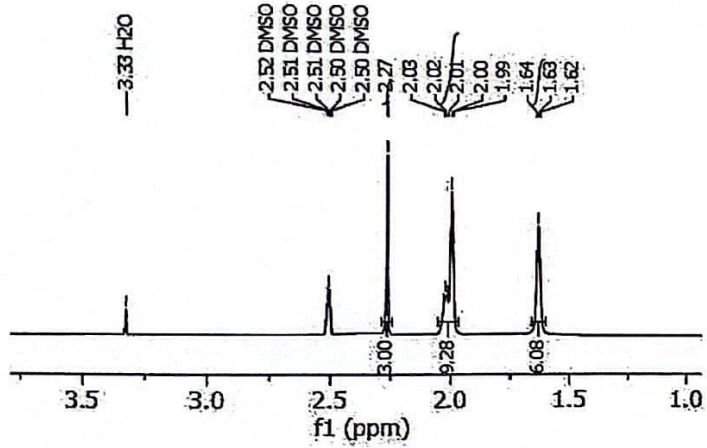
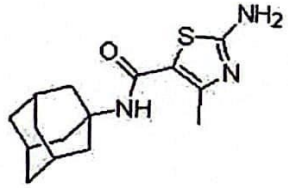
Exact Mass: 313.07

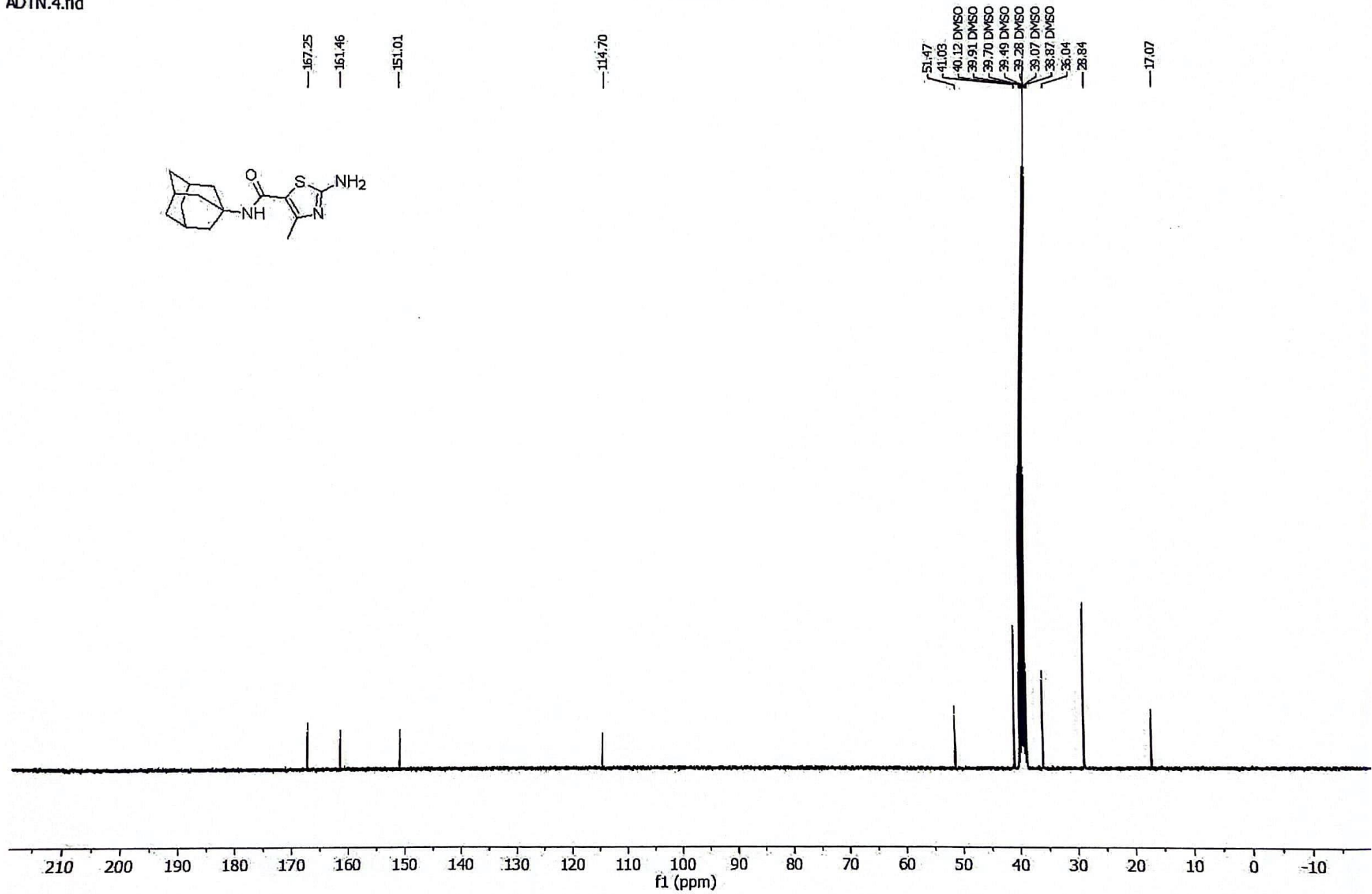
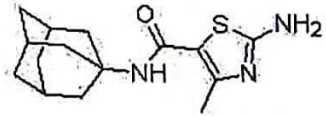
Molecular Weight: 314.22

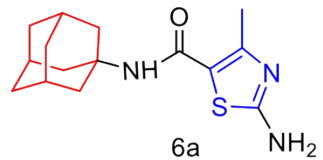
m/z: 313.07 (100.0%), 315.07 (97.3%), 314.07 (15.1%), 316.07 (14.7%), 315.07 (1.1%), 317.07 (1.0%)

Elemental Analysis: C, 53.51; H, 6.42; Br, 25.43; N, 4.46; O, 10.18

Fragmentation pattern:







Chemical Formula: C₁₅H₂₁N₃OS

Exact Mass: 291.14

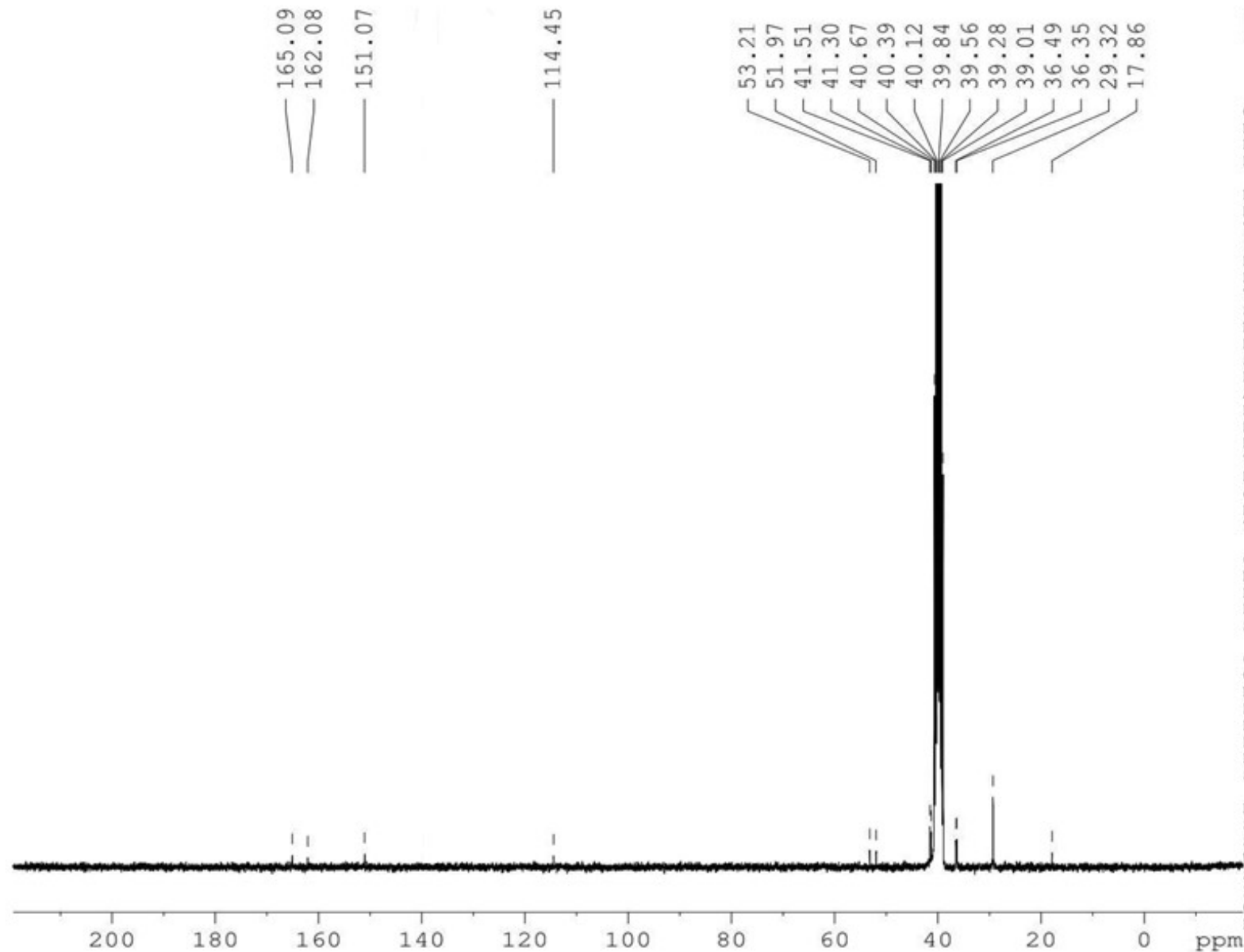
Molecular Weight: 291.41

m/z: **291.14** (100.0%), 292.14 (16.2%), 293.14 (4.5%), 293.15 (1.2%), 292.14 (1.1%)

Elemental Analysis: C, 61.82; H, 7.26; N, 14.42; O, 5.49; S, 11.00

Fragmentation pattern:

DR.AAMER SAEED/FATIMA/TH-11_13CNMR_DMSO



Current Data Parameters
NAME TH-11_13CNMR_DMSO
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters
Date_ 20230619
Time 9.30
INSTRUM spect
PROBHD 5 mm BBO BB-1H
PULPROG zgpg30
TD 35968
SOLVENT DMSO
NS 18432
DS 0
SWH 17985.611 Hz
FIDRES 0.500045 Hz
AQ 0.9999604 sec
RG 32768
DW 27.800 usec
DE 6.00 usec
TE 298.2 K
D1 2.00000000 sec
d11 0.03000000 sec
DELTA 1.89999998 sec
TD0 1

----- CHANNEL f1 -----
NUC1 13C
P1 6.00 usec
PL1 -5.00 dB
SFO1 75.4752953 MHz

----- CHANNEL f2 -----
CPDPRG2 waltz16
NUC2 1H
PCPD2 80.00 usec
PL2 2.00 dB
PL12 20.98 dB
PL13 20.00 dB
SFO2 300.1312005 MHz

F2 - Processing parameters
SI 32768
SF 75.4677490 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40