

SUPPORTING INFORMATION

Substituent Effects on EI-MS Fragmentation Patterns of 1-aryl 5-allyloxytetrazoles and 1-aryl 4- allyltetrazole-5-ones; Correlation with UV-Induced Fragmentation Channels

Alina Secrieru ^{1,2} and Rabah Oumeddour ^{1,3} and Maria L. S. Cristiano ^{1,*}

¹ CCMAR and Department of Chemistry and Pharmacy, FCT, Campus de Gambelas, University of Algarve, 8005-039 Faro, Portugal;

² Department of Chemistry, University of Liverpool, Liverpool L69 7ZD, UK

³ Laboratory of Industrial Analysis and Materials Science, Faculty MISM, University 8 Mai1945, Guelma 24000, Algeria

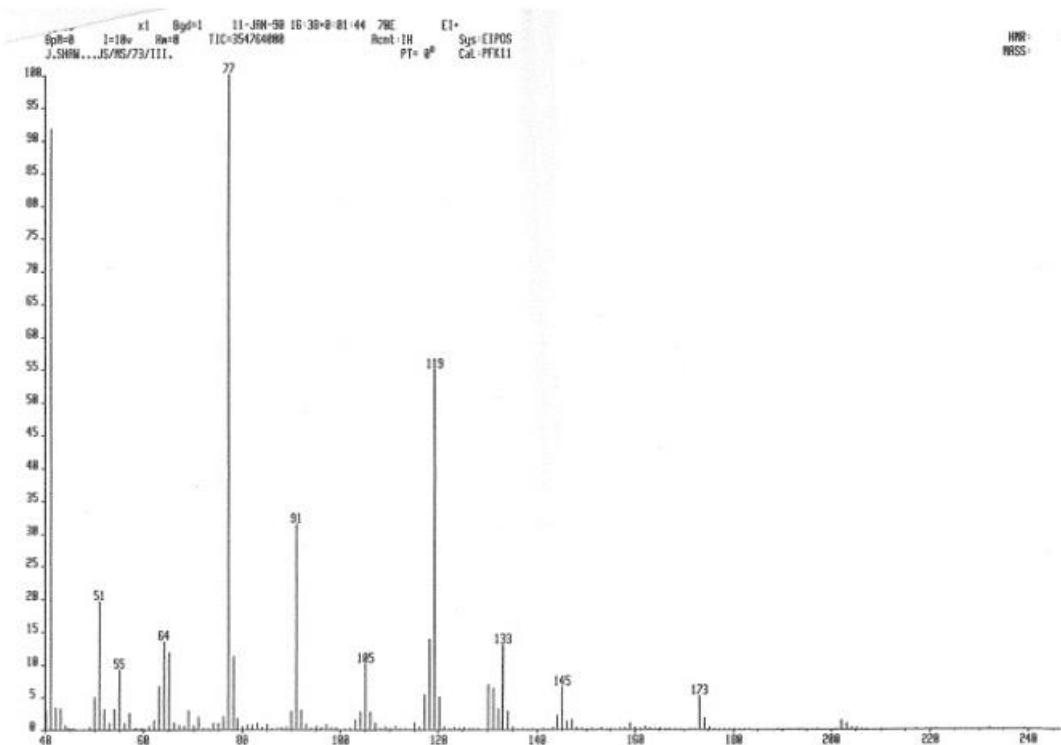


Figure S1. - EI-MS spectrum of 5-(allyloxy)-1-phenyl-1H-tetrazole (2a).

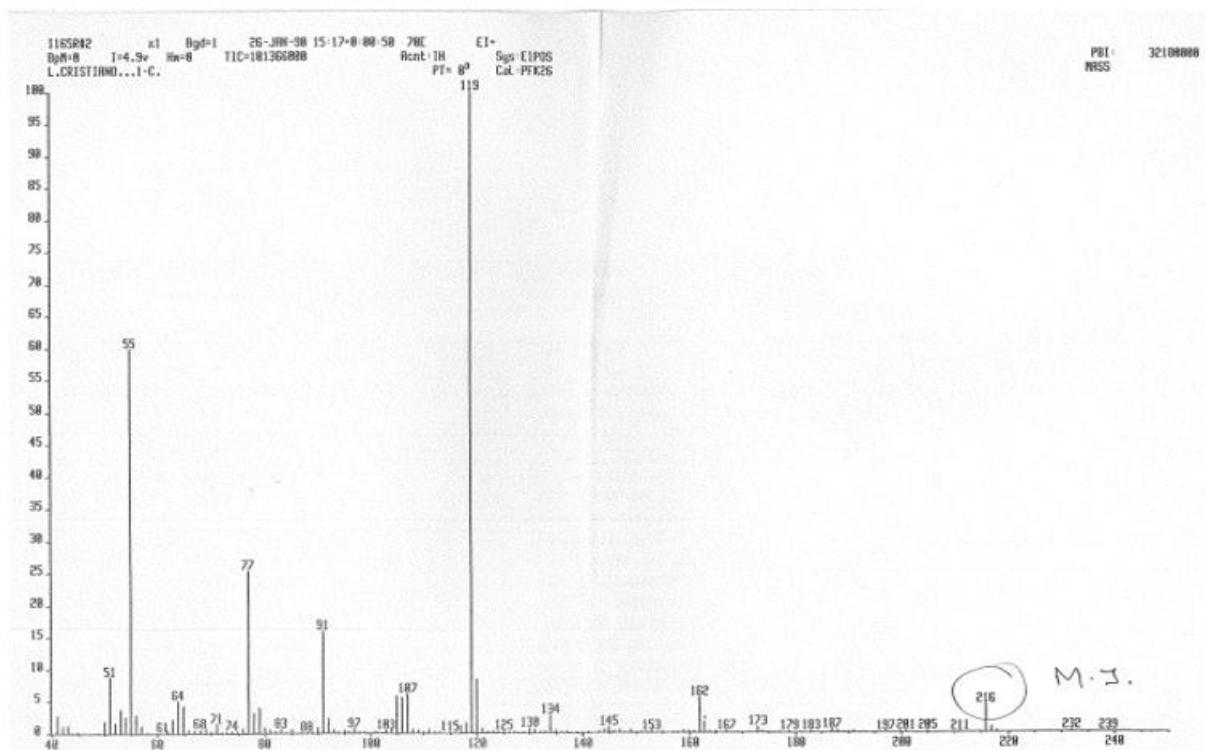


Figure S2. - EI-MS spectrum of 5-(but-2-en-1-yloxy)-1-phenyl-1H-tetrazole (2b).

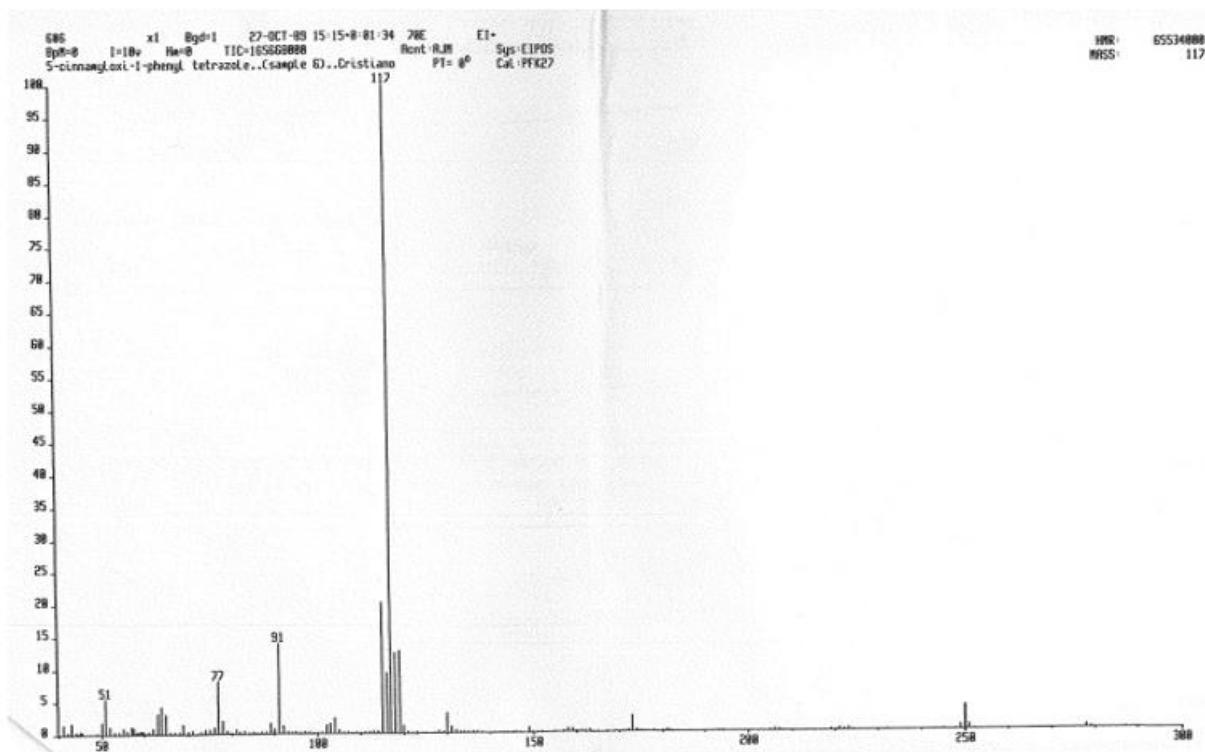


Figure S3. - EI-MS spectrum of 5-(cinnamyoxy)-1-phenyl-1H-tetrazole (2c).

H978 MM V.G LAB-BASE Chem. Dept. L'pool Univ. TRIO-1000 18-Aug-08
 Sample: Nam3.. L.CRIStIANO.. 11:22

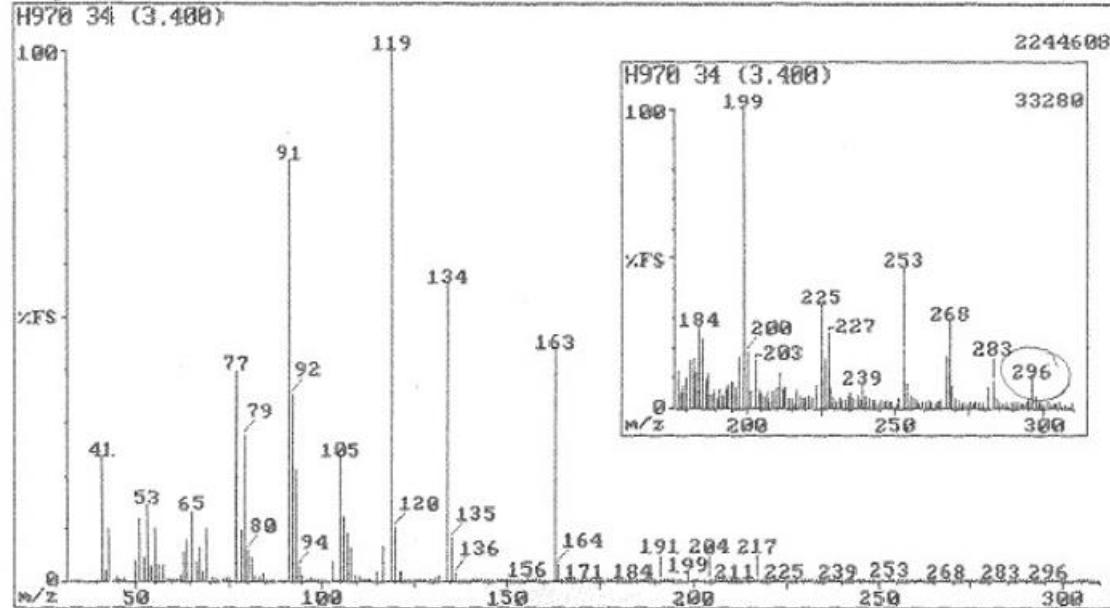


Figure S4. - EI-MS spectrum of 5-((6,6-dimethylbicyclo[3.1.1]hept-2-en-3-yl)methoxy)-1-phenyl-1H-tetrazole (2d).

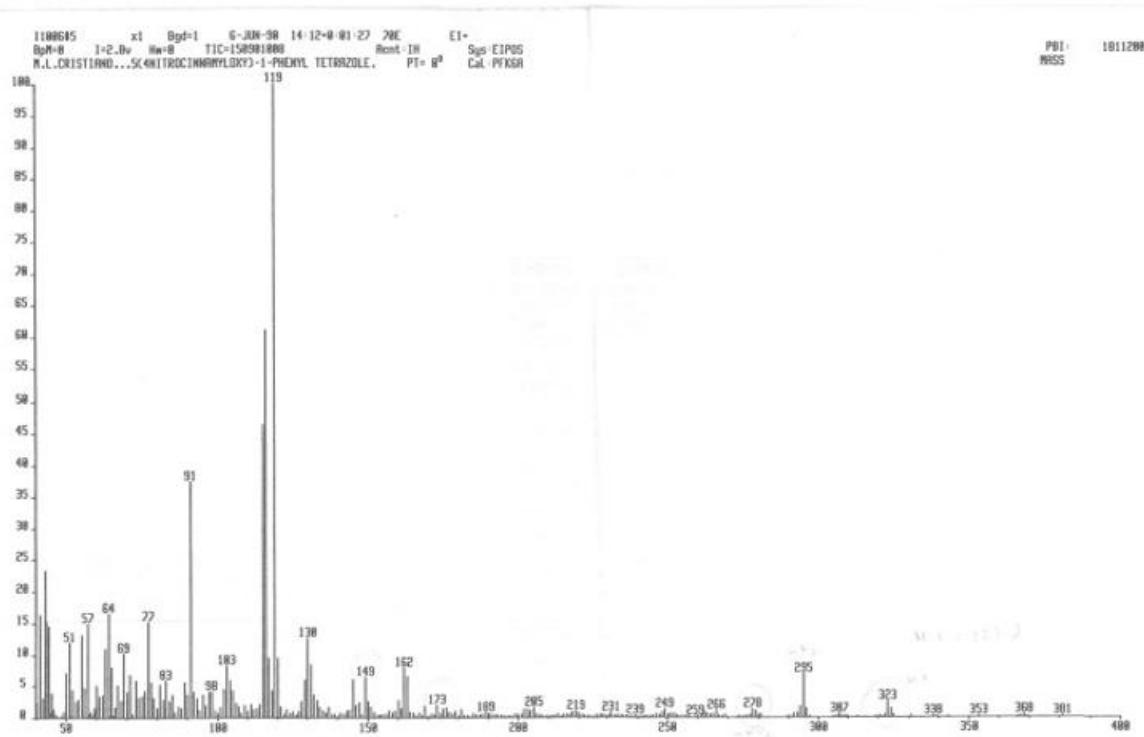


Figure S5. - EI-MS spectrum of 5-((3-(4-nitrophenyl)allyl)oxy)-1-phenyl-1H-tetrazole (2e).

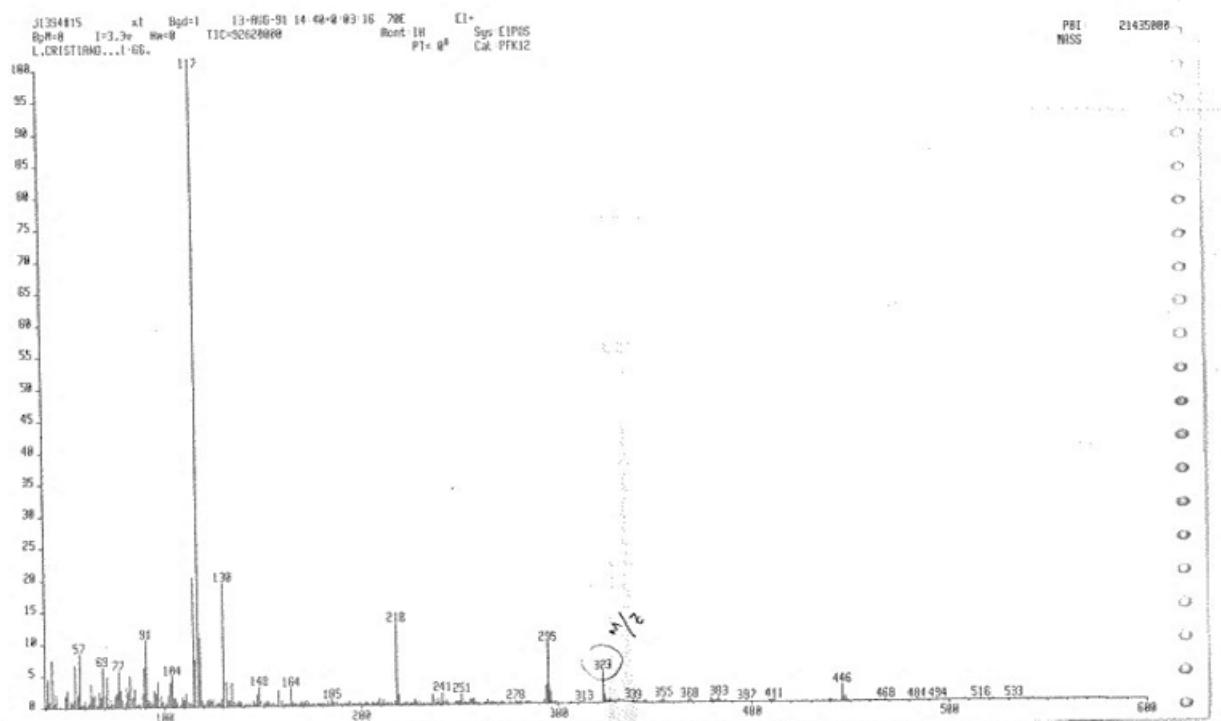


Figure S6. - EI-MS spectrum of 5-(cinnamylloxy)-1-(4-nitrophenyl)-1H-tetrazole (2f).

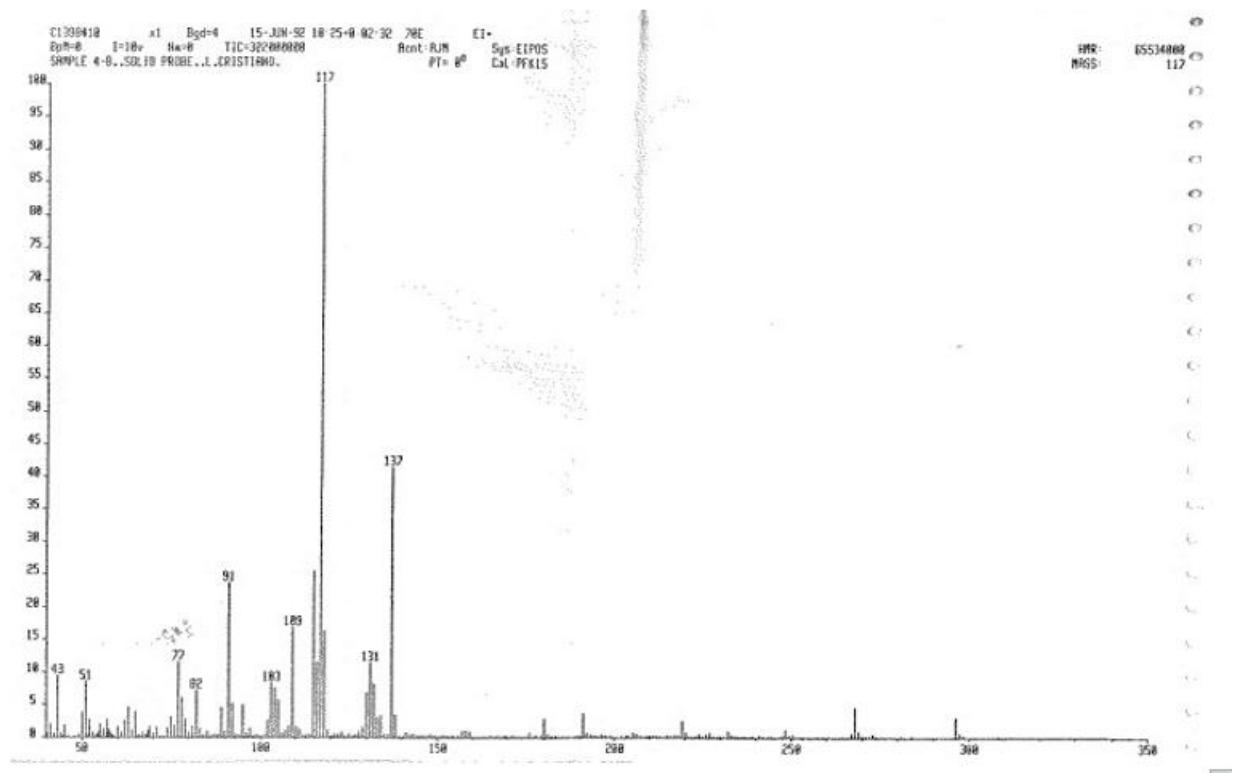


Figure S7. - EI-MS spectrum of 5-(cinnamylloxy)-1-(4-fluorophenyl)-1H-tetrazole (2g).

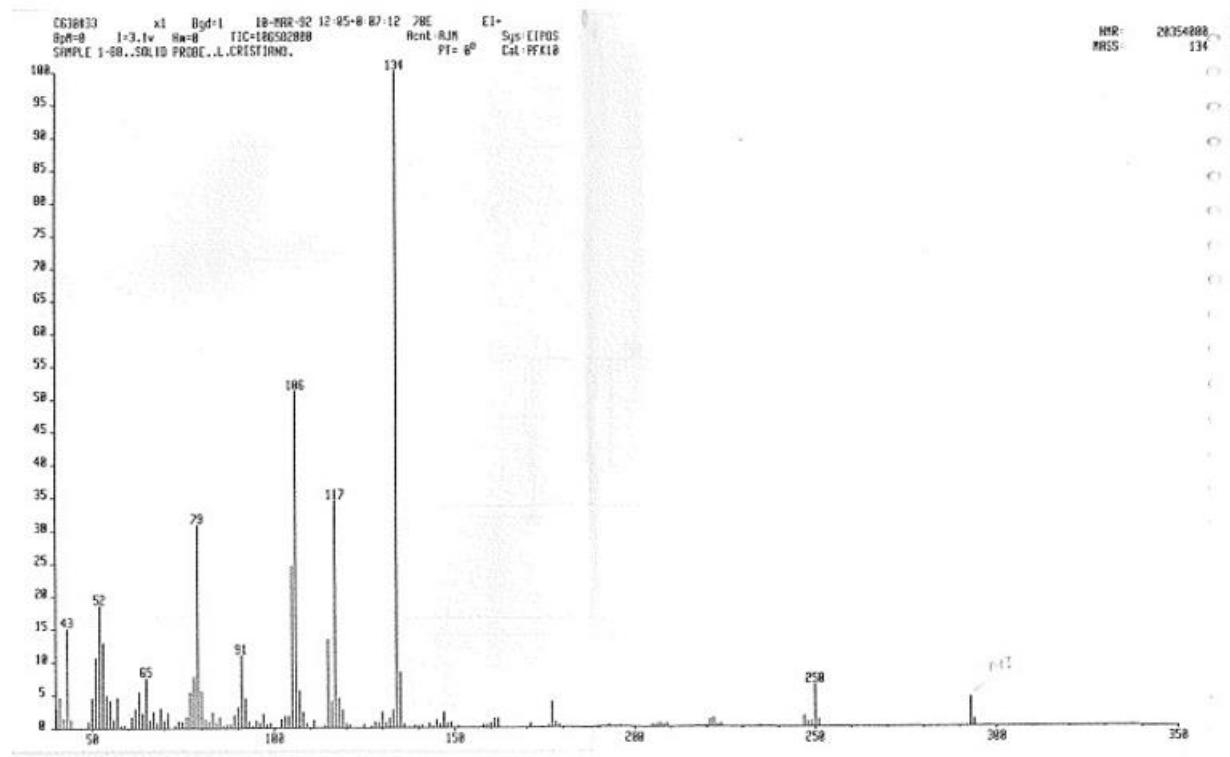


Figure S8. - EI-MS spectrum of 4-(5-(cinnamylloxy)-1H-tetrazol-1-yl)aniline (2h).

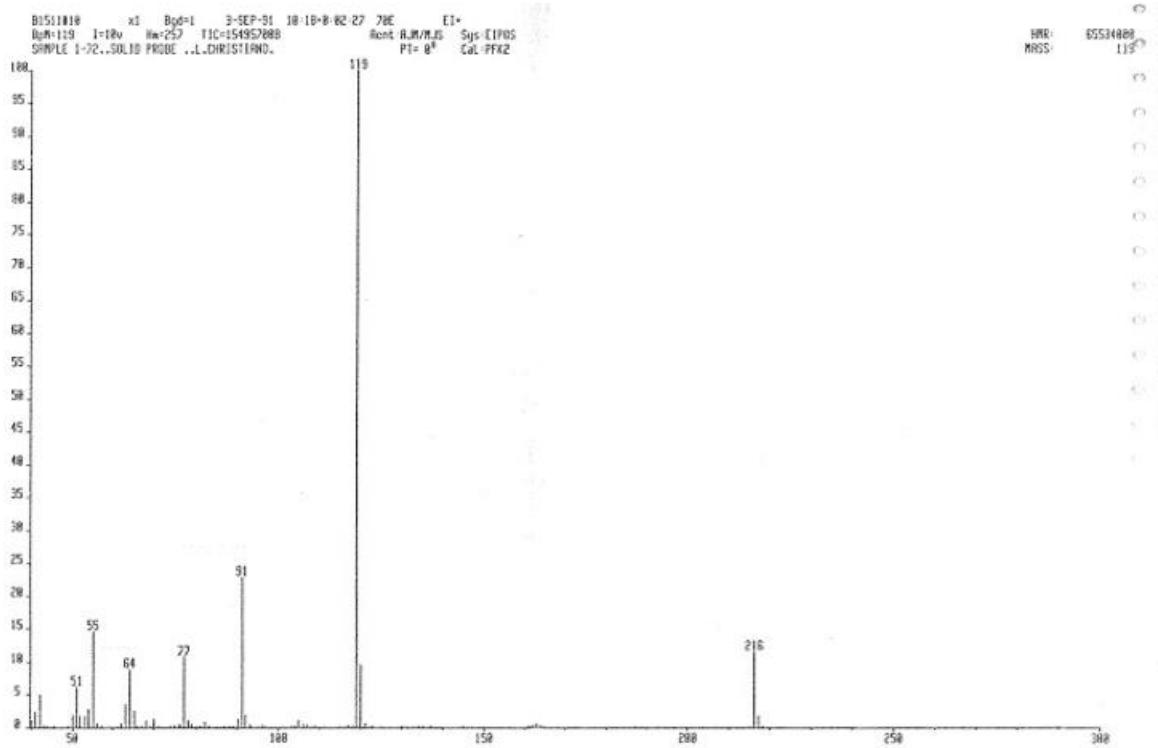


Figure S9. - EI-MS spectrum of 1-(but-3-en-2-yl)-4-phenyl-1H-tetrazol-5(4H)-one (3b).

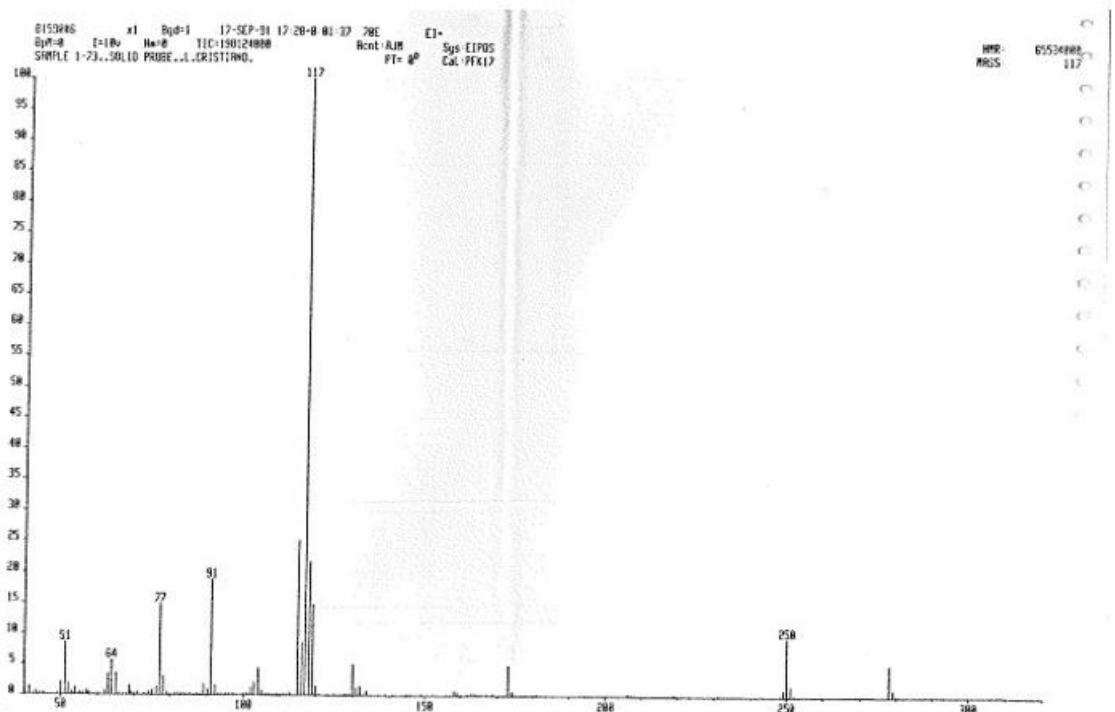


Figure S10. - EI-MS spectrum of 1-phenyl-4-(1-phenylallyl)-1H-tetrazol-5(4H)-one (3c).

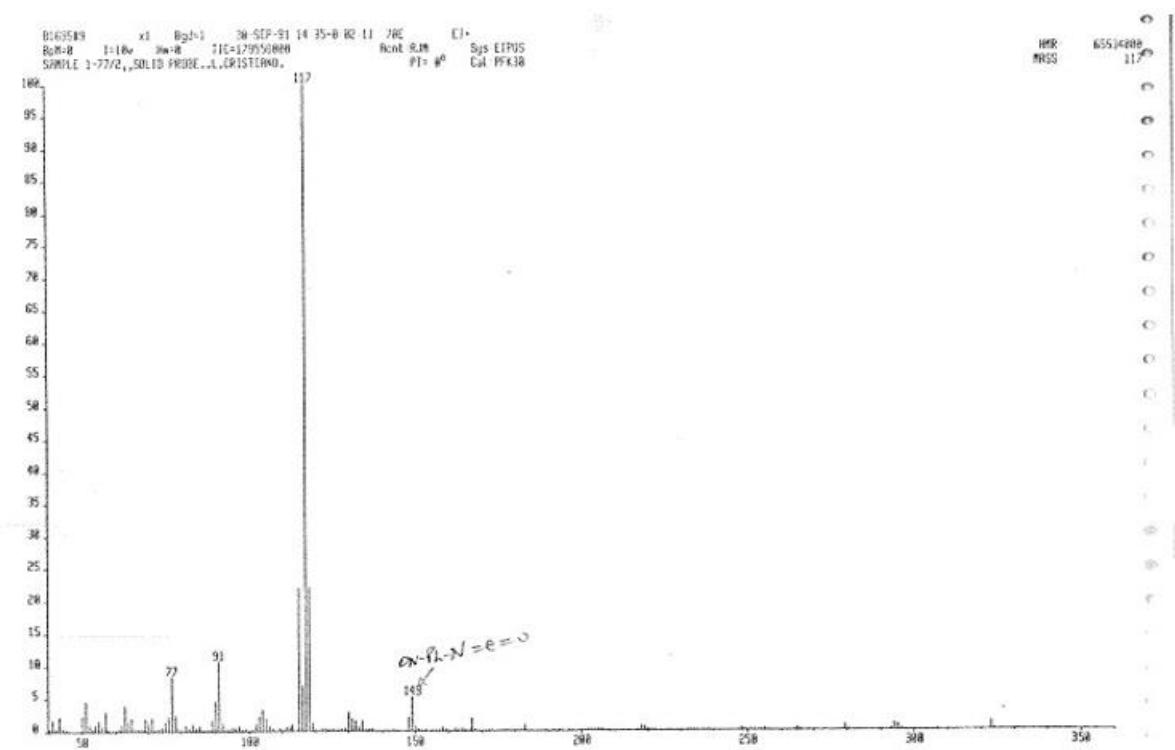


Figure S11. - EI-MS spectrum of 1-(4-nitrophenyl)-4-(1-phenylallyl)-1H-tetrazol-5(4H)-one (3f).

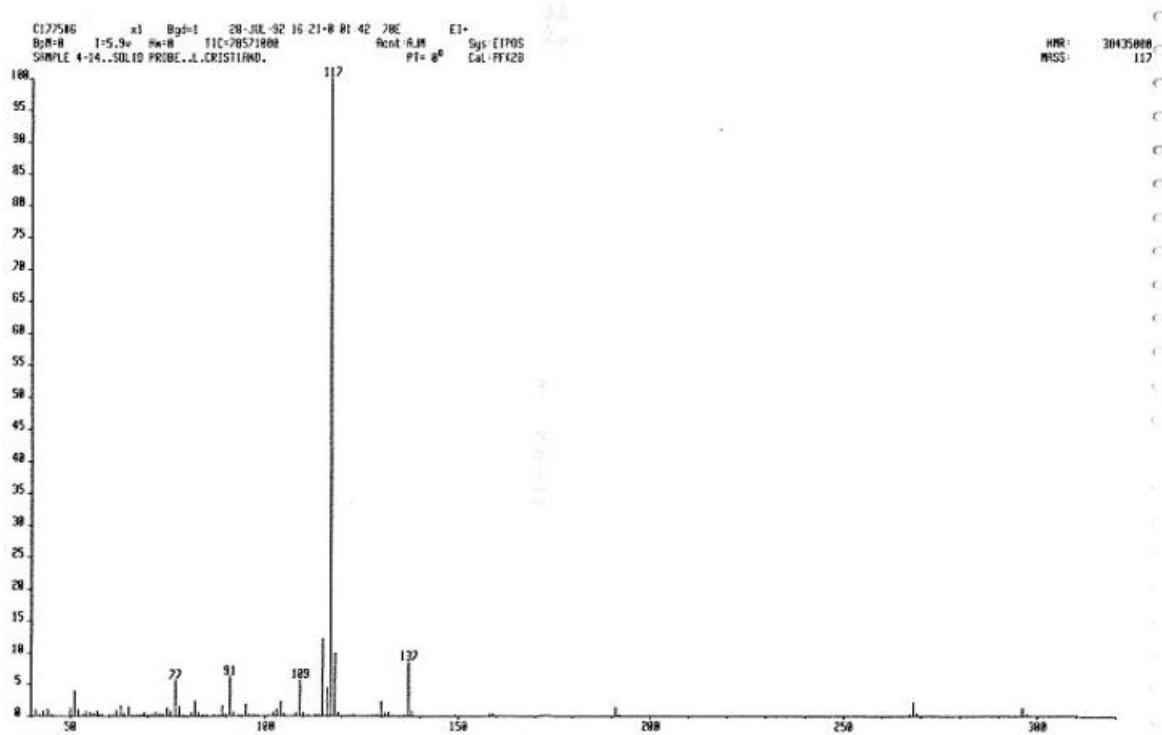


Figure S12. - EI-MS spectrum of 1-(4-fluorophenyl)-4-(1-phenylallyl)-1H-tetrazol-5(4H)-one (3g).

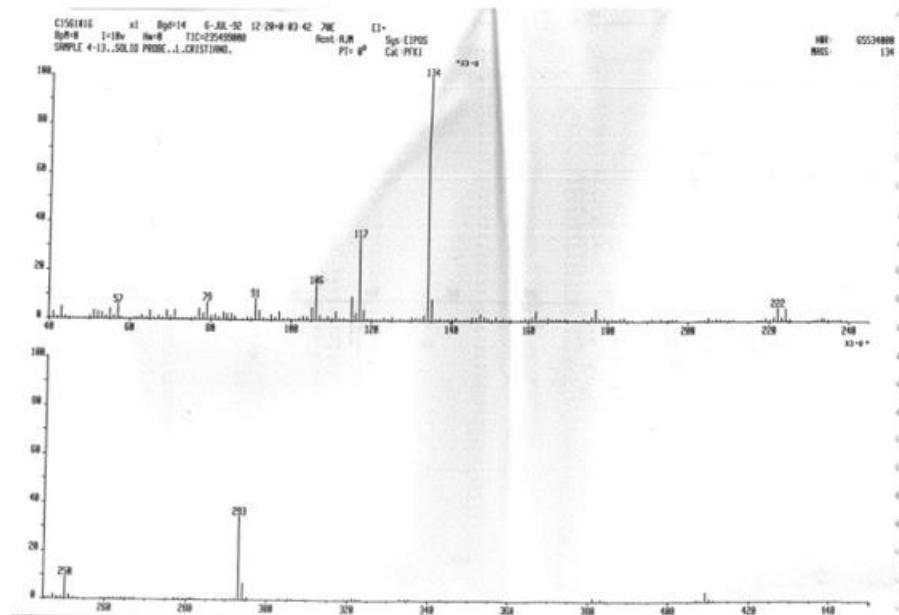


Figure S13. - EI-MS spectrum of 1-(4-aminophenyl)-4-(1-phenylallyl)-1H-tetrazol-5(4H)-one (3h) (amplified 3x).