

Supplementary Materials: Ethyl 5-methyl-7-(4-morpholinophenyl)-4,7- dihydrotetrazolo[1,5-a]pyrimidine-6- carboxylate

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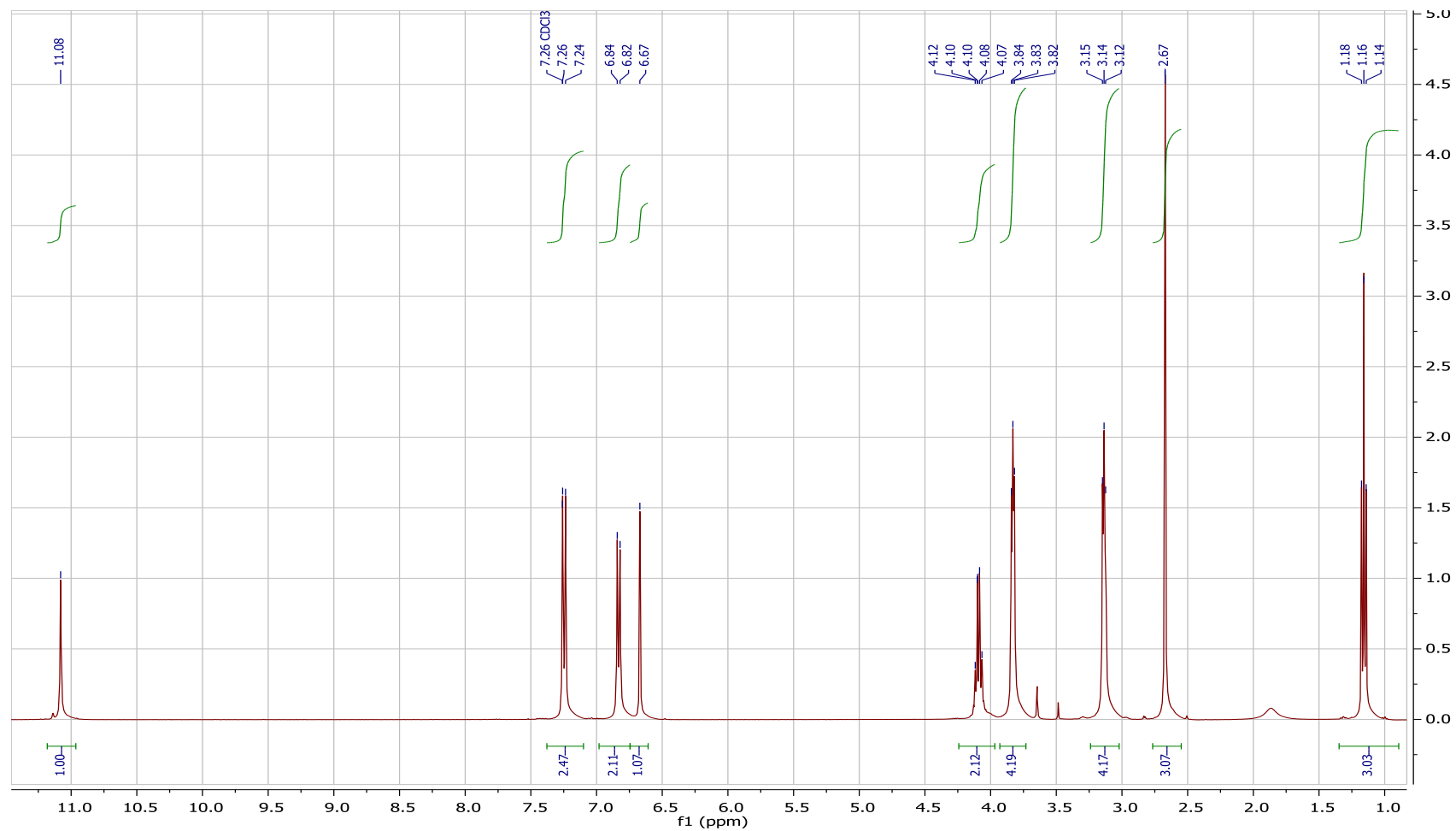


Figure S1. ¹H-NMR spectra of Ethyl 5-methyl-7-(4-morpholinophenyl)-4,7-dihydrotetrazolo[1,5-a]pyrimidine-6-carboxylate.

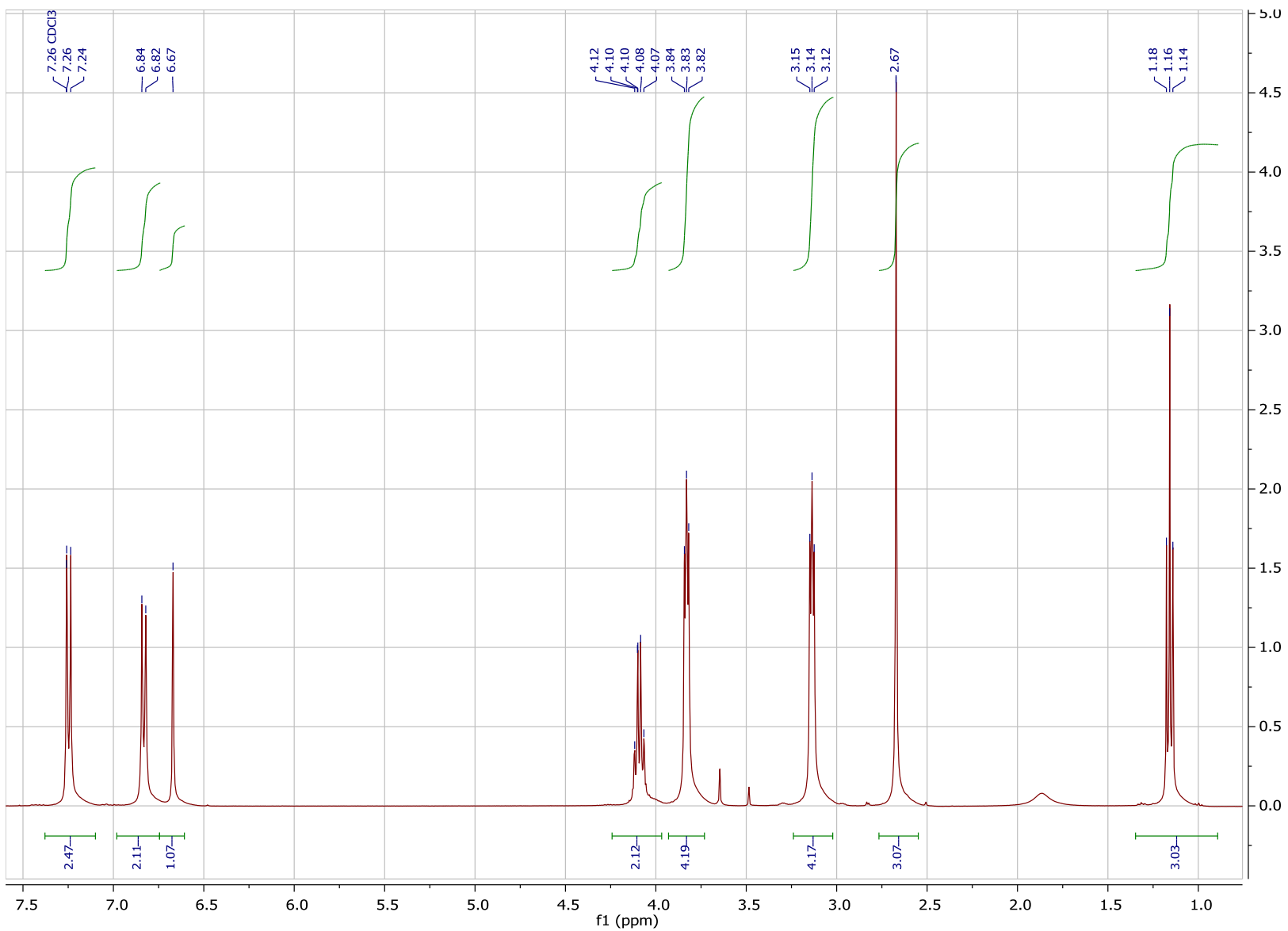


Figure S2. ¹H-NMR spectra of Ethyl 5-methyl-7-(4-morpholinophenyl)-4,7-dihydro-1,5-benzodiazepine-6-carboxylate.

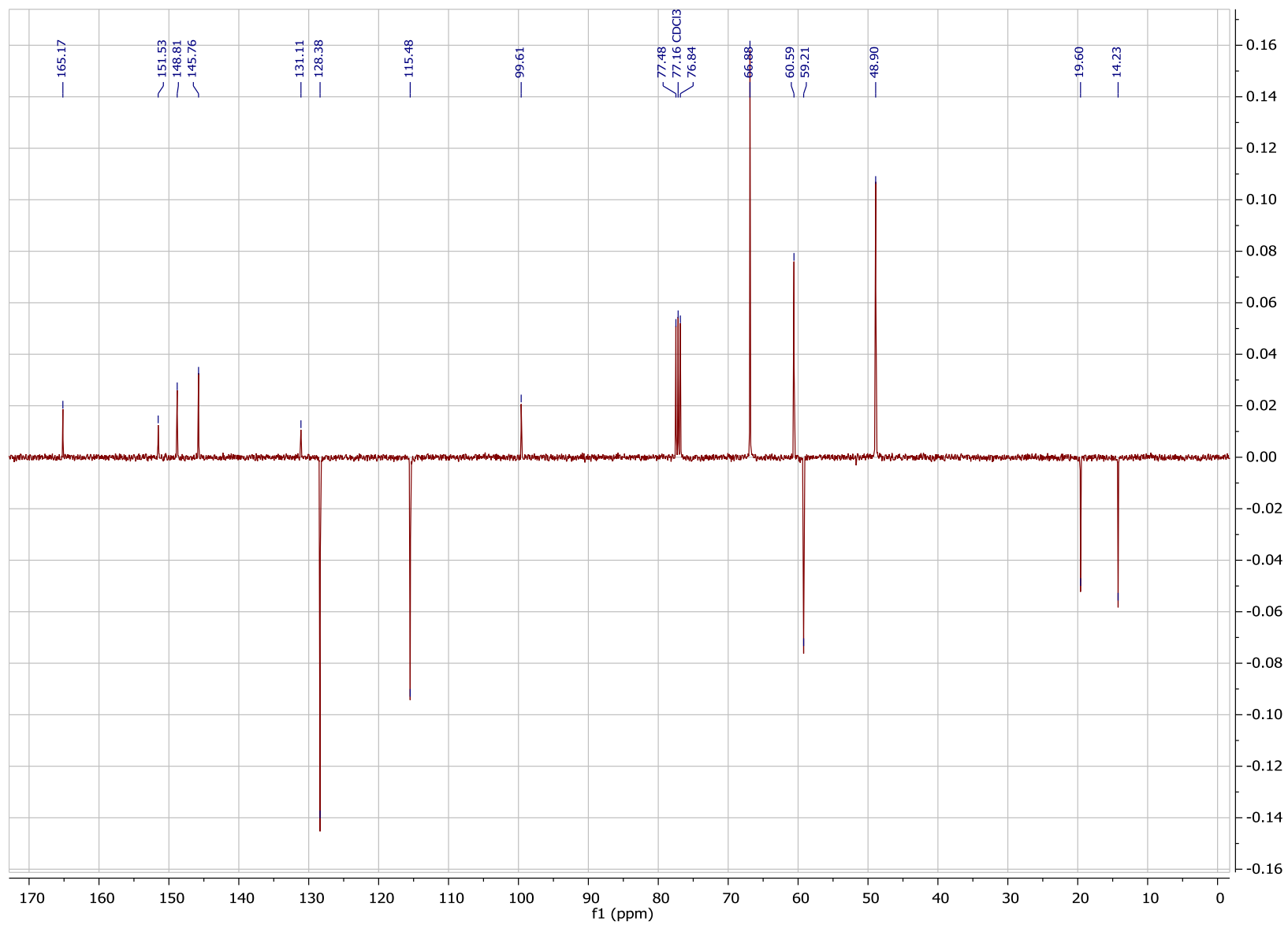


Figure S3. ^{13}C -NMR spectra of Ethyl 5-methyl-7-(4-morpholinophenyl)-4,7-dihydro-1H-tetrazolo[1,5-a]pyrimidine-6-carboxylate.

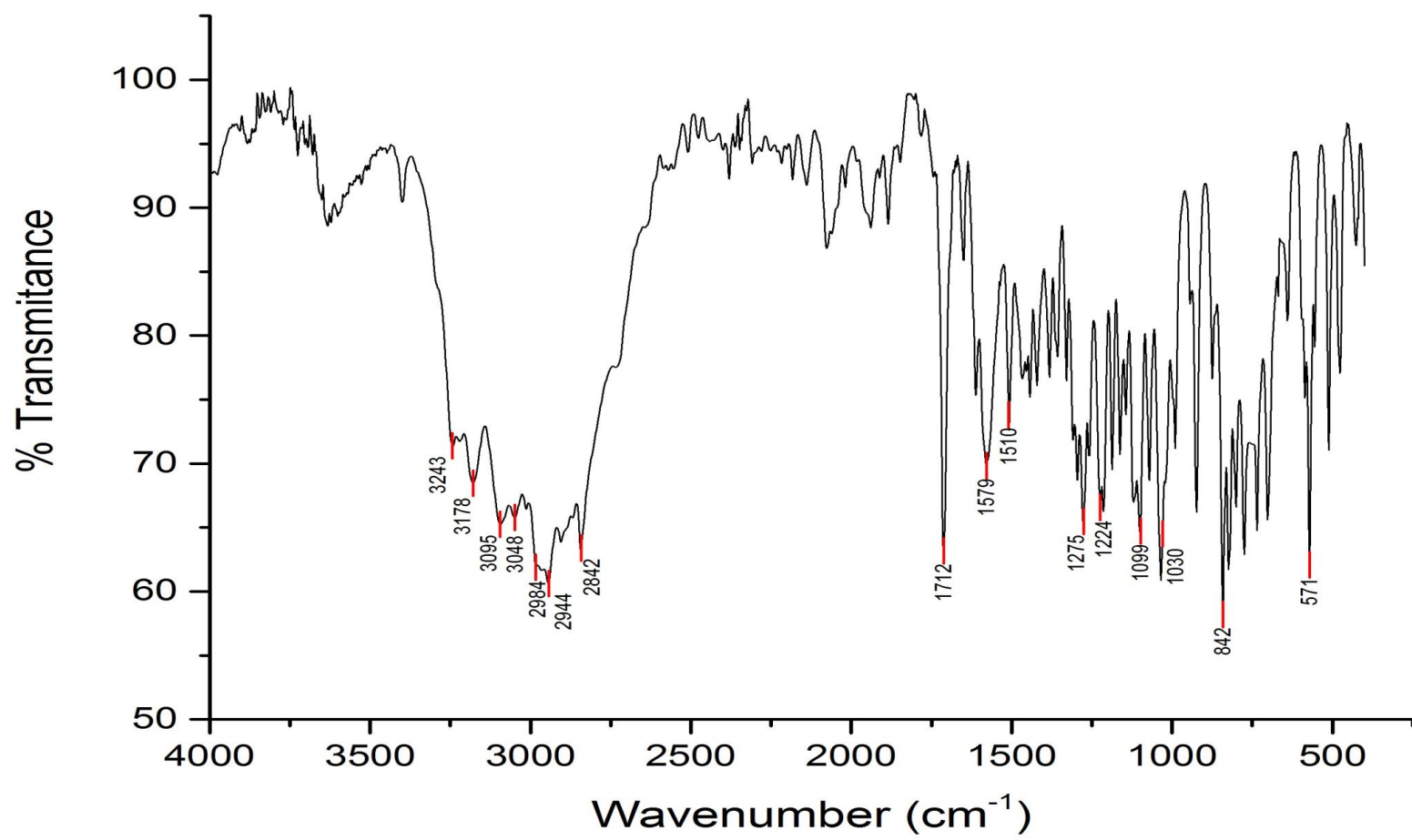


Figure S4. FTIR spectra of Ethyl 5-methyl-7-(4-morpholinophenyl)-4,7-dihydro-1H-tetrazolo[1,5-a]pyrimidine-6-carboxylate.

Elemental Composition Report

Page 1

Single Mass Analysis

Tolerance = 50.0 PPM / DBE: min = -1.5, max = 50.0

Element prediction: Off

Number of isotope peaks used for i-FIT = 5

Monoisotopic Mass, Even Electron Ions

1042 formula(e) evaluated with 43 results within limits (up to 1 closest results for each mass)

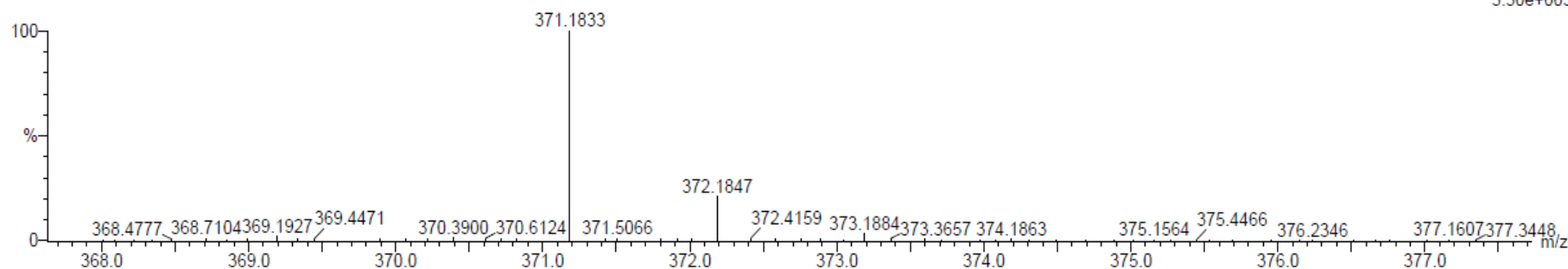
Elements Used:

C: 0-1000 H: 0-1000 N: 0-500 O: 0-500

standard

Unair_Salma_MTDPT2-pos 18 (0.563)

TOF MS ES+
3.30e+003



Minimum: -1.5
Maximum: 30.0 50.0 50.0

Mass	Calc. Mass	mDa	PPM	DBE	i-FIT	i-FIT (Norm)	Formula
371.1833	371.1832	0.1	0.3	10.5	123.6	0.0	C18 H23 N6 O3

Figure S5. HRESI-MS spectra of Ethyl 5-methyl-7-(4-morpholinophenyl)-4,7-dihydro-tetrazolo[1,5-a]pyrimidine-6-carboxylate.

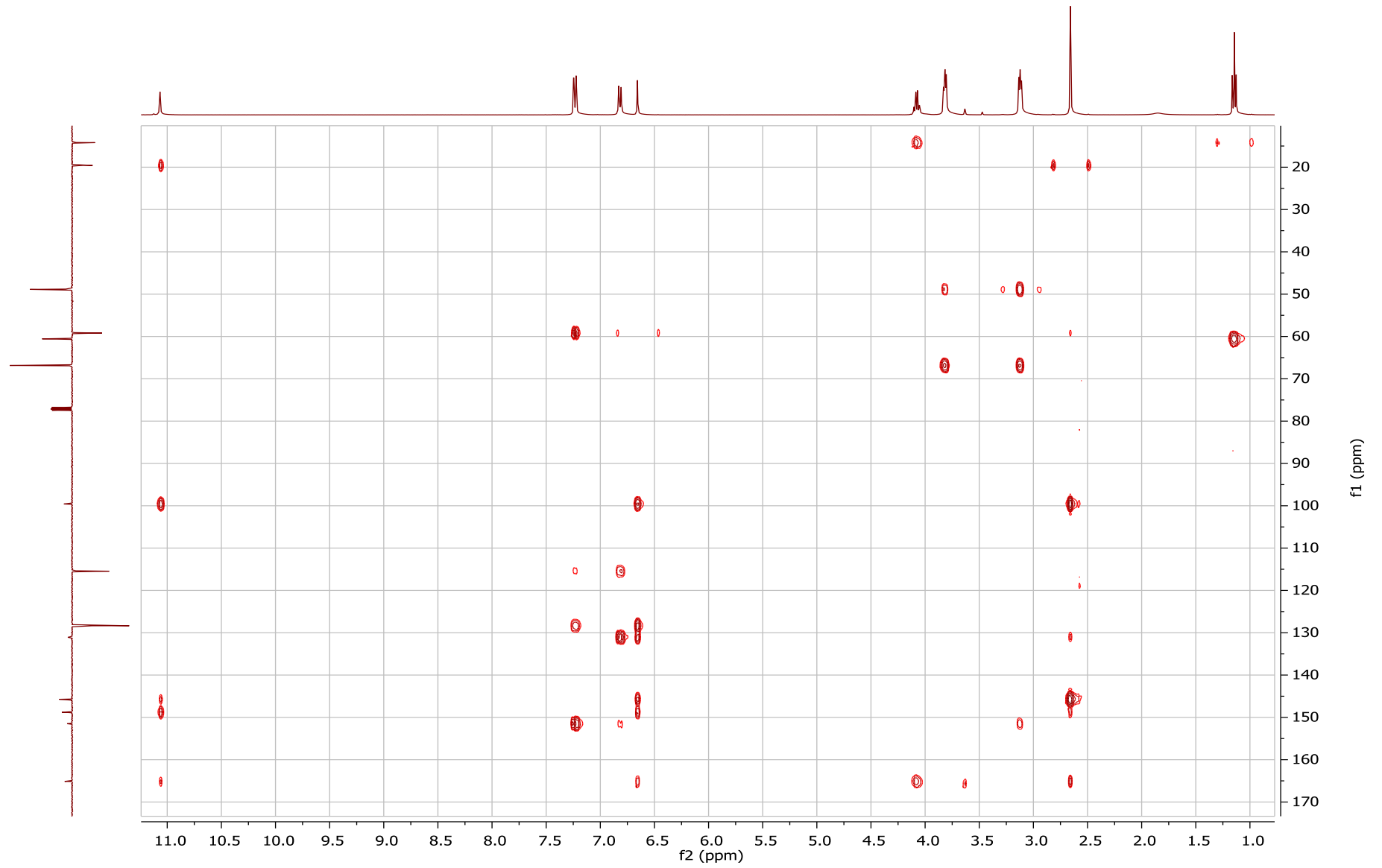


Figure S6. HMBC spectra of Ethyl 5-methyl-7-(4-morpholinophenyl)-4,7-dihydro-1,5-benzodiazepin-6-carboxylate.