

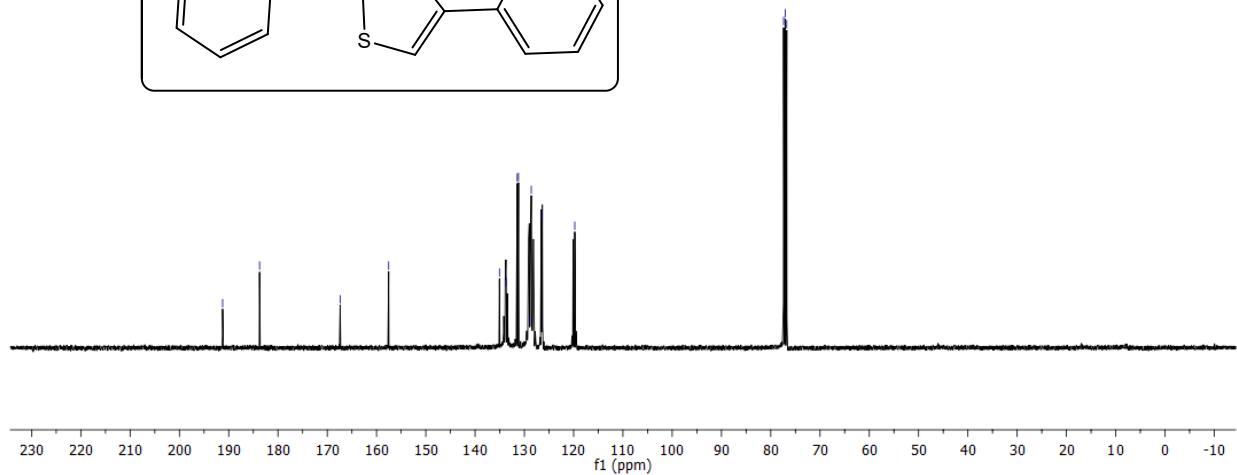
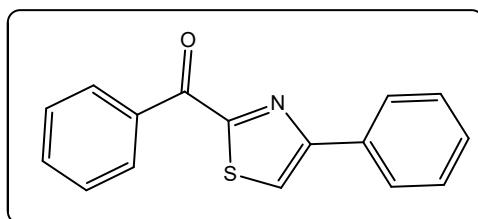
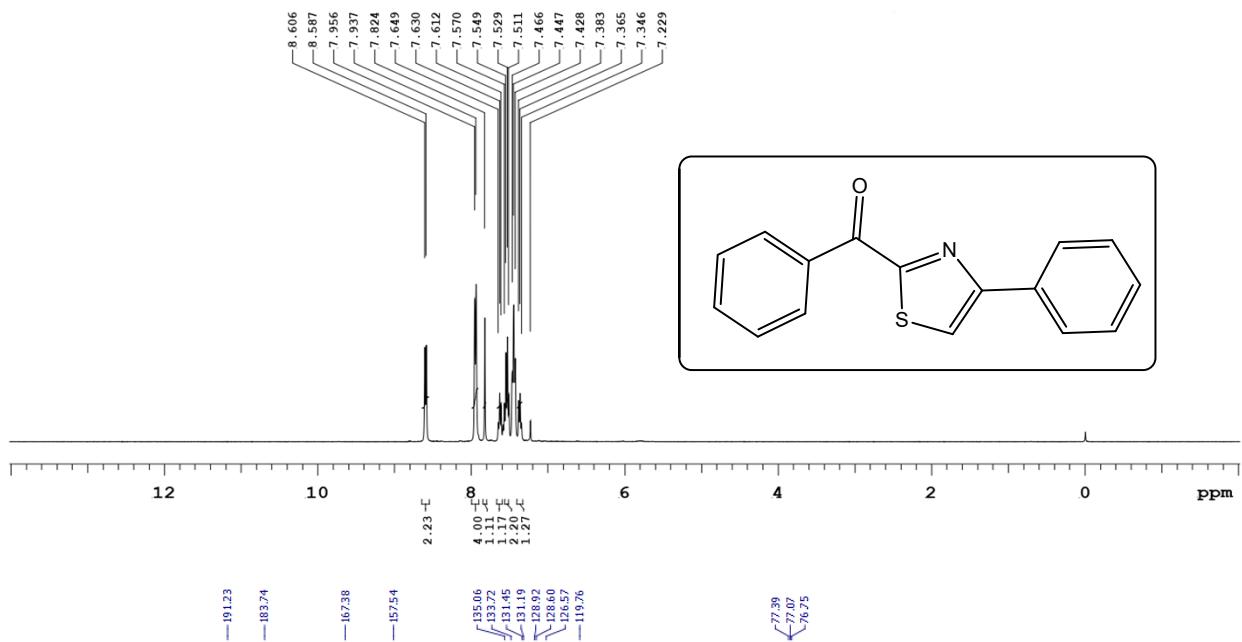
A panoramic view on synthetic applications of α -oxothioamides: A highly regioselective synthesis of 2-acyl-4-(het)arylthiazoles and thioethers

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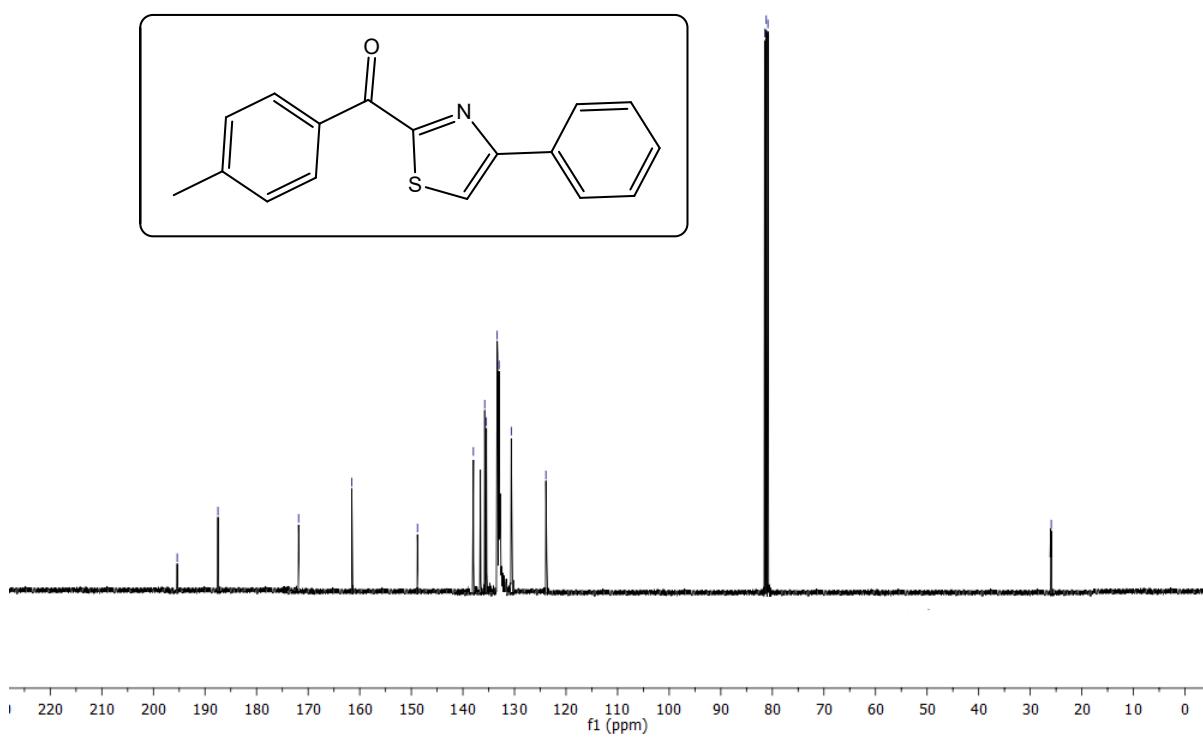
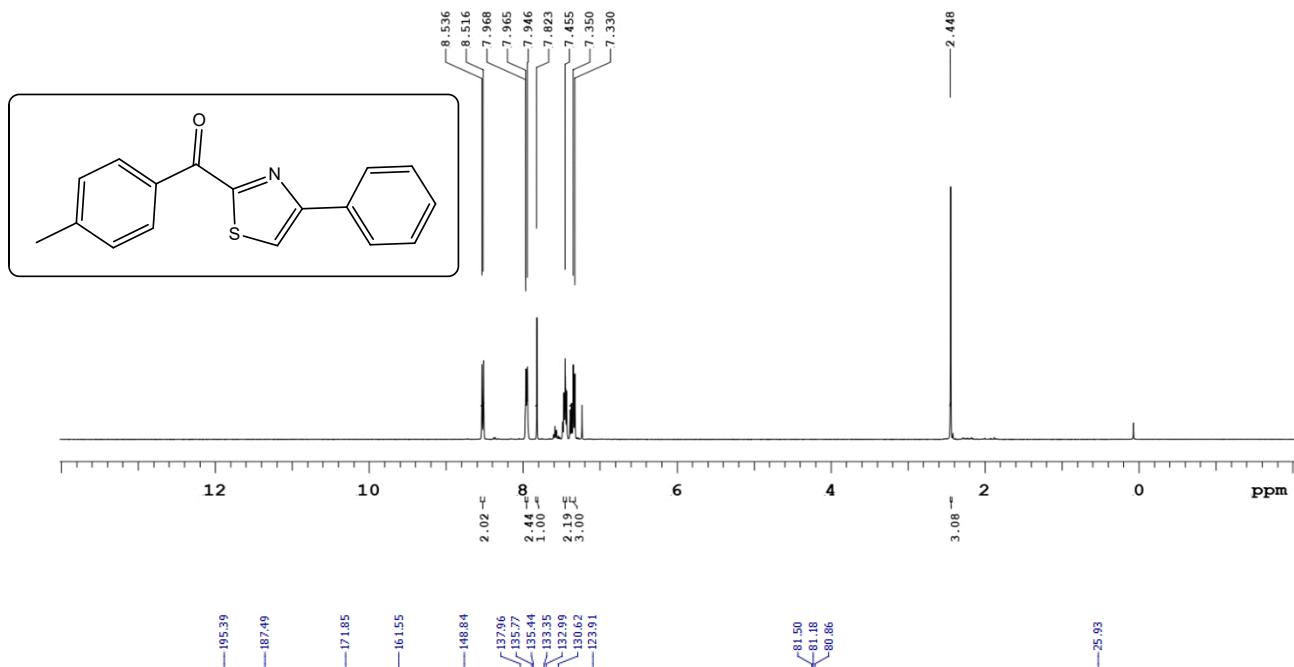
^a Department of Studies in Chemistry, University of Mysore, Manasagangotri, Mysuru - 570 006, Karnataka, India

^b Department of Studies in Organic Chemistry, University of Mysore, Manasagangotri, Mysuru - 570 006, Karnataka, India.

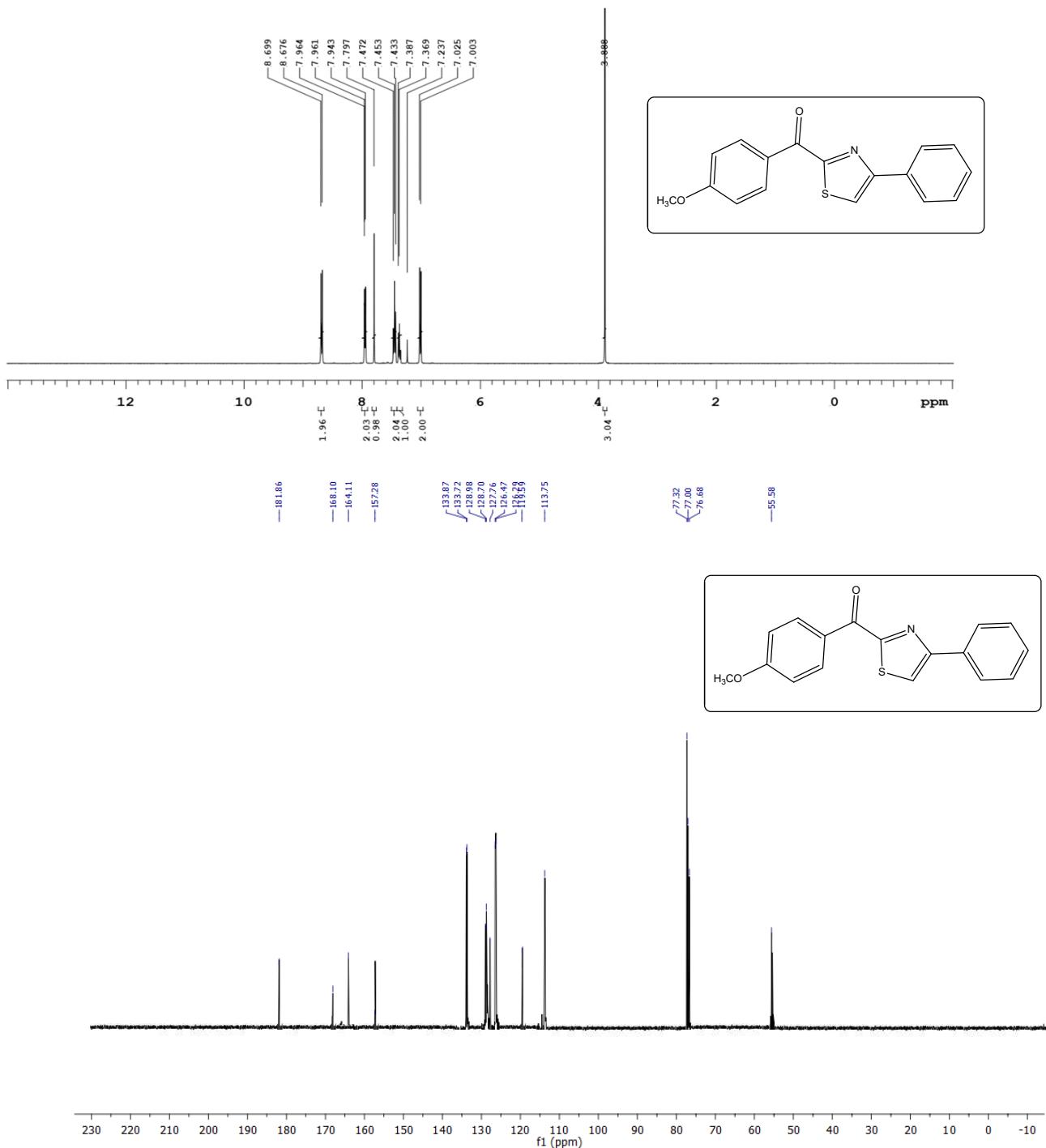
^c Vijnana Bhavana, University of Mysore, Manasagangotri, Mysuru - 570 006, Karnataka, India.



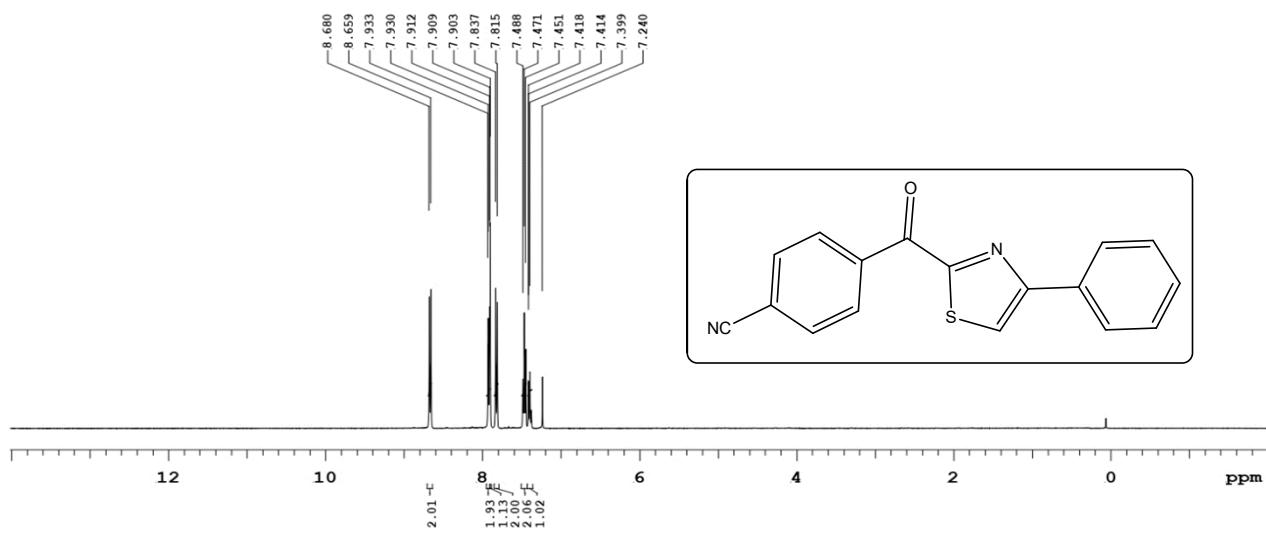
¹H & ¹³C NMR of Phenyl(4-phenylthiazol-2-yl)methanone (4a)



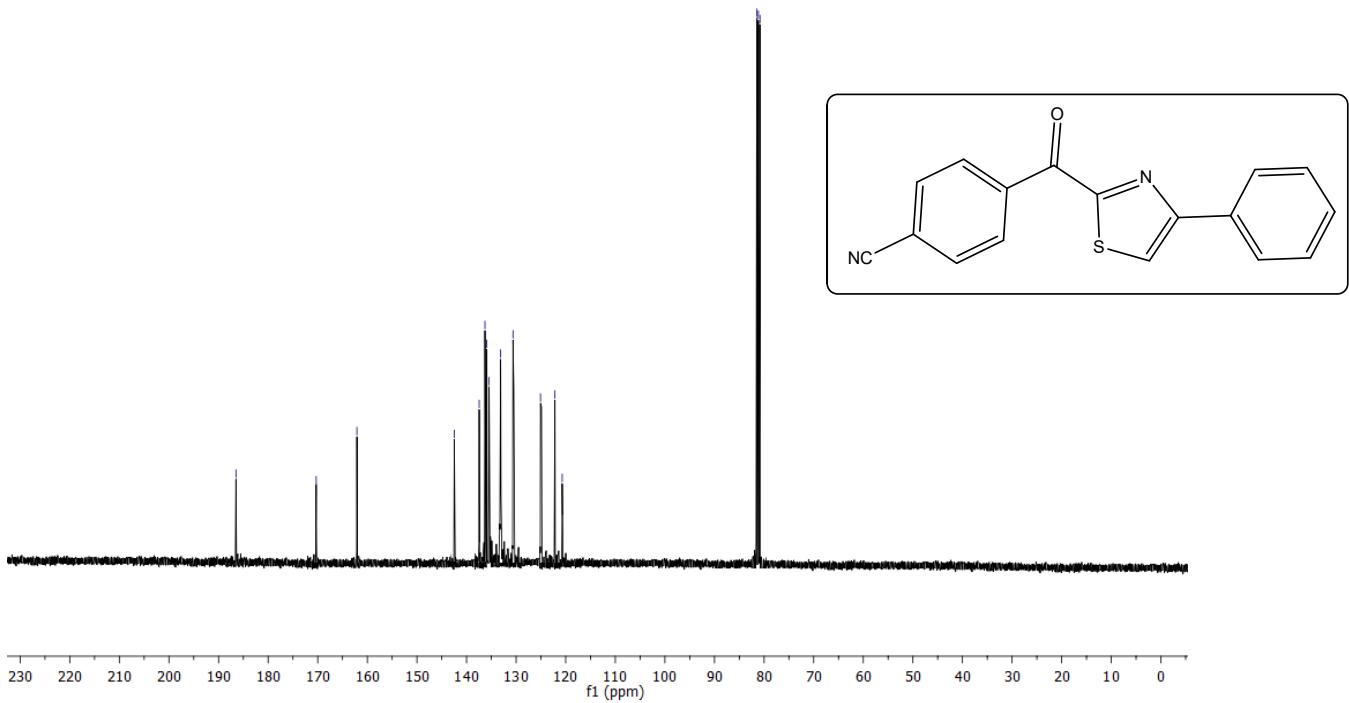
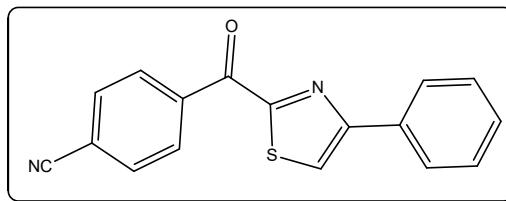
¹H & ¹³C NMR of (4-Phenylthiazol-2-yl)(p-tolyl)methanone (4b)



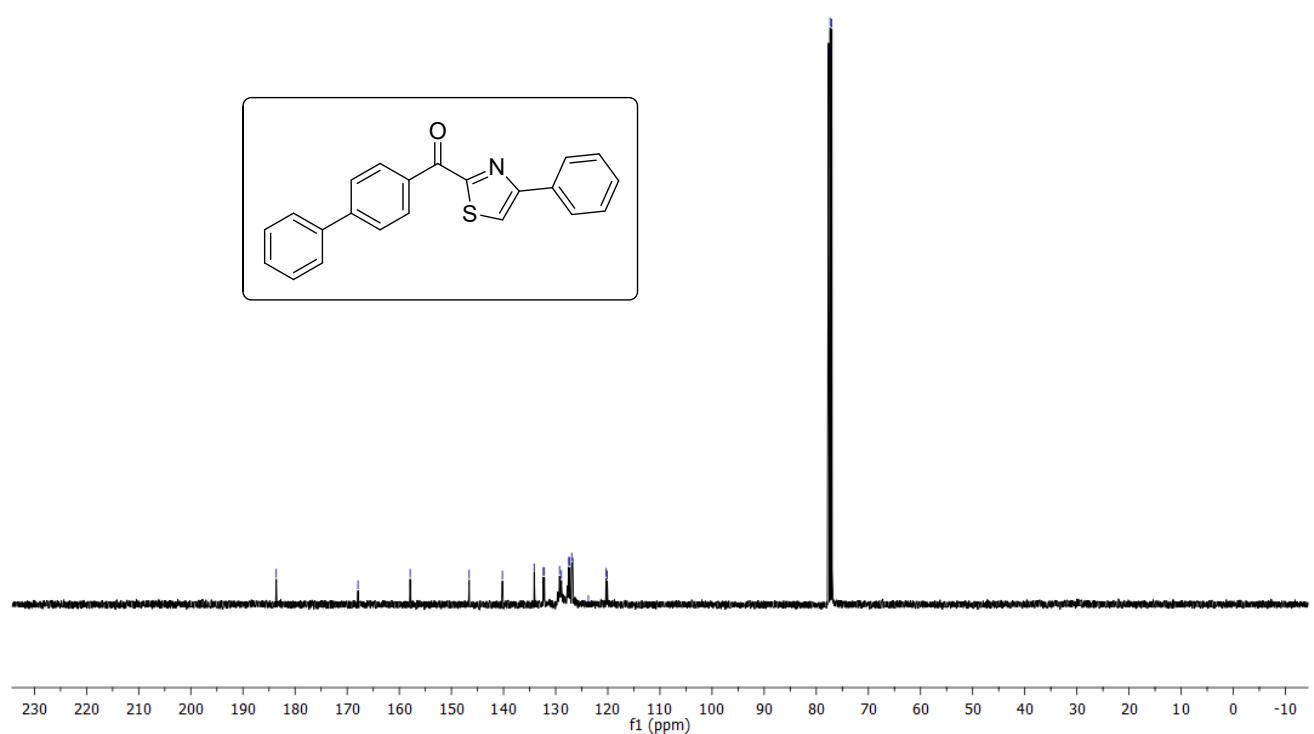
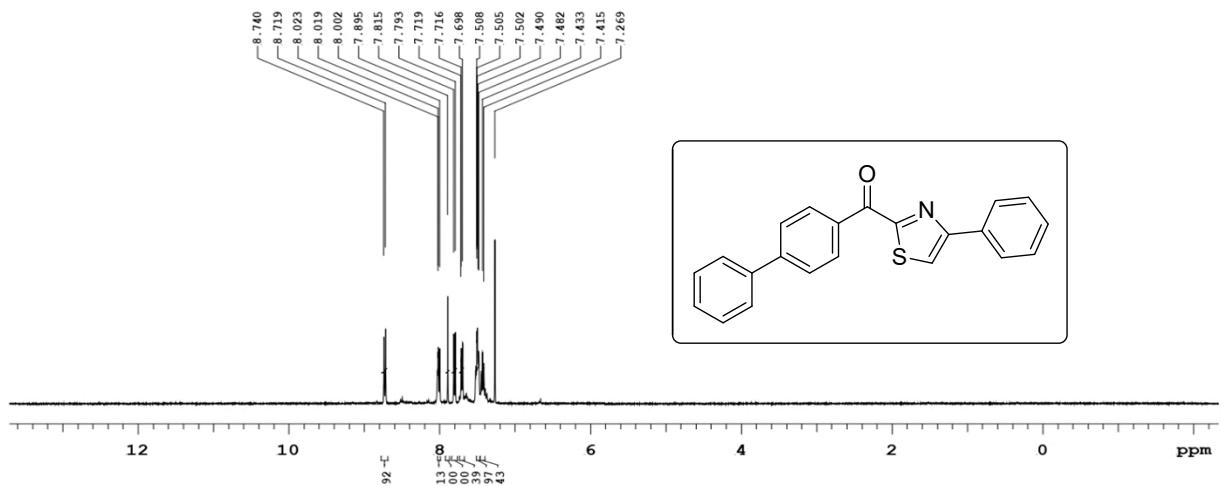
¹H & ¹³C NMR of (4-Methoxyphenyl)(4-phenylthiazol-2-yl)methanone (4c)



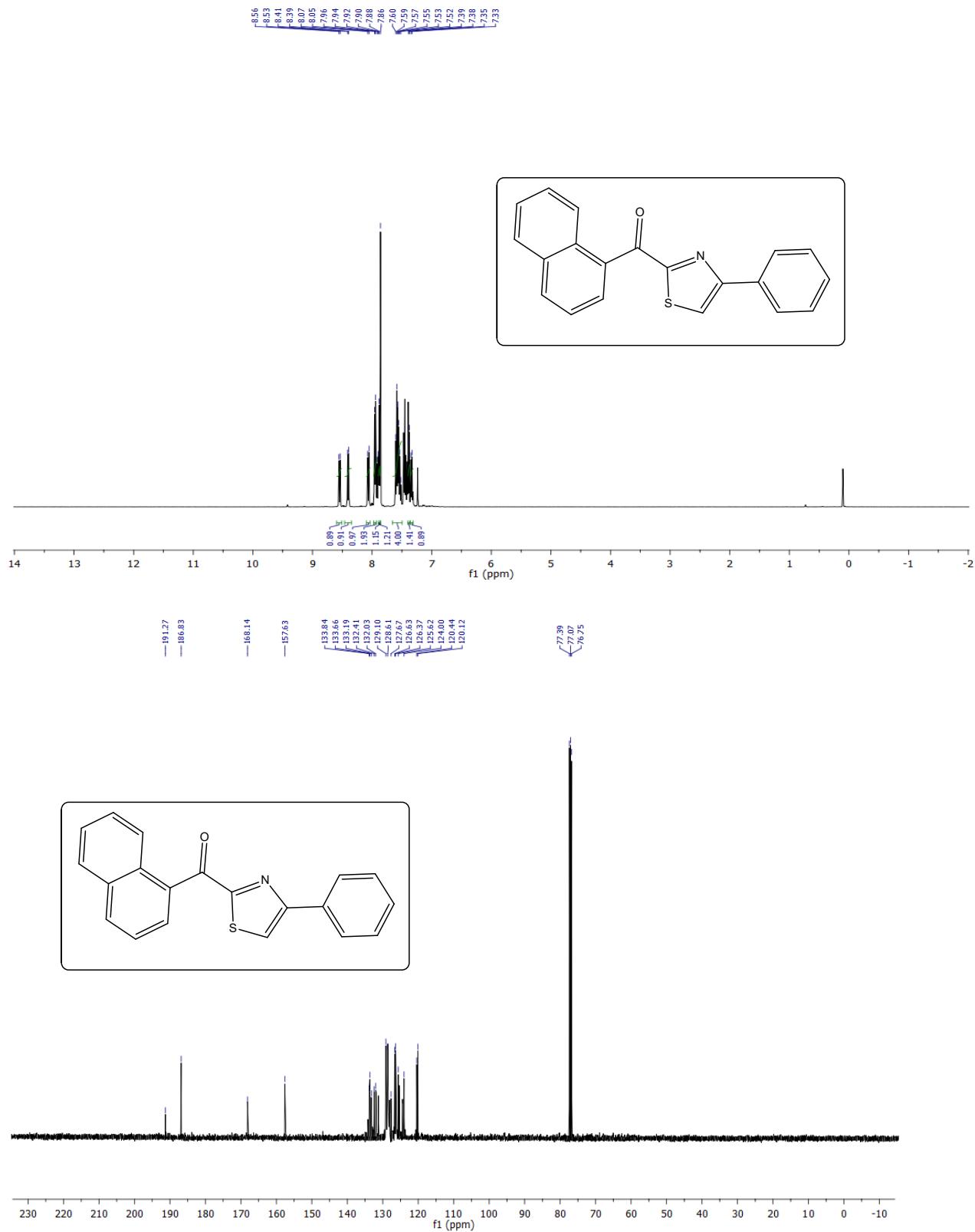
186.49
170.31
162.10
142.46
137.44
136.58
135.52
135.45
133.12
130.63
125.69
122.18
120.71



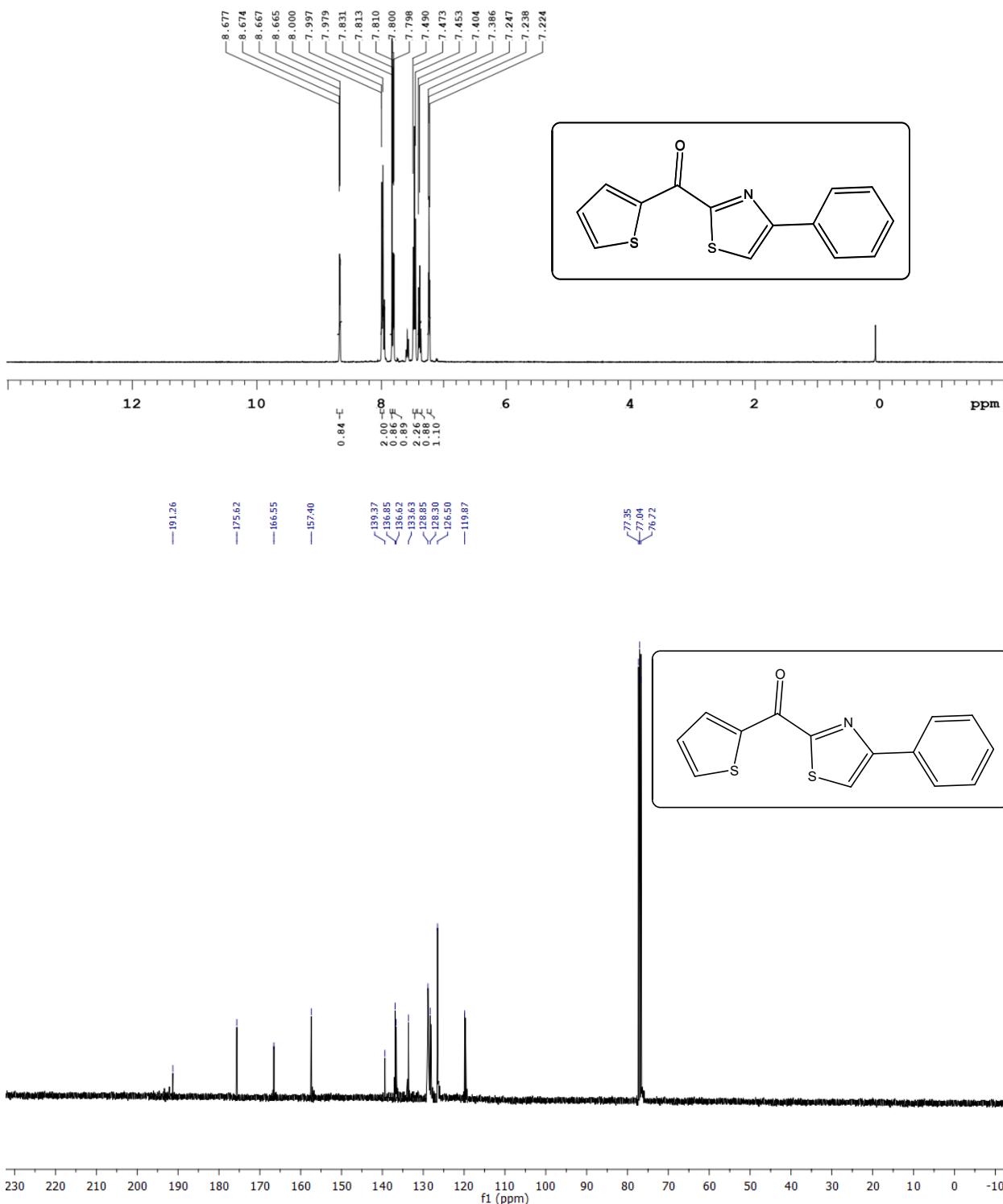
¹H & ¹³C NMR of 4-(4-Phenylthiazole-2-carbonyl)benzonitrile (4d)

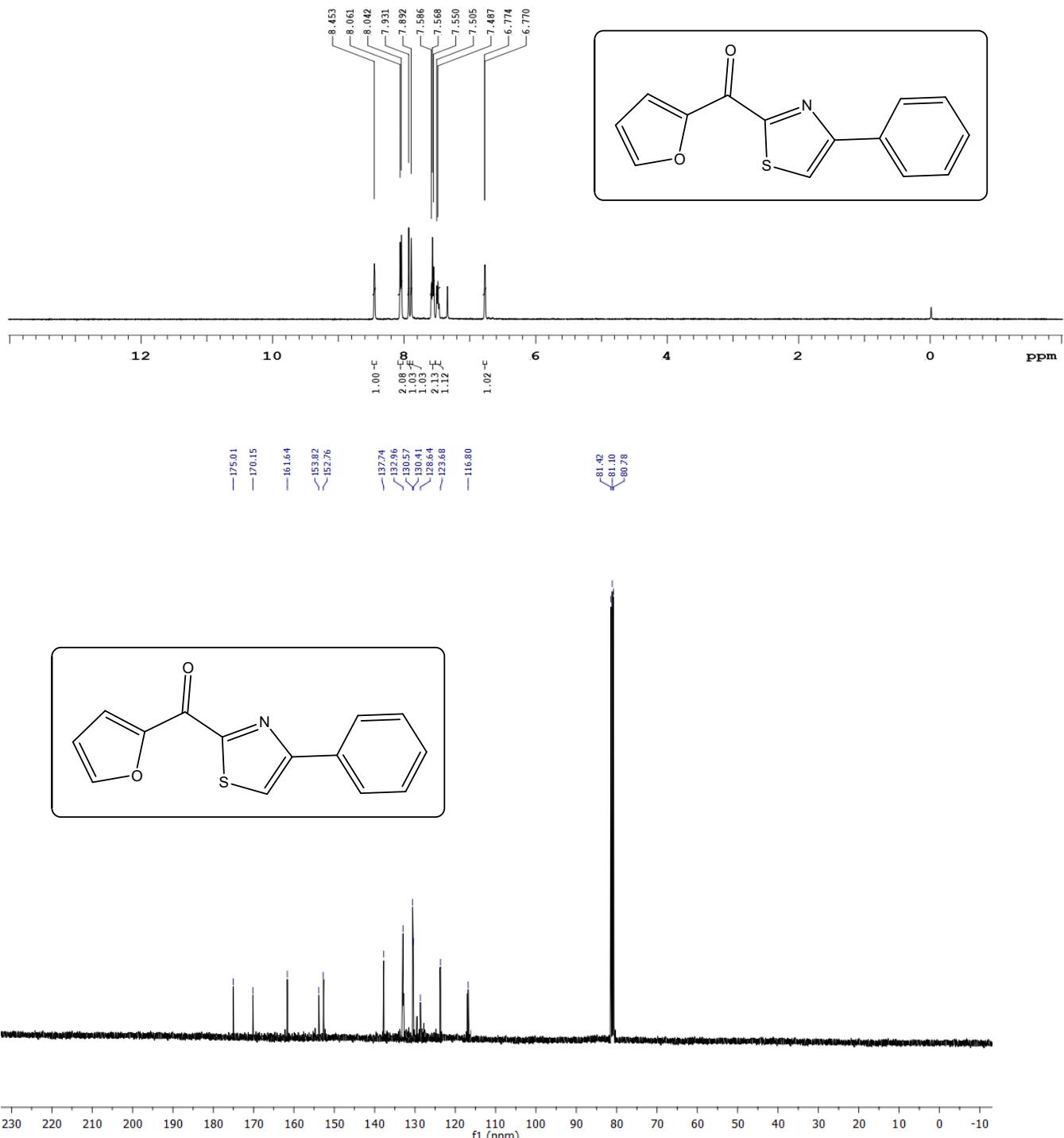


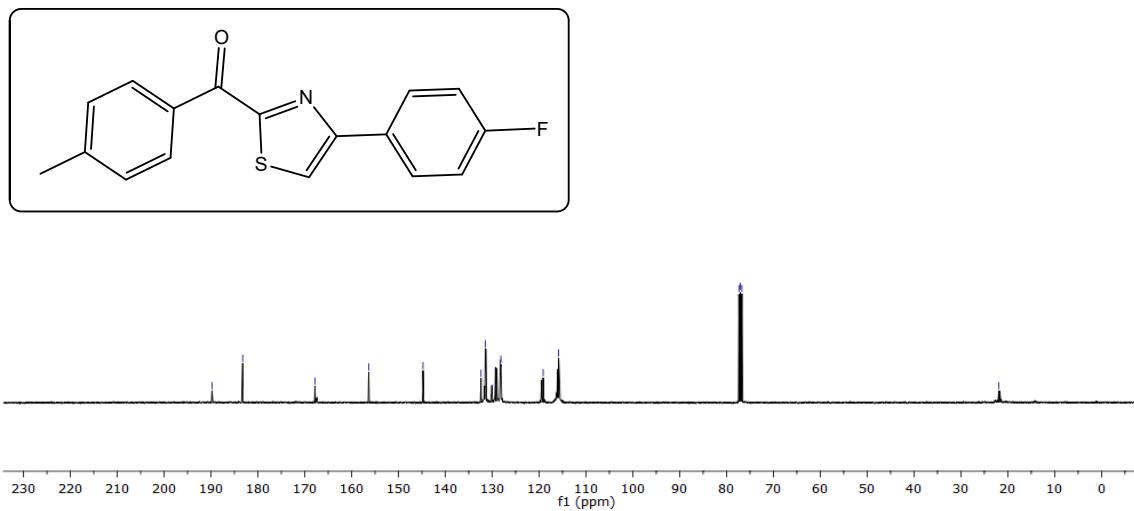
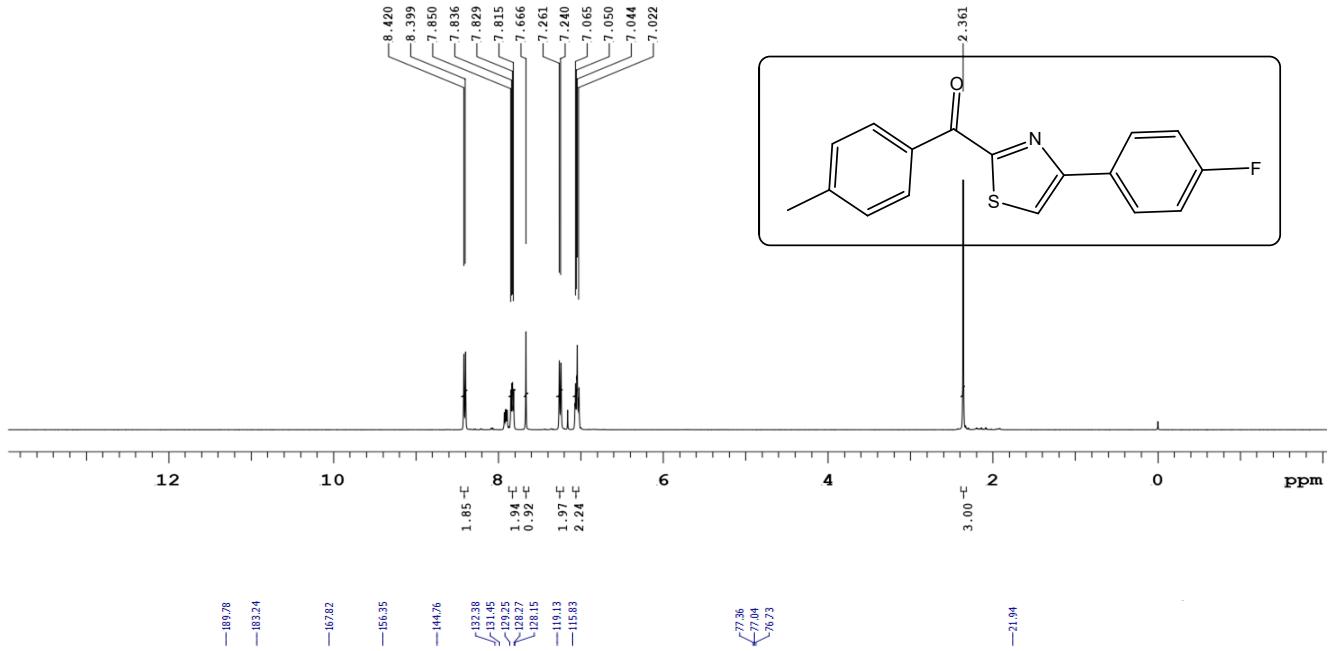
¹H & ¹³C NMR of [1,1'-biphenyl]-4-yl(4-phenylthiazol-2-yl)methanone (4e)



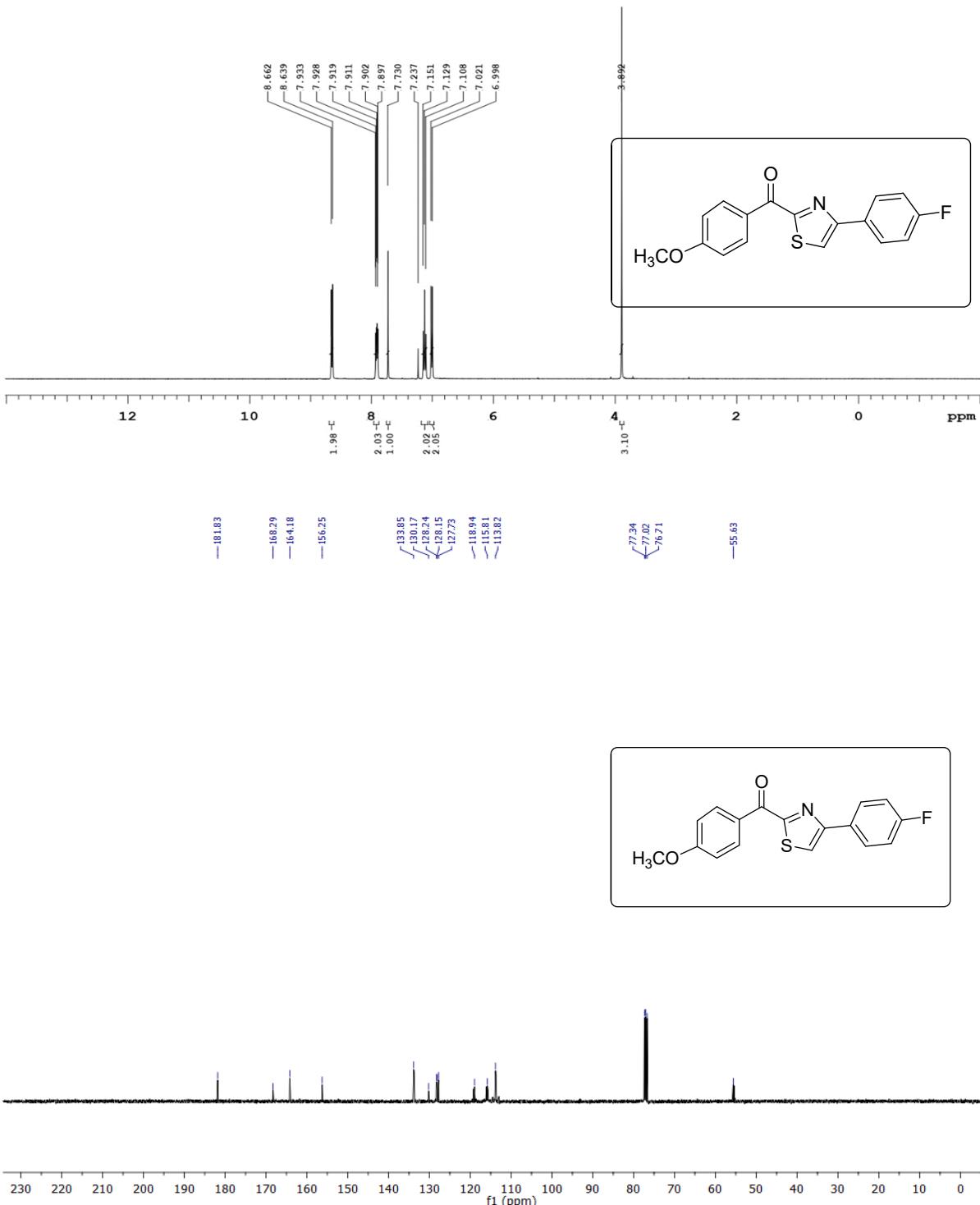
¹H & ¹³C NMR of Naphthalen-1-yl(4-phenylthiazol-2-yl)methanone (4f)



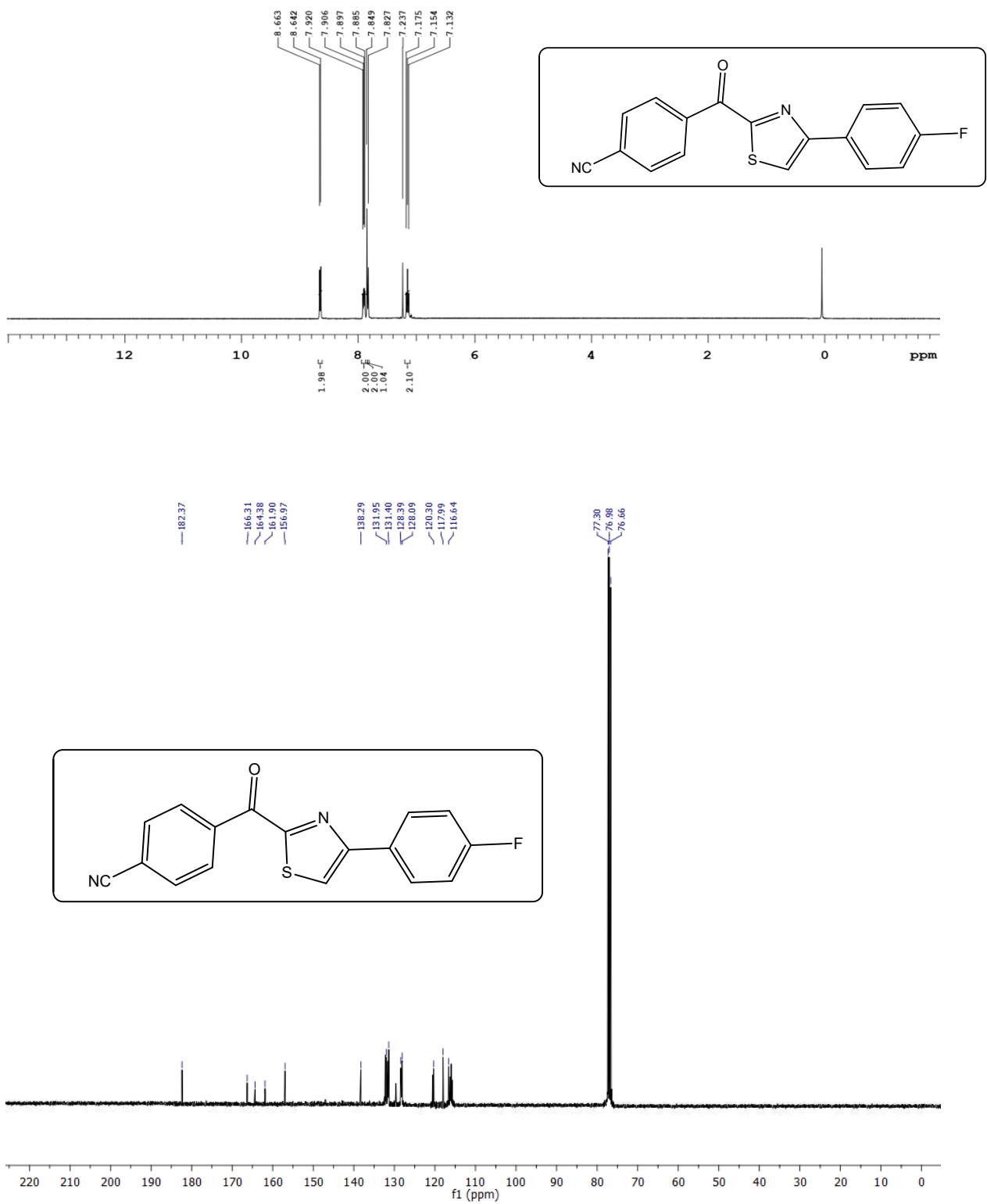




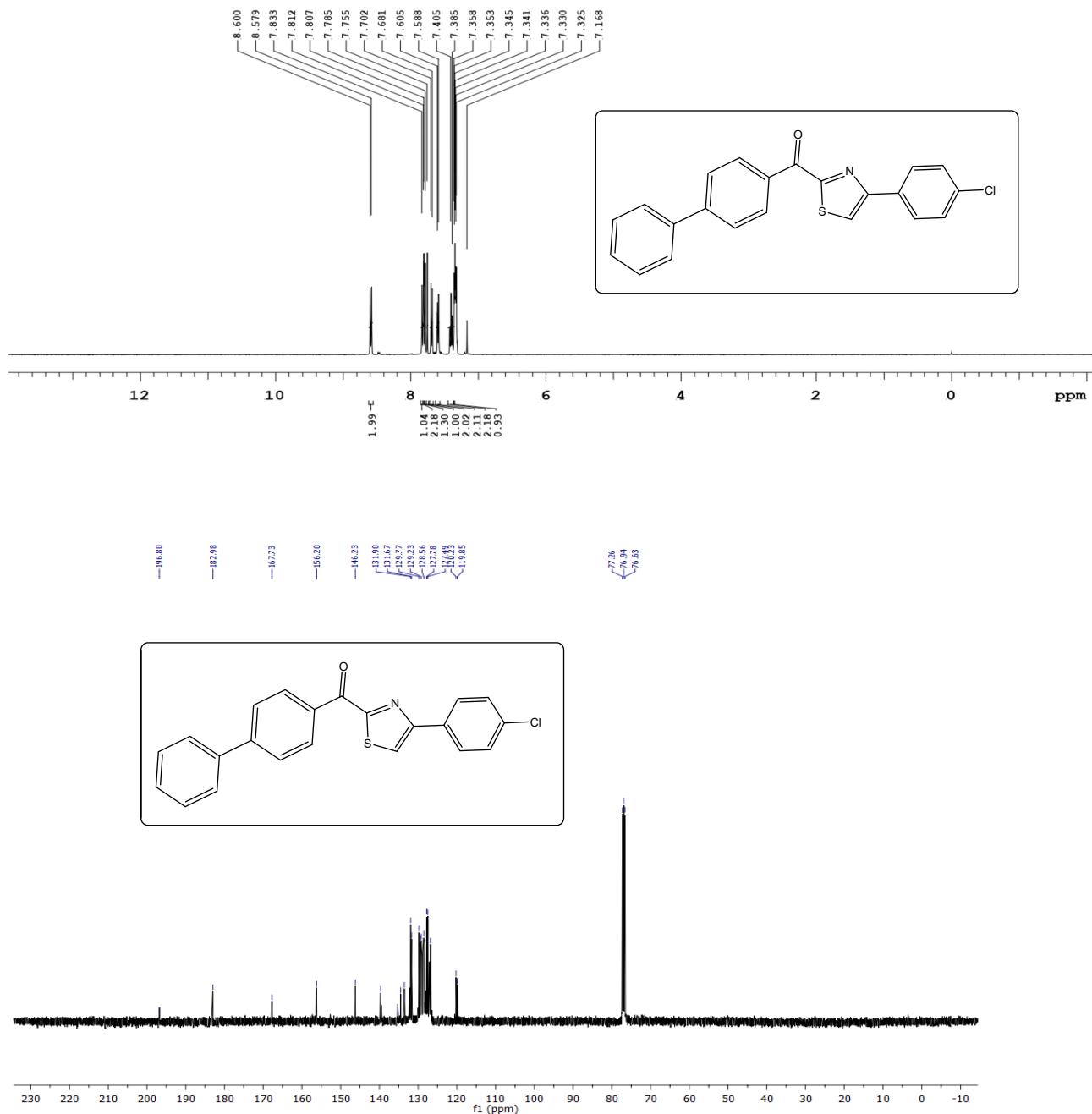
¹H & ¹³C NMR of (4-(4-Fluorophenyl)thiazol-2-yl)(p-tolyl)methanone (4i)



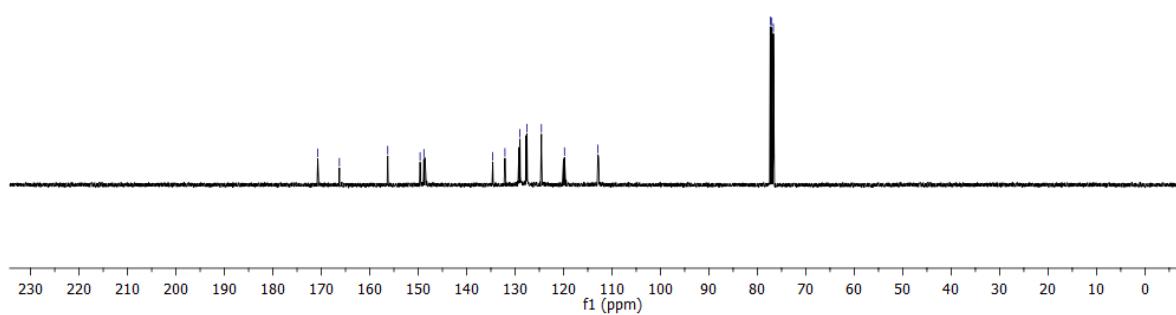
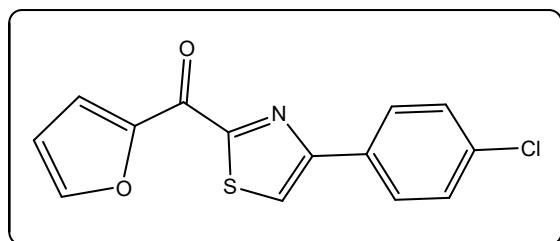
¹H & ¹³C NMR of (4-(4-Fluorophenyl)thiazol-2-yl)(4-methoxyphenyl)methanone (4j)



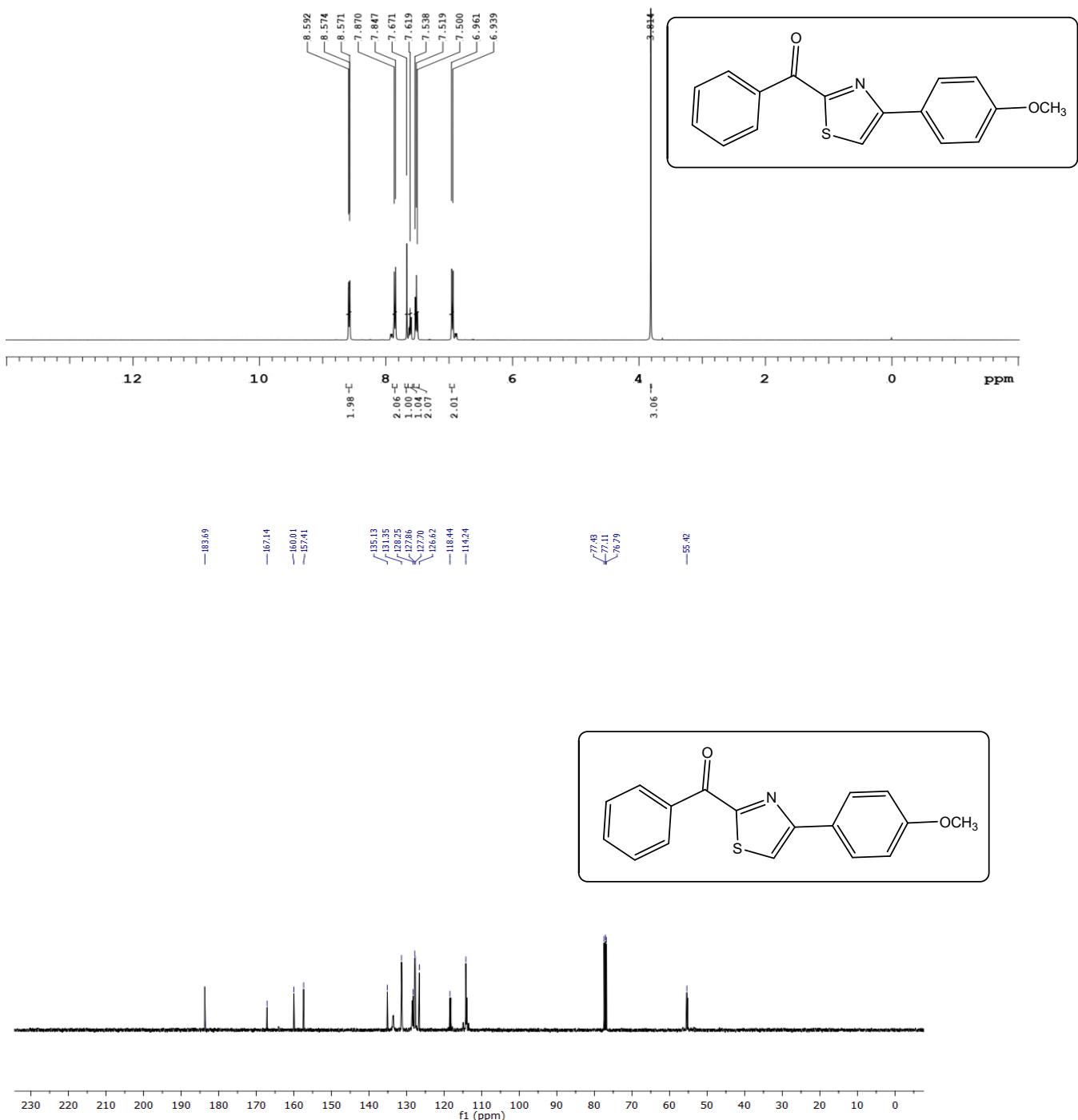
¹H & ¹³C NMR of 4-(4-(4-Fluorophenyl)thiazole-2-carbonyl)benzonitrile (4k)



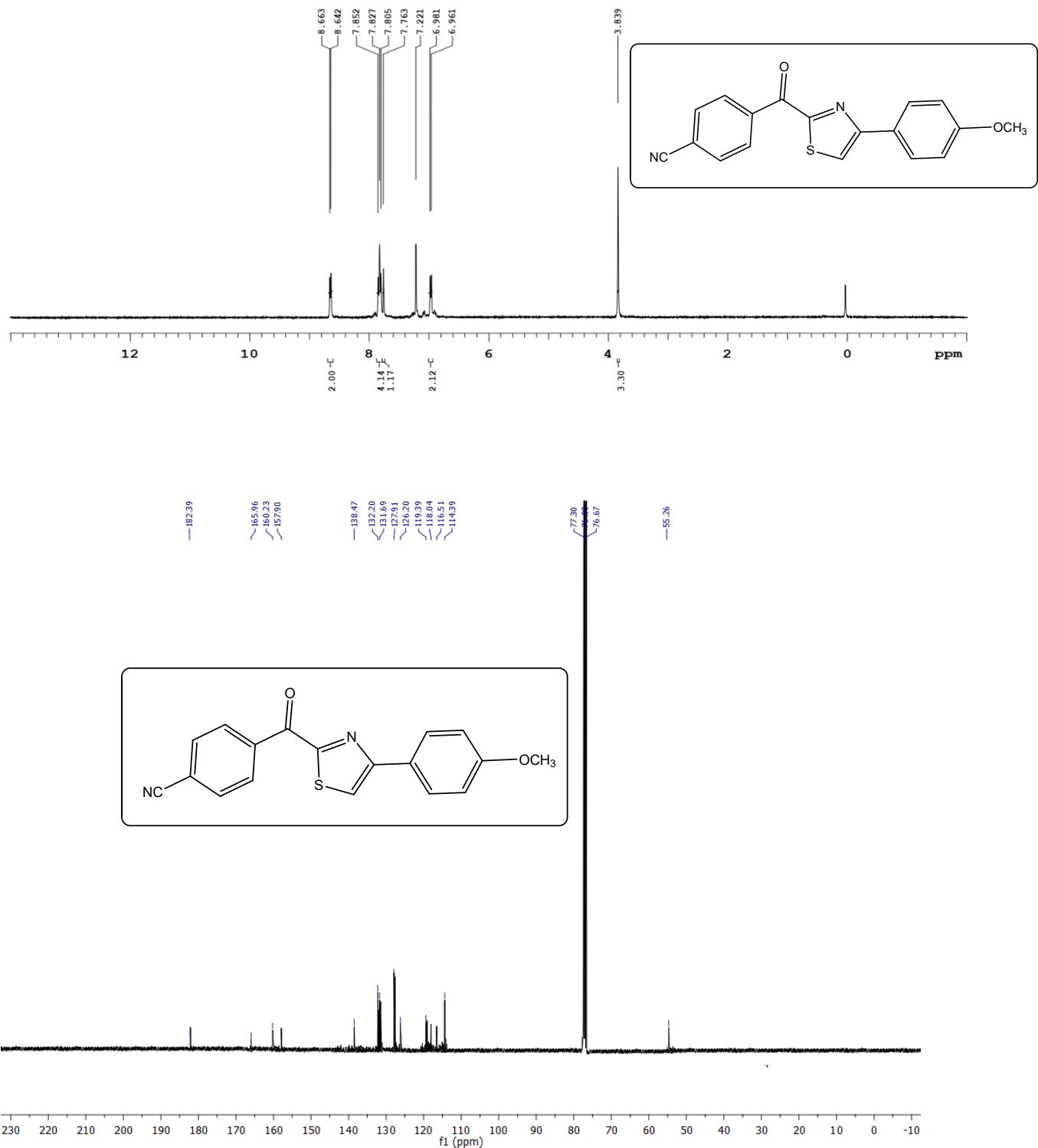
¹H & ¹³C NMR of [1,1'-Biphenyl]-4-yl(4-(4-chlorophenyl)thiazol-2-yl)methanone (4l)



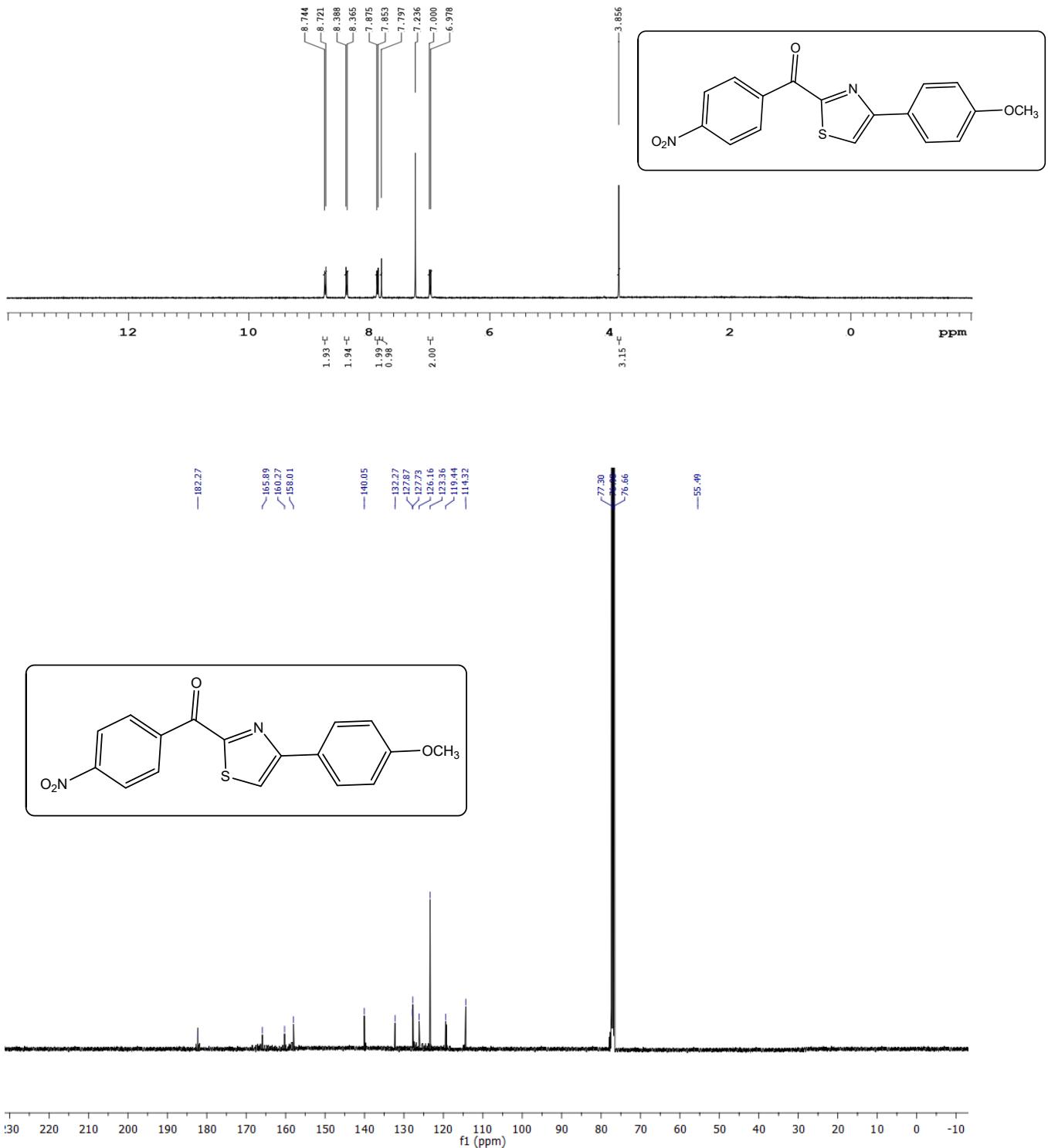
¹H & ¹³C NMR of (4-(4-Chlorophenyl)thiazol-2-yl)(furan-2-yl)methanone (4m)



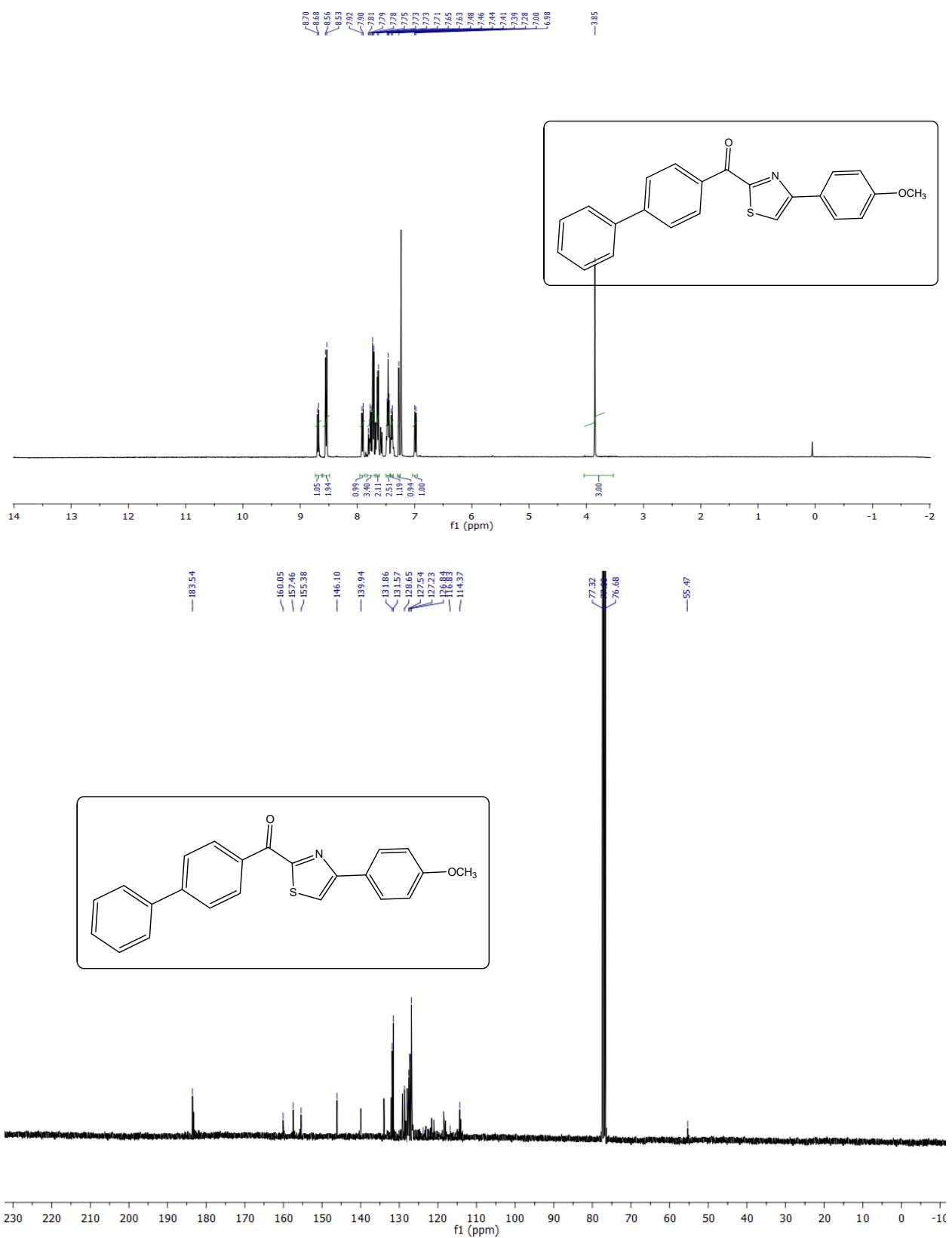
¹H & ¹³C NMR of (4-(4-Methoxyphenyl)thiazol-2-yl)(phenyl)methanone (4n)



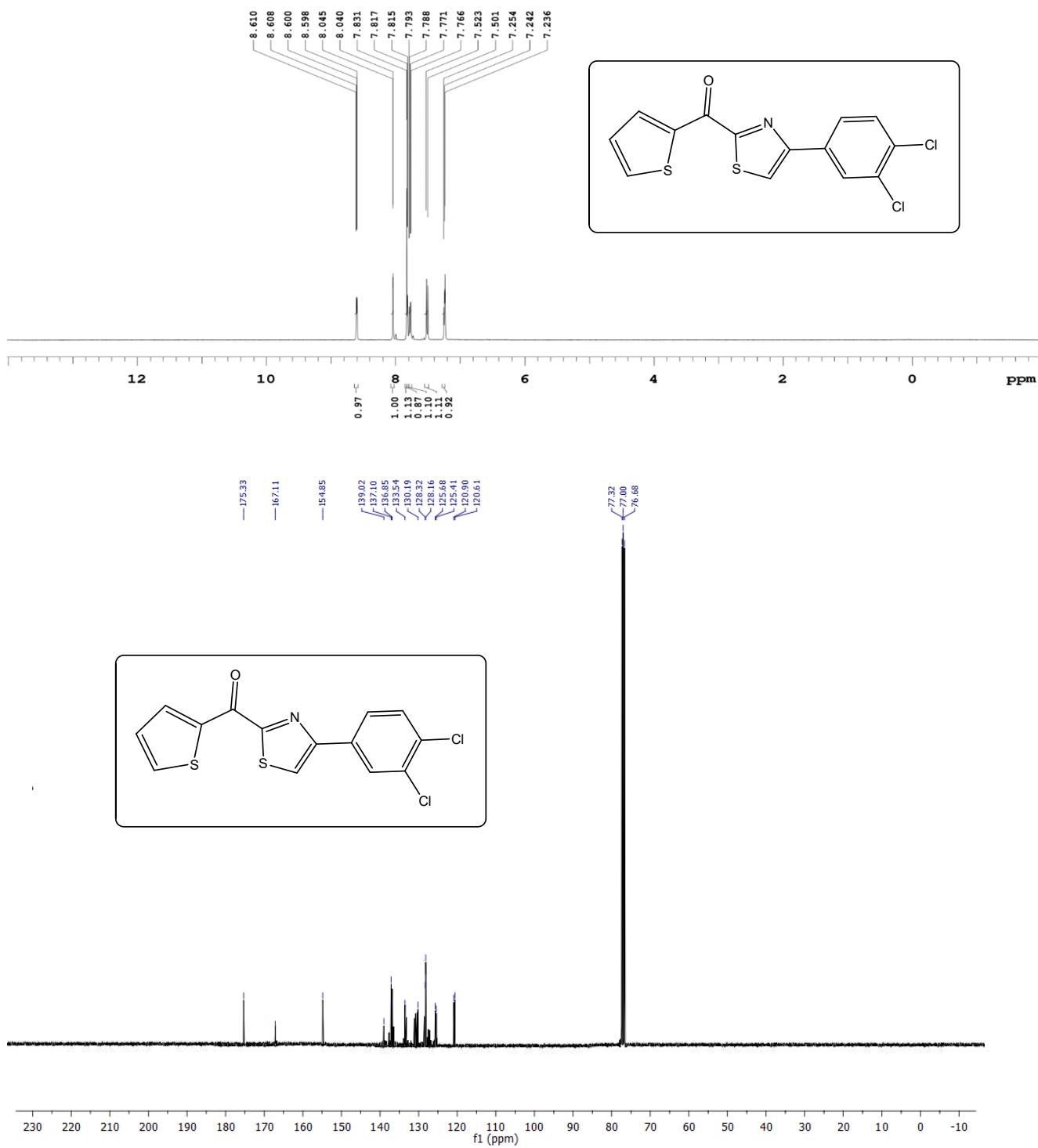
¹H & ¹³C NMR of 4-(4-Methoxyphenyl)thiazole-2-carbonylbenzonitrile (4o**)**



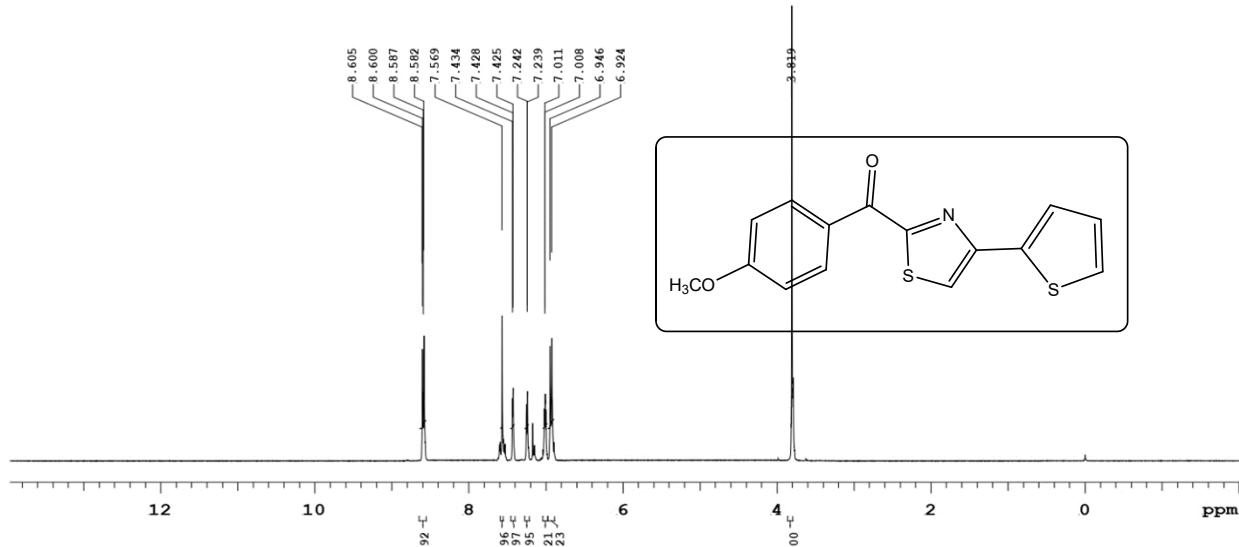
¹H & ¹³C NMR of (4-(4-Methoxyphenyl)thiazol-2-yl)(4-nitrophenyl)methanone (4p)



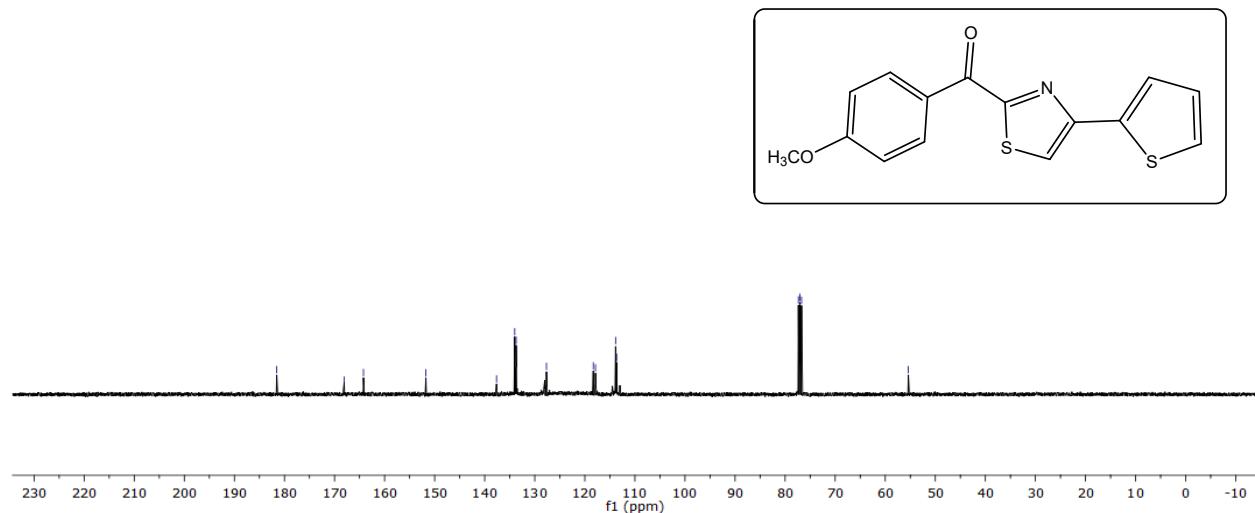
¹H & ¹³C NMR of [1,1'-Biphenyl]-4-yl(4-(4-methoxyphenyl)thiazol-2-yl)methanone (4q)



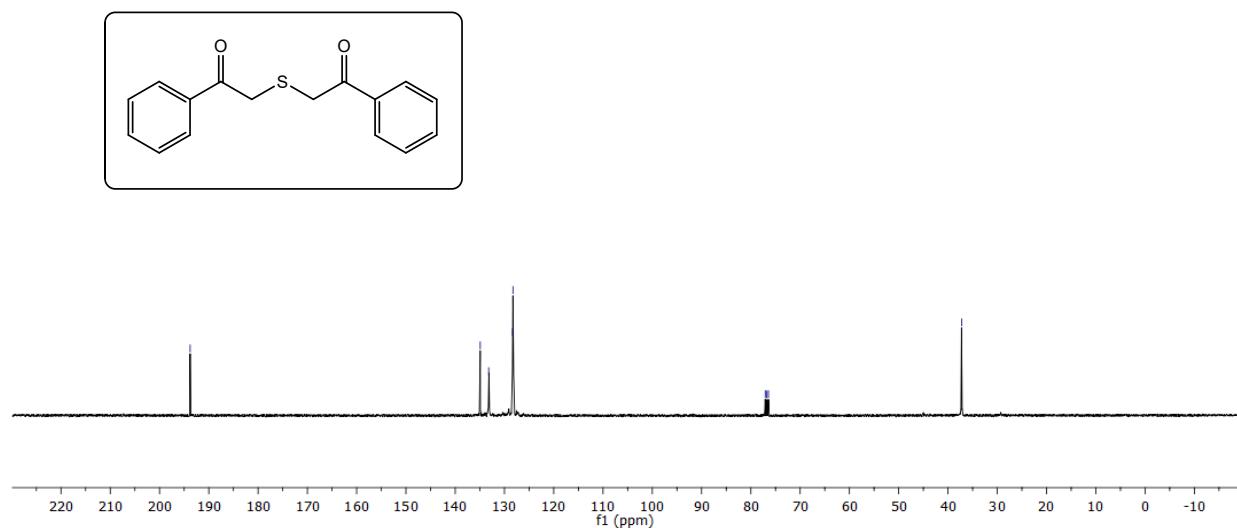
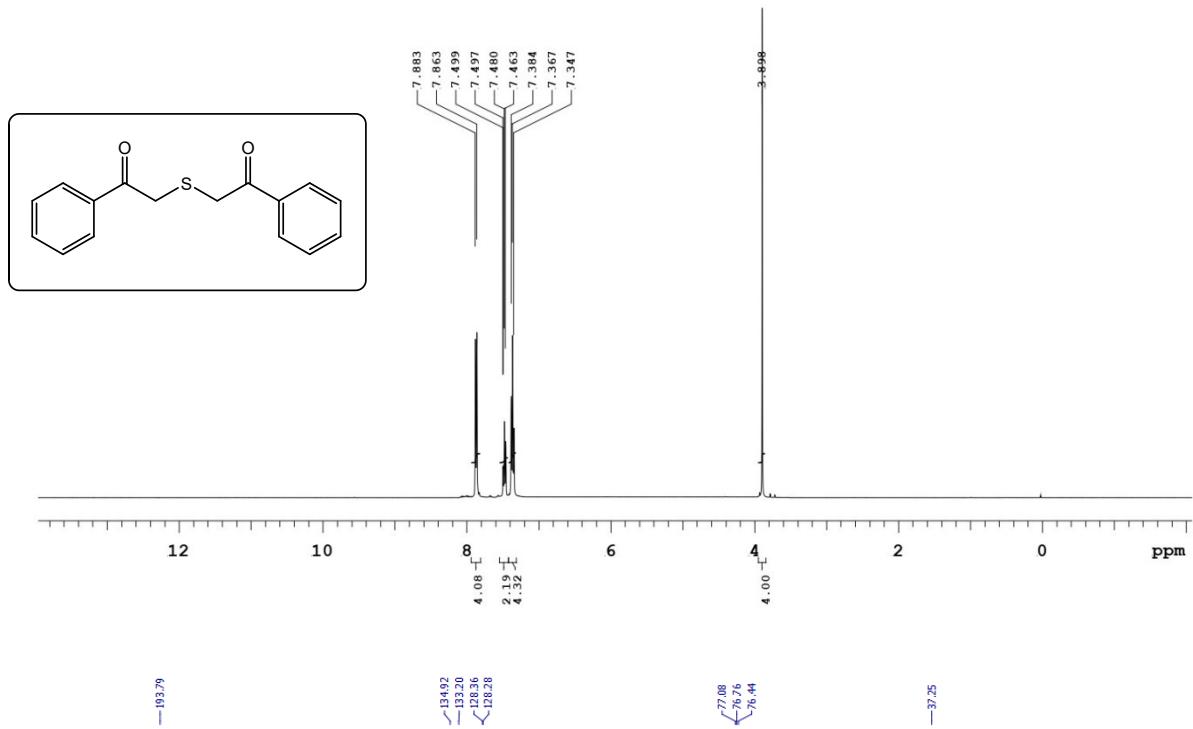
¹H & ¹³C NMR of (4-(3,4-Dichlorophenyl)thiazol-2-yl)(thiophen-2-yl)methanone (4r)



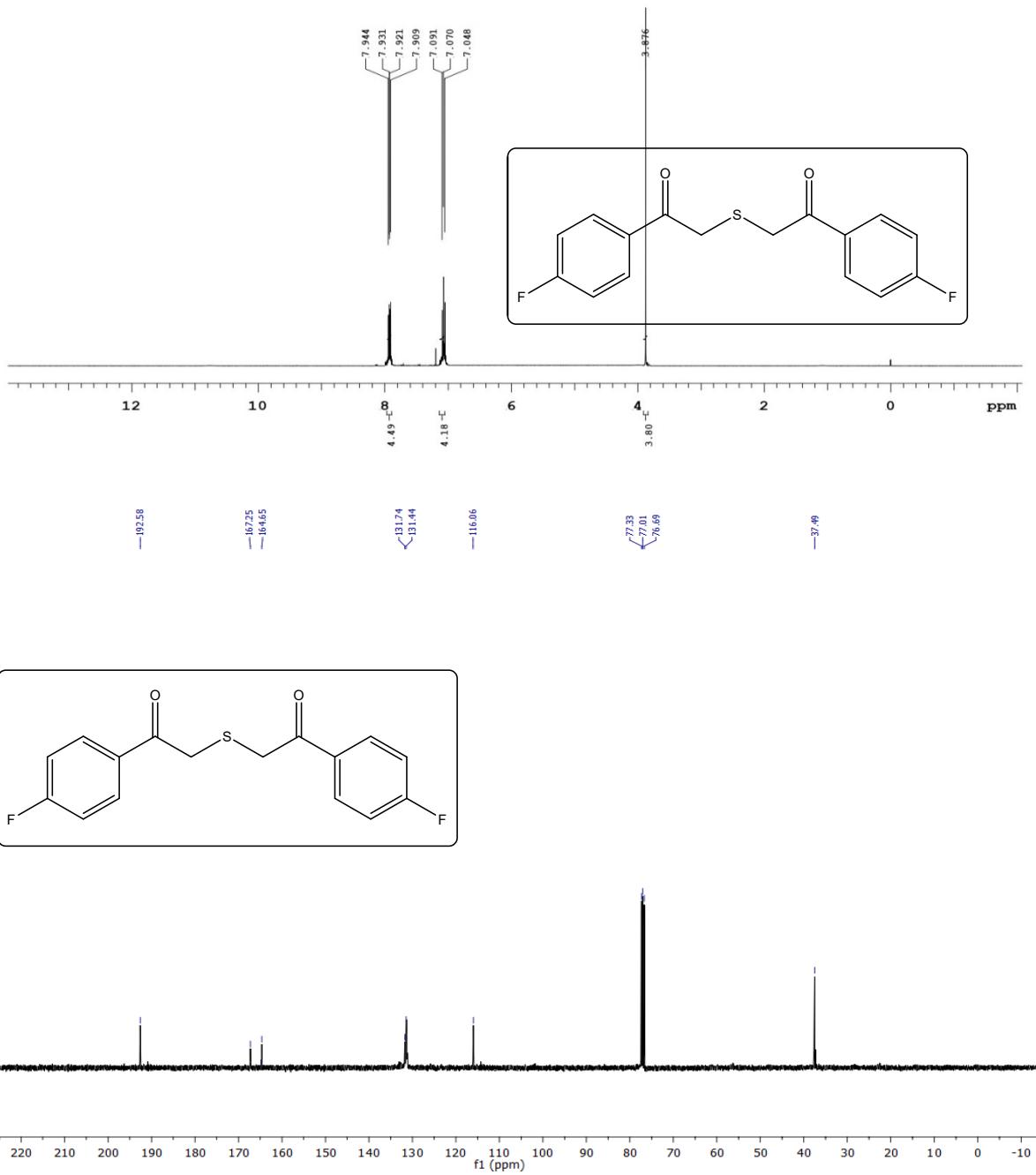
—181.54
—168.07
—165.22
—151.75
—137.64
—135.00
—132.23
—127.67
—118.32
—117.88
—113.86
—113.66
—55.36



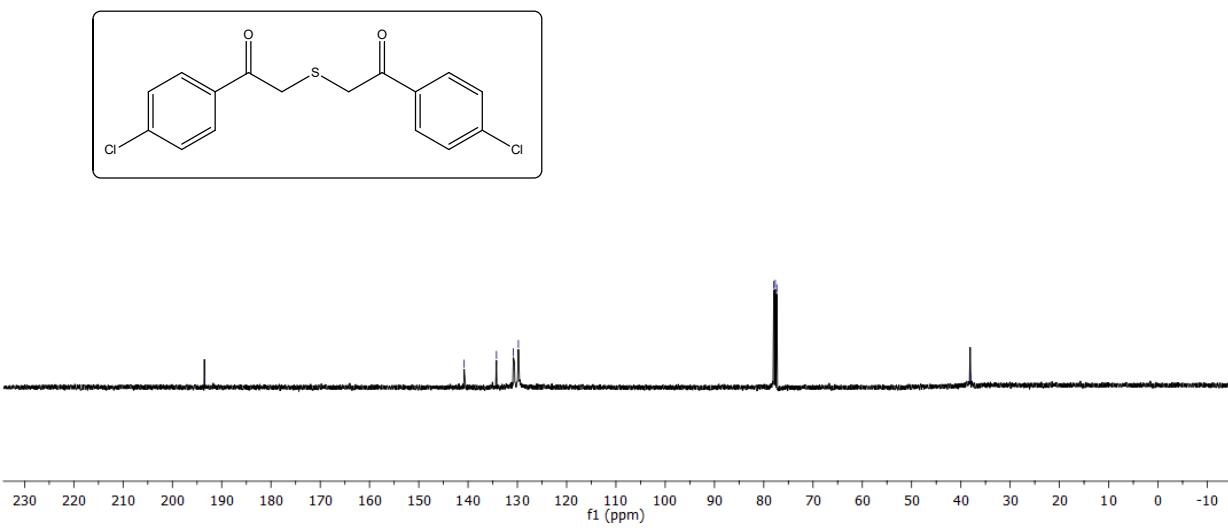
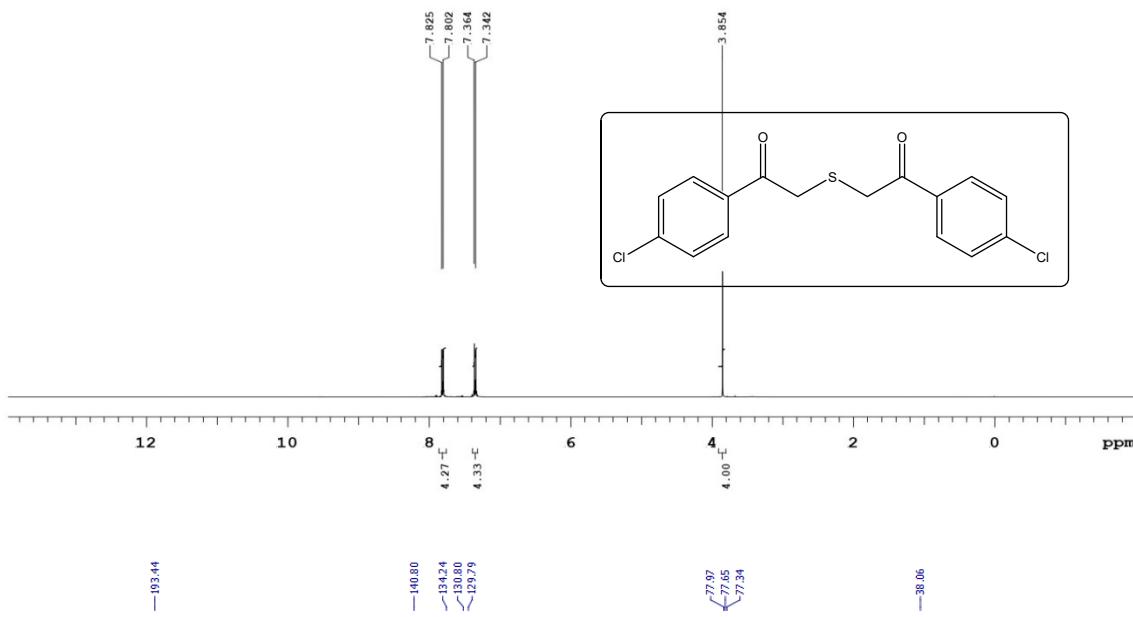
¹H & ¹³C NMR of (4-methoxyphenyl)(4-(thiophen-2-yl)thiazol-2-yl)methanone (4s)



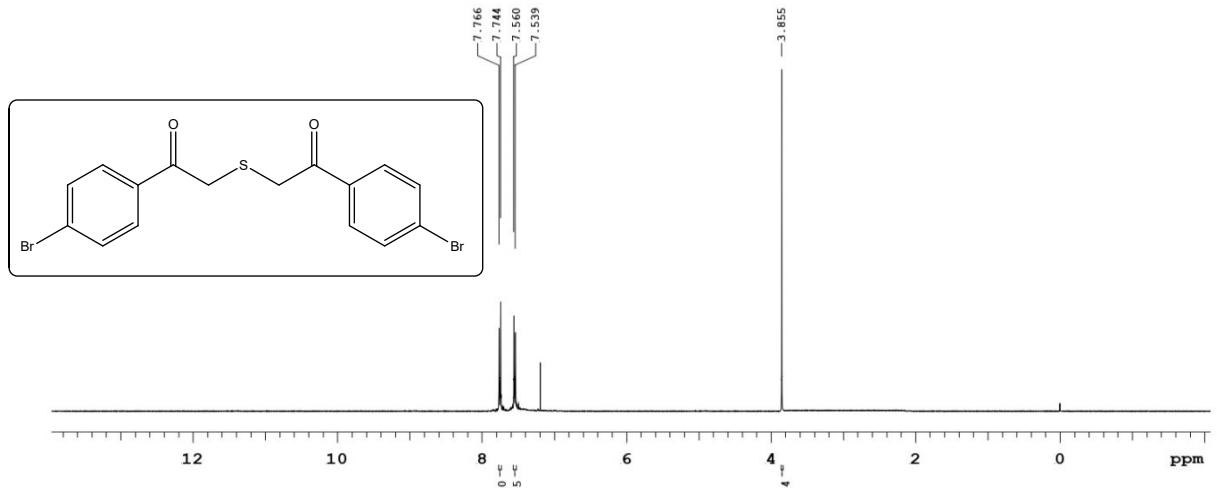
^1H & ^{13}C NMR of 2,2'-thiobis(1-phenylethanone) (5a)



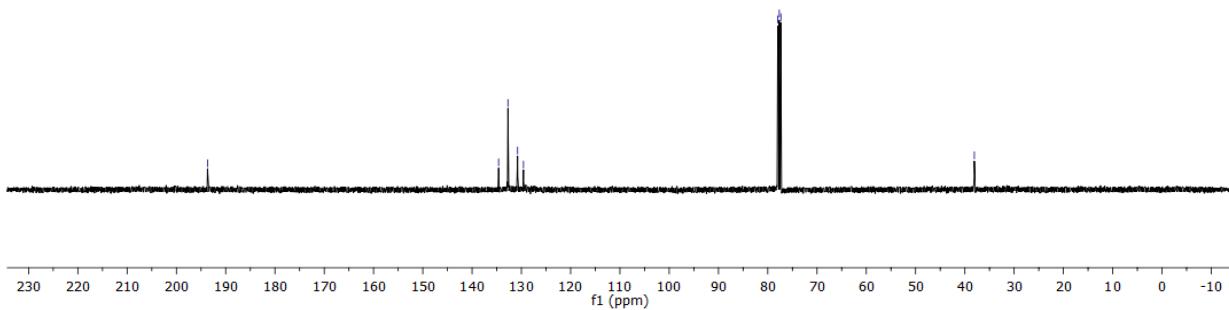
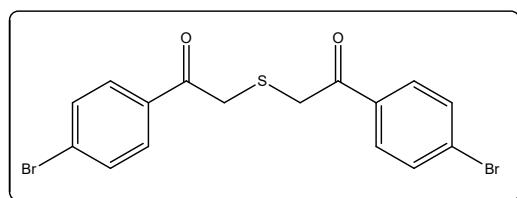
¹H & ¹³C NMR of 2, 2'-thiobis(1-(4-fluorophenyl)ethanone) (5b)



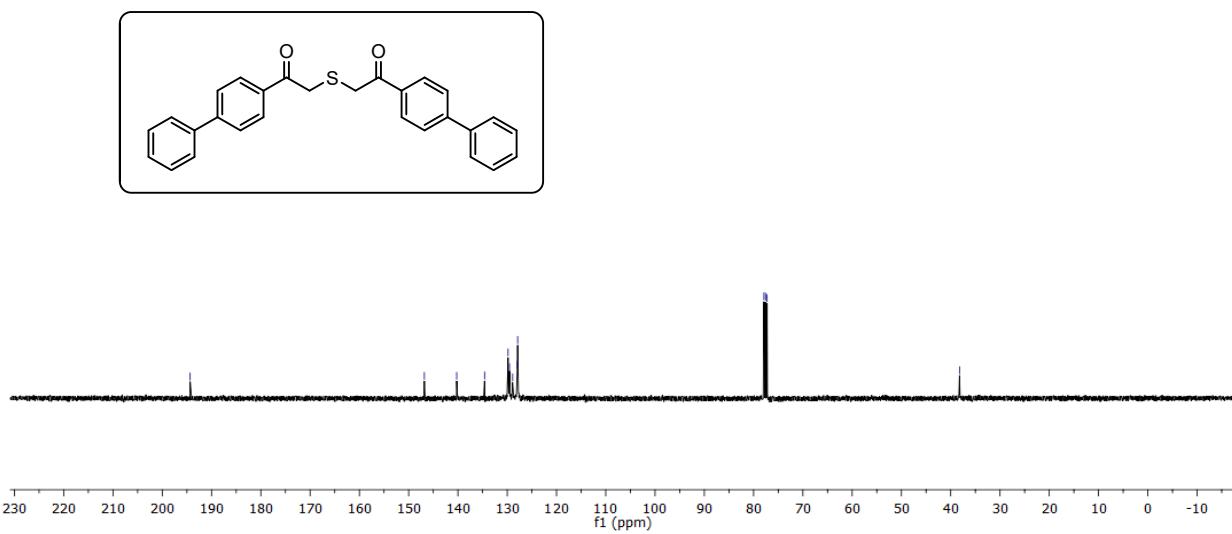
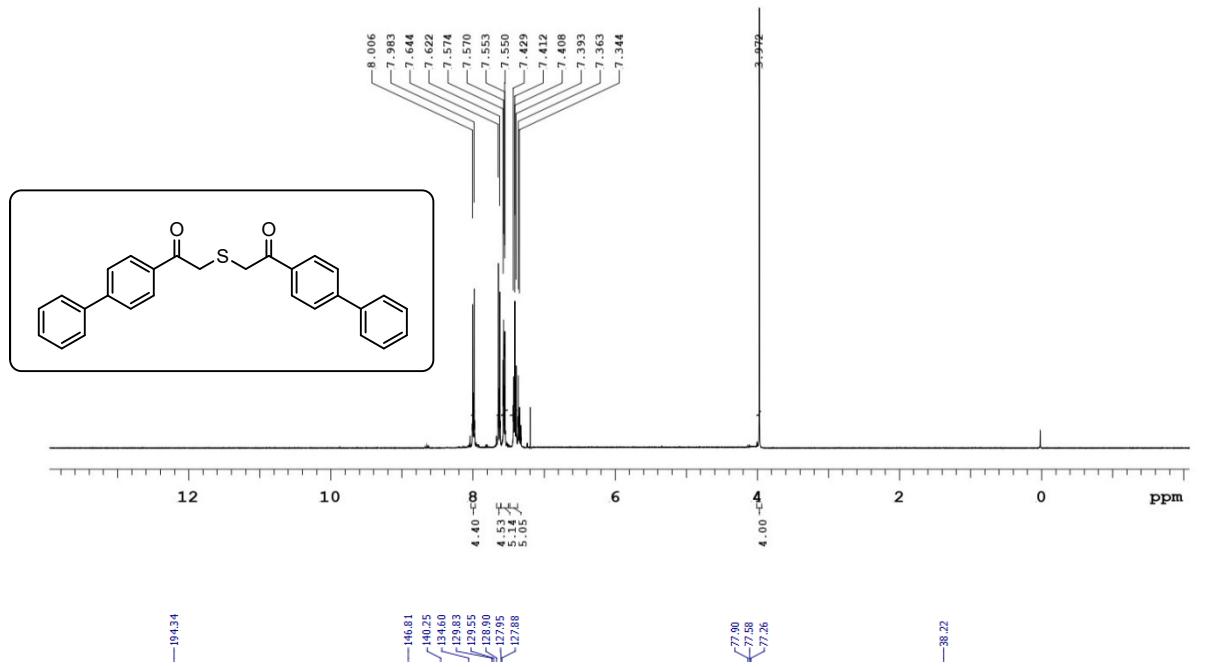
¹H & ¹³C NMR of 2,2'-thiobis(1-(4-chlorophenyl)ethanone) (5c)



13C NMR chemical shifts (δ) in ppm: 134.62, 133.73, 130.76, 129.57, 77.98, 77.66, 77.34, 38.08.



¹H & ¹³C NMR of 2,2'-thiobis(1-(4-bromophenyl)ethanone) (5d)



¹H & ¹³C NMR of 2,2'-thiobis(1-([1,1'-biphenyl]-4-yl)ethanone) (5e)



¹H & ¹³C NMR of 2,2'-thiobis(1-(thiophen-2-yl)ethanone) (5f)