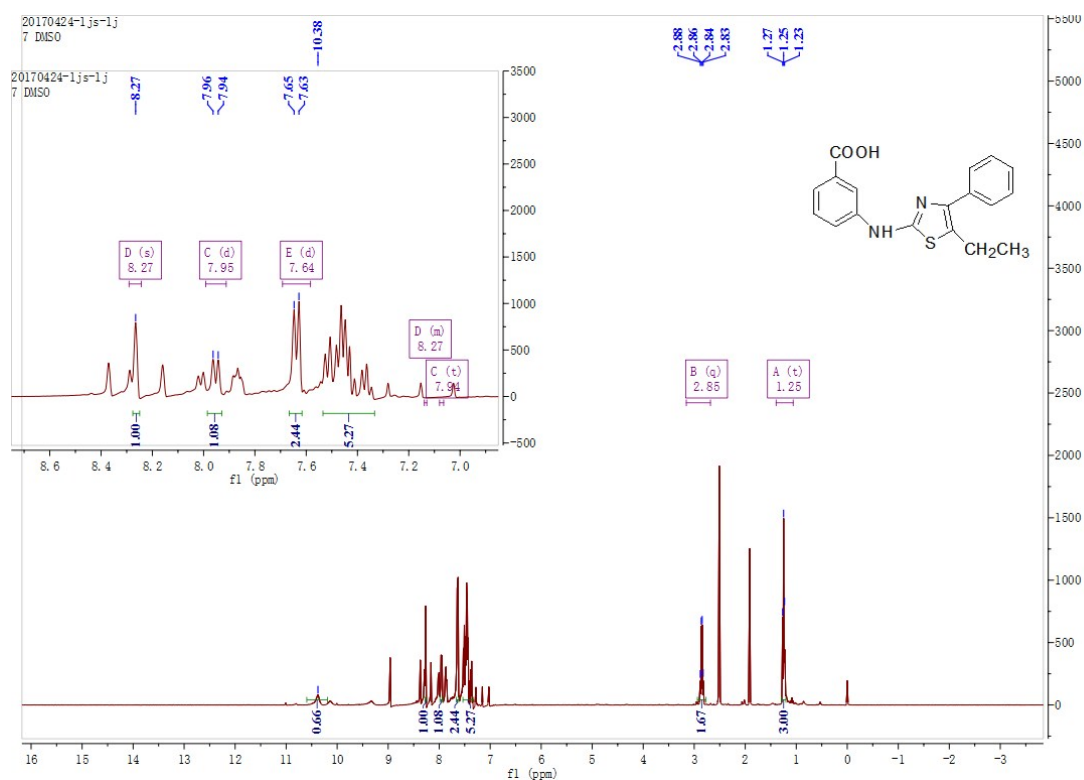
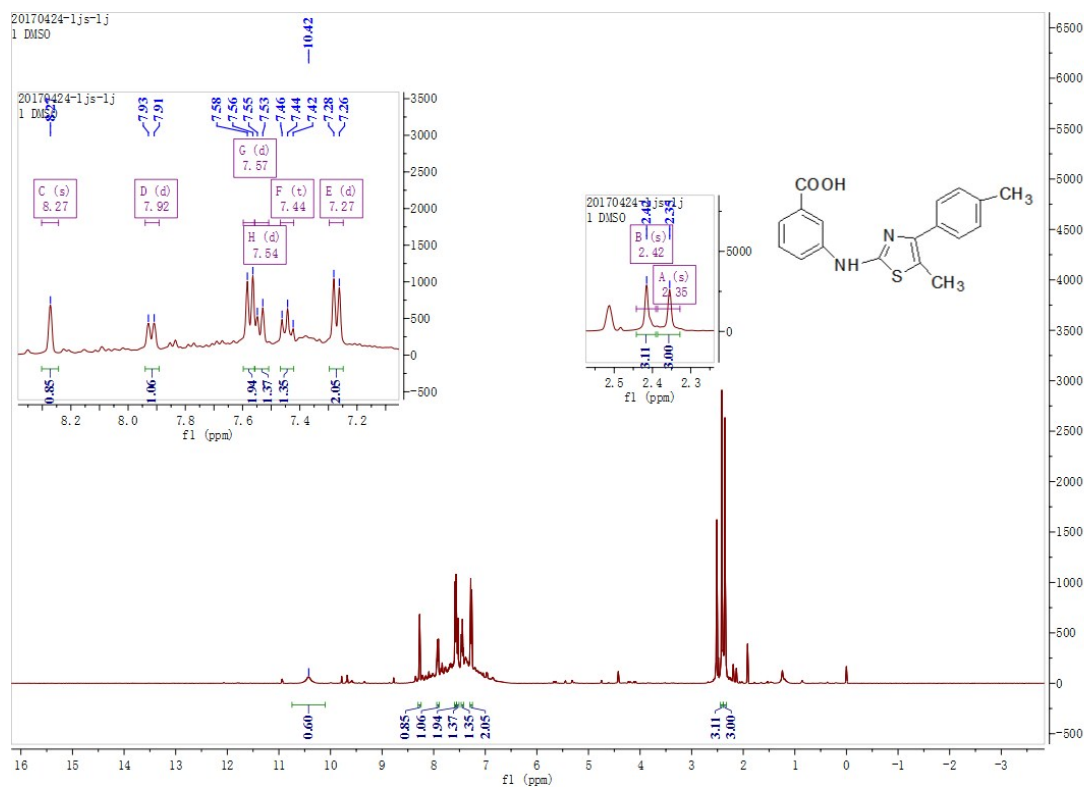


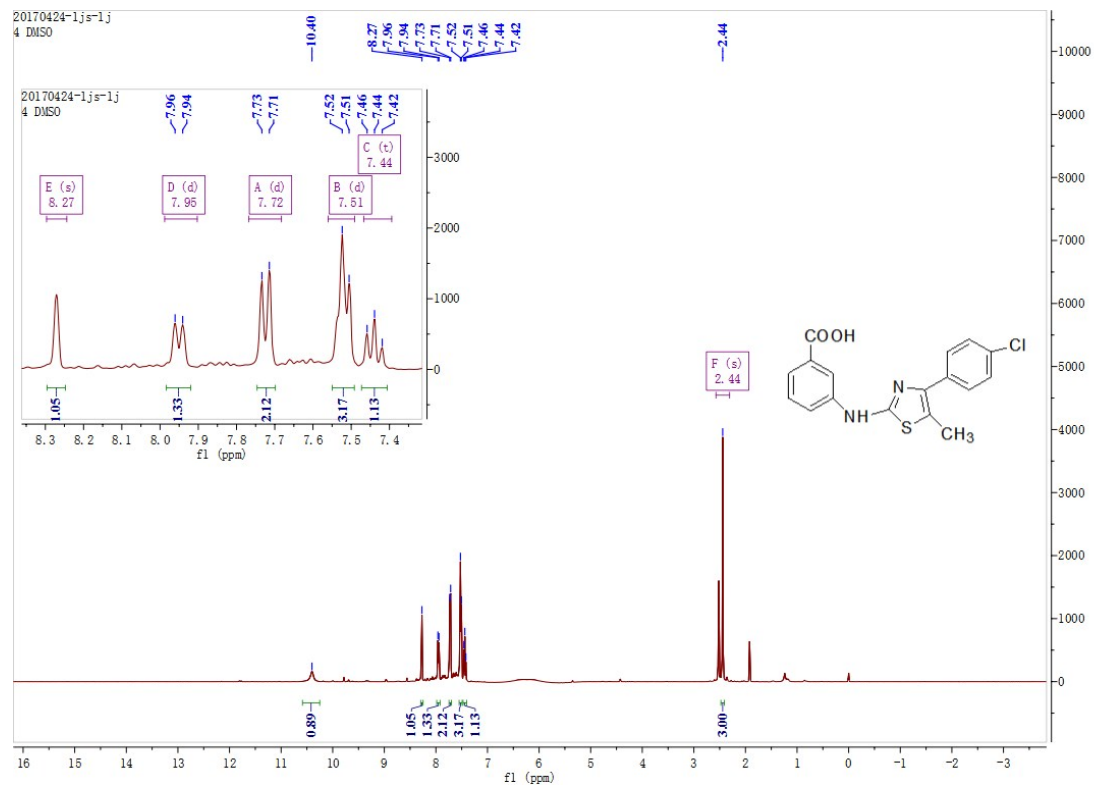
2a



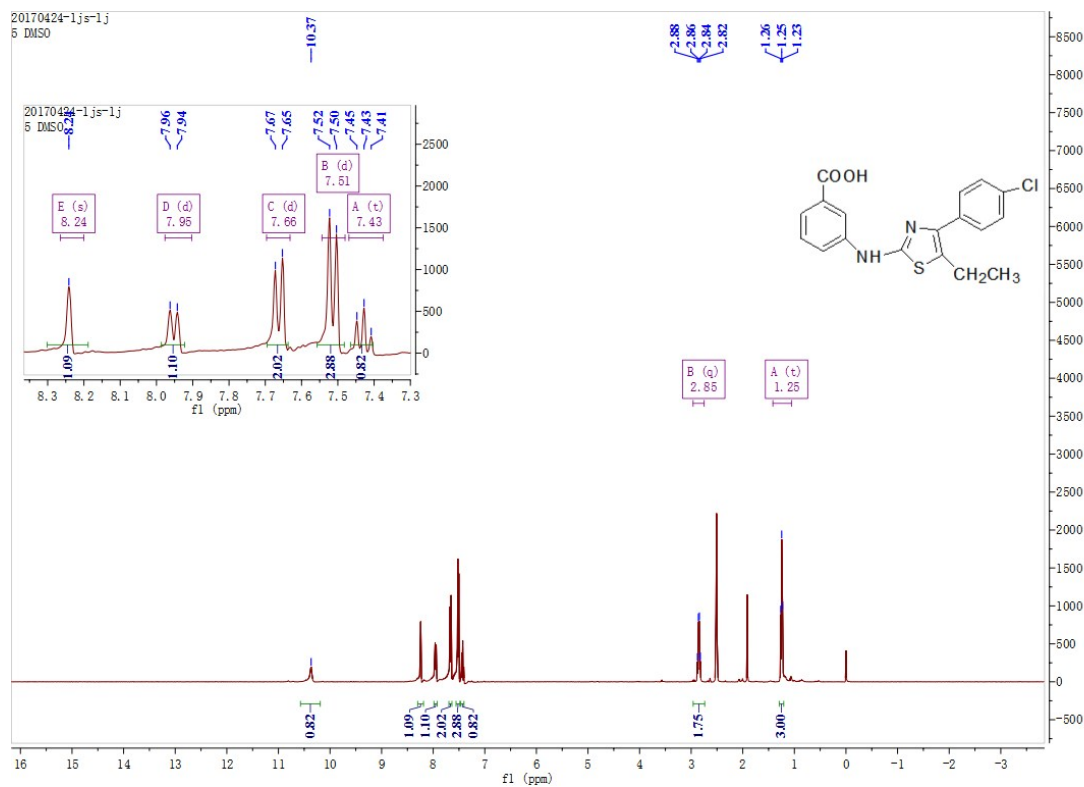
2b



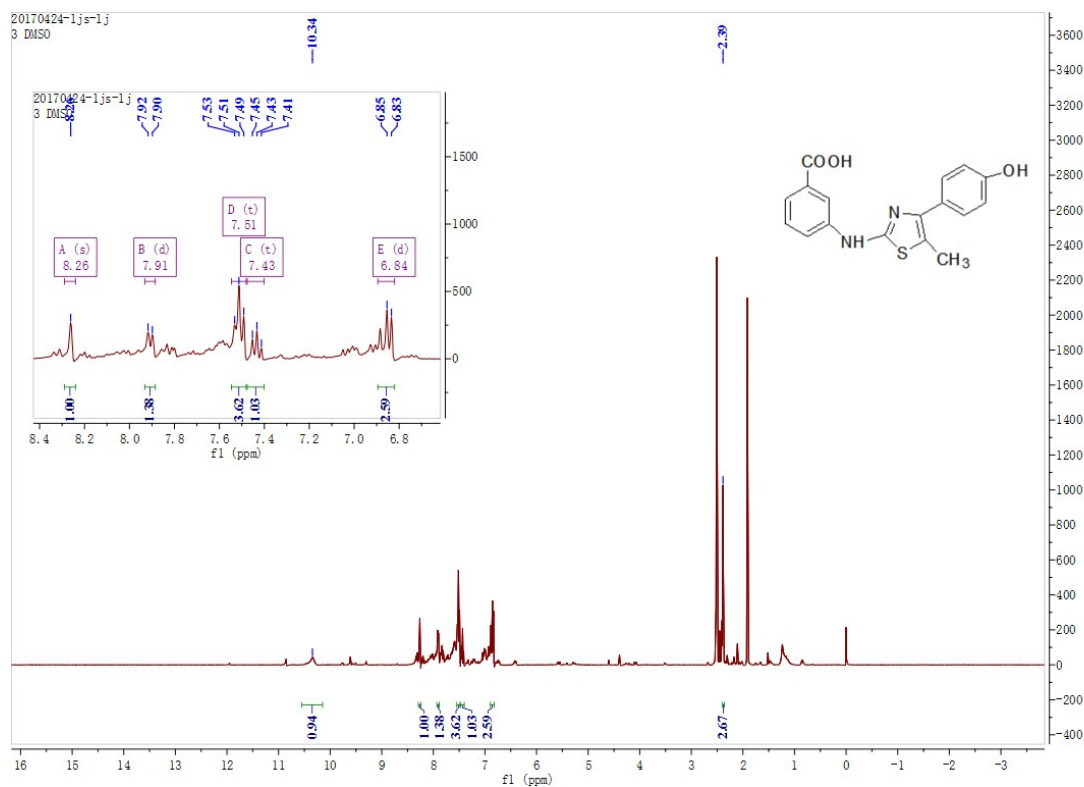
2c



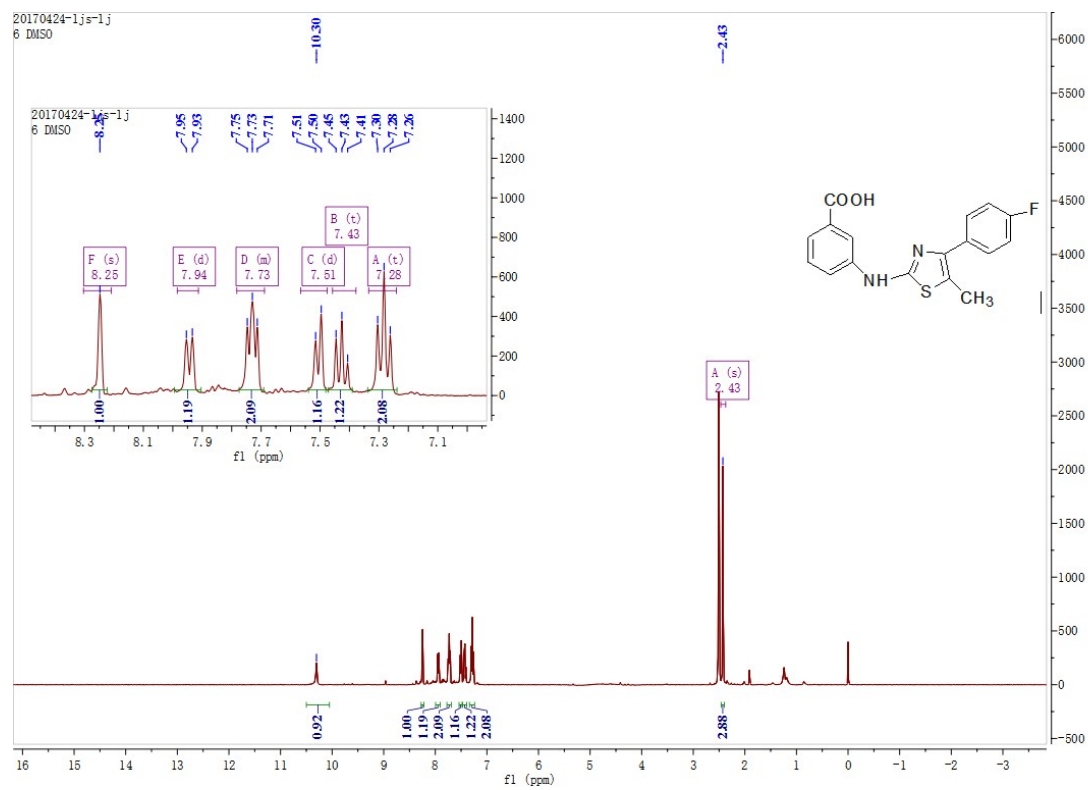
2d



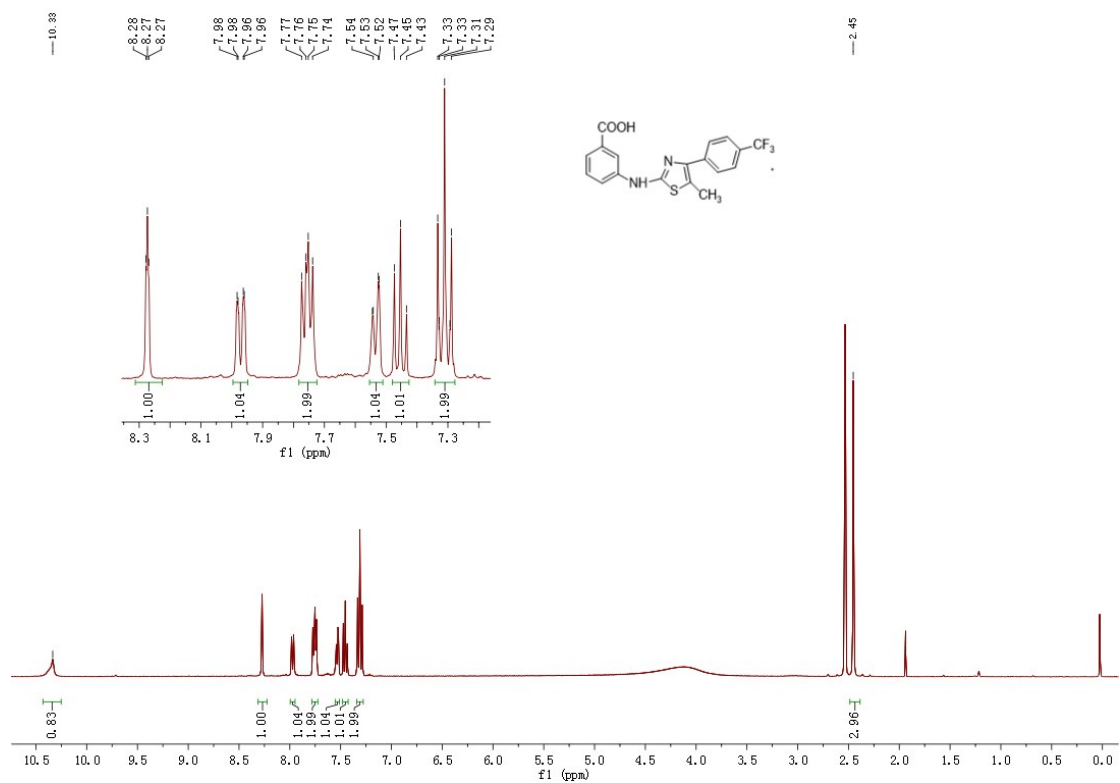
2e



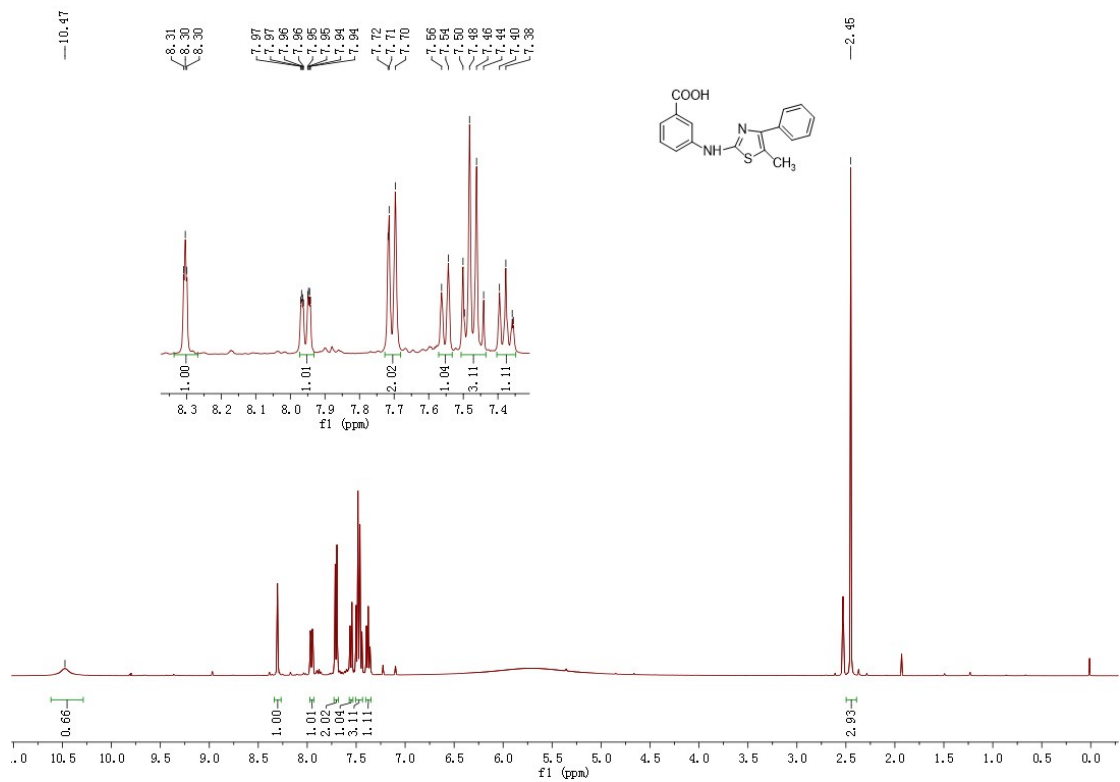
2f



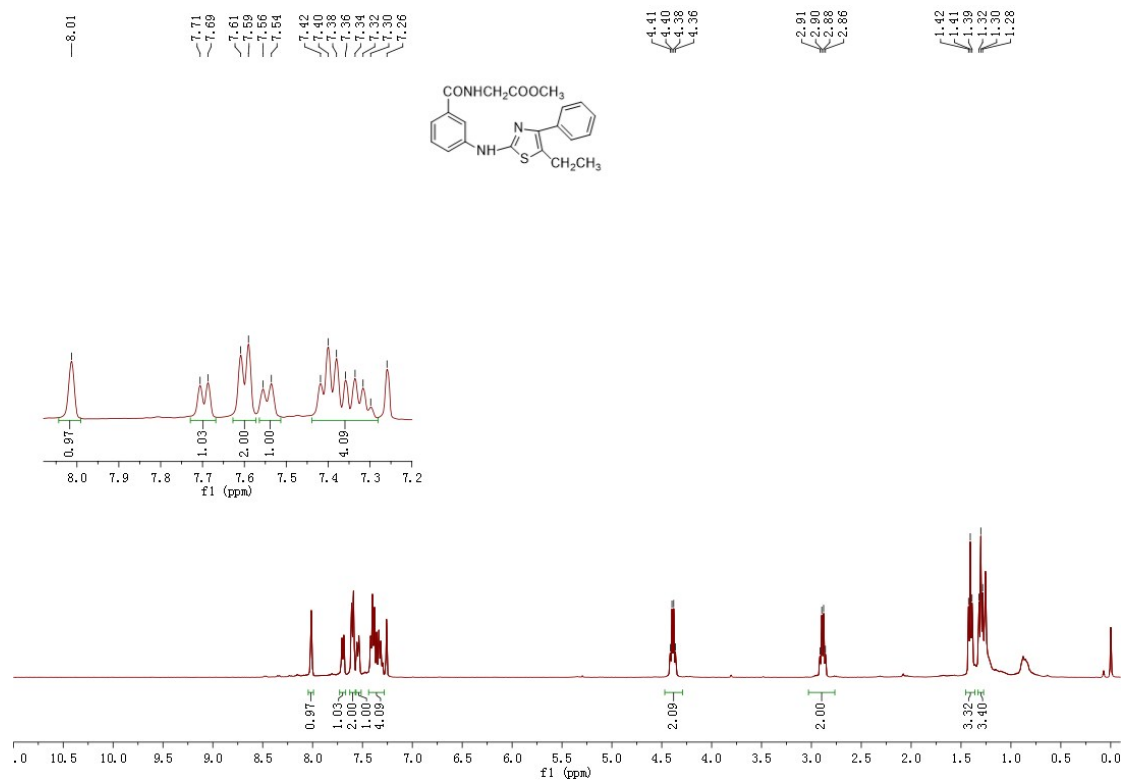
2g



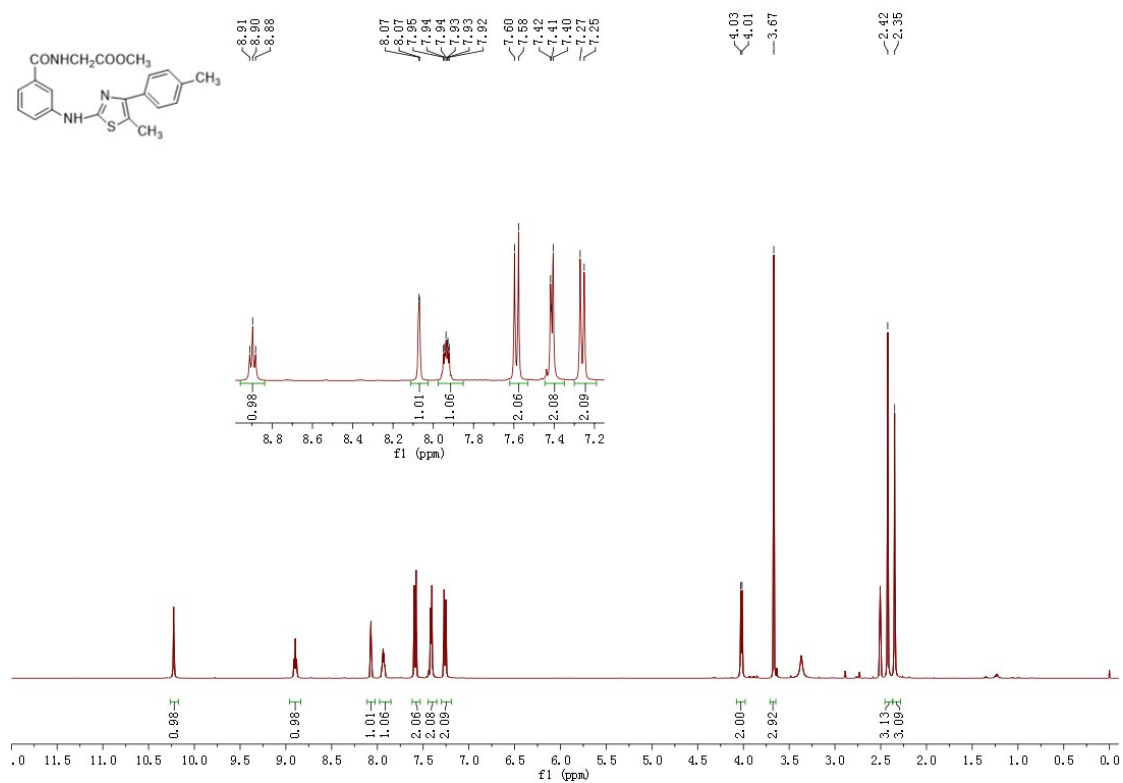
2h



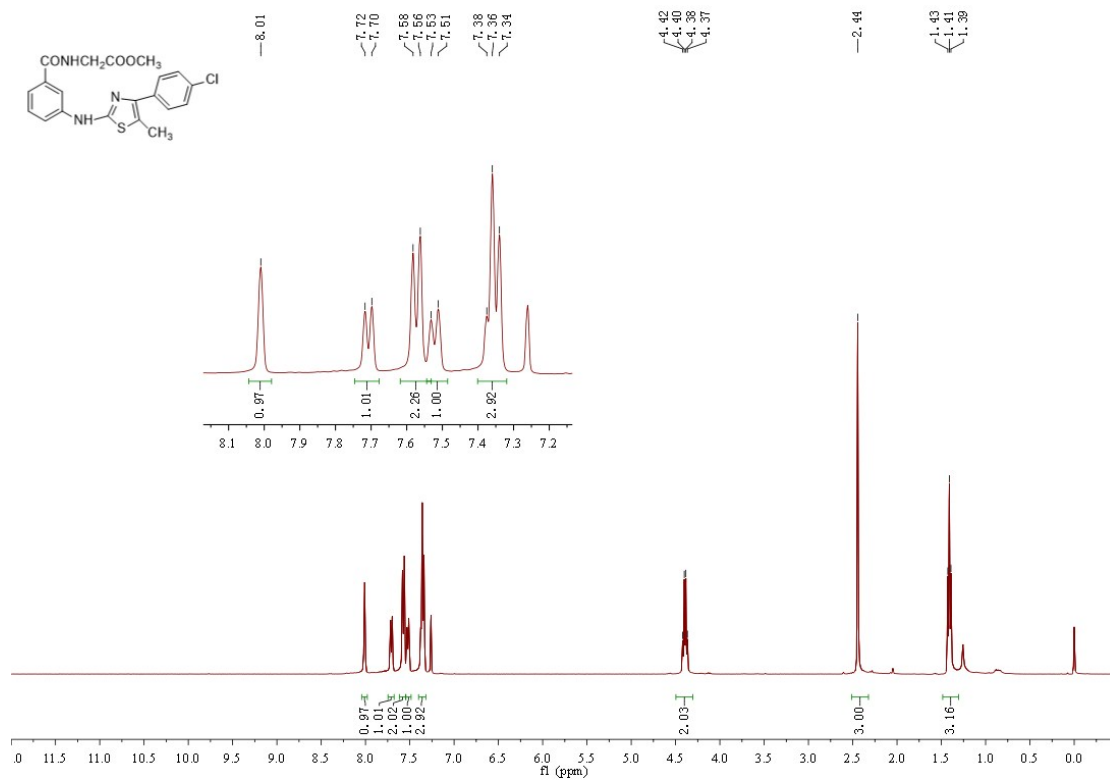
3a



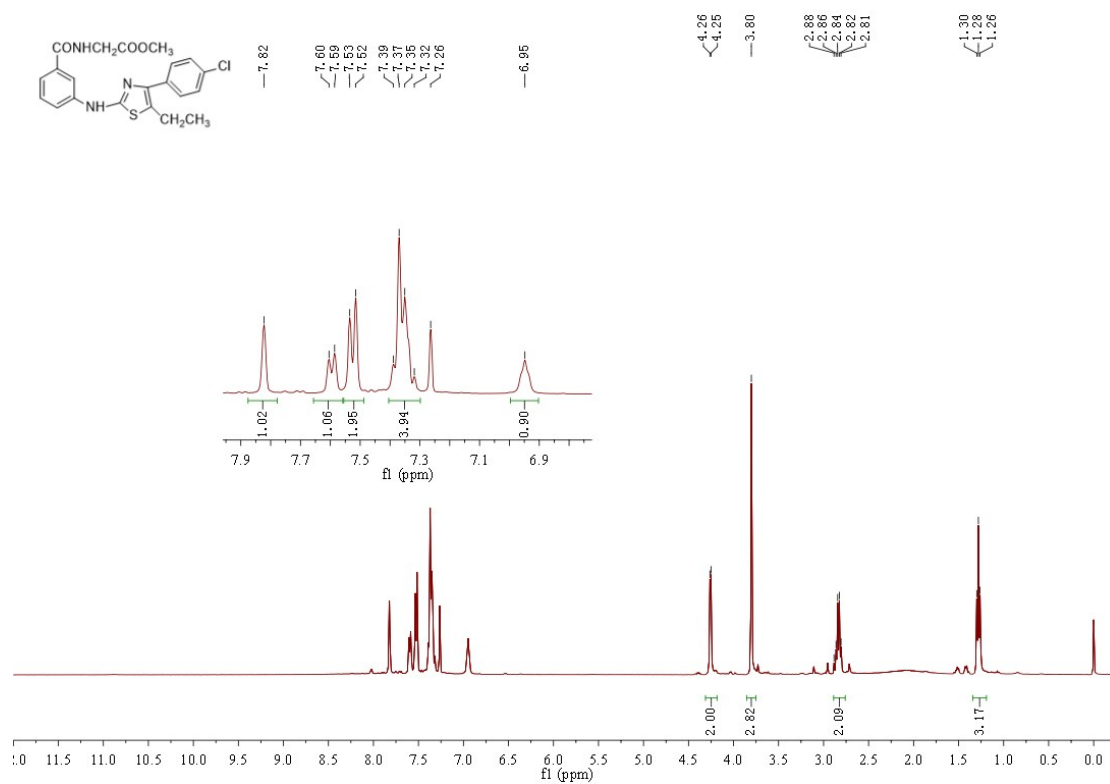
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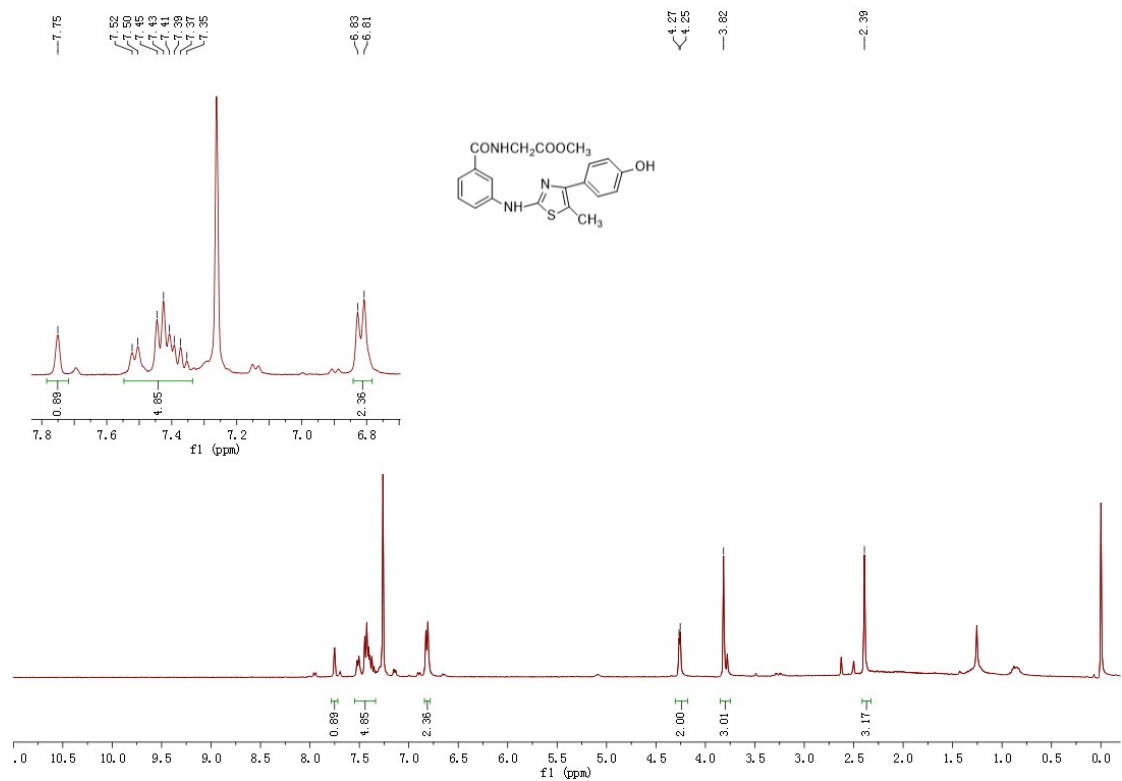
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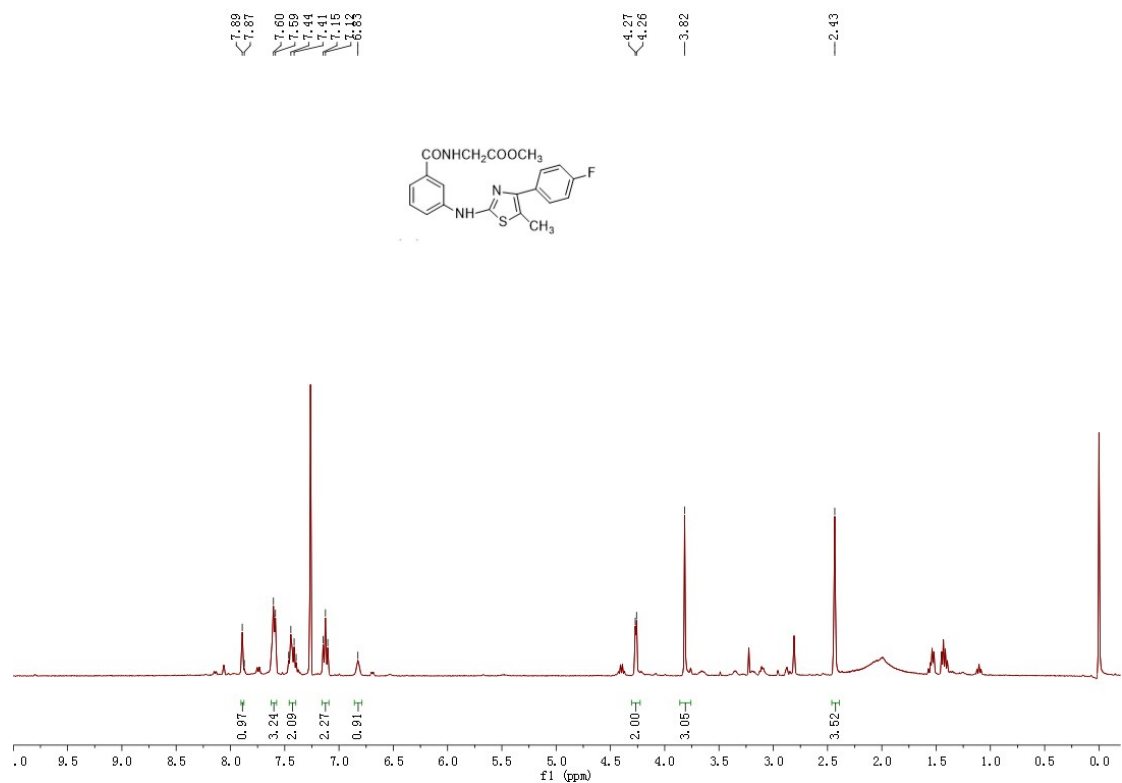
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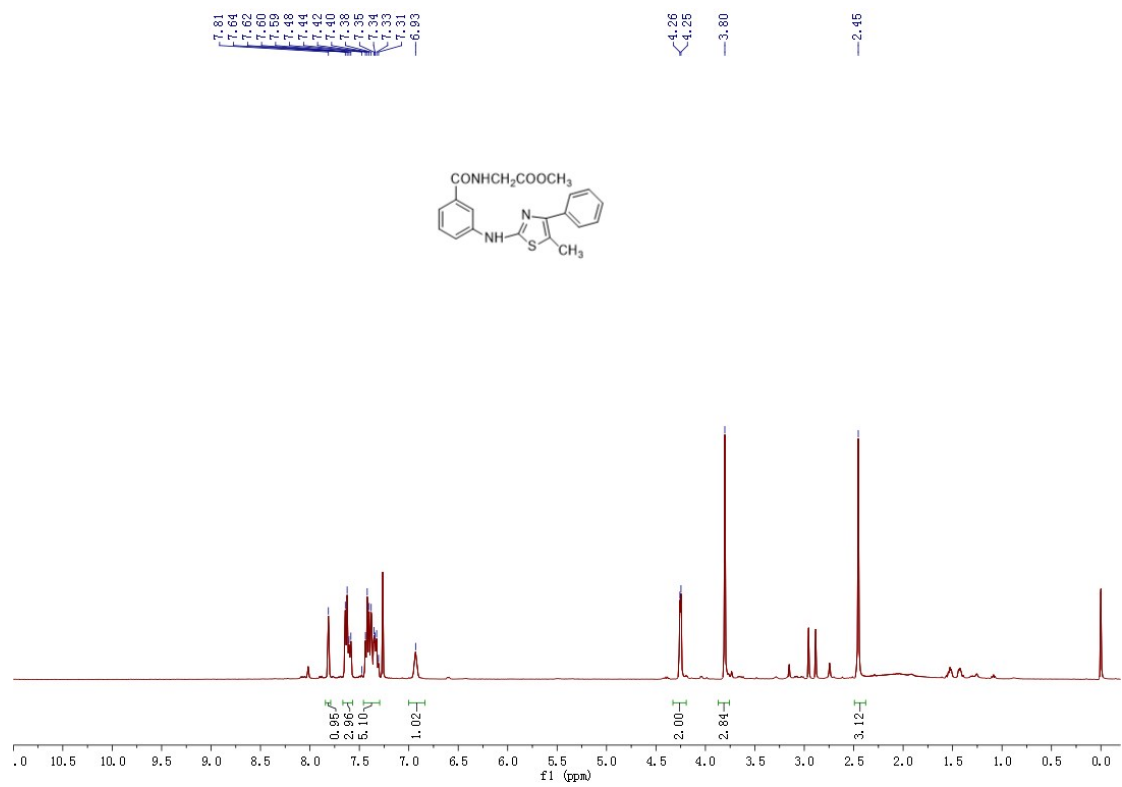
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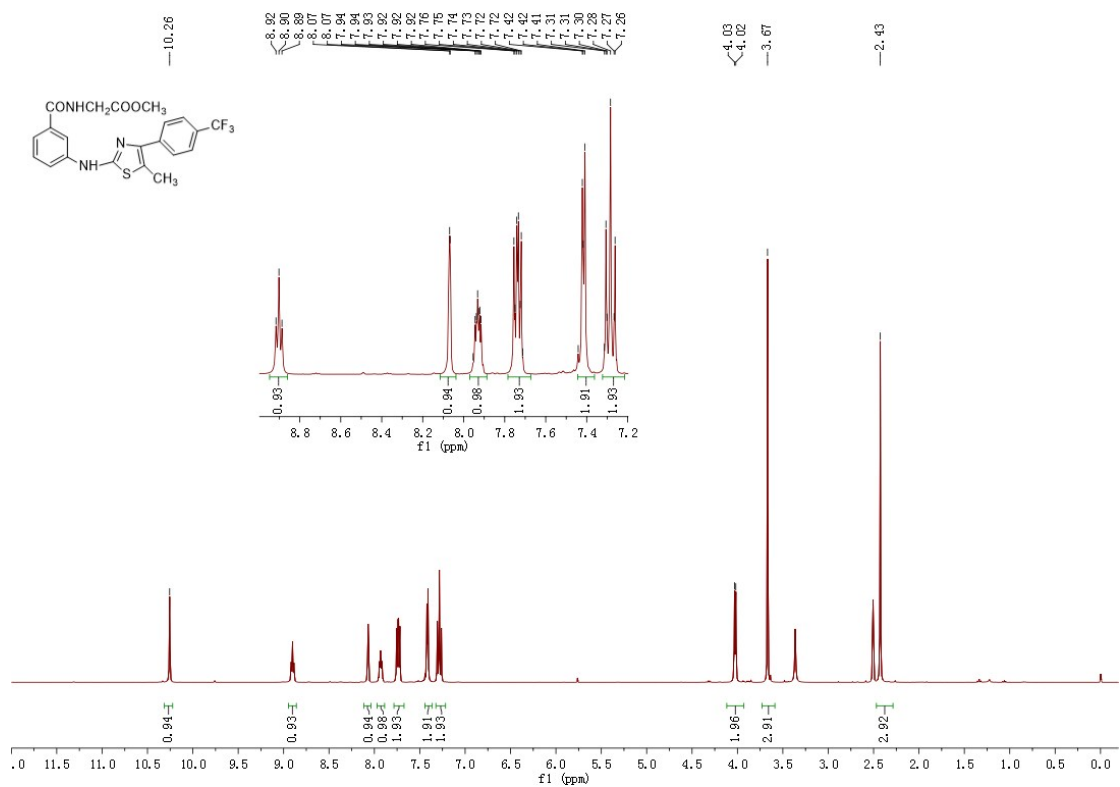
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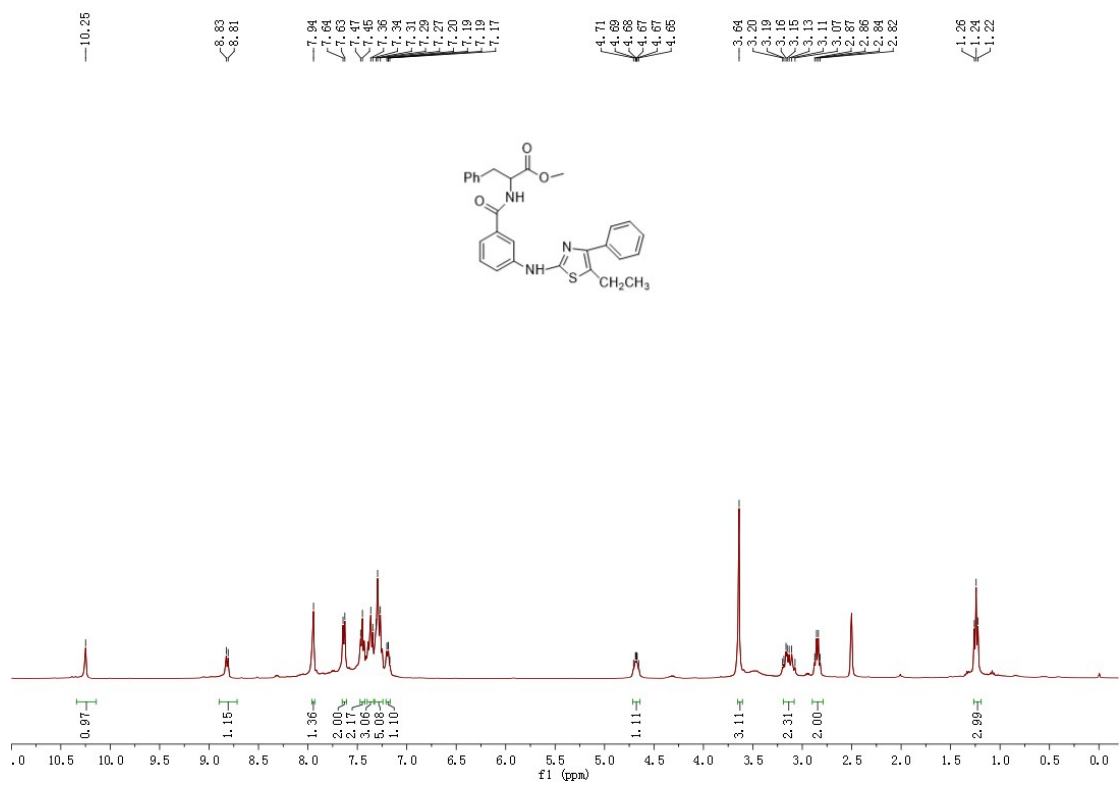
3g



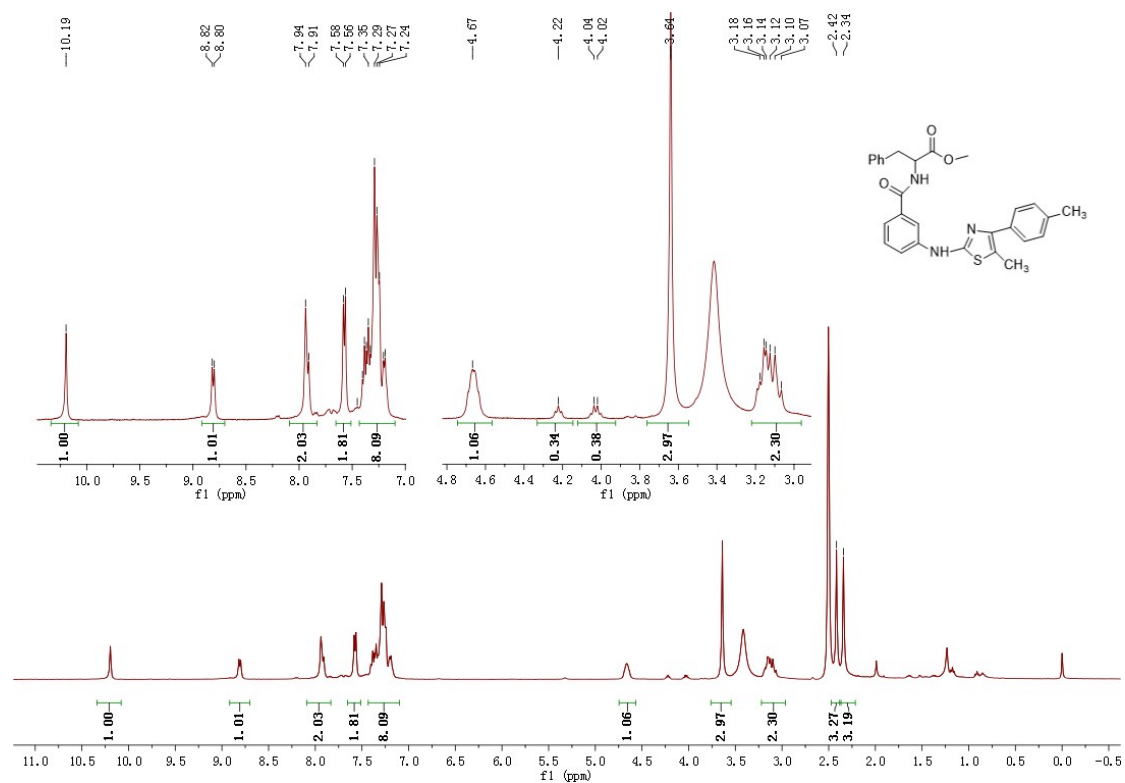
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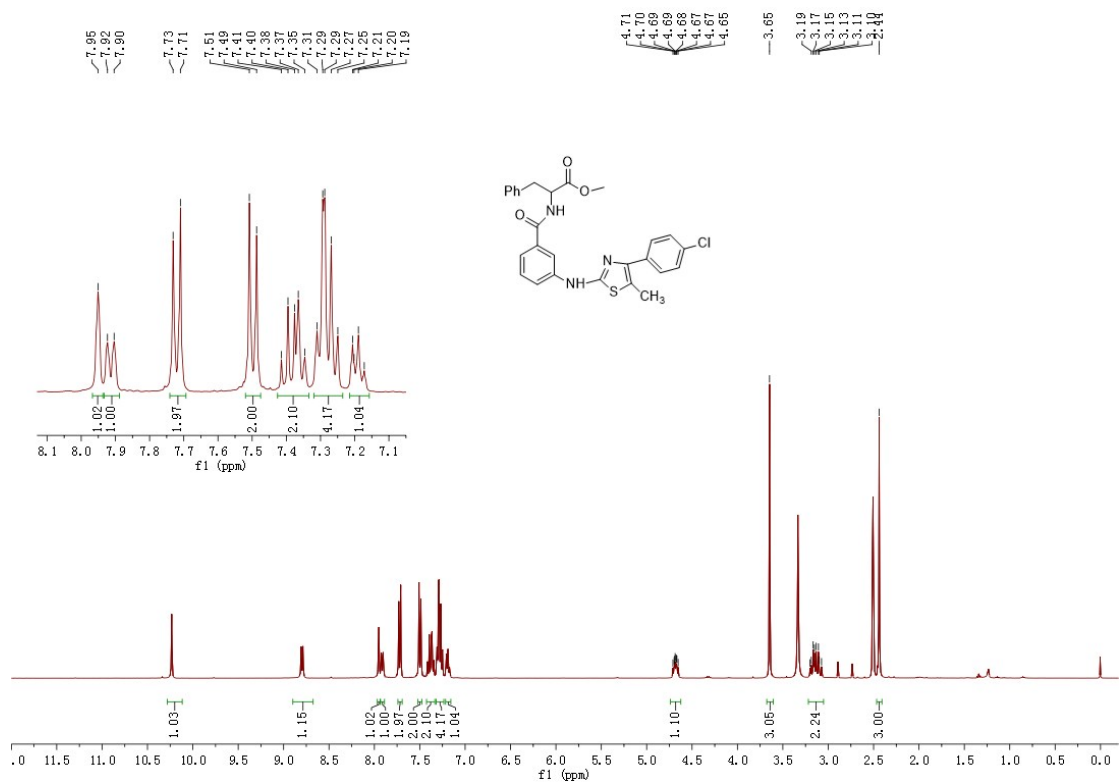
3i



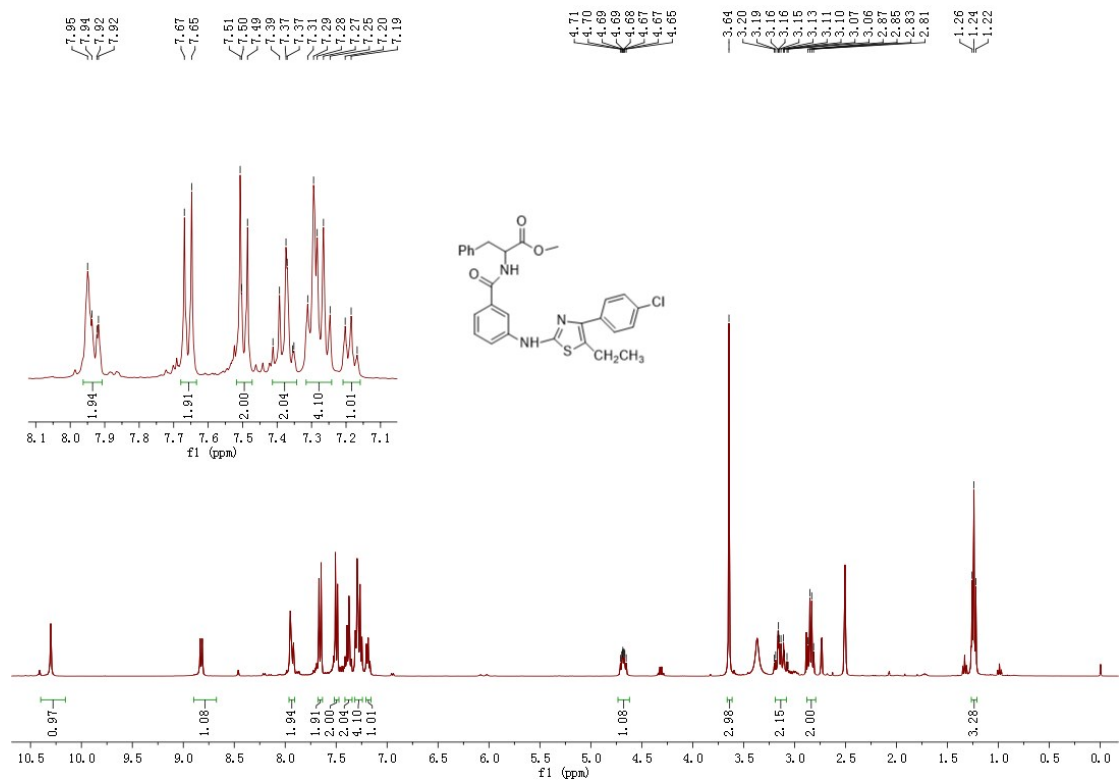
3j



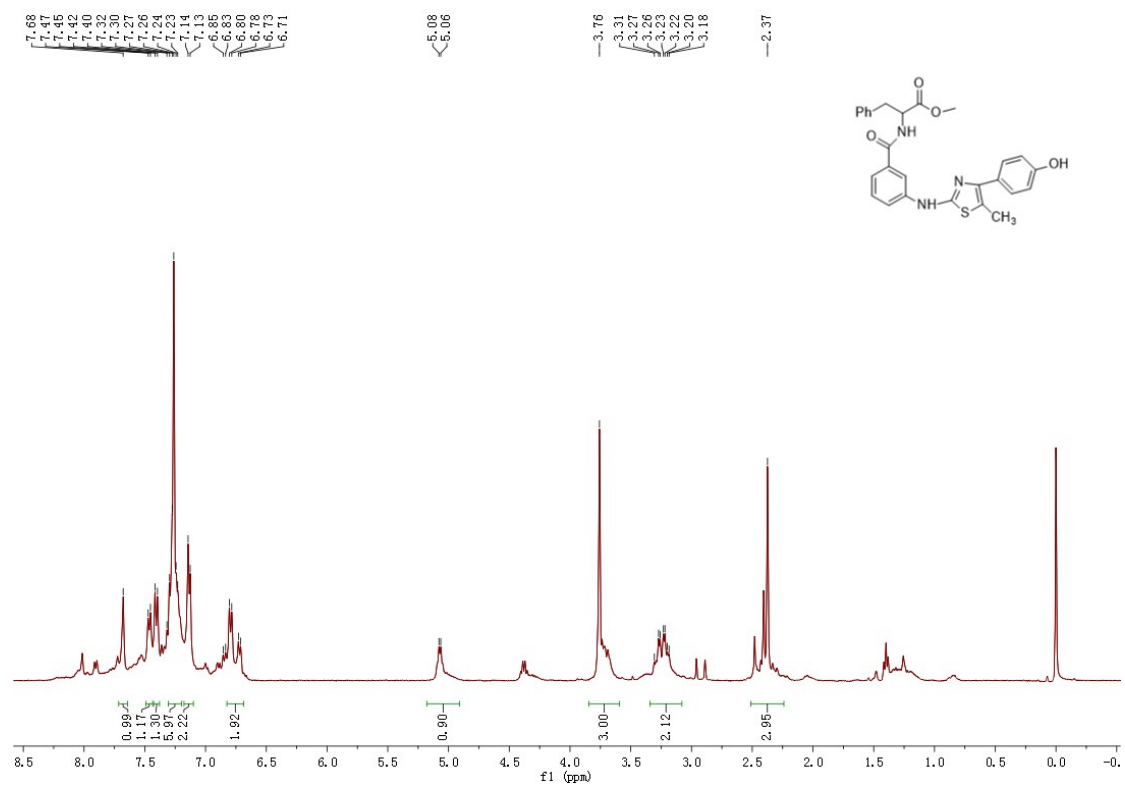
3k



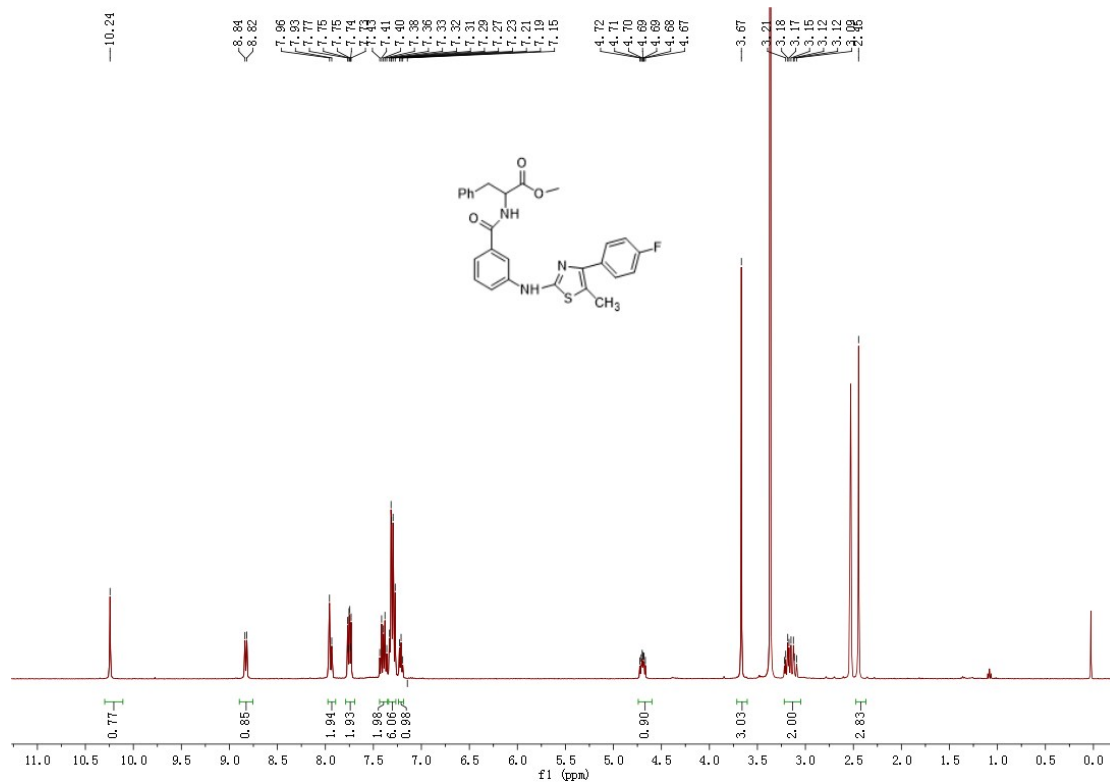
31



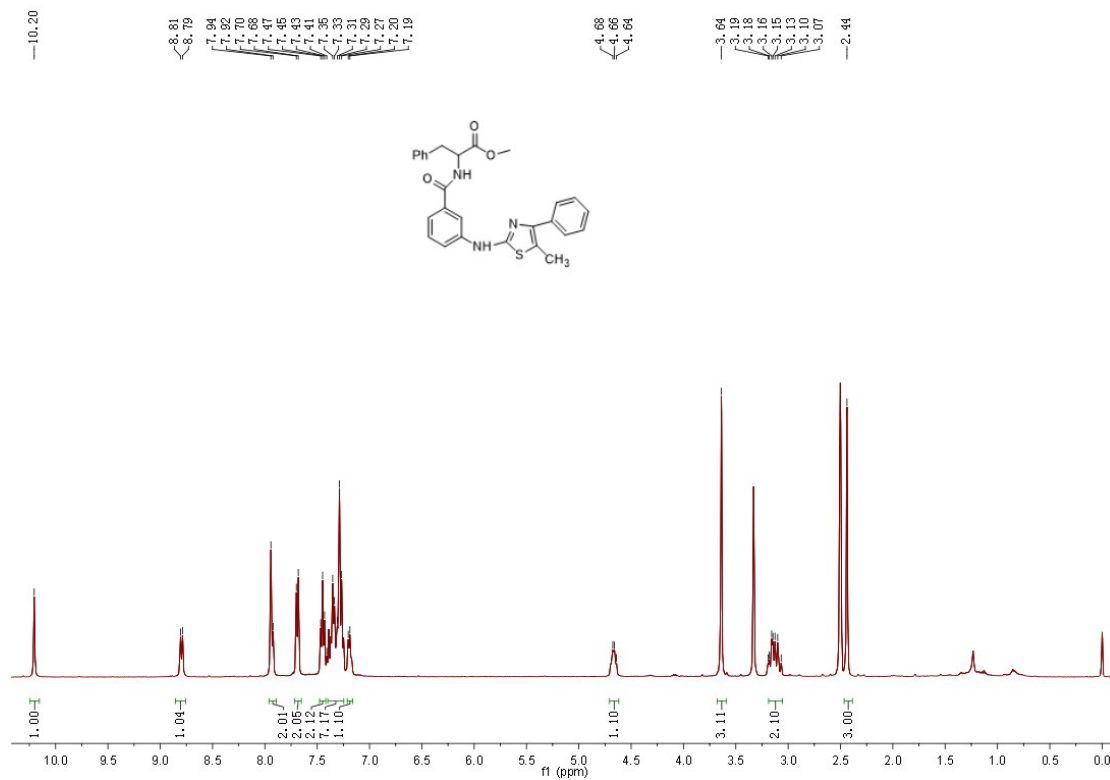
3m



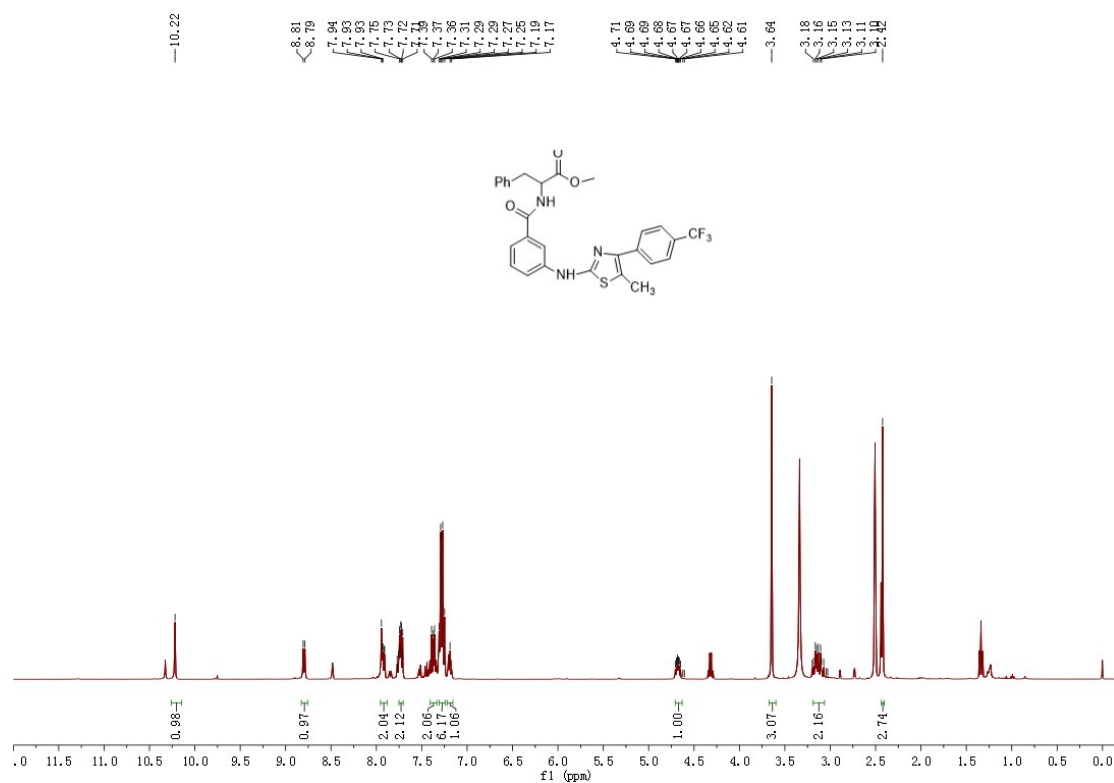
3n



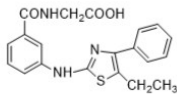
30



3p



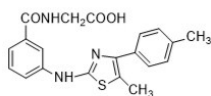
4a



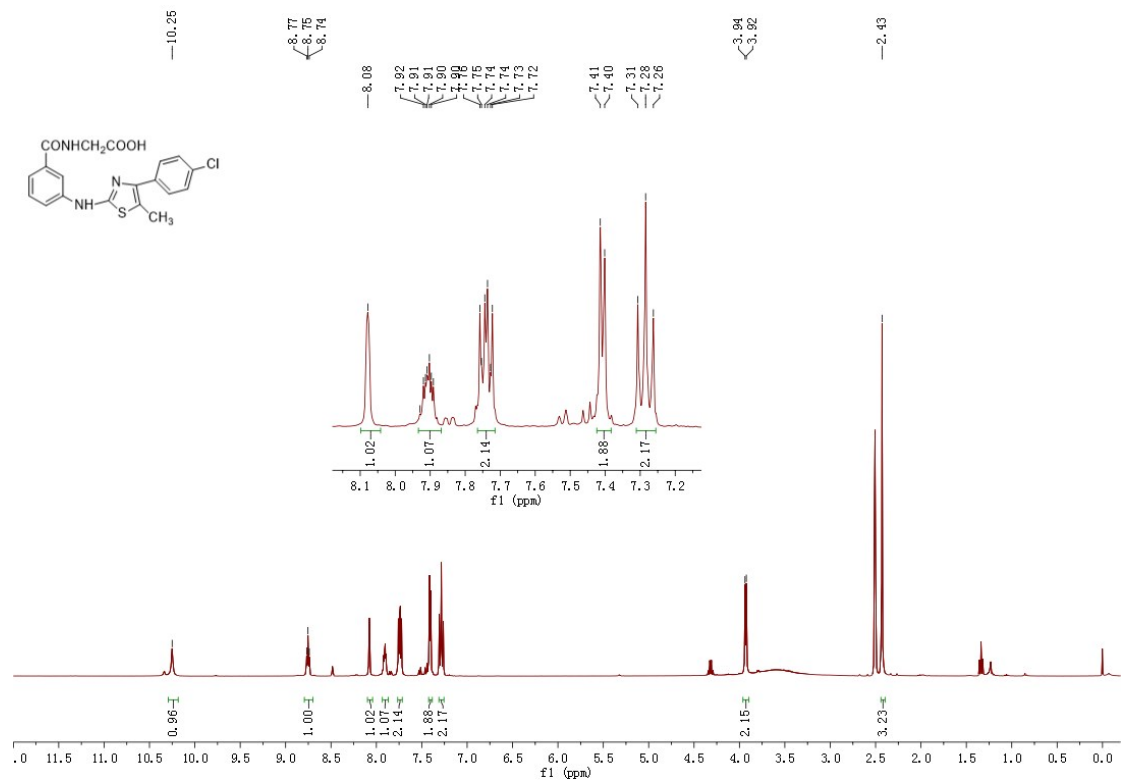
Chemical structure of 4-(4-methylphenyl)-2-((3-(4-oxo-4-oxopropyl)phenyl)hydrazono)thiazole is shown in the top left corner.

The ^1H NMR spectrum (DMSO- d_6) displays the following peaks and integrations:

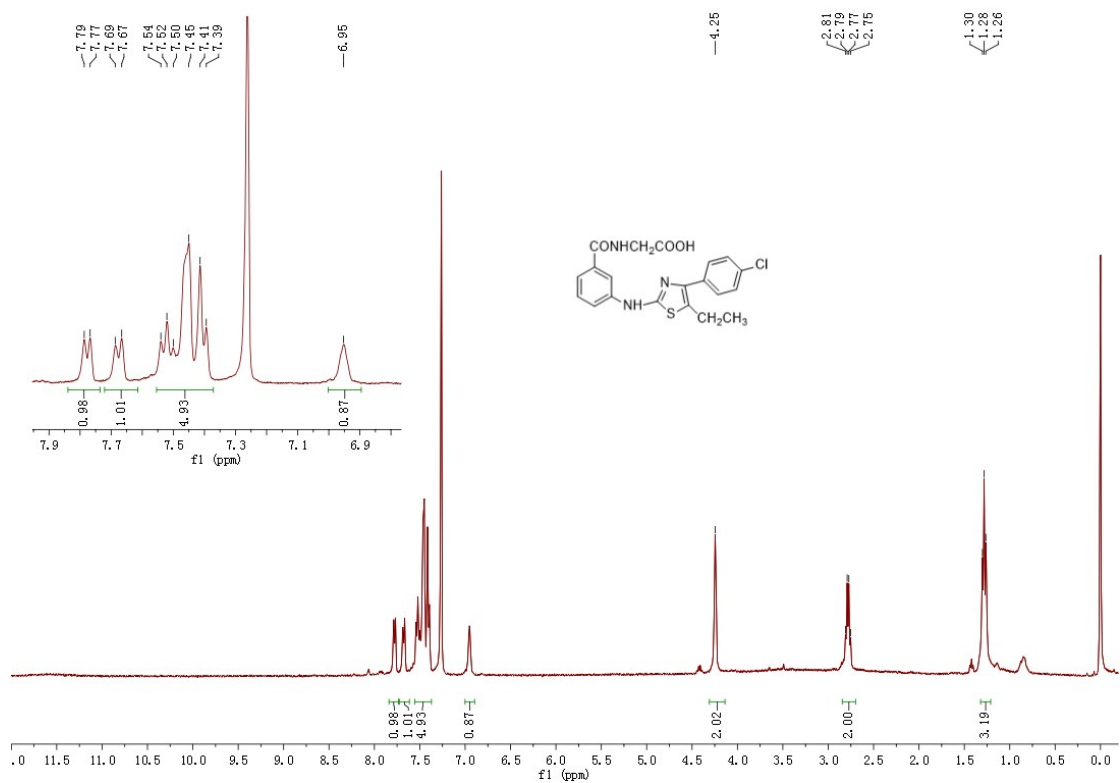
- 10.49 (s, 1H, integration 0.95)
- 8.84, 8.82, 8.81 (m, 3H, integration 0.92)
- 8.08, 8.07, 7.92, 7.91, 7.90, 7.90, 7.89 (m, 6H, integration 2.08)
- 7.58, 7.56, 7.47, 7.45, 7.44, 7.43, 7.42, 7.38, 7.26 (m, 8H, integration 2.15)
- 3.94, 3.92 (s, 2H, integration 2.06)
- 2.41, 2.36 (s, 3H, integration 3.00, 2.96)



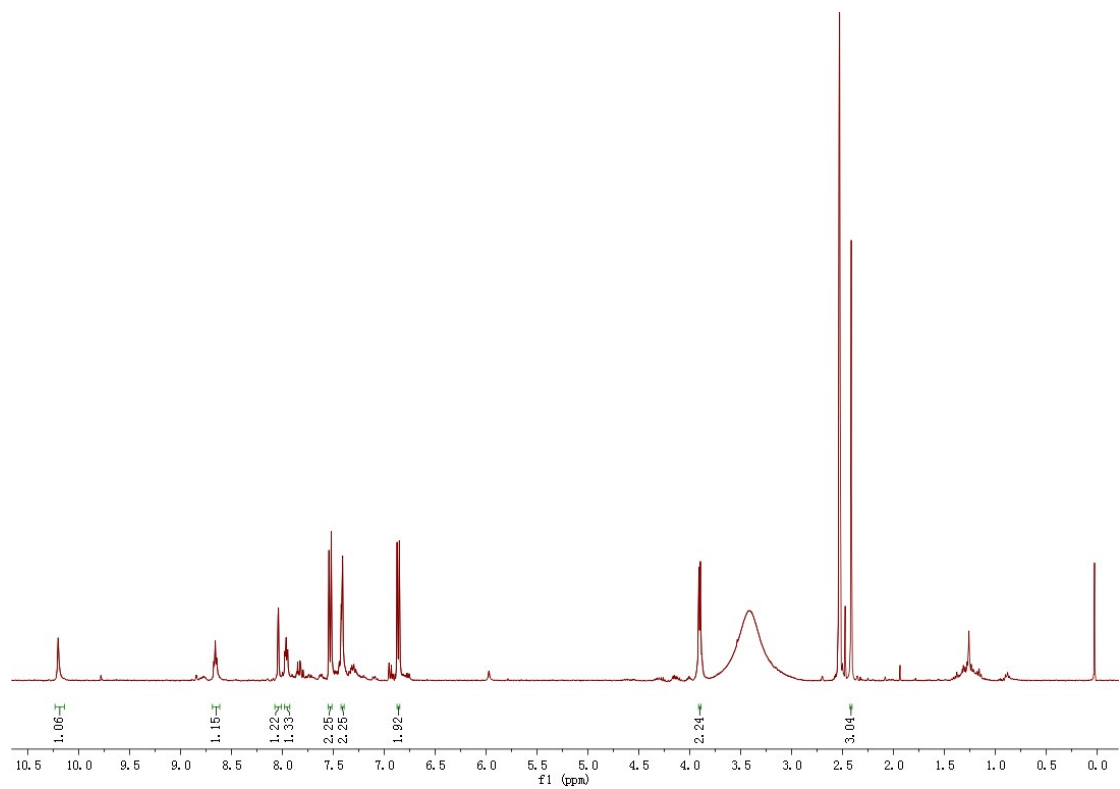
4c



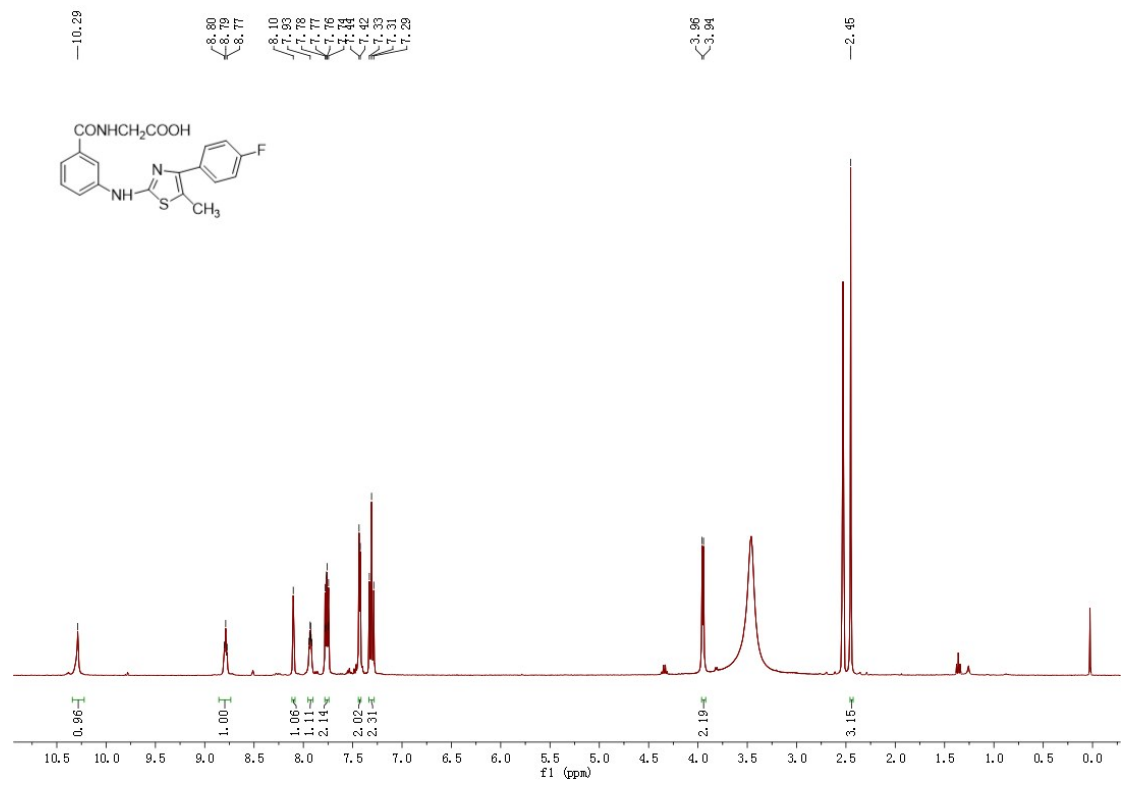
4d



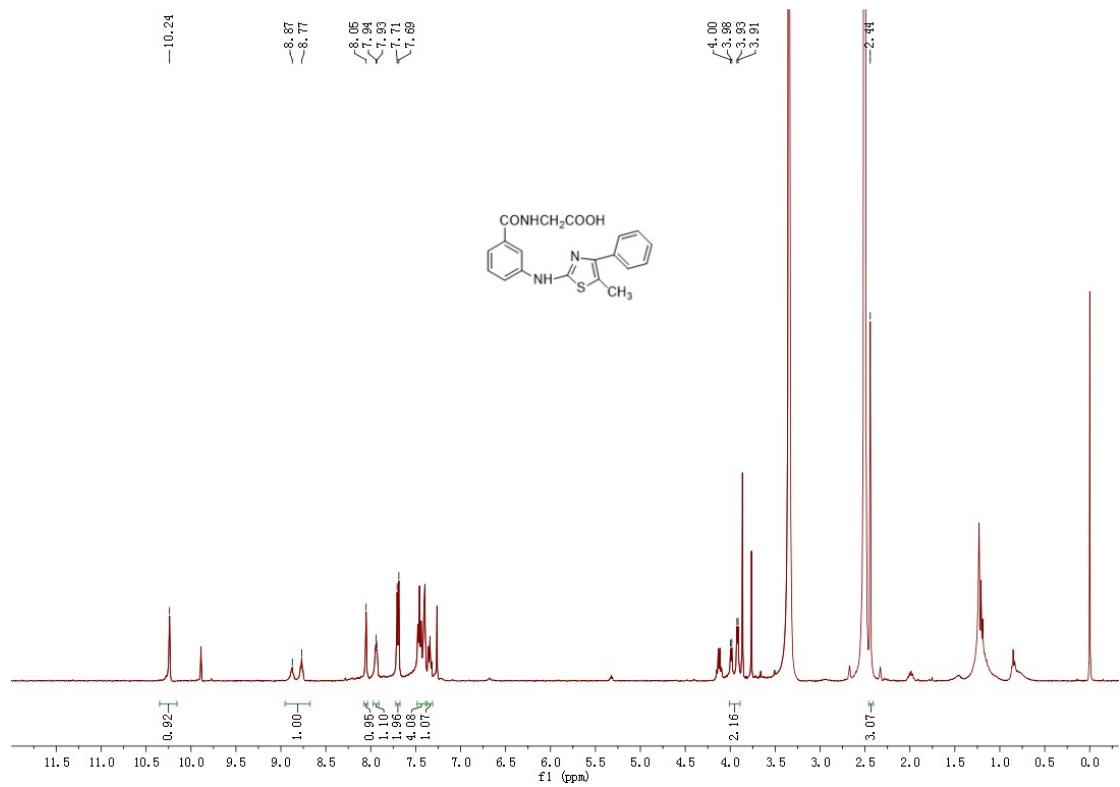
4e



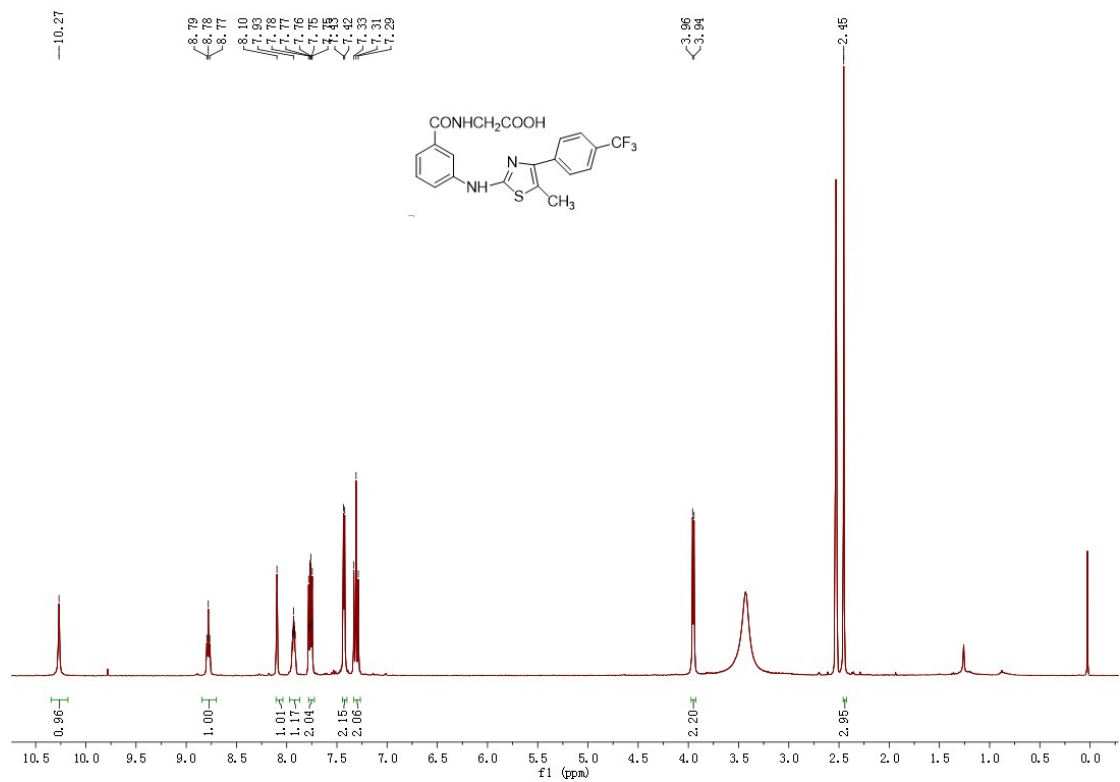
4f



4g

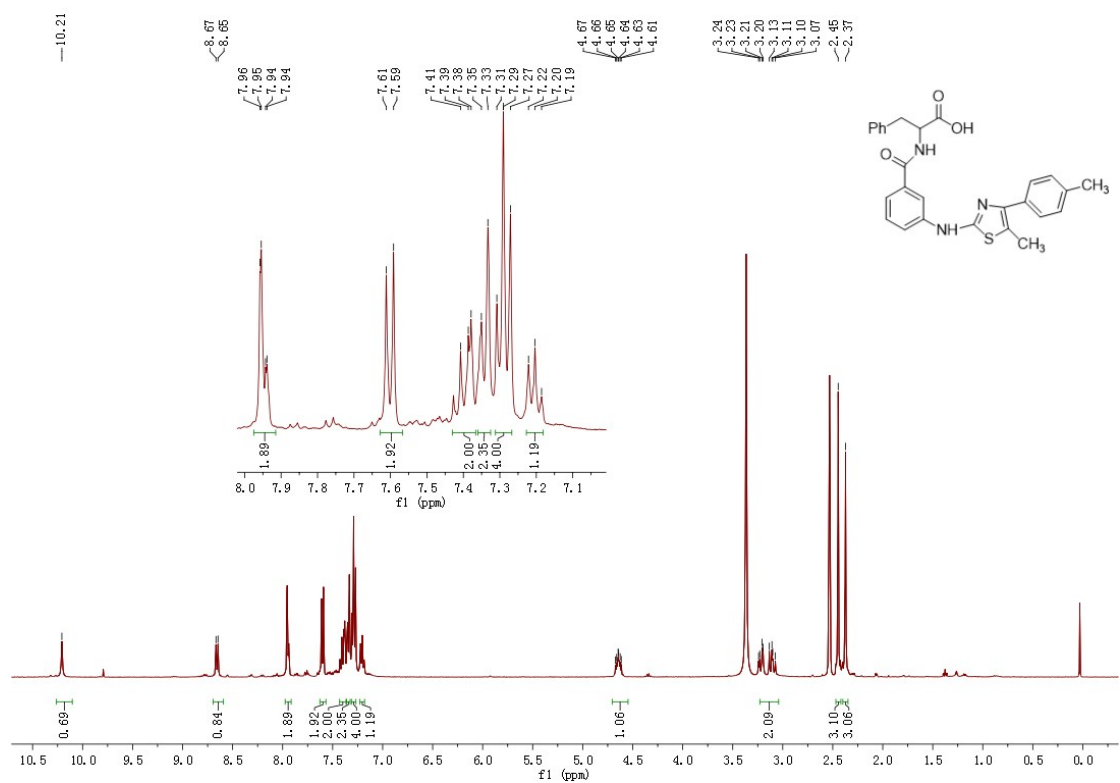


4h

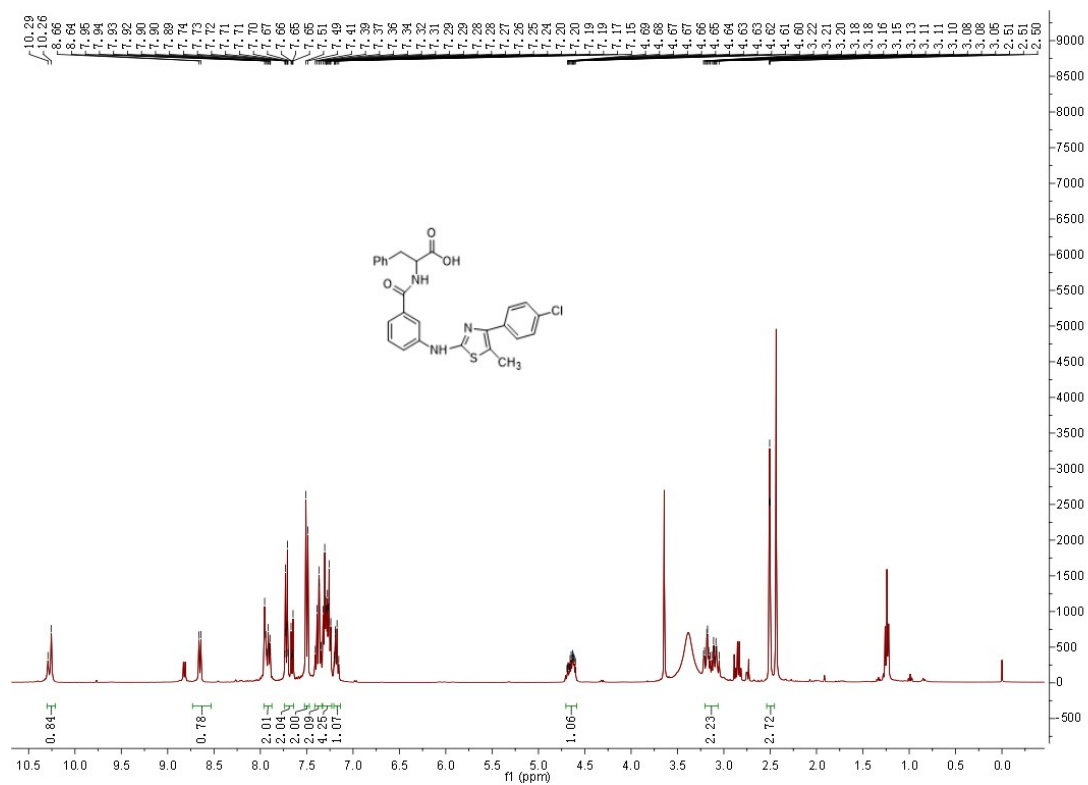


Chemical structure of compound 10: CCc1cc2sc(NC3=CC=C(C=C3)C(=O)NC(Cc4ccccc4)C(=O)O)c2cc1

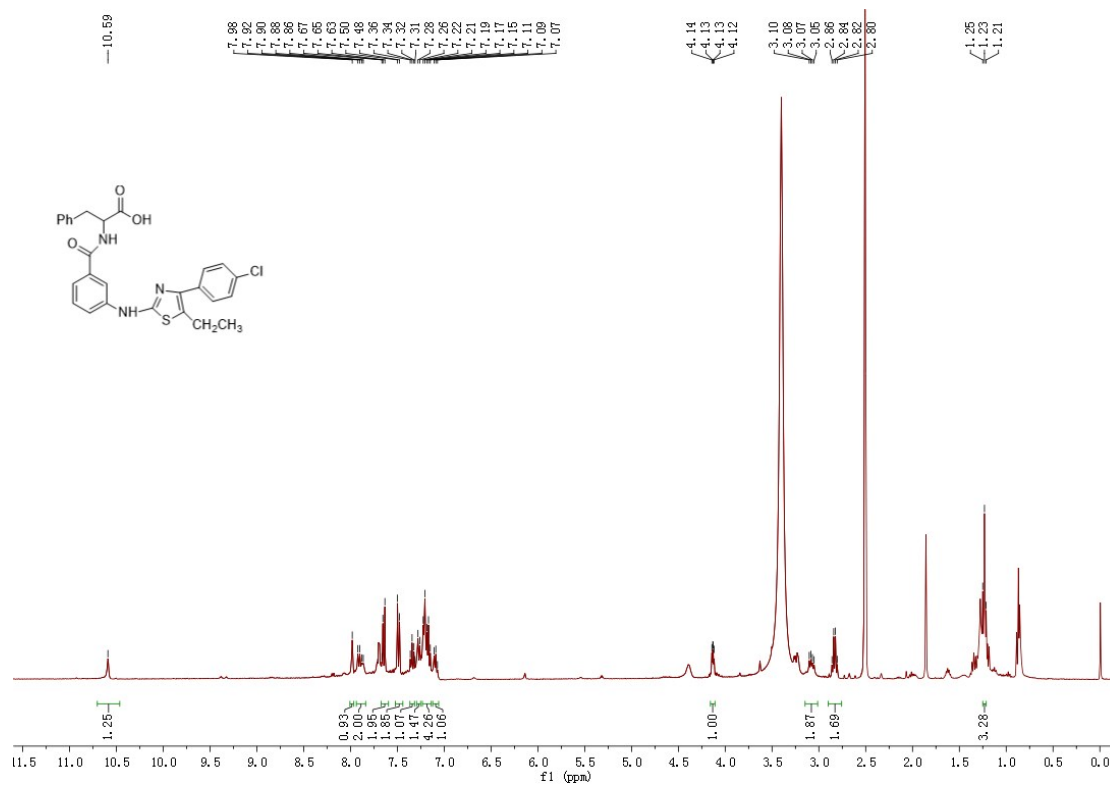
¹H NMR spectrum (CDCl₃) of compound 10. The spectrum shows peaks from 0.0 to 10.5 ppm. Key features include a broad peak at ~10.2 ppm (NH), aromatic signals between 7.1-8.0 ppm, a thiazole ring signal at ~7.3 ppm, a CH₂ group at ~2.5 ppm, and a CH₃ group at ~1.2 ppm. Integration values are shown below the peaks.



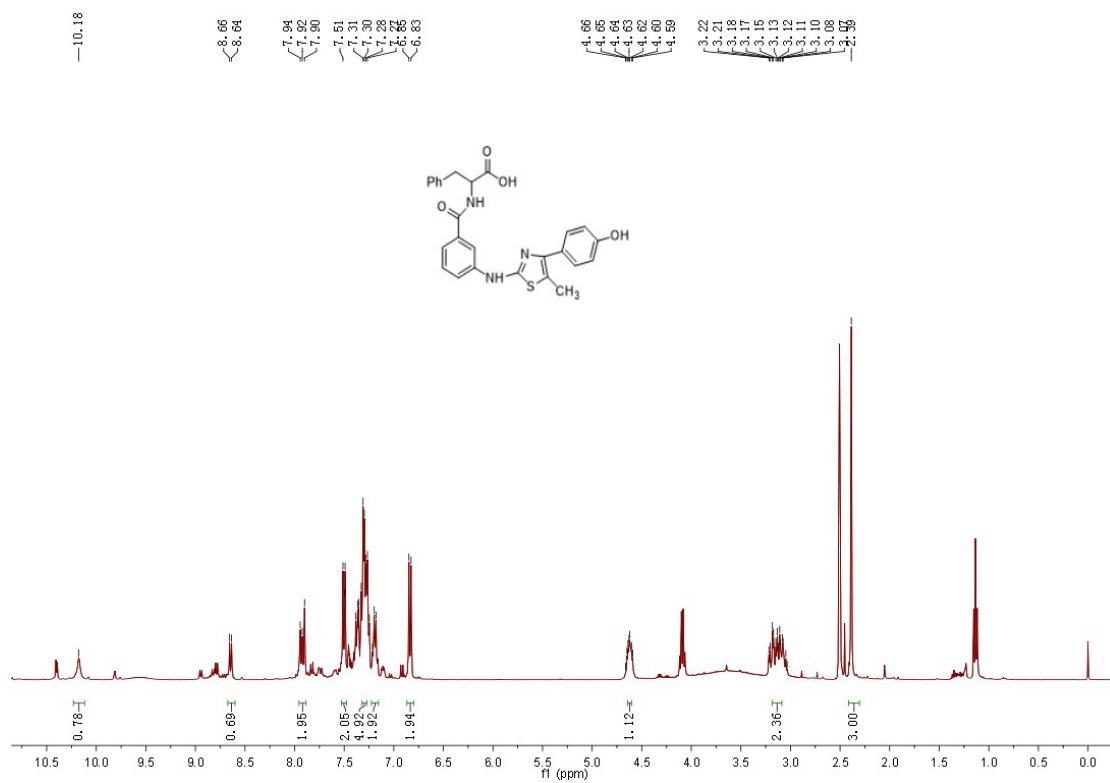
4k



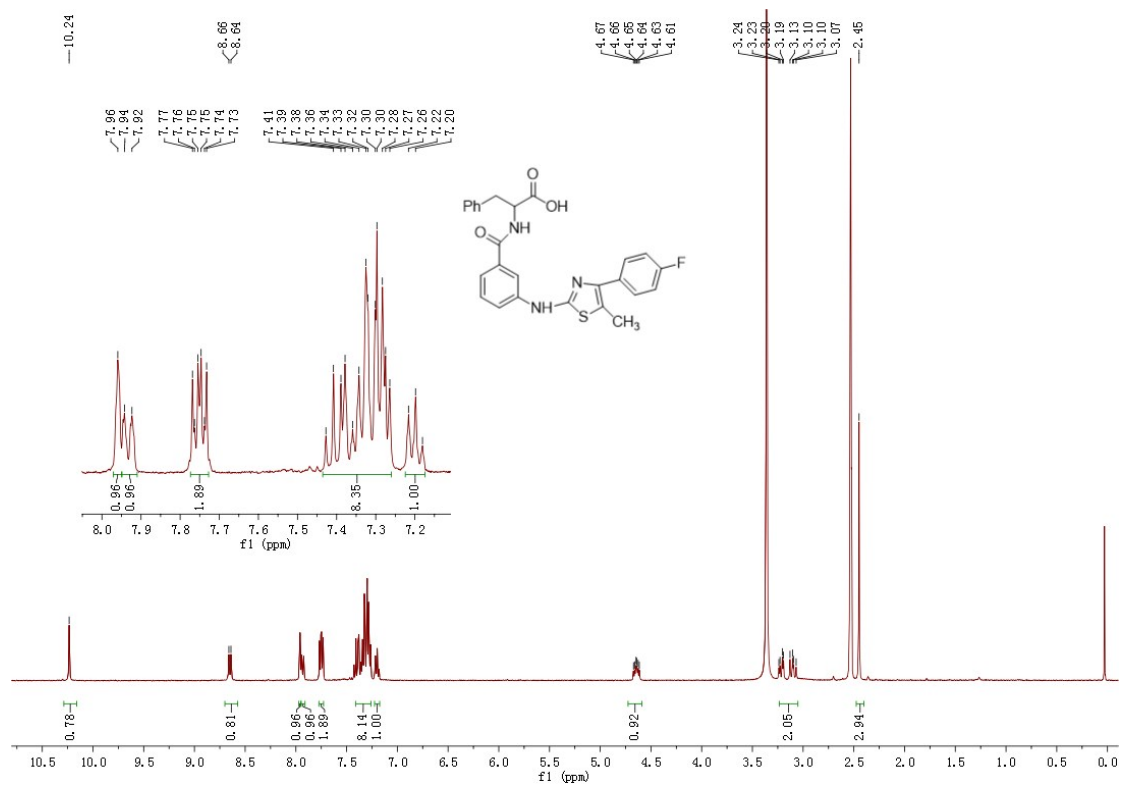
4l



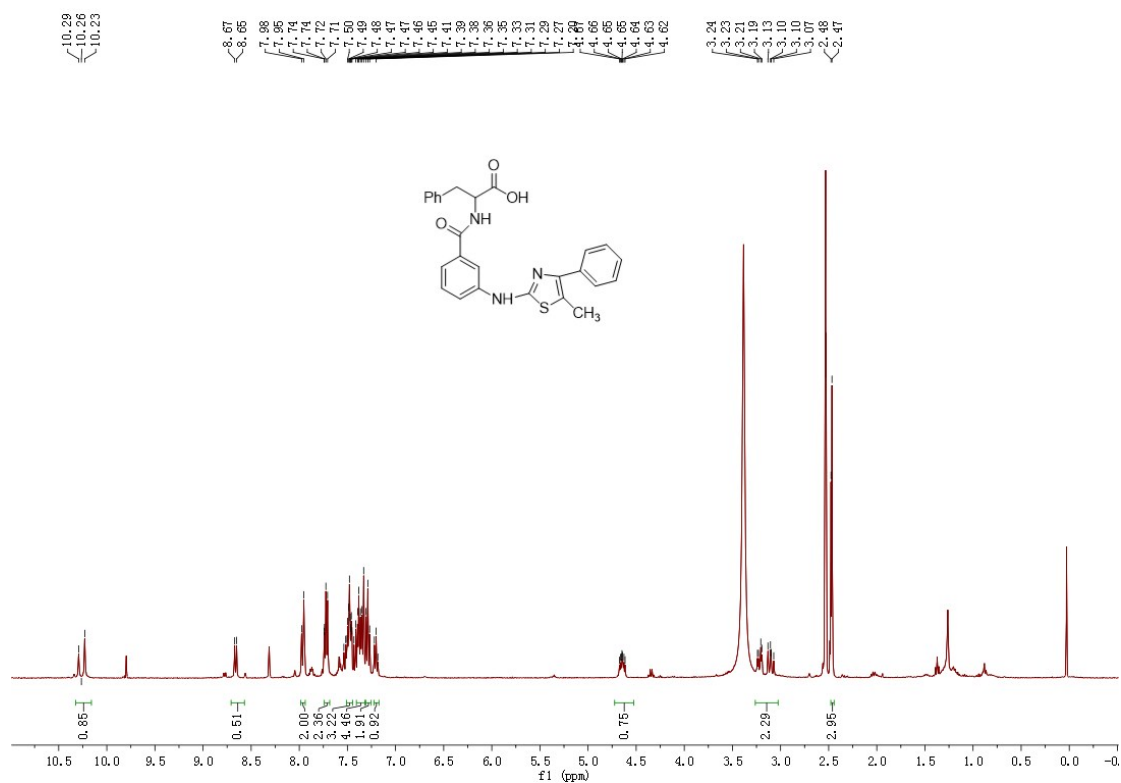
4m



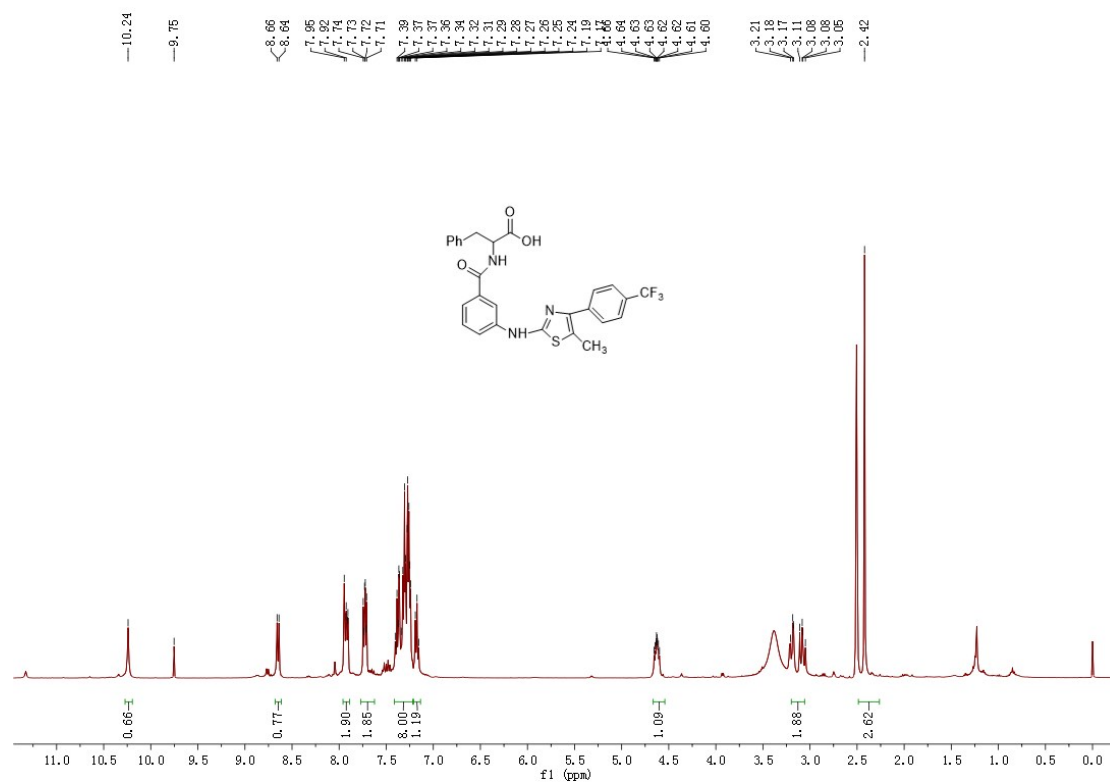
4n



4o

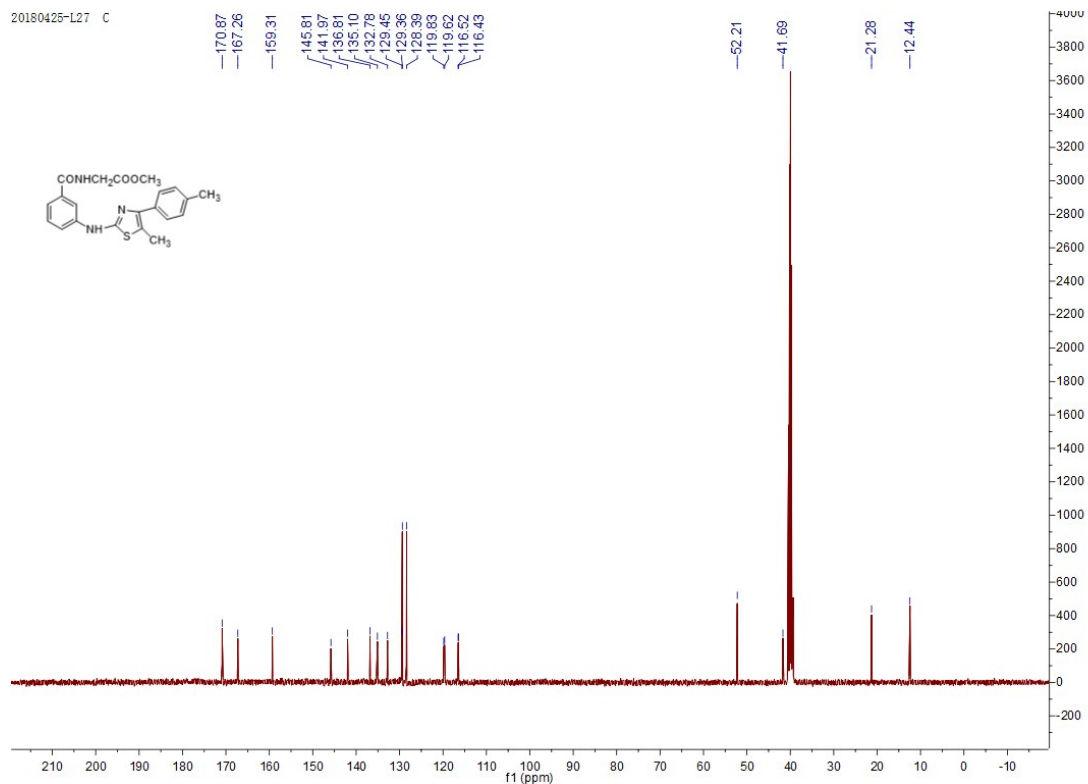


4p



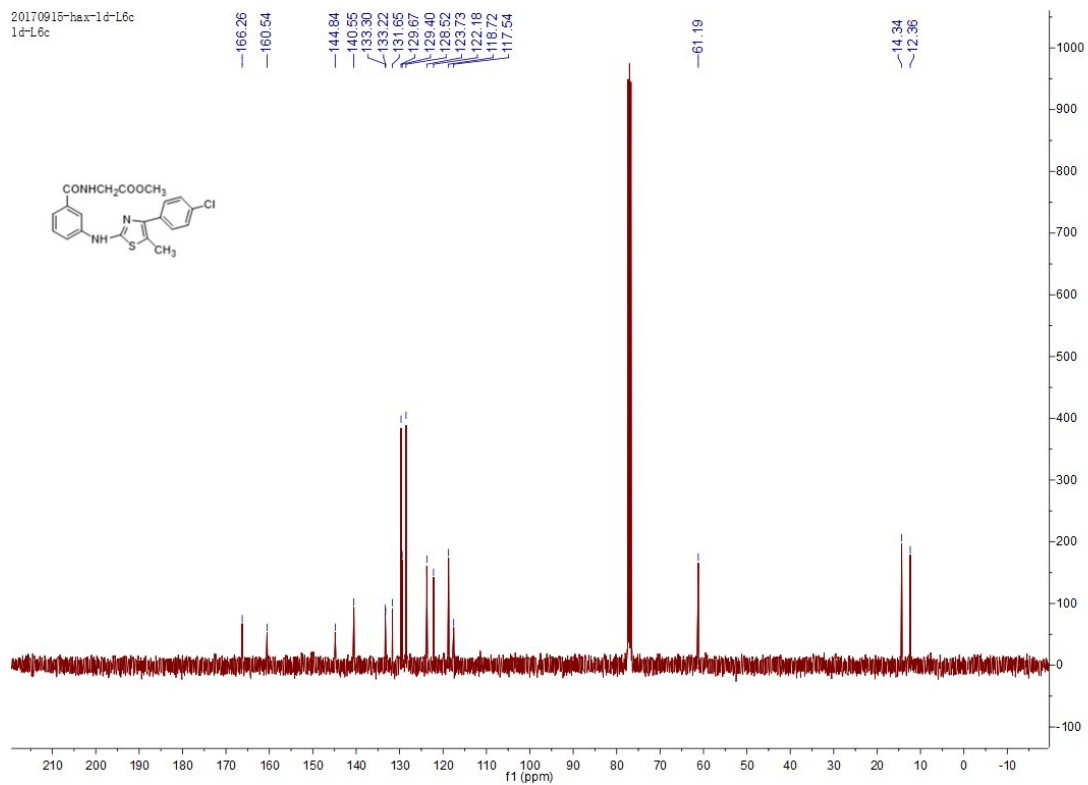
3b

20180425-L27 C

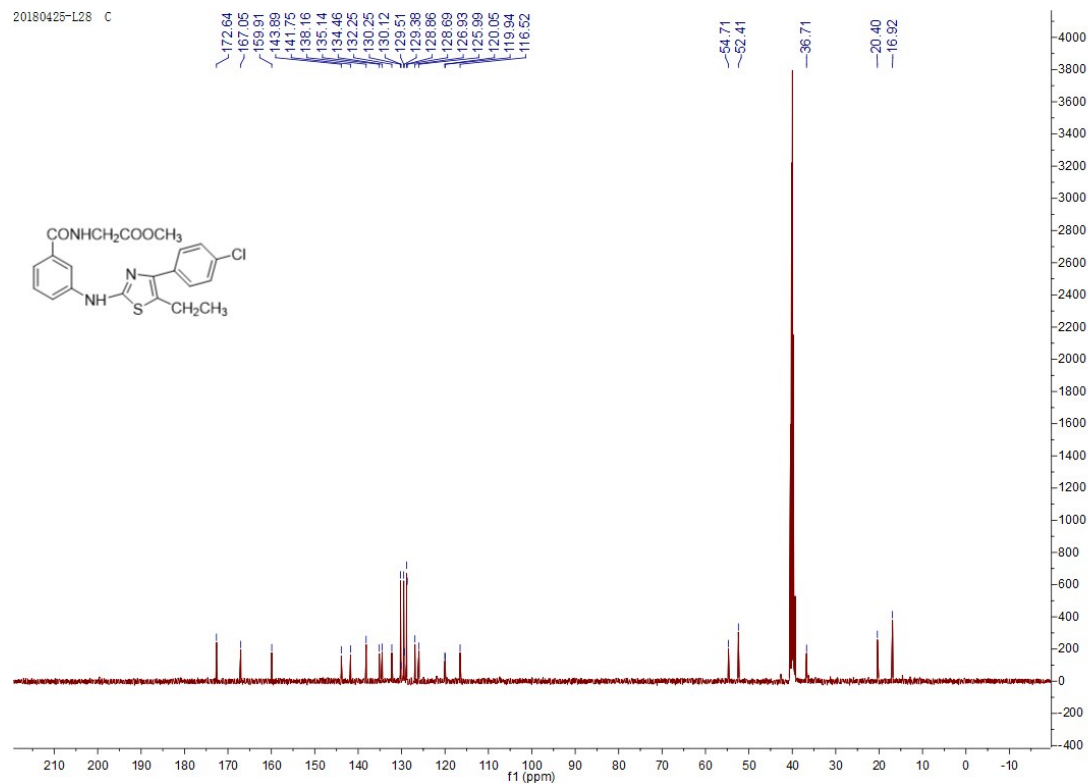


3c

20170915-hax-ld-L6c
1d-L6c



3d



3l

20180425-L31 C

