

Synthesis of novel 2-methyl-3-furyl sulfide flavor derivatives as efficient preservatives

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(A) Antimicrobial activity of 2-methyl-3-furyl sulfide flavor derivatives

Table 3. Inhibition zone of 2-methyl-3-furyl sulfide flavor derivatives on foodborne microorganisms. Inhibition zone values are represented as mean $\pm$ standard error (including the disc diameter, 6 mm). The means with different letters (^{a-i}) are significantly different ($p < 0.05$).

Compounds	Inhibition zone (mm)					
	Bacteria					
	EC	BS	SA	SP	LM	VP
3a	0	12.57 $\pm$ 0.29 ^{abc}	11.57 $\pm$ 0.29 ^c	11.03 $\pm$ 0.47 ^c	0	0
3b	8.77 $\pm$ 0.33 ^g	17.57 $\pm$ 0.29 ^{bcd}	8.57 $\pm$ 0.29 ^{ef}	19.90 $\pm$ 0.42 ^a	18.57 $\pm$ 0.29 ^a	0
3c	0	0	0	0	0	0
3d	8.90 $\pm$ 0.20 ^g	0	7.37 $\pm$ 0.13 ^g	8.43 $\pm$ 0.33 ^d	0	0
3e	12.03 $\pm$ 0.29 ^d	0	0	0	0	0
3f	0	7.23 $\pm$ 0.13 ^e	0	8.37 $\pm$ 0.13 ^d	7.57 $\pm$ 0.47 ^f	10.03 $\pm$ 0.29 ^b
3g	7.57 $\pm$ 0.29 ^h	9.23 $\pm$ 0.47 ^{cde}	10.23 $\pm$ 0.47 ^d	13.70 $\pm$ 0.42 ^b	13.57 $\pm$ 0.29 ^c	0
3h	0	8.70 $\pm$ 0.42 ^{cde}	7.57 $\pm$ 0.29 ^g	11.23 $\pm$ 0.47 ^c	9.70 $\pm$ 0.42 ^d	9.70 $\pm$ 0.42 ^{bc}
3i	11.23 $\pm$ 0.47 ^e	16.77 $\pm$ 0.33 ^a	17.77 $\pm$ 0.33 ^a	8.70 $\pm$ 0.42 ^d	8.57 $\pm$ 0.29 ^e	0
3j	17.70 $\pm$ 0.20 ^b	13.43 $\pm$ 0.33 ^{ab}	18.17 $\pm$ 0.33 ^a	0	0	0
3k	0	0	7.57 $\pm$ 0.29 ^g	0	0	0
3l	19.57 $\pm$ 0.29 ^a	14.57 $\pm$ 0.29 ^{ab}	17.57 $\pm$ 0.29 ^a	11.90 $\pm$ 0.20 ^{bc}	10.03 $\pm$ 0.29 ^d	9.57 $\pm$ 0.29 ^c
3m	0	0	8.90 $\pm$ 0.20 ^e	0	0	0
5a	9.37 $\pm$ 0.13 ^g	0	7.83 $\pm$ 0.33 ^g	0	0	0
5b	0	0	0	0	0	0
5c	0	0	0	0	0	0
5d	0	0	0	0	0	0
5e	0	0	0	0	0	0
5f	9.77 $\pm$ 0.33 ^f	7.83 $\pm$ 0.33 ^{de}	8.57 $\pm$ 0.29 ^{ef}	7.57 $\pm$ 0.29 ^d	7.23 $\pm$ 0.13 ^f	0
Pcn	15.83 $\pm$ 0.33 ^c	15.37 $\pm$ 0.13 ^{ab}	16.17 $\pm$ 0.33 ^b	17.83 $\pm$ 0.33 ^a	14.57 $\pm$ 0.29 ^b	12.37 $\pm$ 0.13 ^a

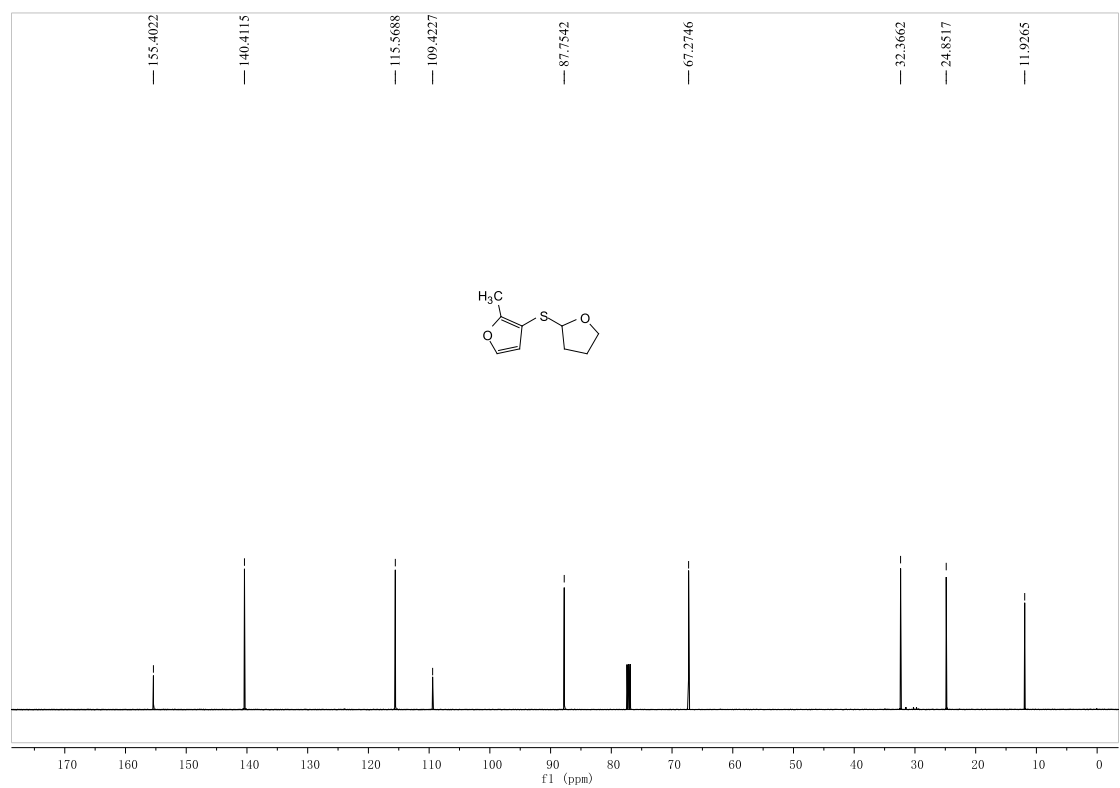
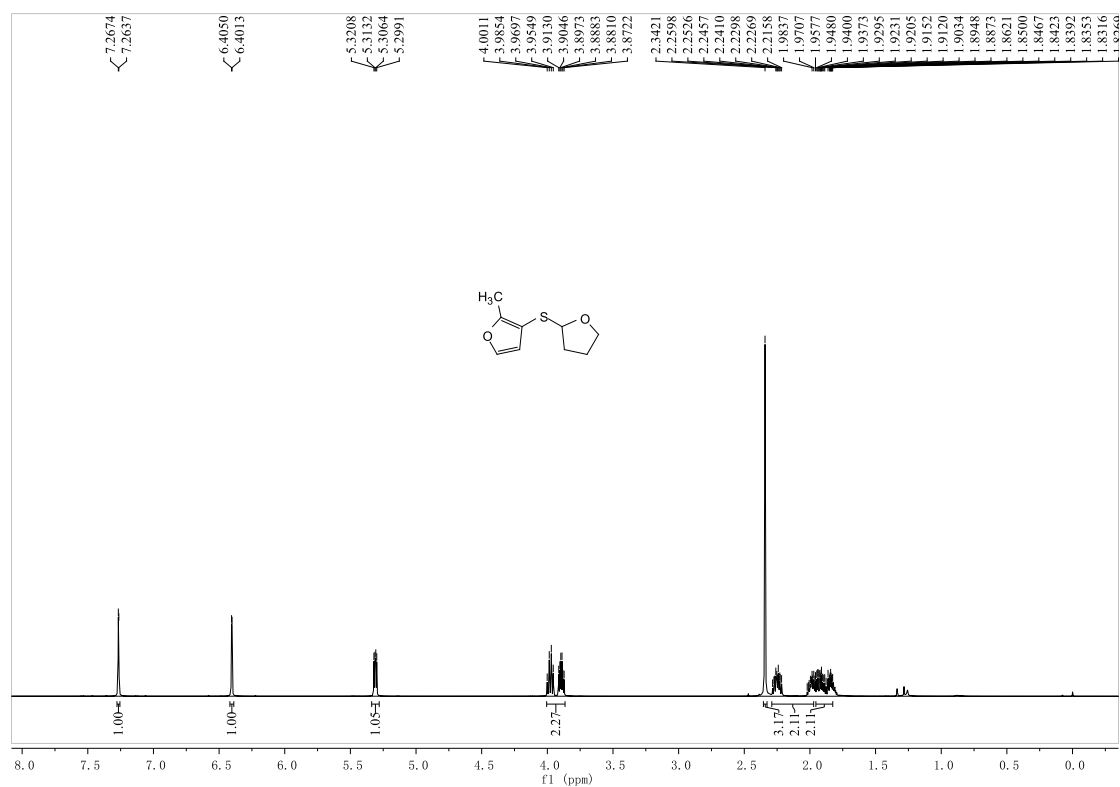
EC = *E. coli*. BS = *B. subtilis*. SA = *S. aureus*. SP = *S. paratyphi*. LM = *L. monocytogenes*. VP = *V. parahemolyticus*. Pcn = penicillin. The means with different letters (^{a-h}) are significantly different ($p < 0.05$).

Compounds	Inhibition zone (mm)			
	Fungi			
	PI	AN	MR	RO
3a	12.37 ± 0.13 ^f	13.70 ± 0.42 ^d	15.83 ± 0.33 ^f	11.03 ± 0.29 ^{cd}
3b	7.37 ± 0.13 ^h	0	9.17 ± 0.33 ⁱ	0
3c	0	0	0	0
3d	20.23 ± 0.47 ^b	0	20.57 ± 0.29 ^b	8.17 ± 0.33 ^{gh}
3e	21.57 ± 0.29 ^a	8.70 ± 0.42 ^f	0	7.70 ± 0.42 ^h
3f	0	0	12.17 ± 0.33 ^g	0
3g	9.37 ± 0.47 ^g	17.70 ± 0.42 ^b	15.57 ± 0.29 ^f	8.77 ± 0.33 ^g
3h	13.37 ± 0.13 ^e	12.77 ± 0.33 ^e	18.03 ± 0.29 ^d	9.57 ± 0.29 ^f
3i	15.77 ± 0.33 ^d	18.70 ± 0.20 ^a	10.43 ± 0.33 ^h	19.70 ± 0.42 ^a
3j	0	9.23 ± 0.13 ^f	7.37 ± 0.13 ⁱ	10.70 ± 0.20 ^{de}
3k	0	0	7.70 ± 0.20 ⁱ	0
3l	9.23 ± 0.13 ^g	8.70 ± 0.20 ^f	21.43 ± 0.33 ^a	8.57 ± 0.29 ^g
3m	8.70 ± 0.42 ^g	7.83 ± 0.33 ^g	19.43 ± 0.33 ^c	11.57 ± 0.29 ^c
5a	0	0	0	0
5b	7.57 ± 0.29 ^h	0	0	0
5c	0	0	0	0
5d	0	0	0	0
5e	0	0	0	0
5f	14.03 ± 0.29 ^e	0	16.70 ± 0.20 ^e	10.17 ± 0.33 ^{ef}
Amb	17.83 ± 0.33 ^c	18.37 ± 0.13 ^a	18.57 ± 0.29 ^d	17.57 ± 0.29 ^b
TMTD	15.23 ± 0.13 ^d	15.70 ± 0.20 ^c	17.90 ± 0.20 ^d	17.37 ± 0.13 ^b

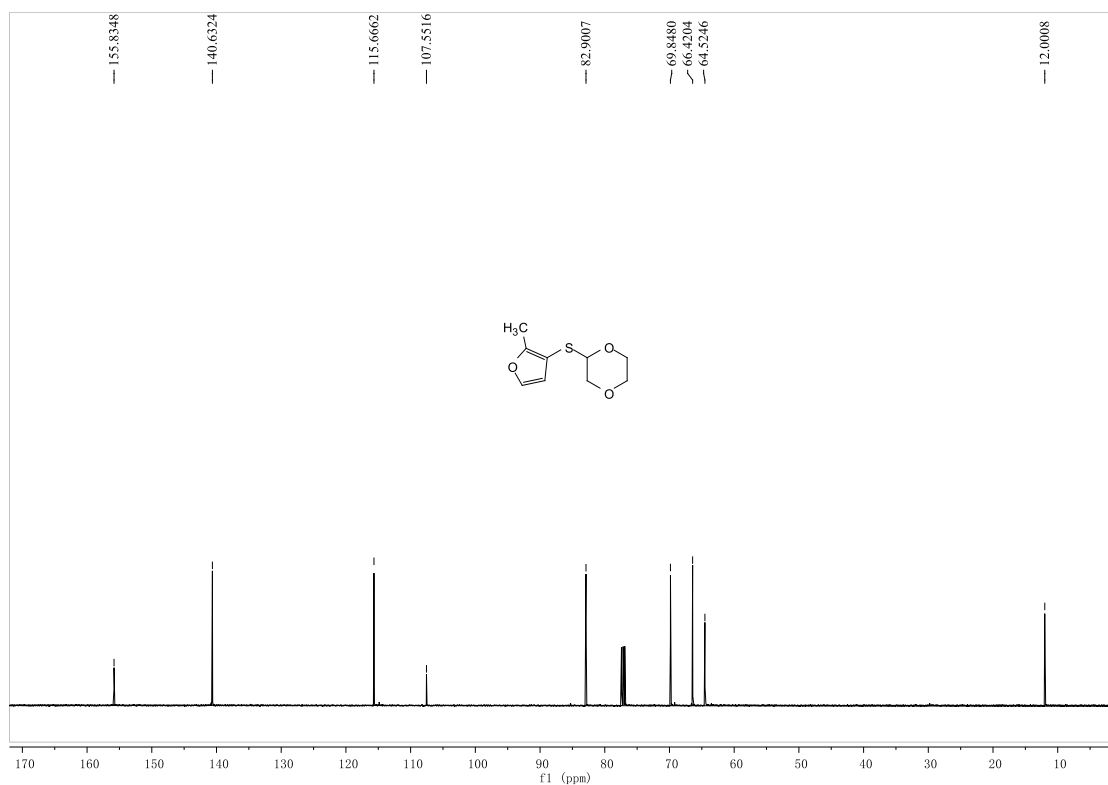
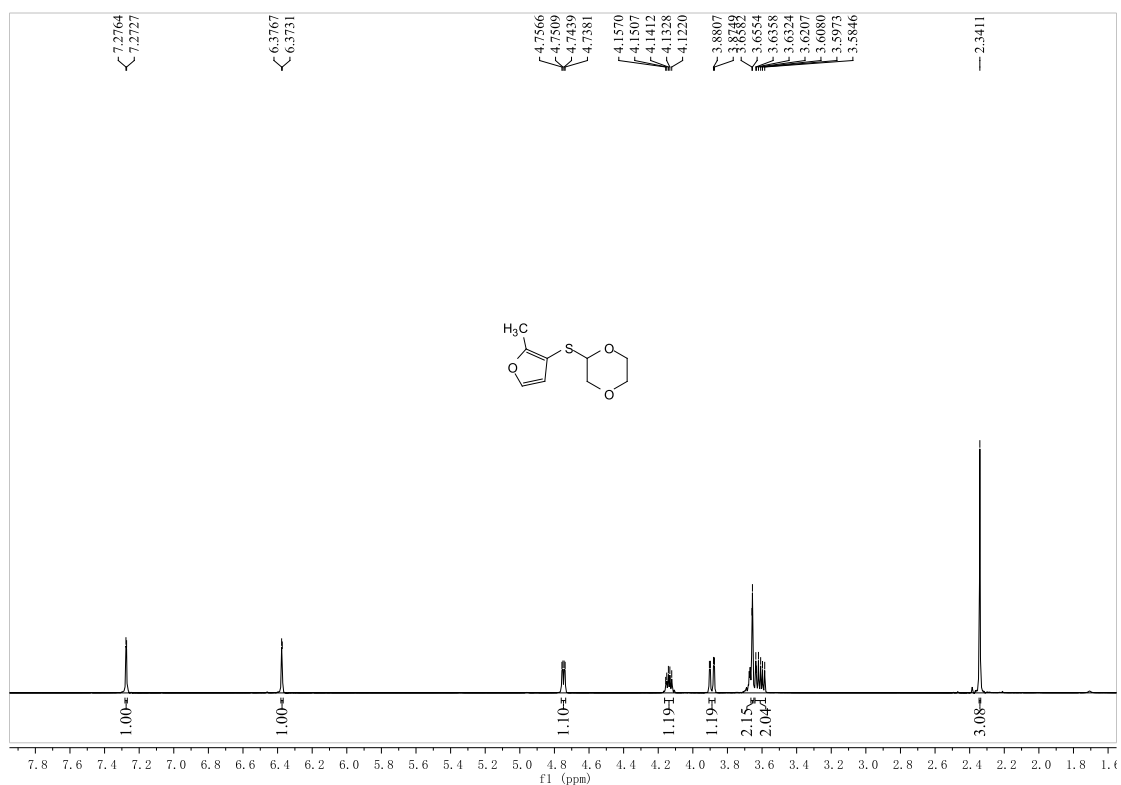
PI = *P. italicum*. **AN** = *A. niger*. **MR** = *M. racemosus*. **RO** = *R. oryzae*. **Pcn** = penicillin. **Amb** = amphotericin B. **TMTD** = thiram. The means with different letters (^{a-j}) are significantly different (p<0.05).

(B) NMR Spectra for products

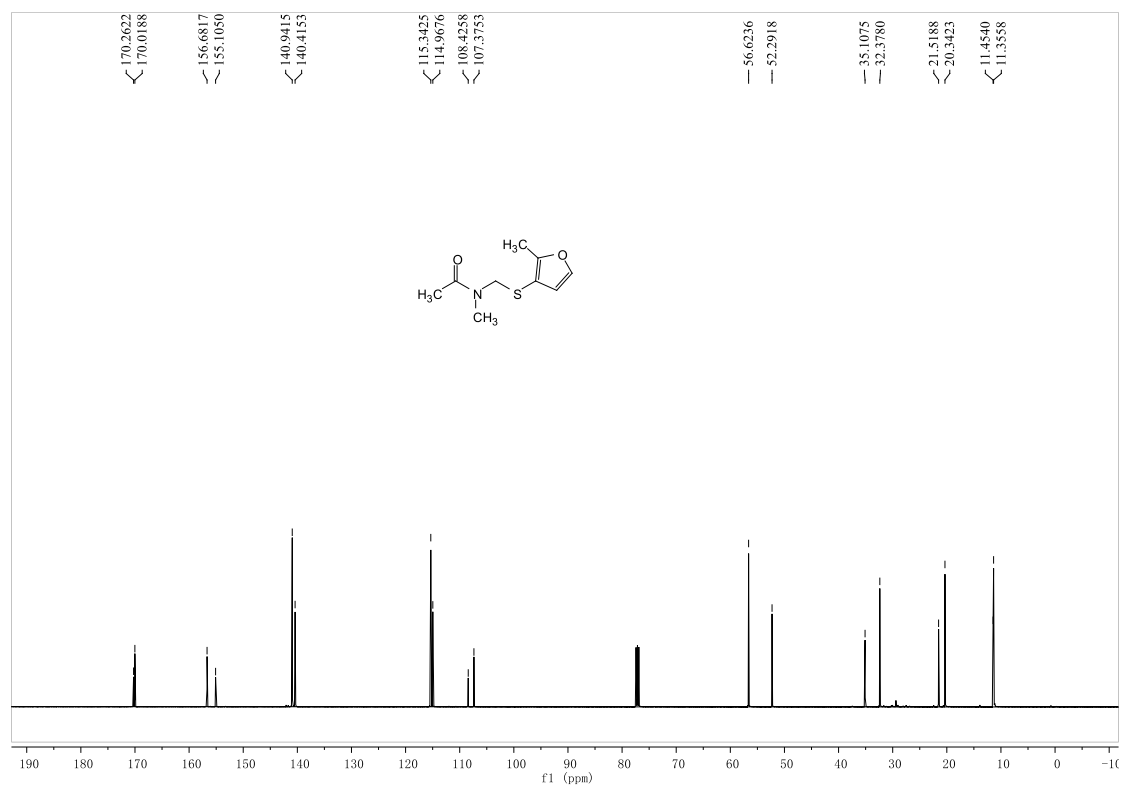
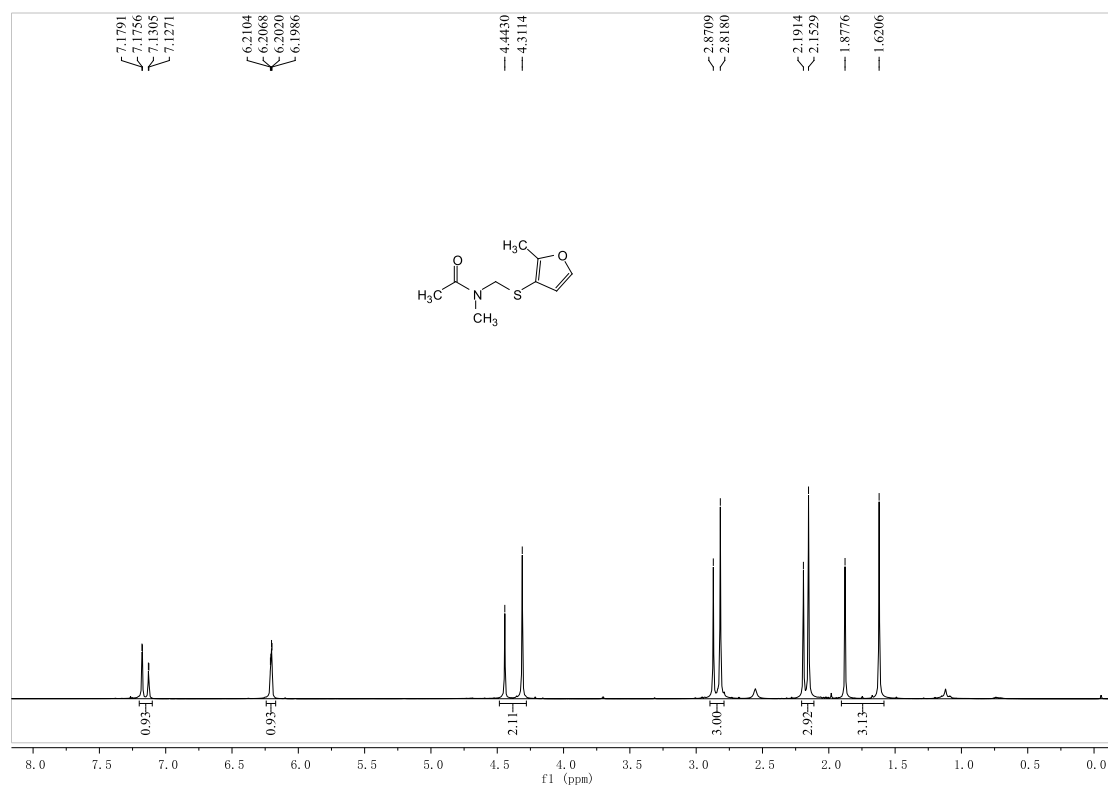
2-methyl-3-((tetrahydrofuran-2-yl)thio)furan (3a)



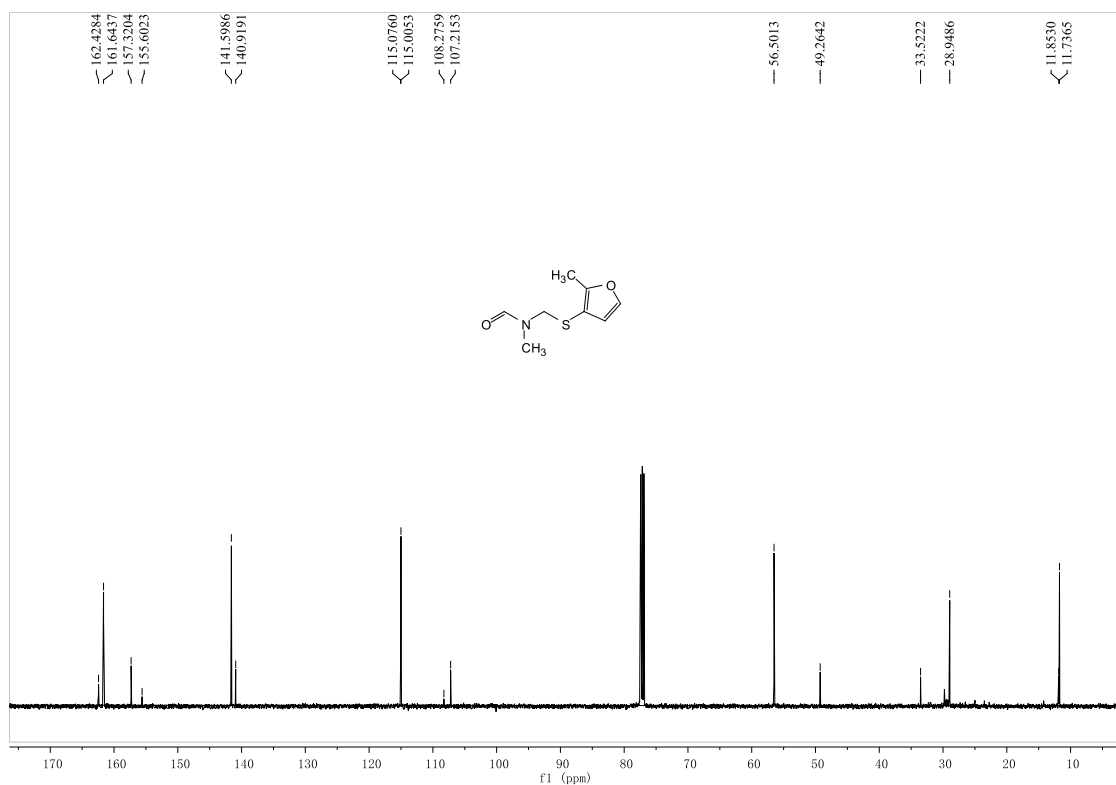
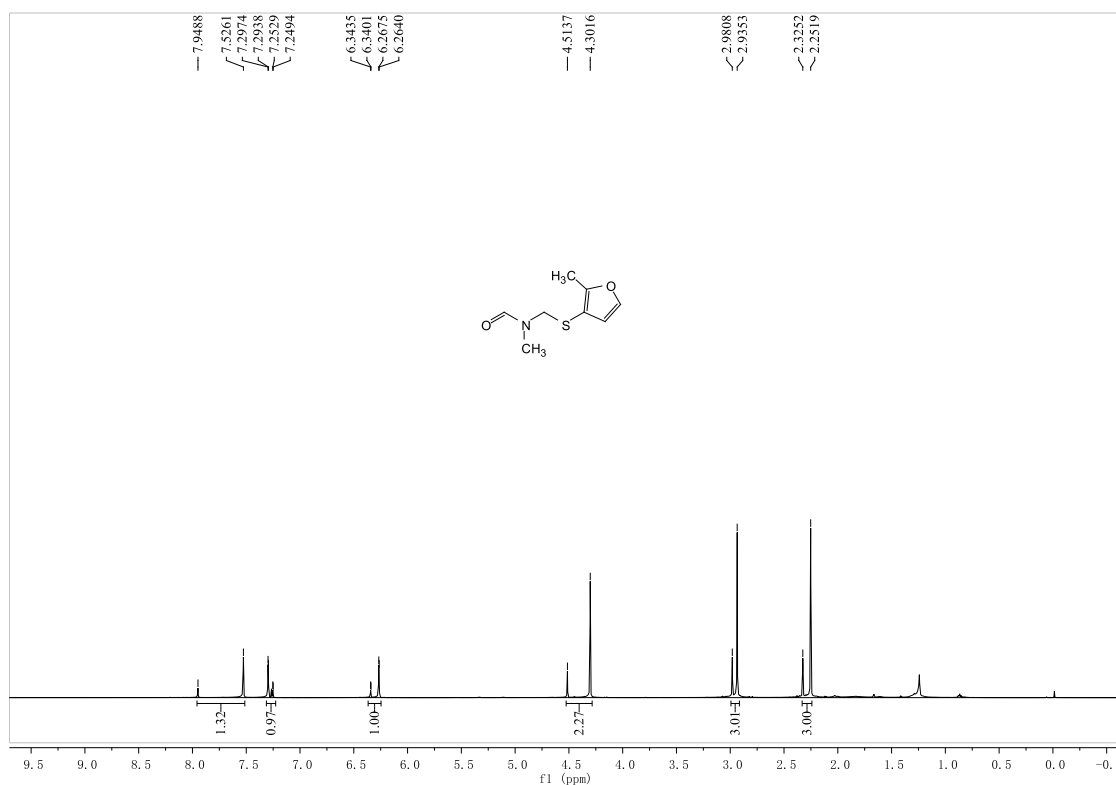
2-((2-methylfuran-3-yl)thio)-1,4-dioxane (3b)



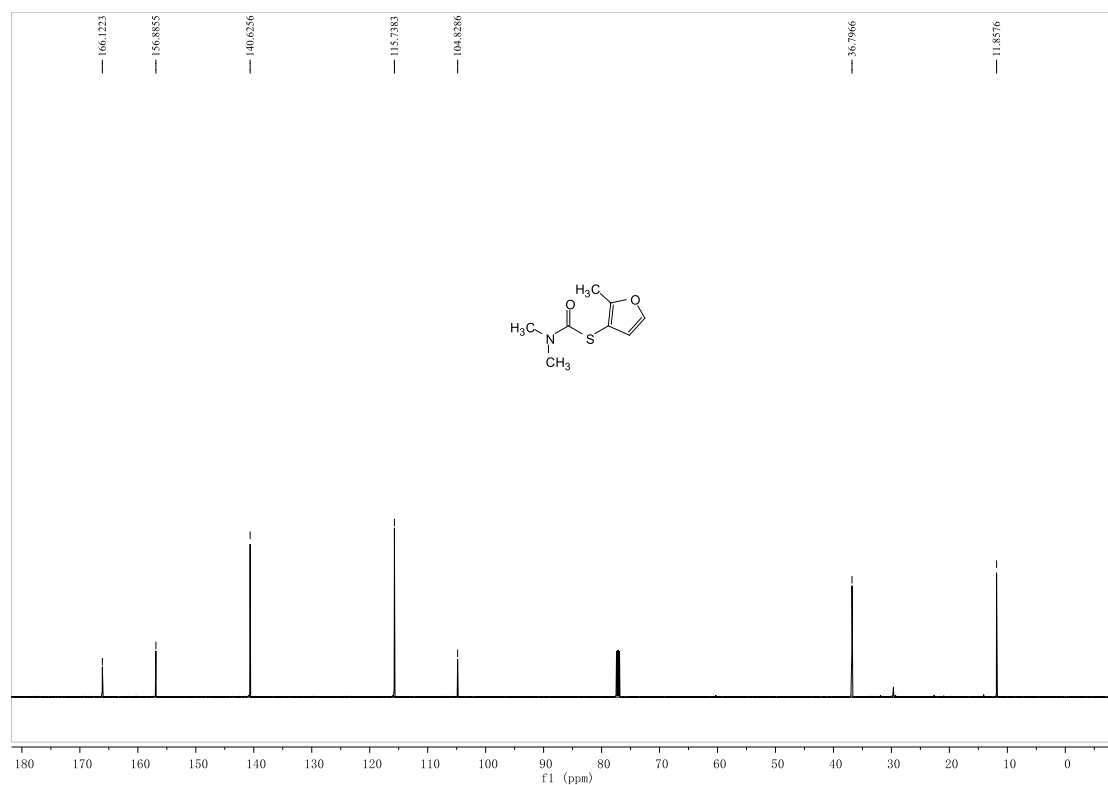
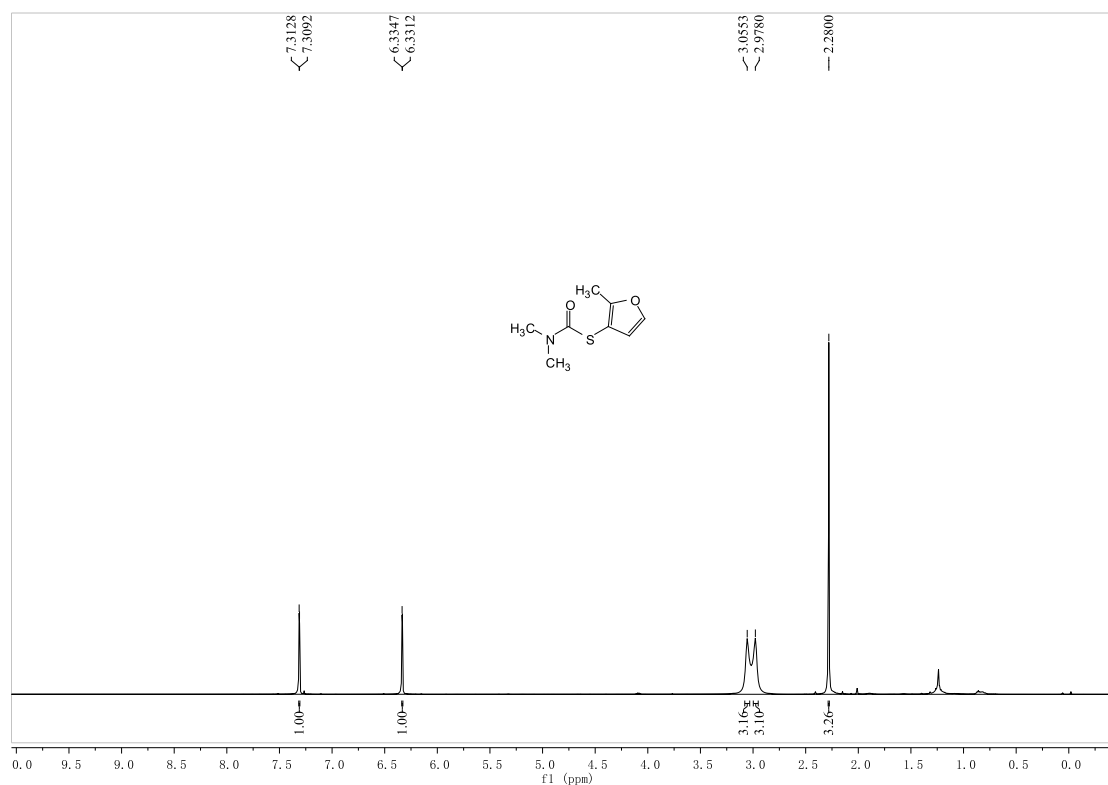
N-methyl-N-(((2-methylfuran-3-yl)thio)methyl)acetamide (3c)



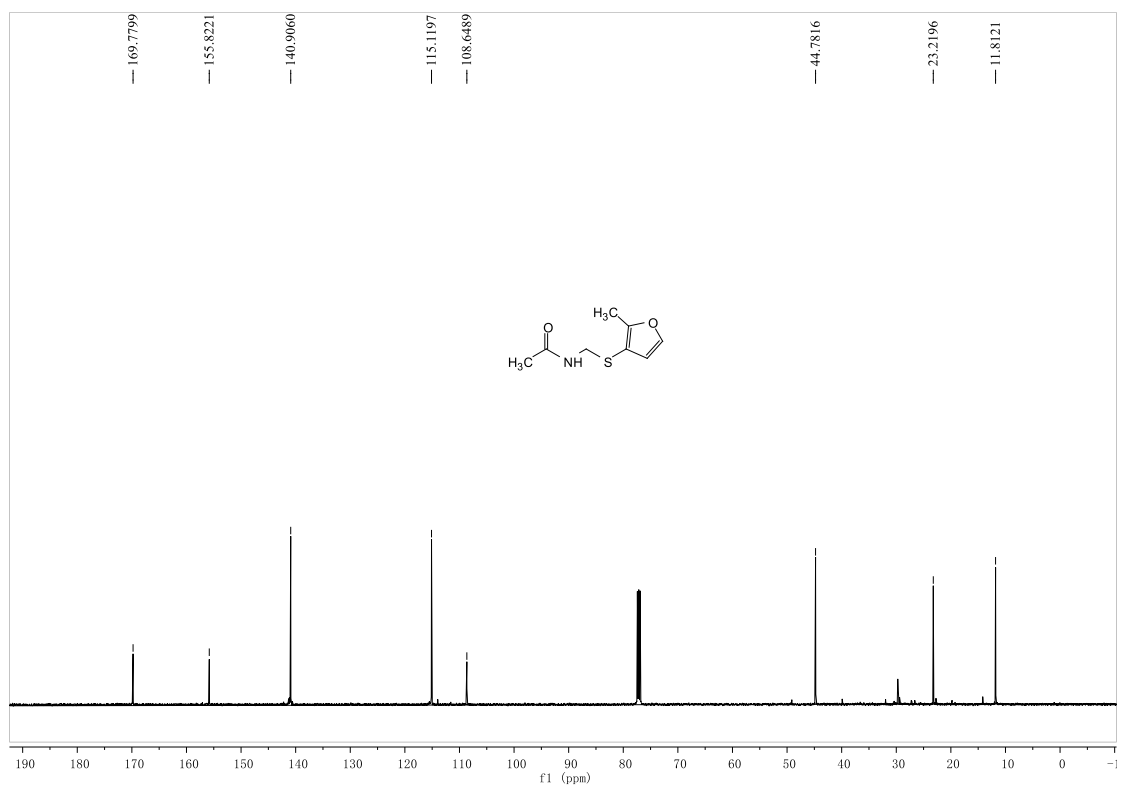
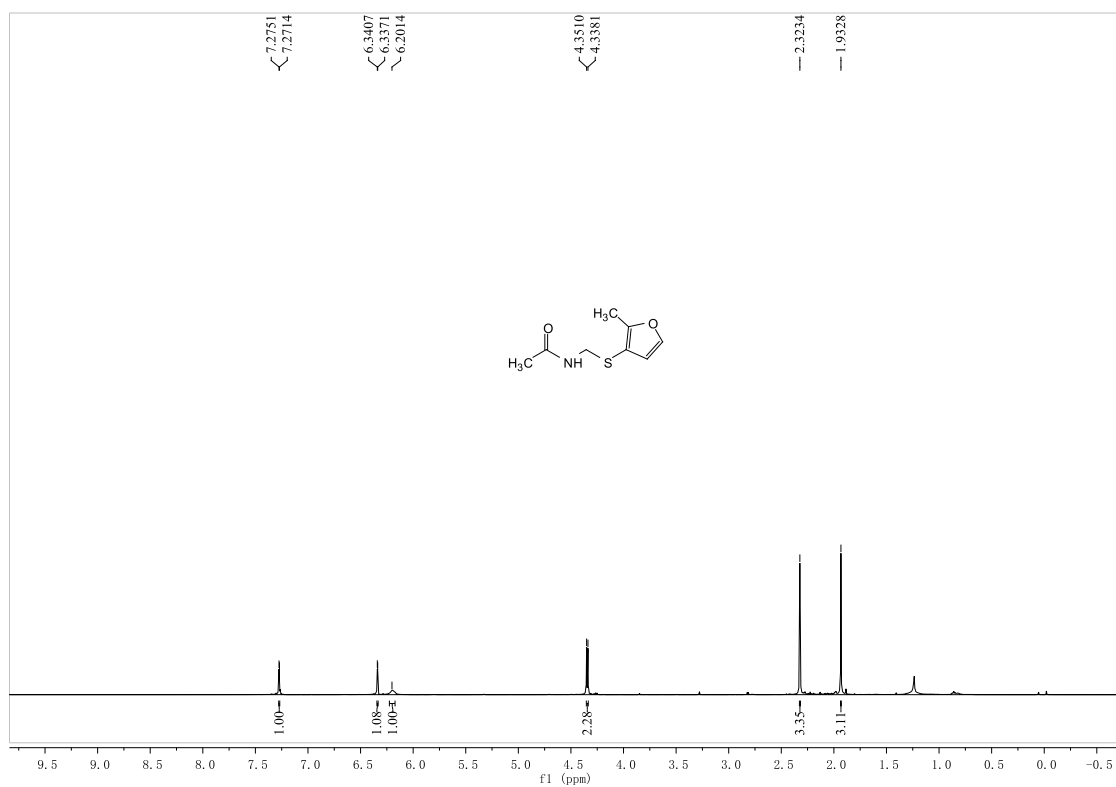
N-methyl-N-(((2-methylfuran-3-yl)thio)methyl)formamide (3d)



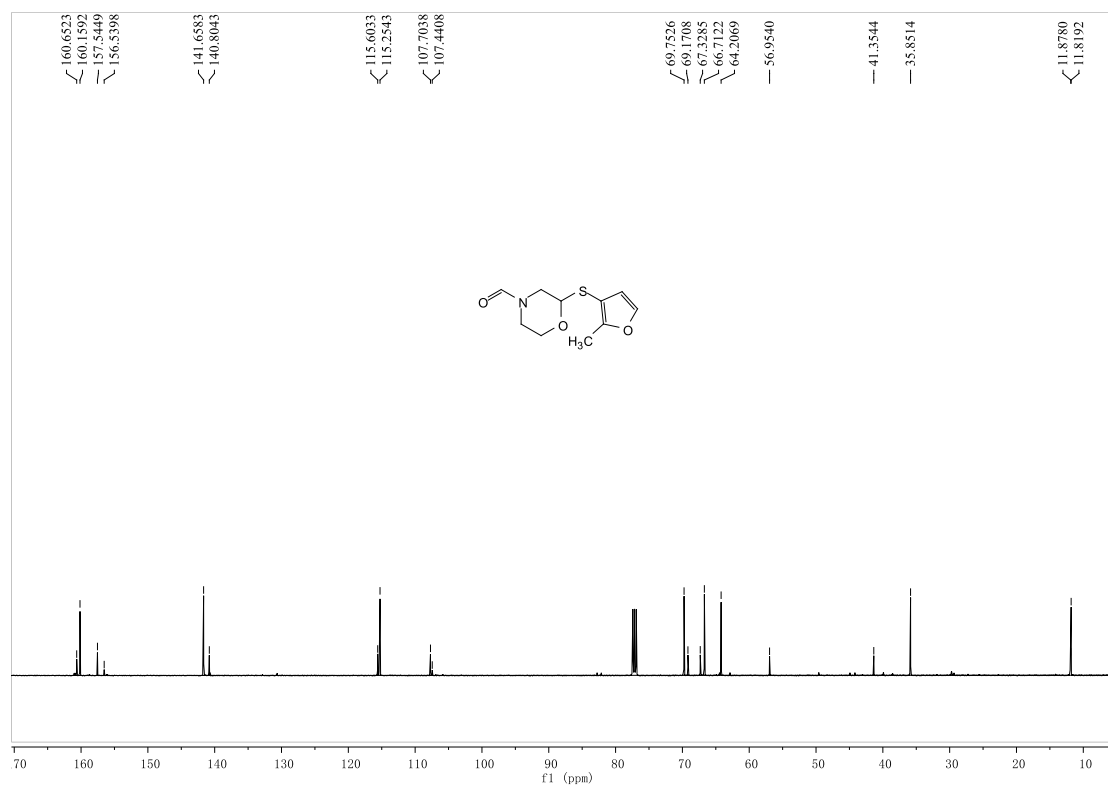
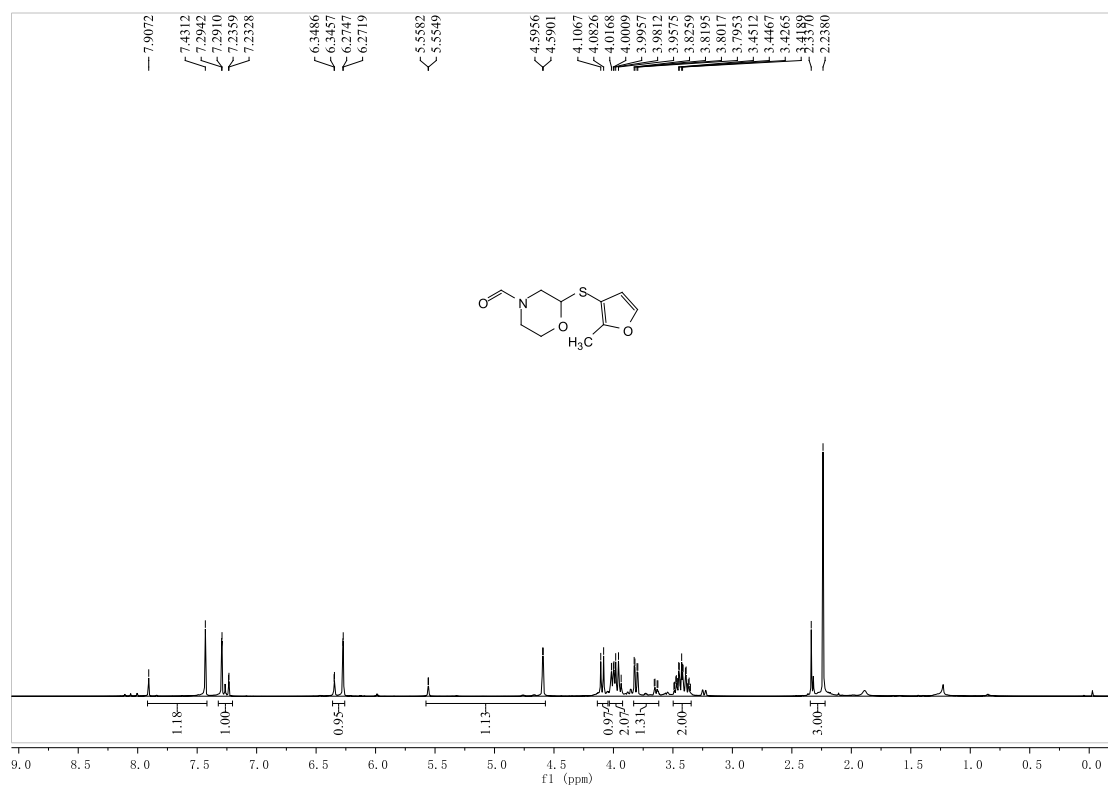
S-(2-methylfuran-3-yl)dimethylcarbamothioate (3e)



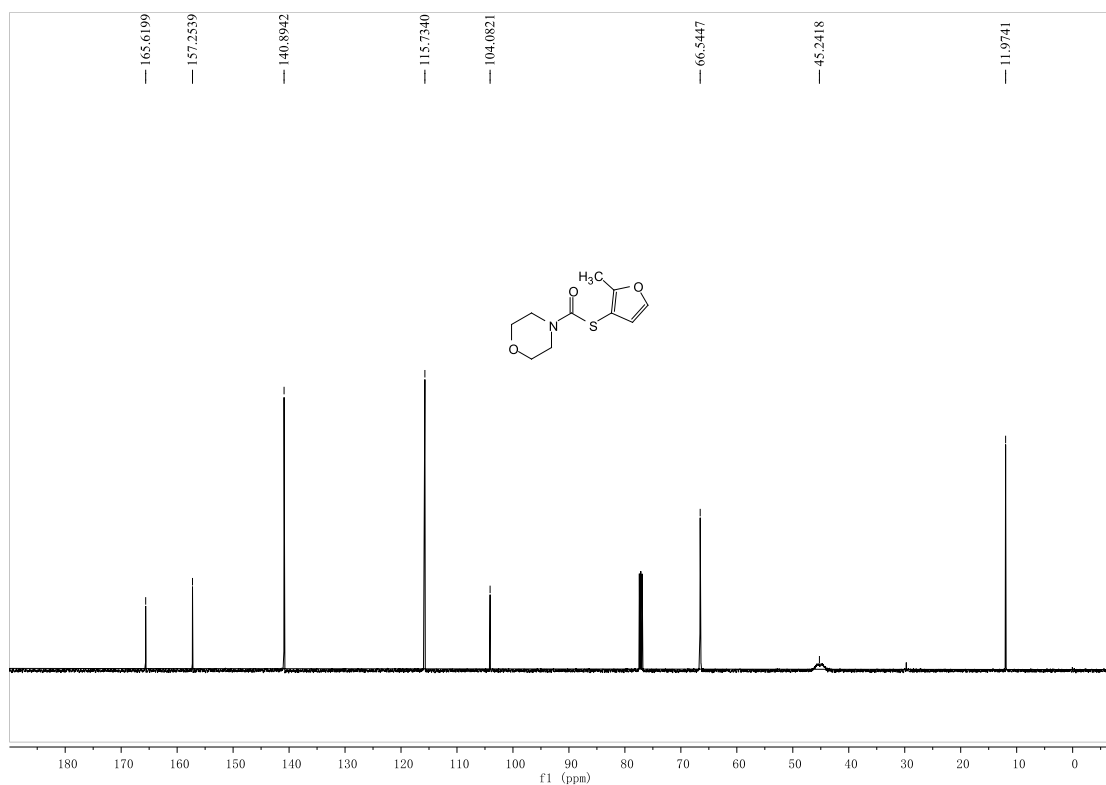
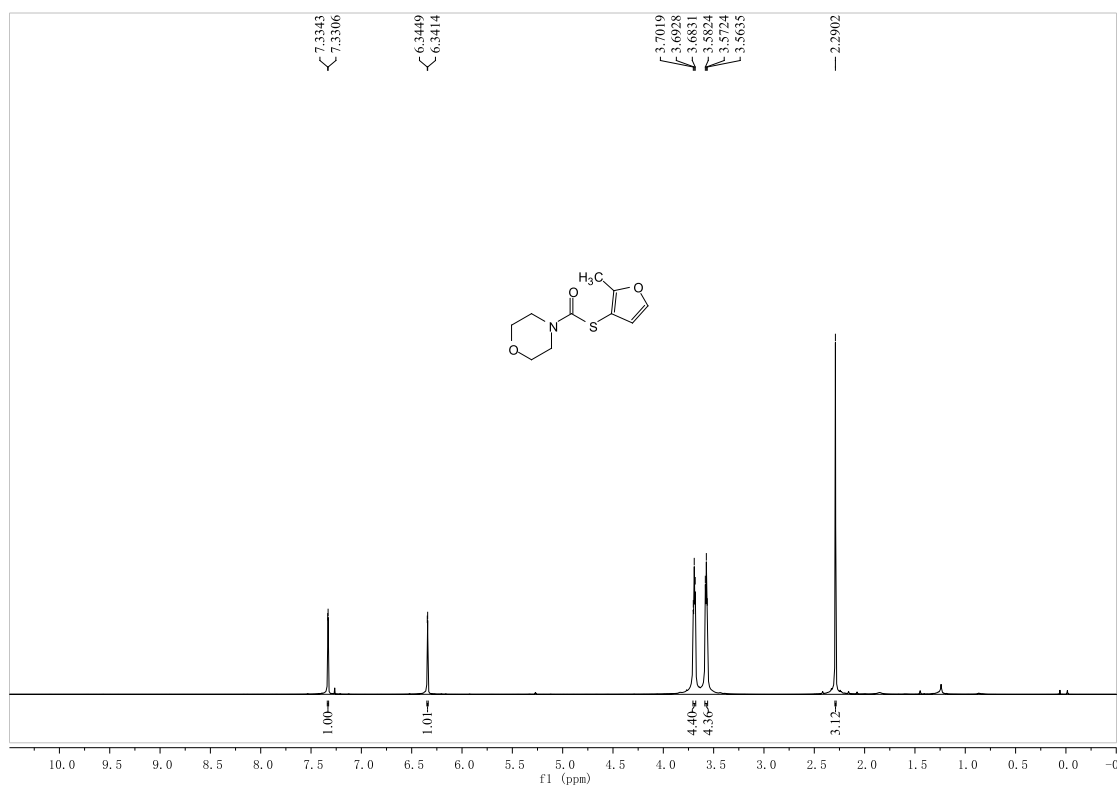
N-(((2-methylfuran-3-yl)thio)methyl)acetamide (3f)



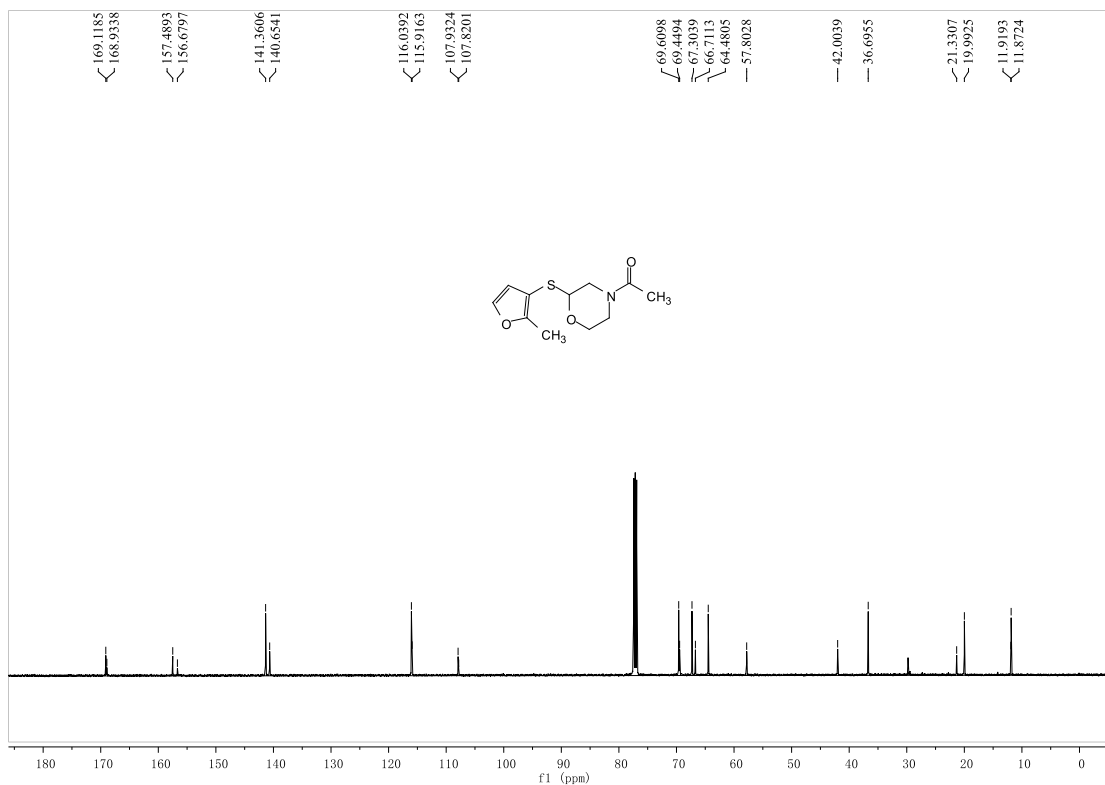
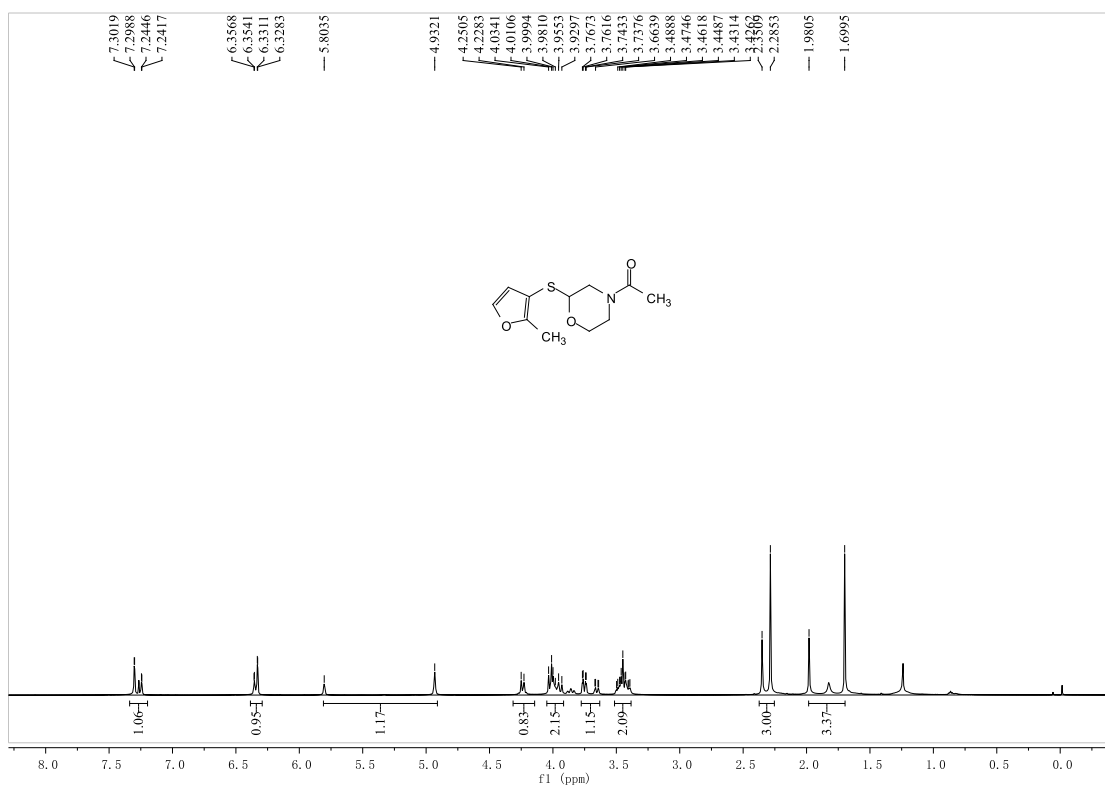
2-((2-methylfuran-3-yl)thio)morpholine-4-carbaldehyde (3g)



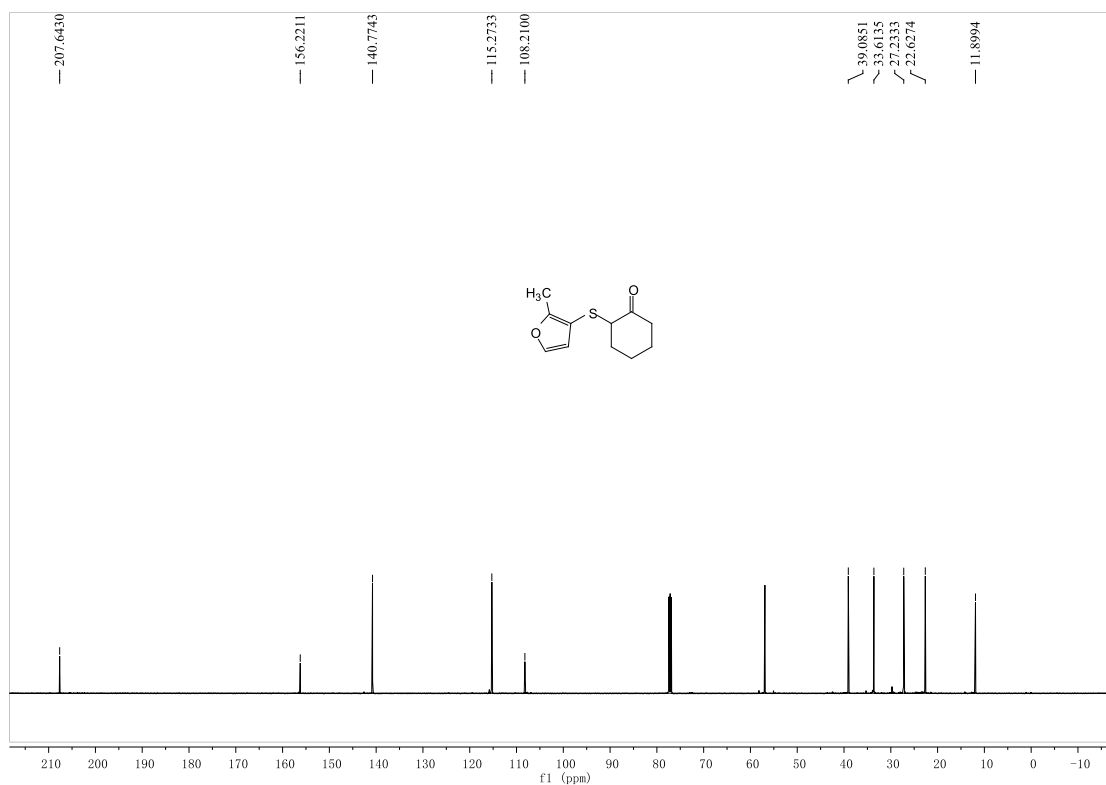
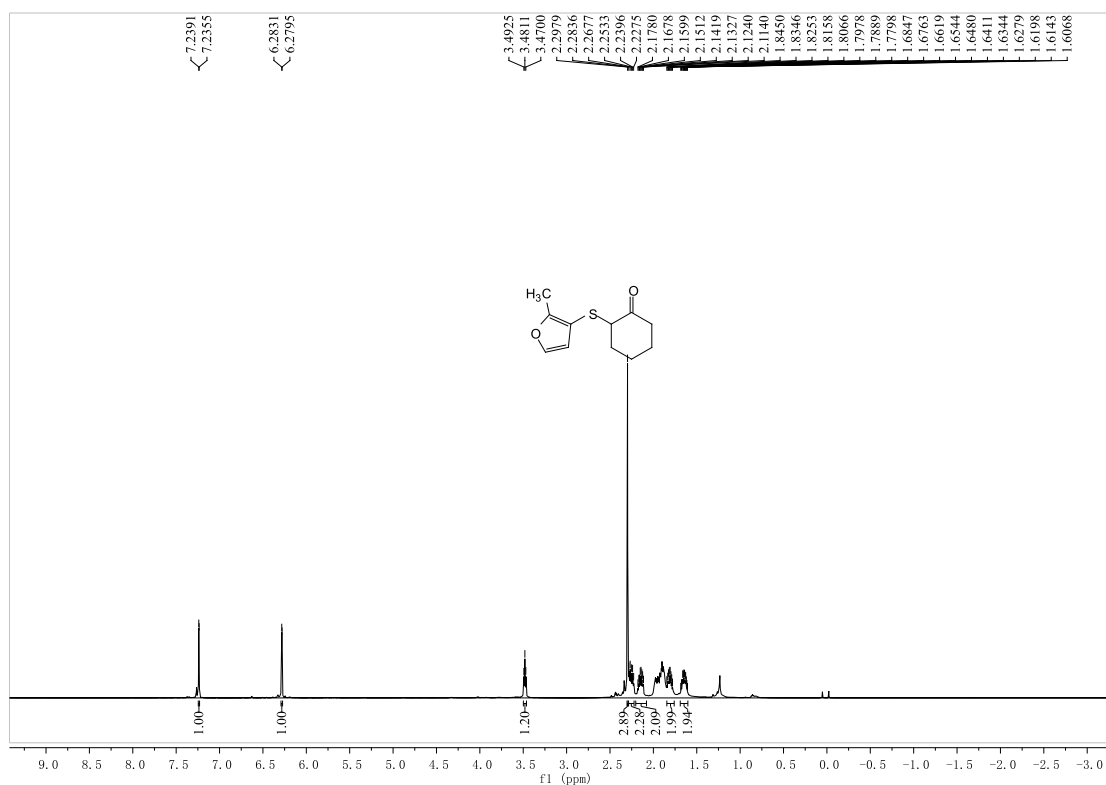
S-(2-methylfuran-3-yl)morpholine-4-carbothioate (3h)



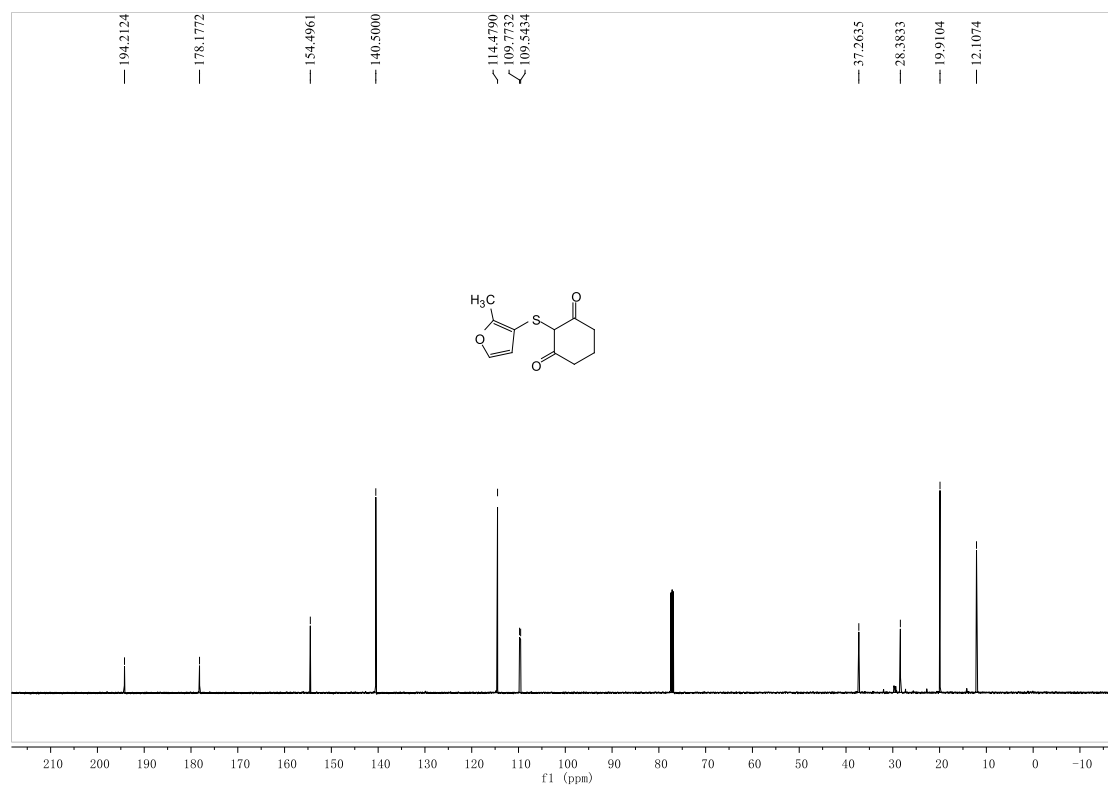
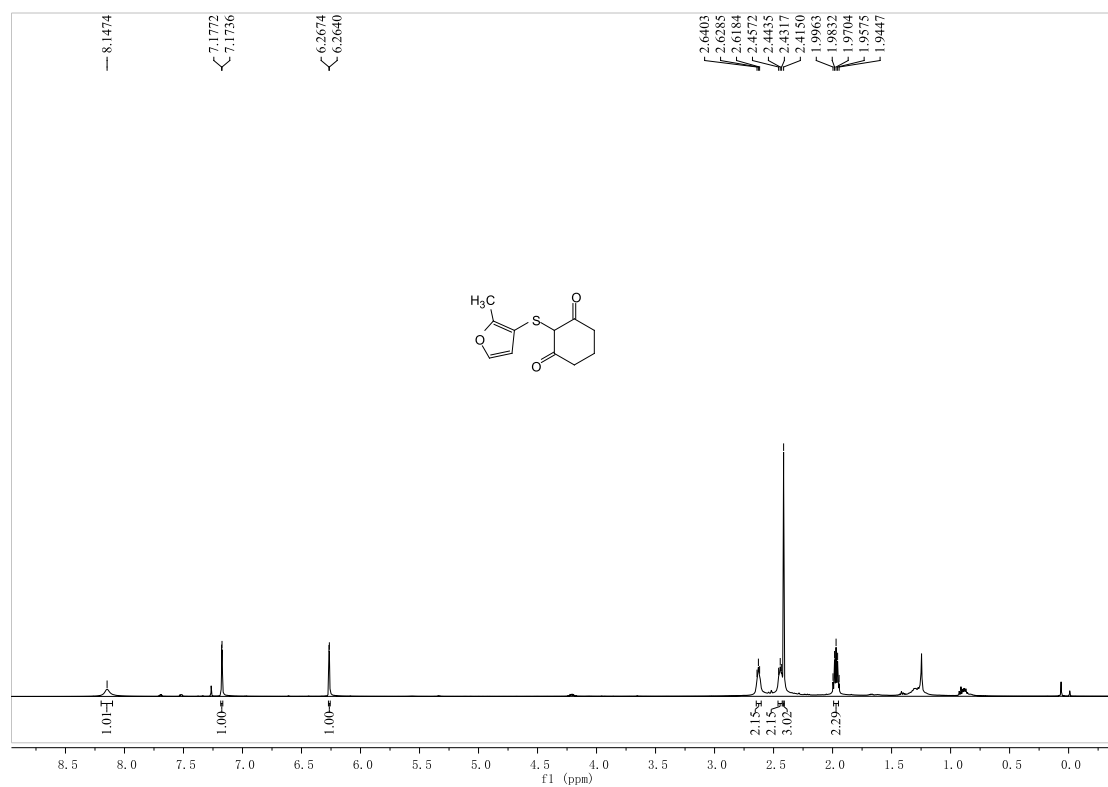
1-(2-((2-methylfuran-3-yl)thio)morpholino)ethan-1-one (3i)



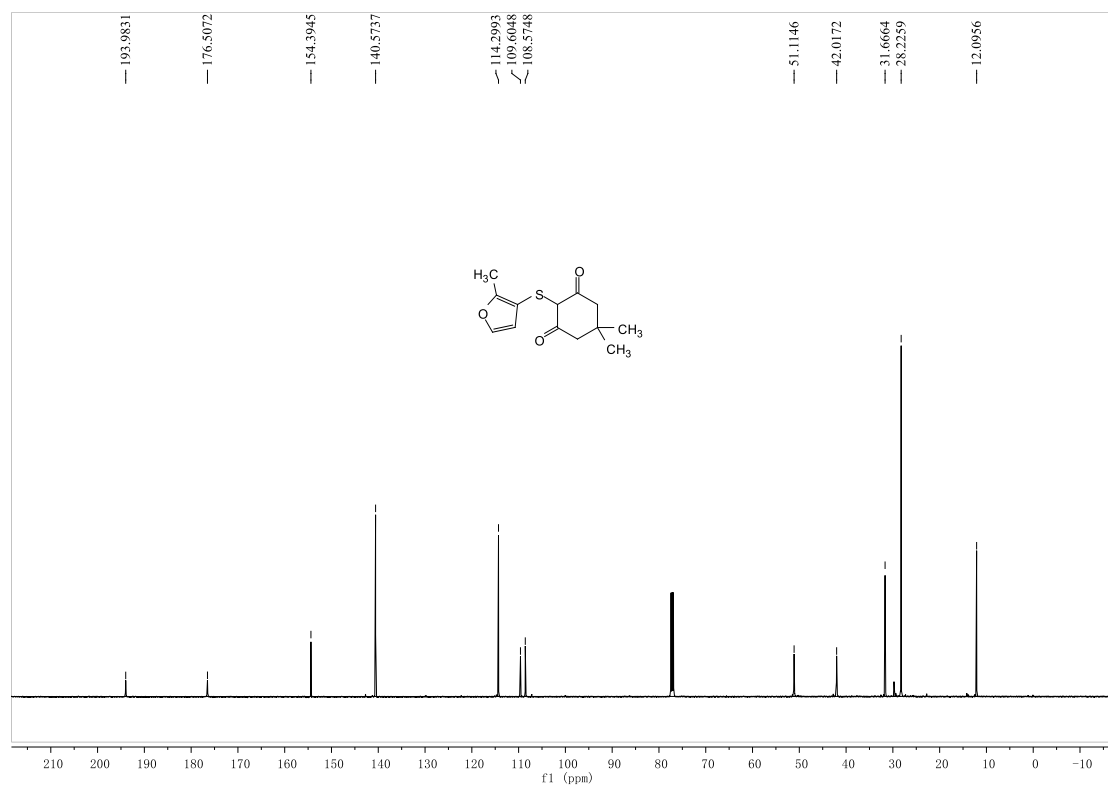
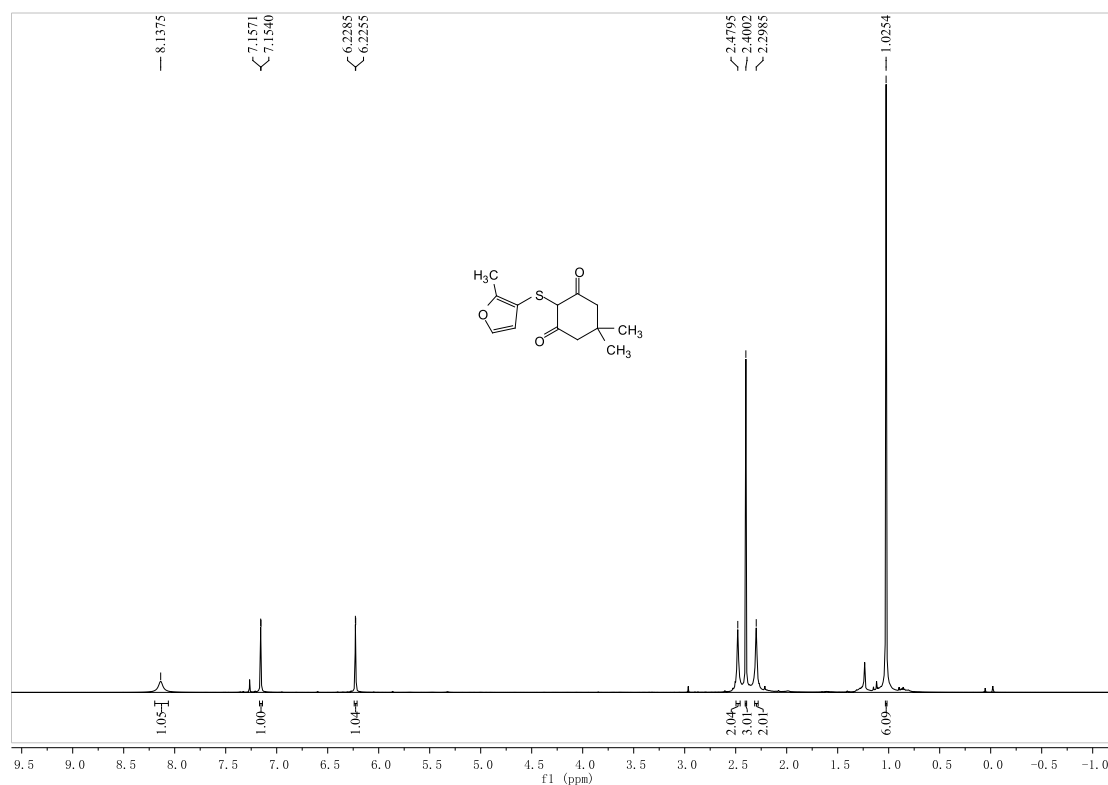
2-((2-methylfuran-3-yl)thio)cyclohexan-1-one (3j)



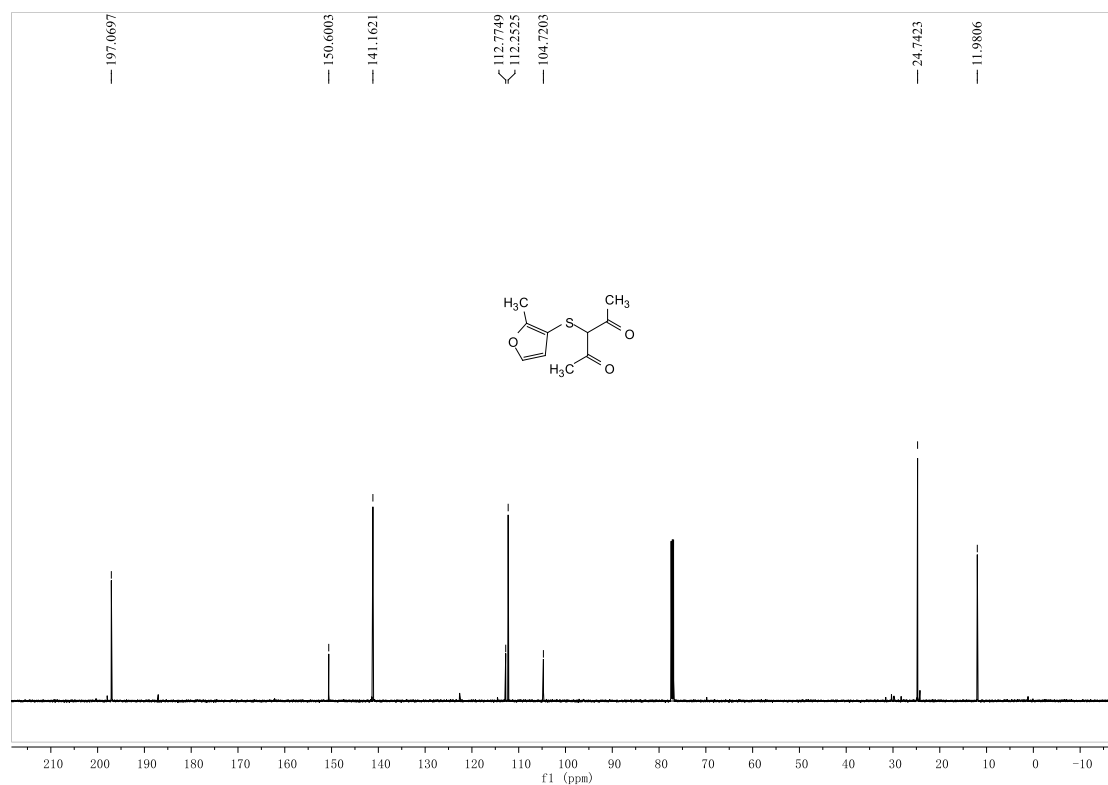
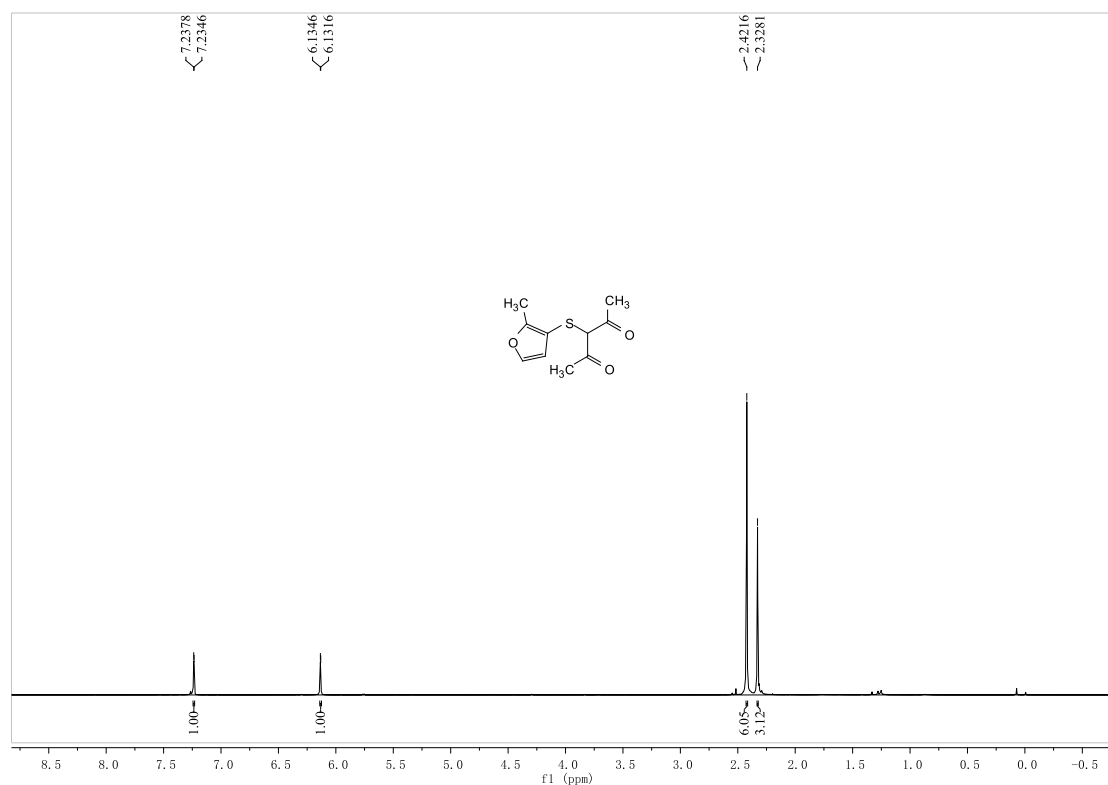
2-((2-methylfuran-3-yl)thio)cyclohexane-1,3-dione (3k)



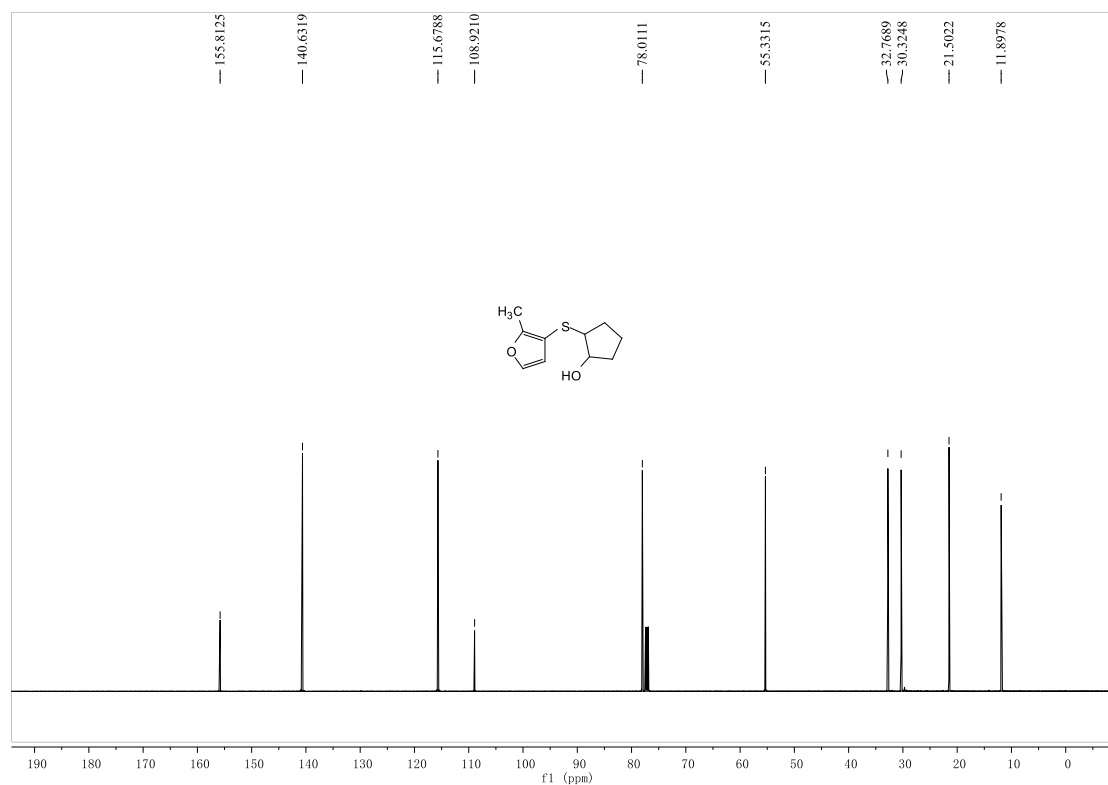
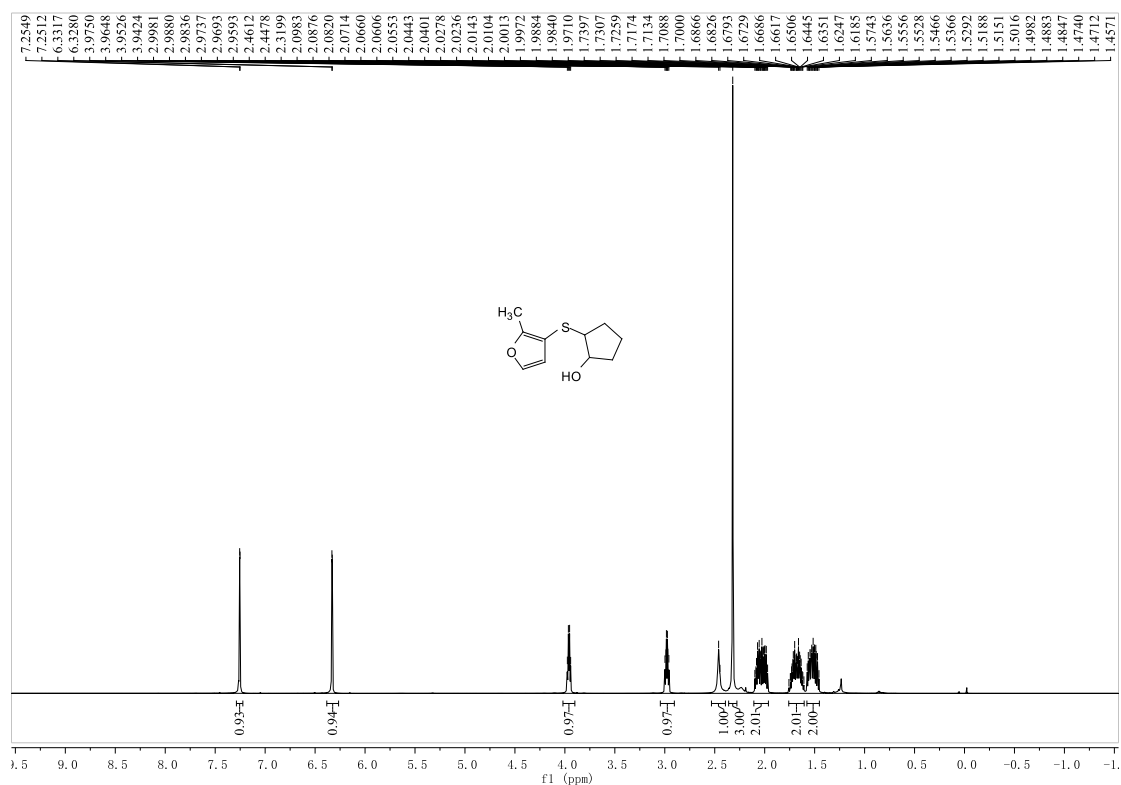
5,5-dimethyl-2-((2-methylfuran-3-yl)thio)cyclohexane-1,3-dione (3l)



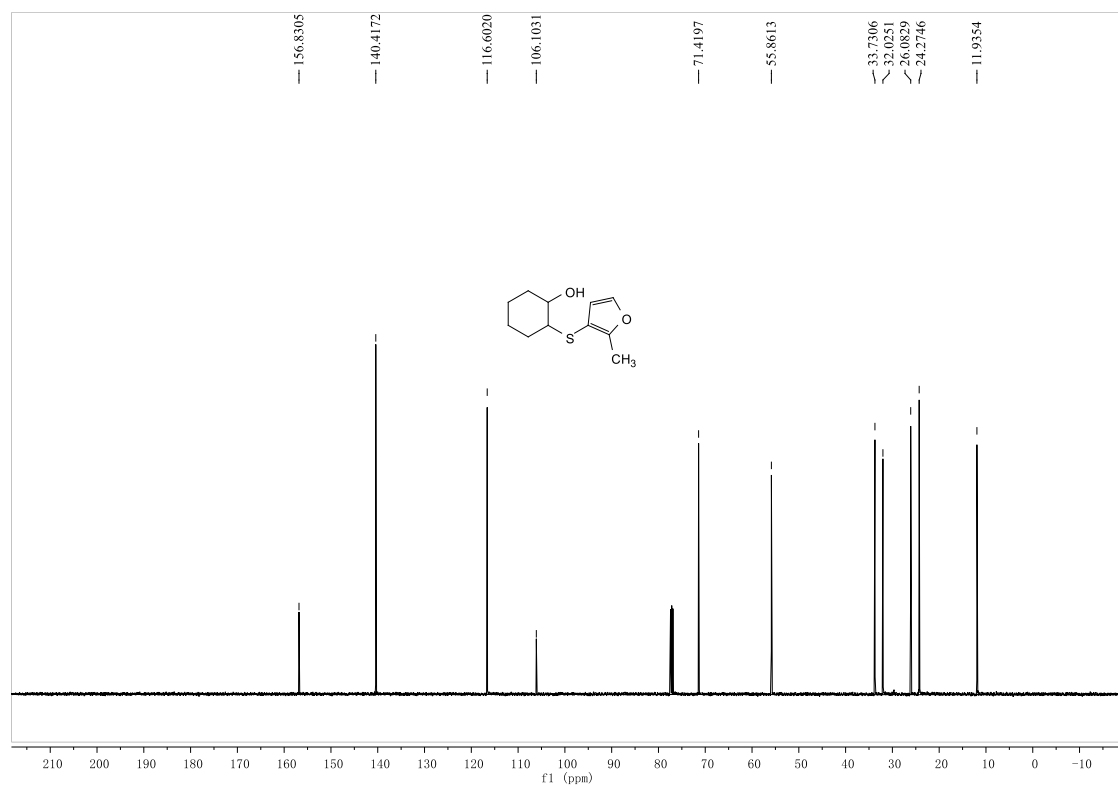
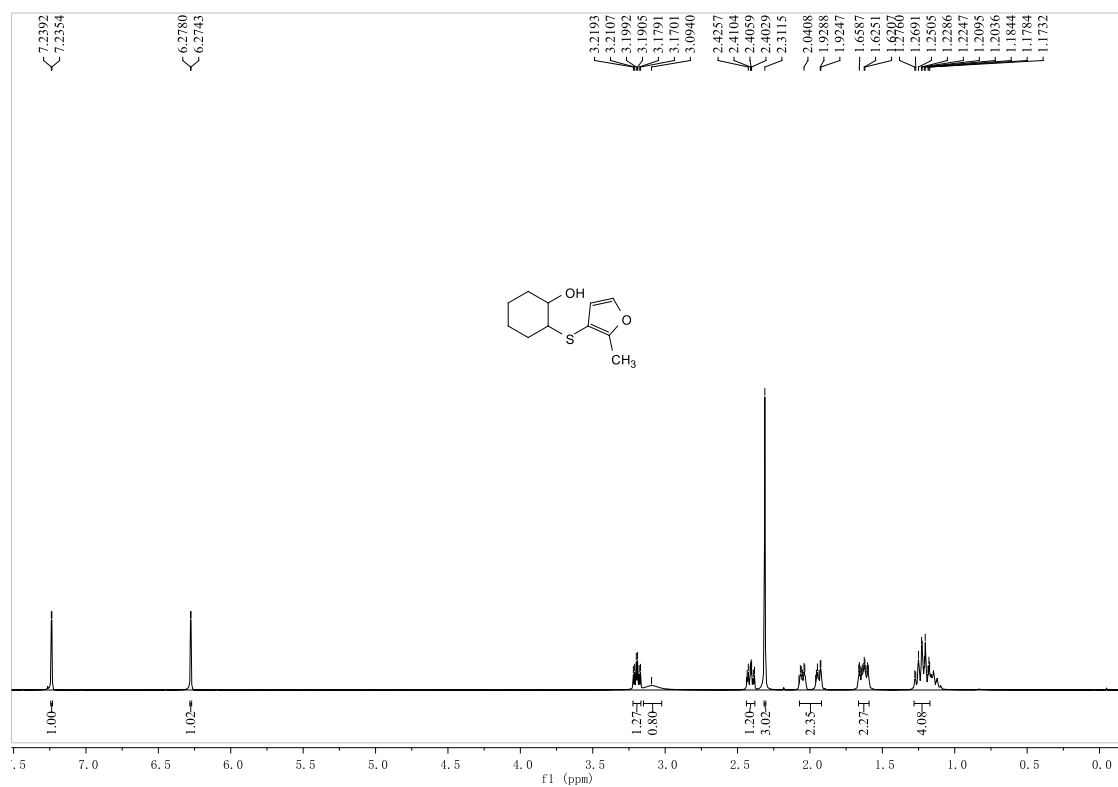
3-((2-methylfuran-3-yl)thio)pentane-2,4-dione (3m)



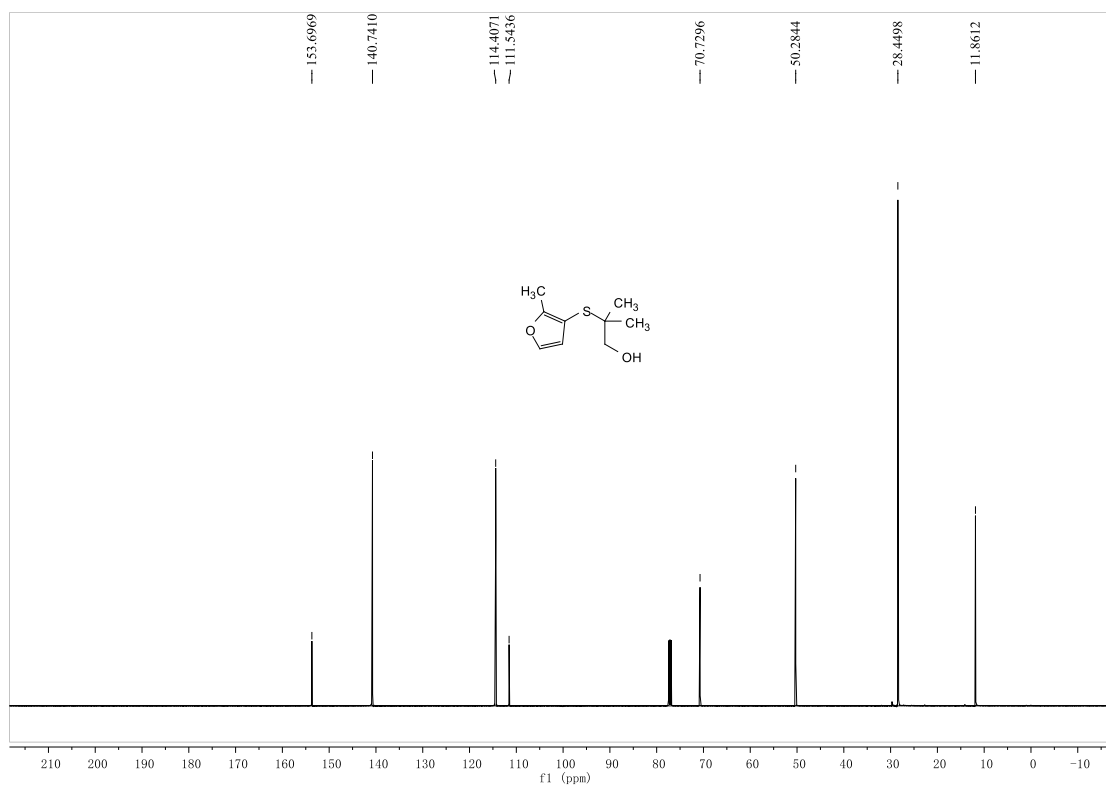
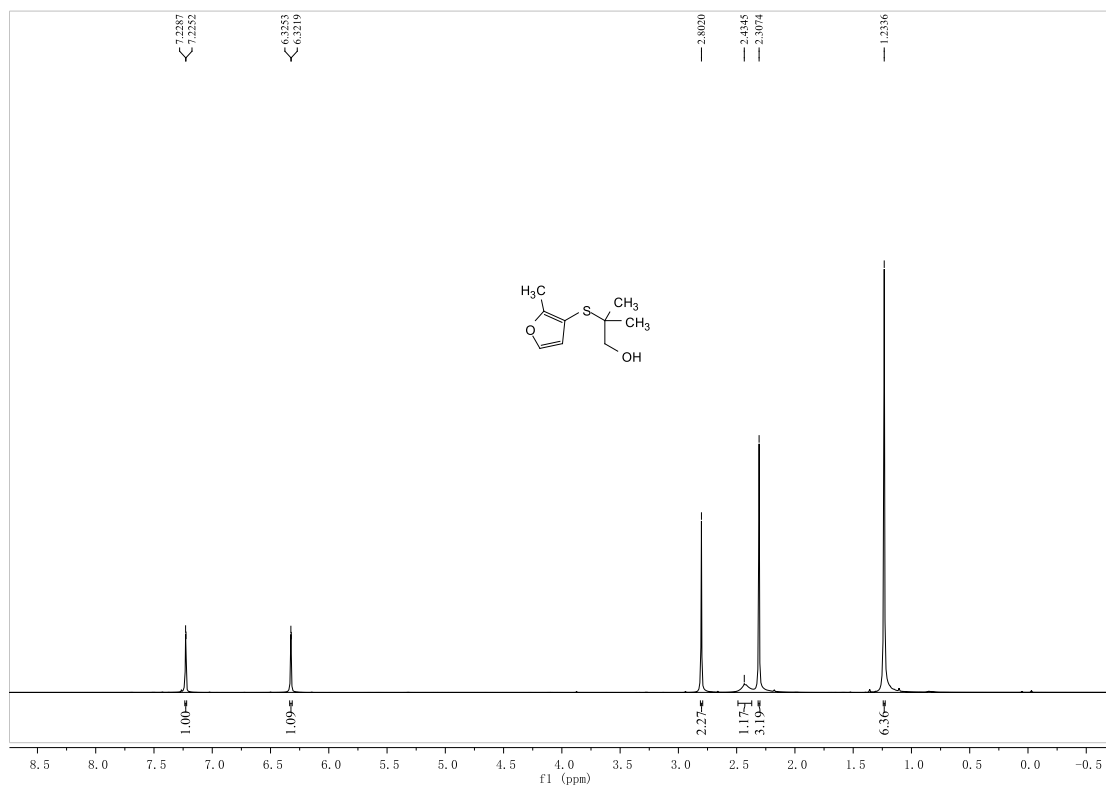
2-((2-methylfuran-3-yl)thio)cyclopentan-1-ol (5a)



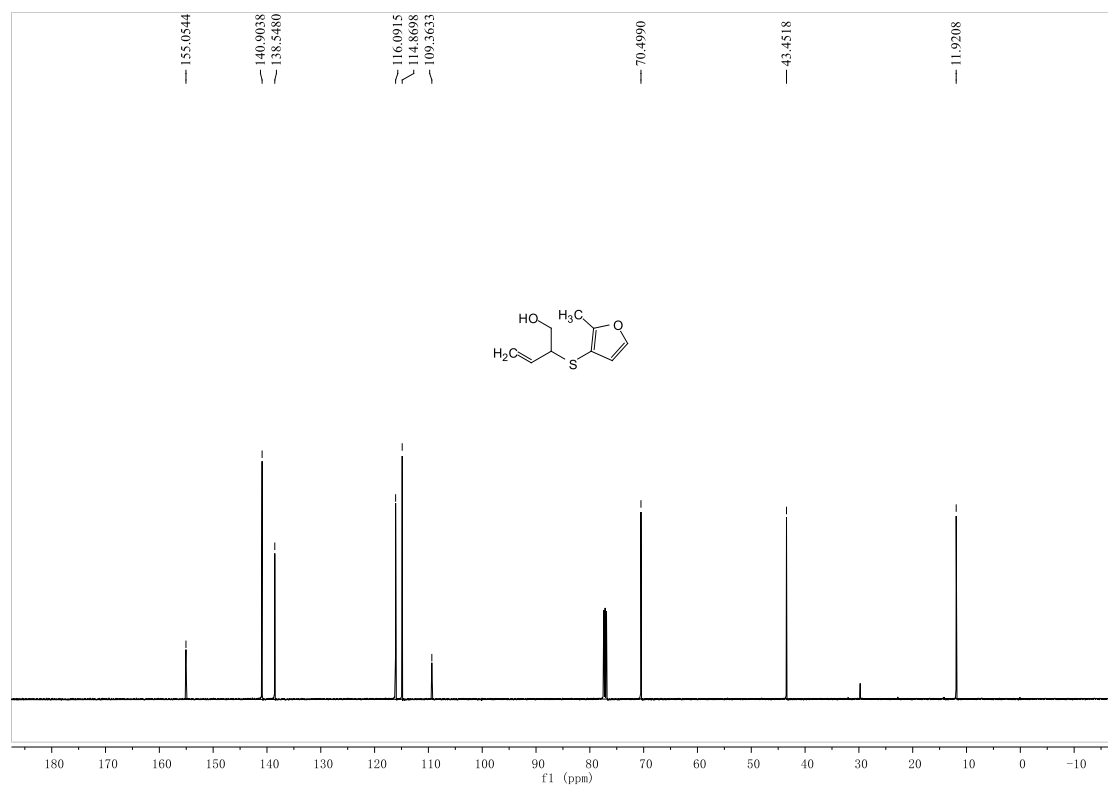
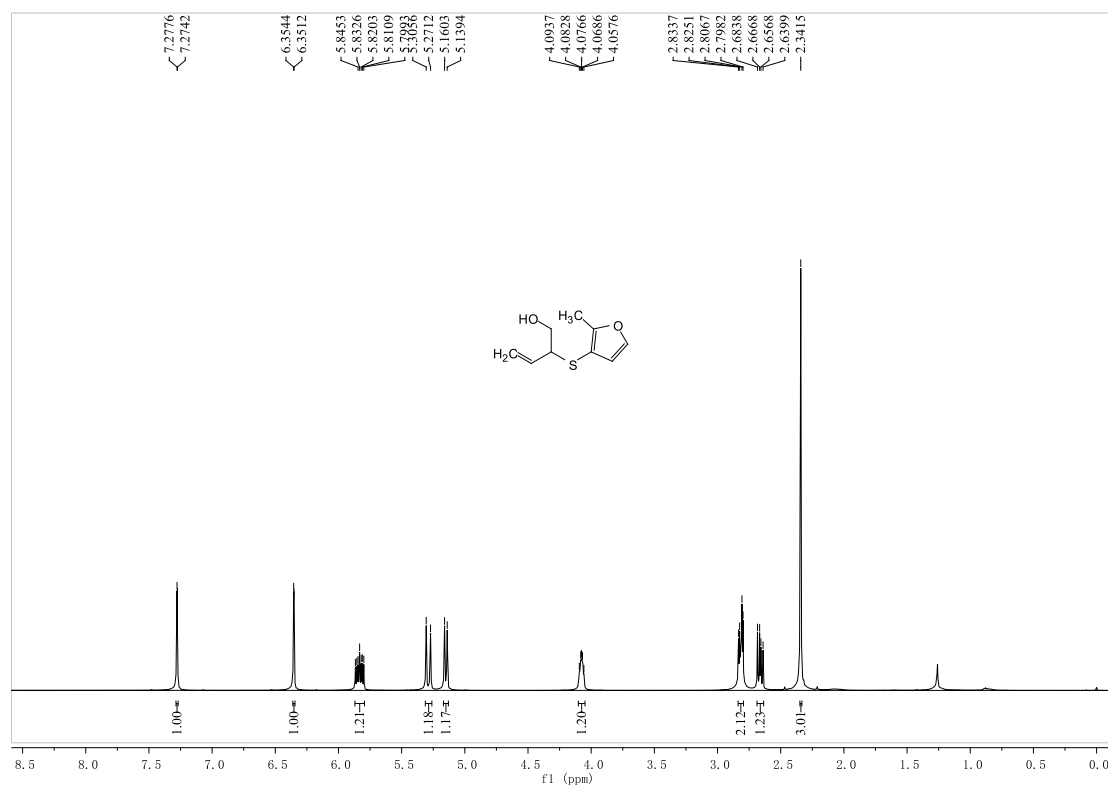
2-((2-methylfuran-3-yl)thio)cyclohexan-1-ol (5b)



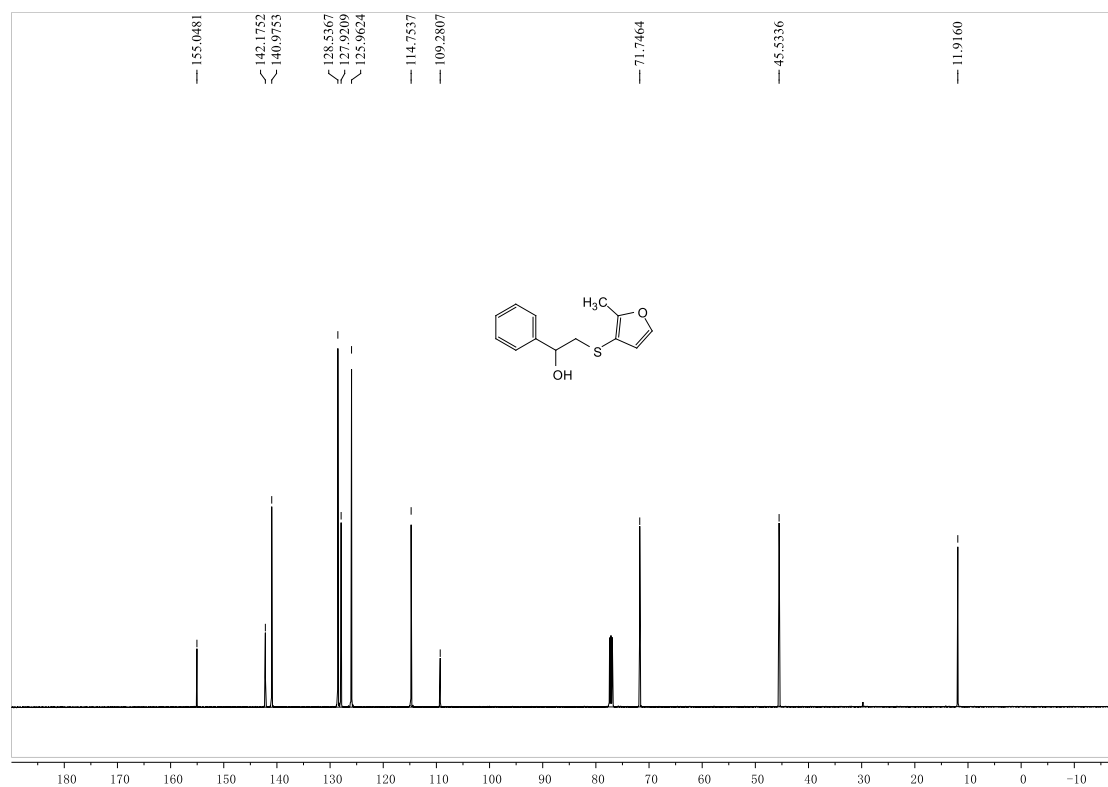
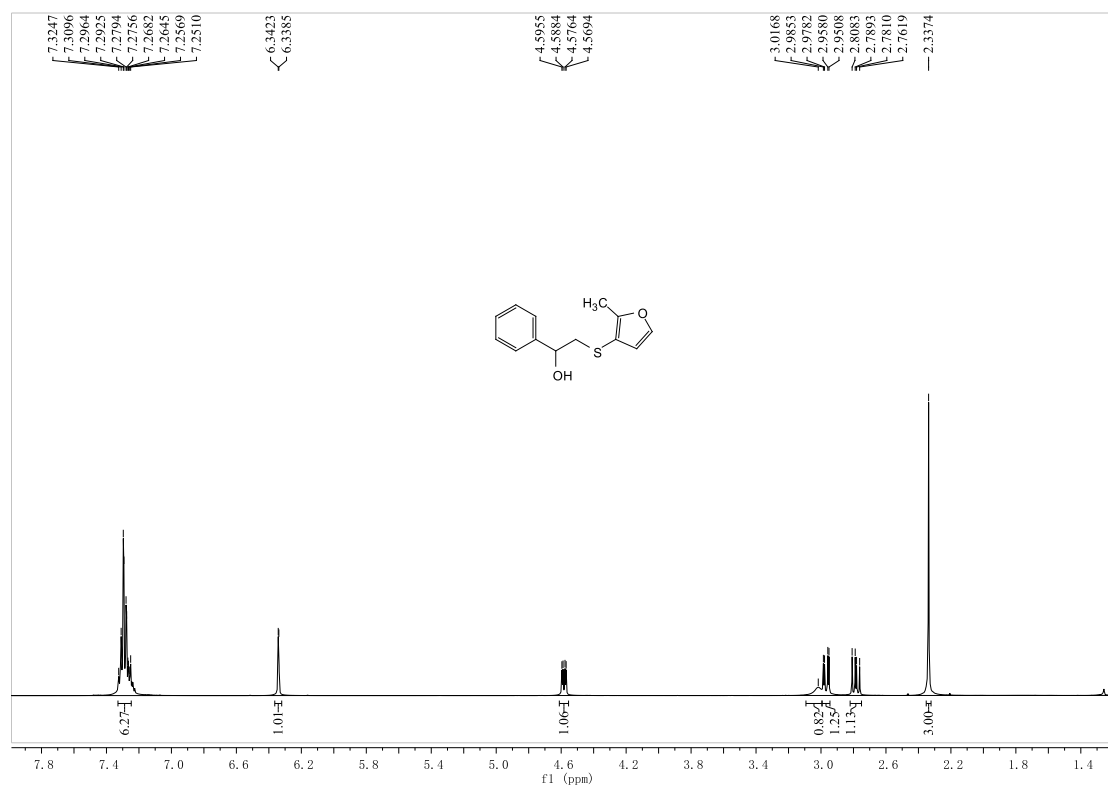
2-methyl-2-((2-methylfuran-3-yl)thio)propan-1-ol (5c)



2-((2-methylfuran-3-yl)thio)but-3-en-1-ol (5d)



2-((2-methylfuran-3-yl)thio)-1-phenylethan-1-ol (5e)



2-((2-methylfuran-3-yl)thio)-2-phenylethan-1-ol (5f)

