

Homoallylic amines as efficient chiral inducing framework in the conjugated addition of amides to α,β -unsaturated esters.

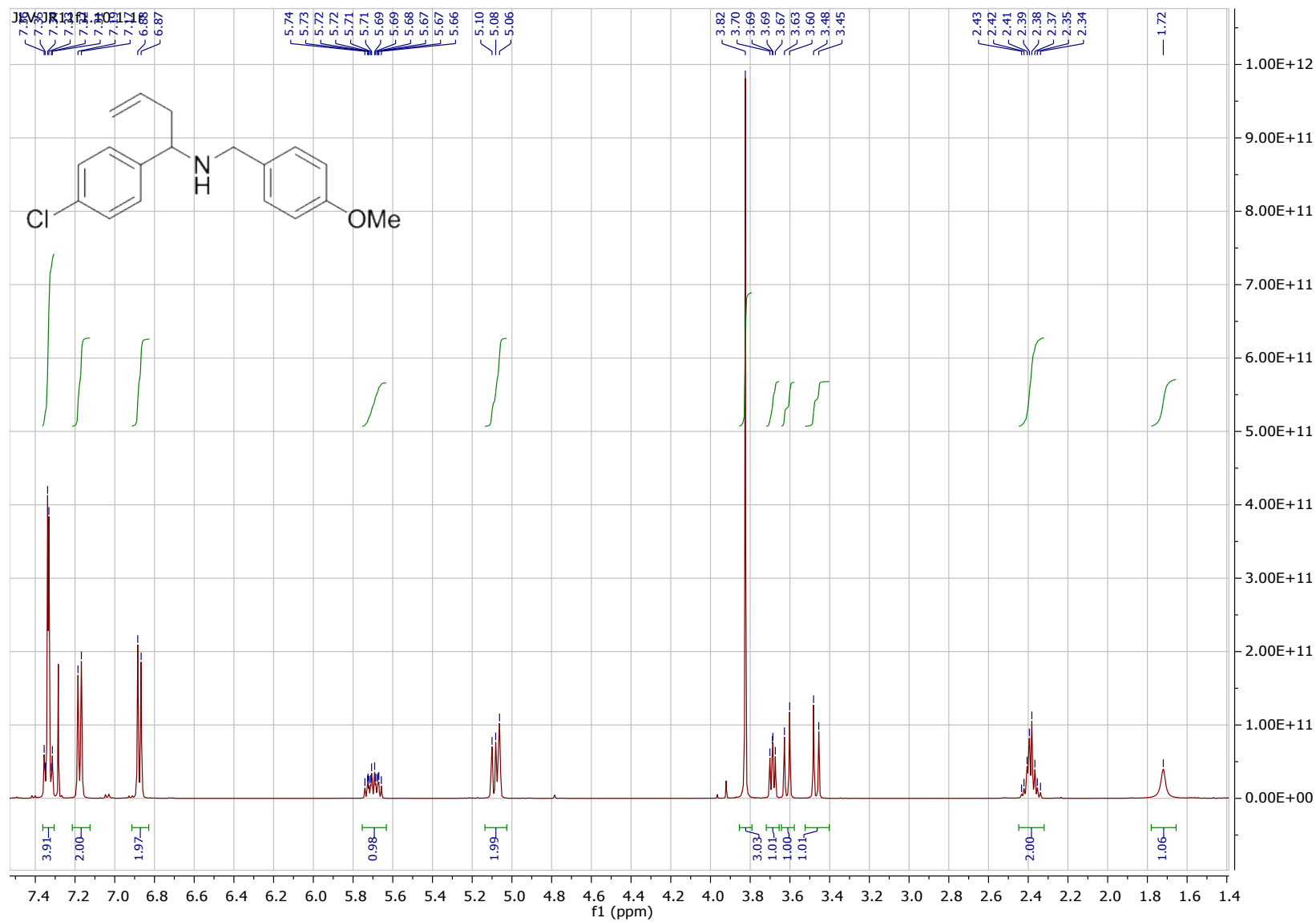
An entry to enantio-enriched diversely substituted amines.

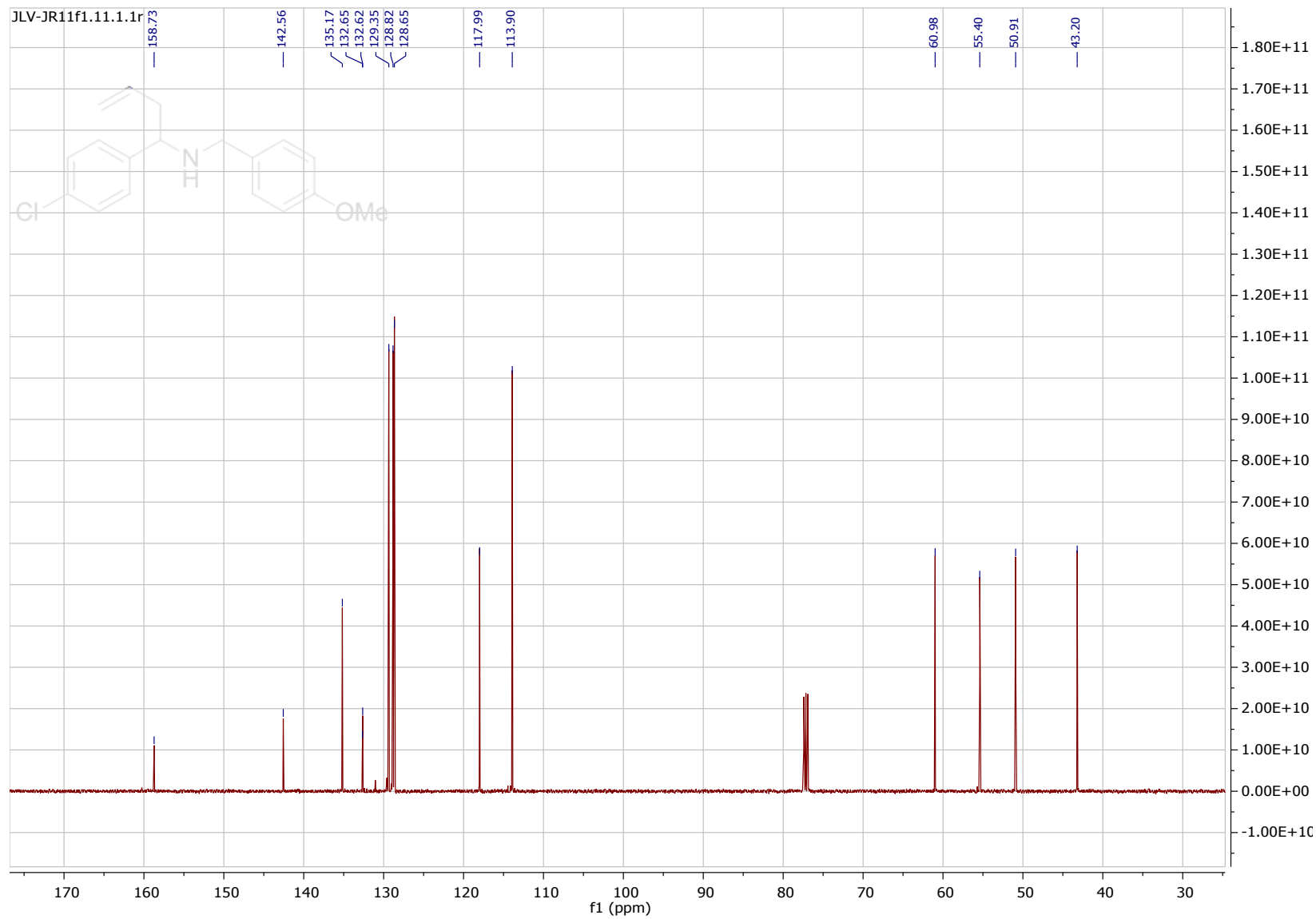
Johann Rogier, Lilia Anani, Aurélien Coelho, Fabien Massicot, Carine Machado-Rodrigues, Jean-Bernard Behr, Jean-Luc Vasse

Institut de Chimie Moléculaire de Reims, CNRS-UMR 7312 and Université de Reims, 51687 Reims Cedex 2, France

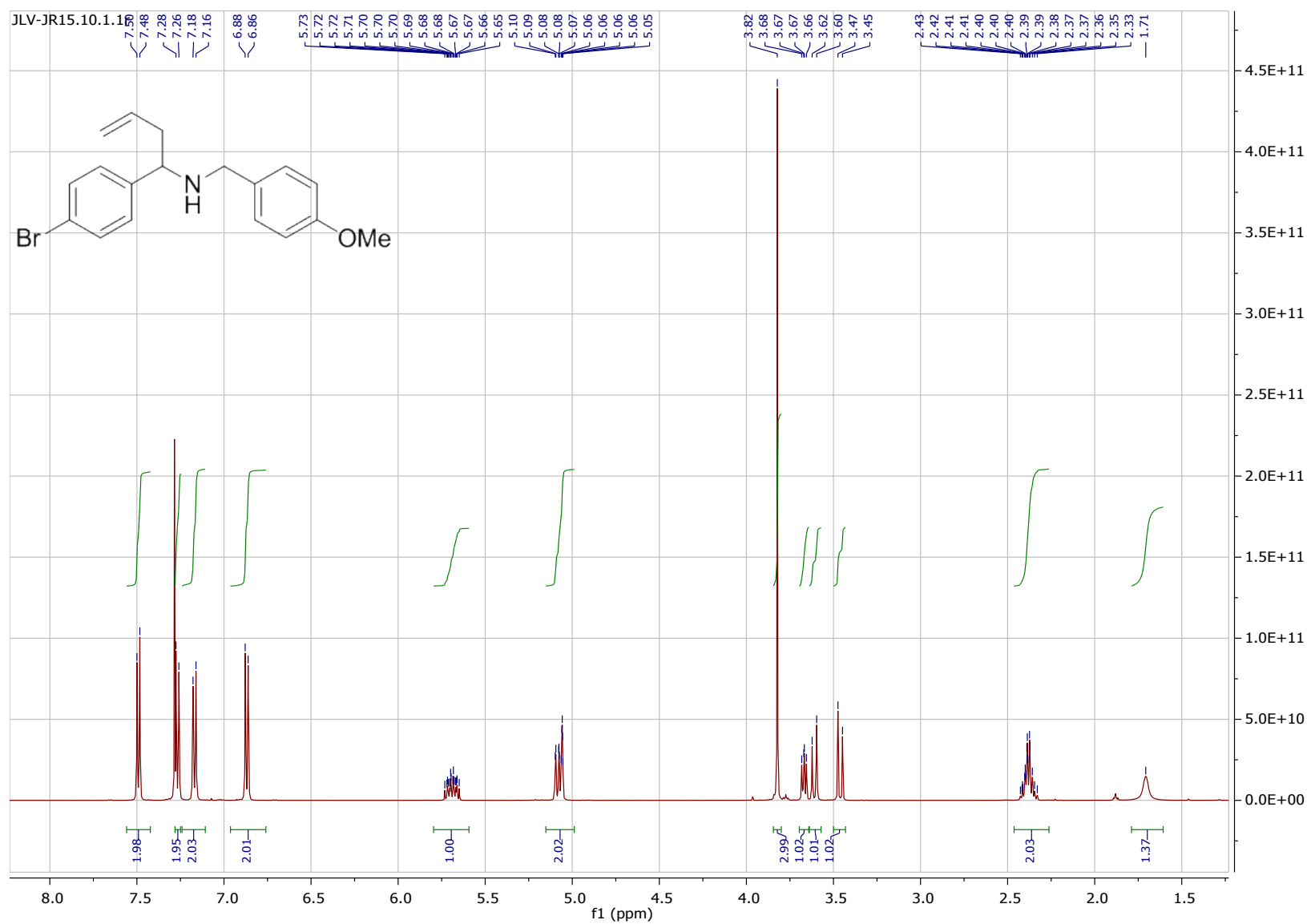
^1H and ^{13}C NRM spectra copies

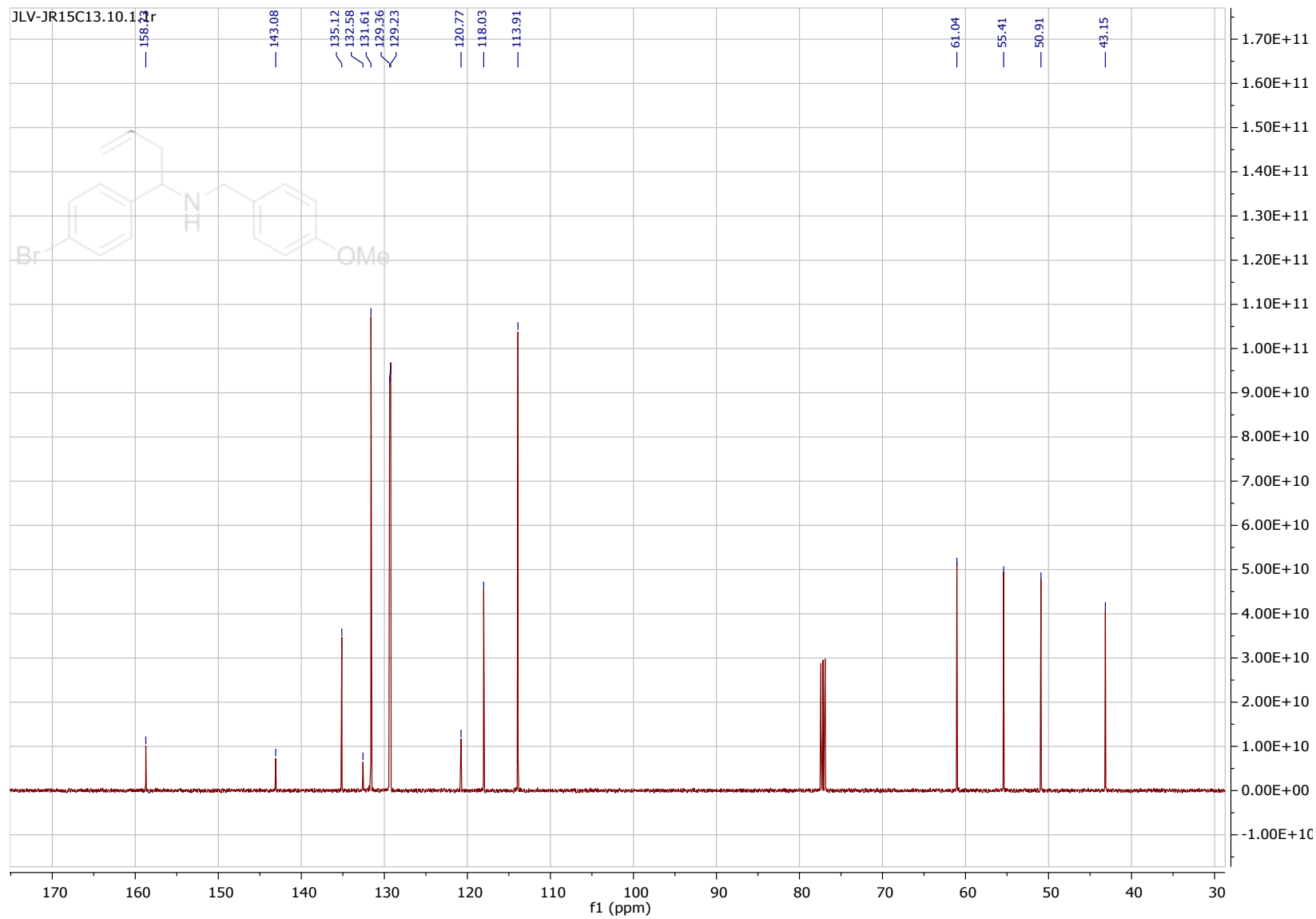
N-Benzyl-1-(4-chlorophenyl)but-3-en-1-amine 1c



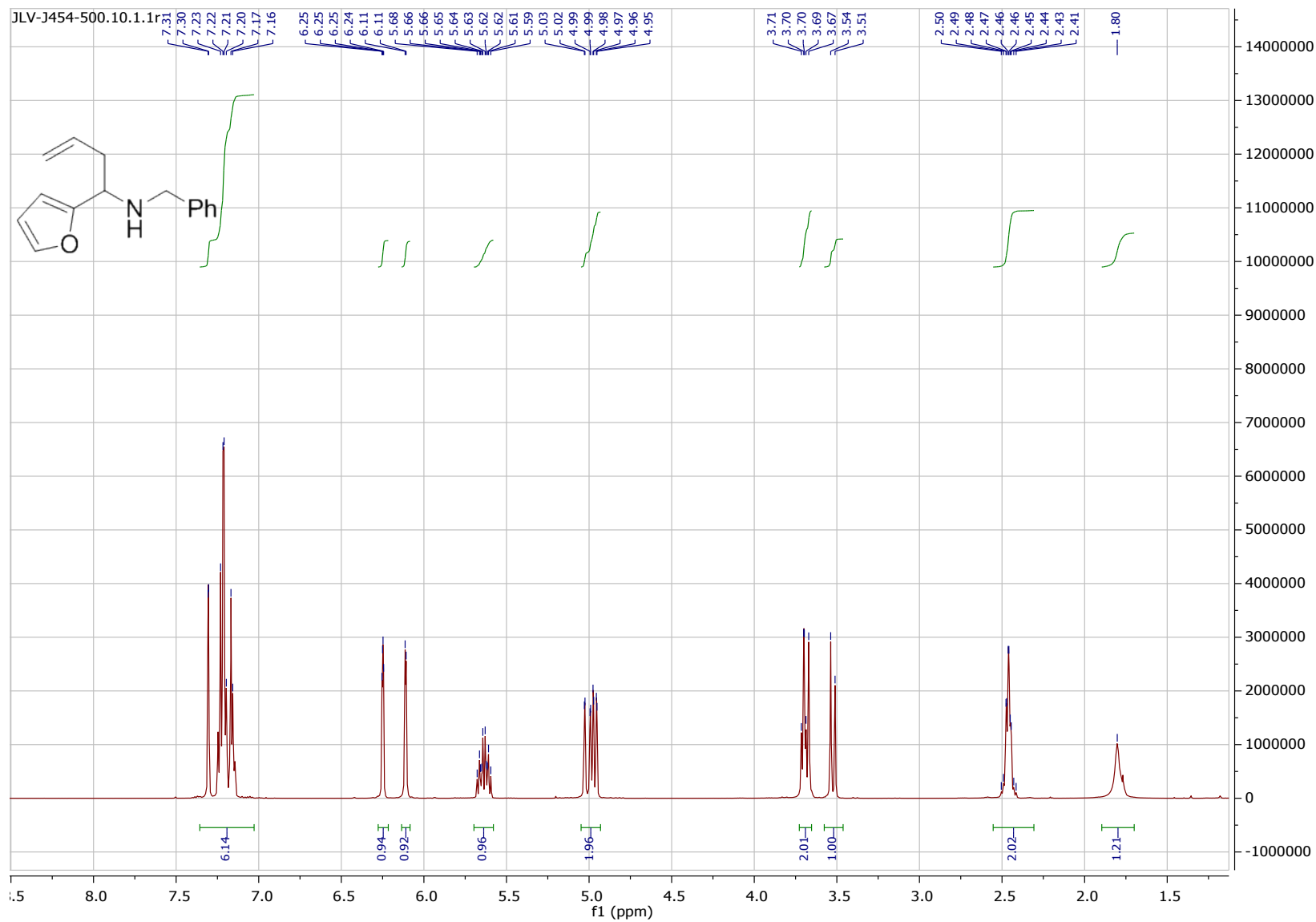


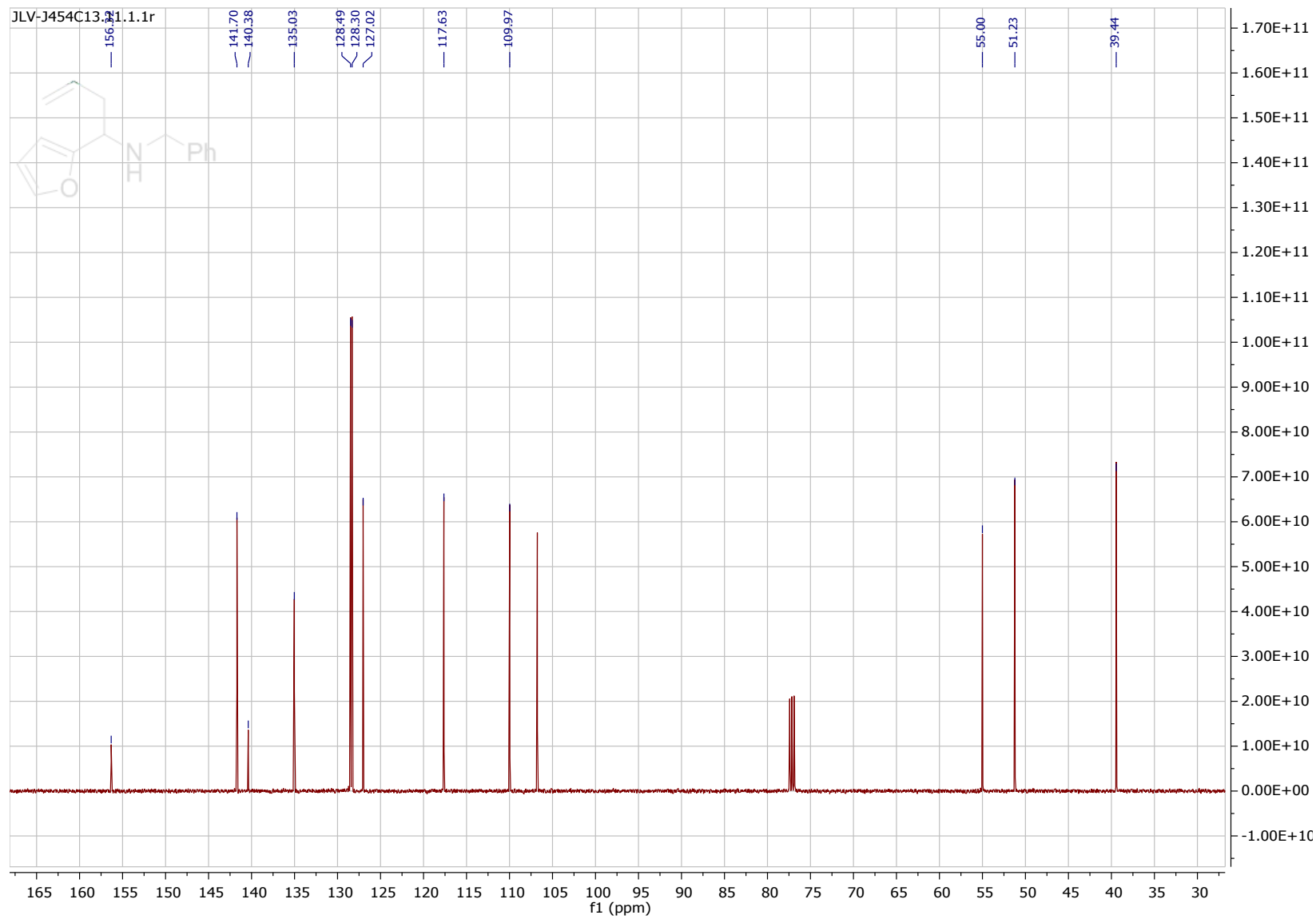
1-(4-Bromophenyl)-N-(4-methoxybenzyl)but-3-en-1-amine 1d



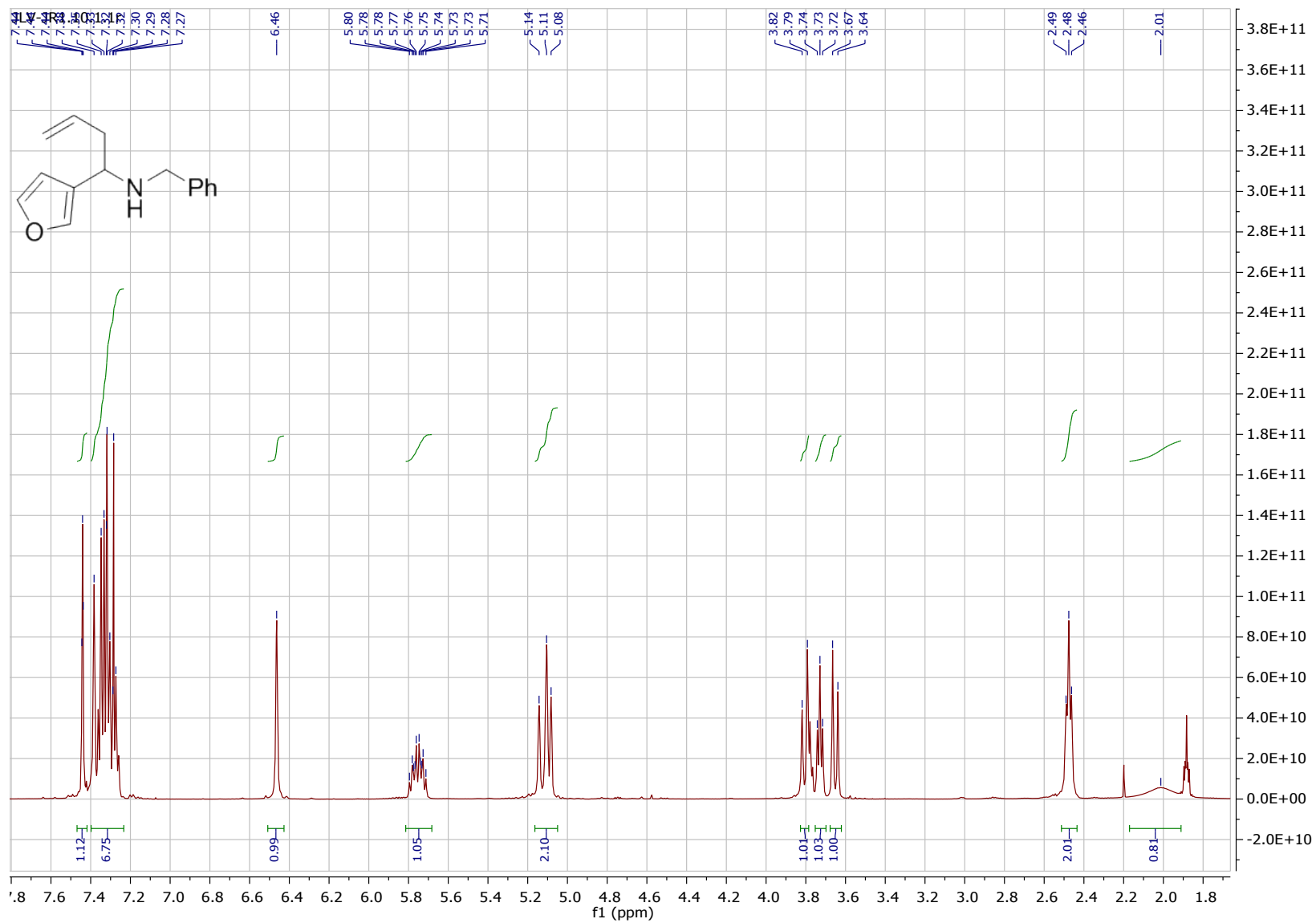


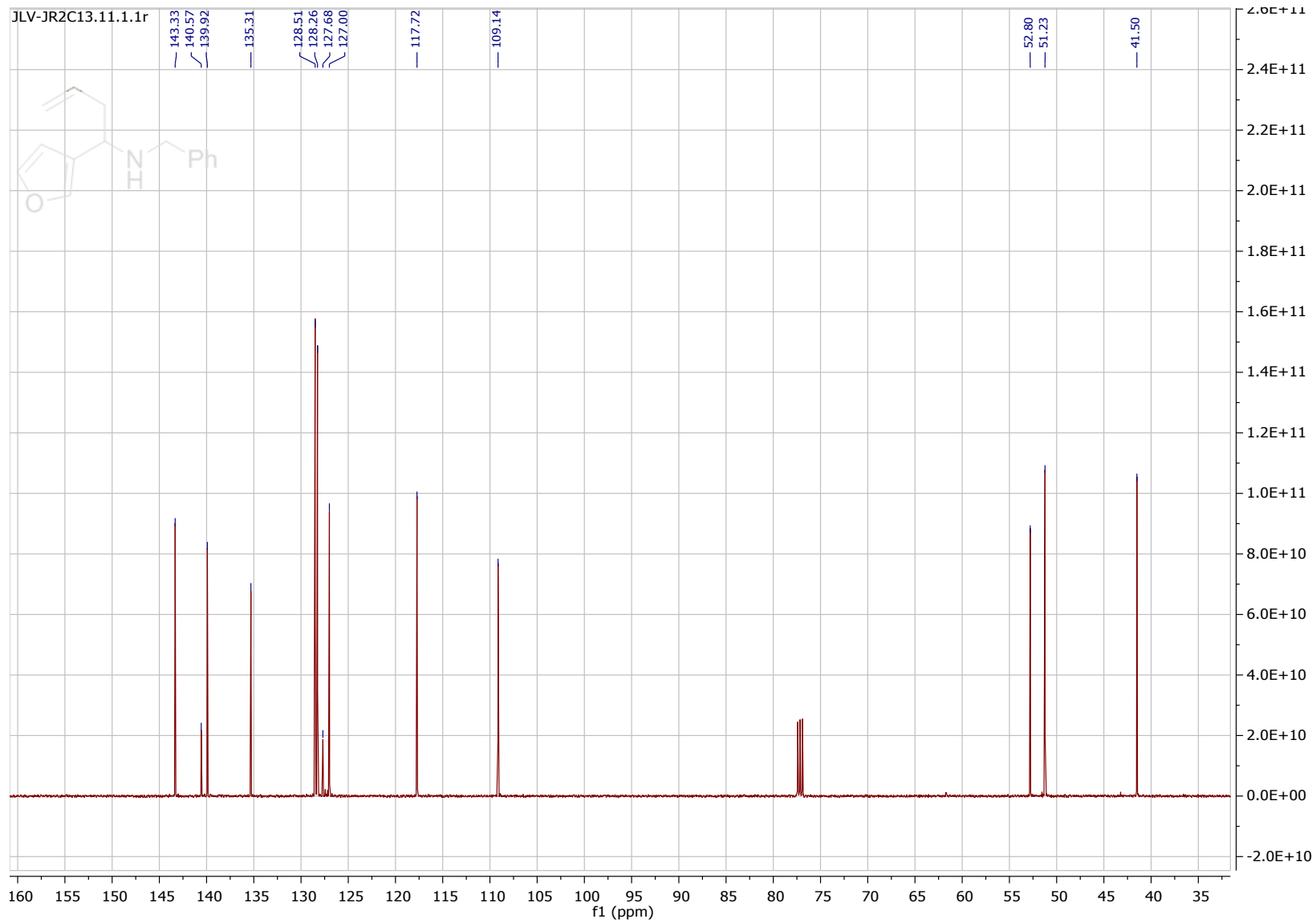
***N*-Benzyl-1-(furan-2-yl)but-3-en-1-amine 1f**





***N*-Benzyl-1-(furan-3-yl)but-3-en-1-amine 1g**





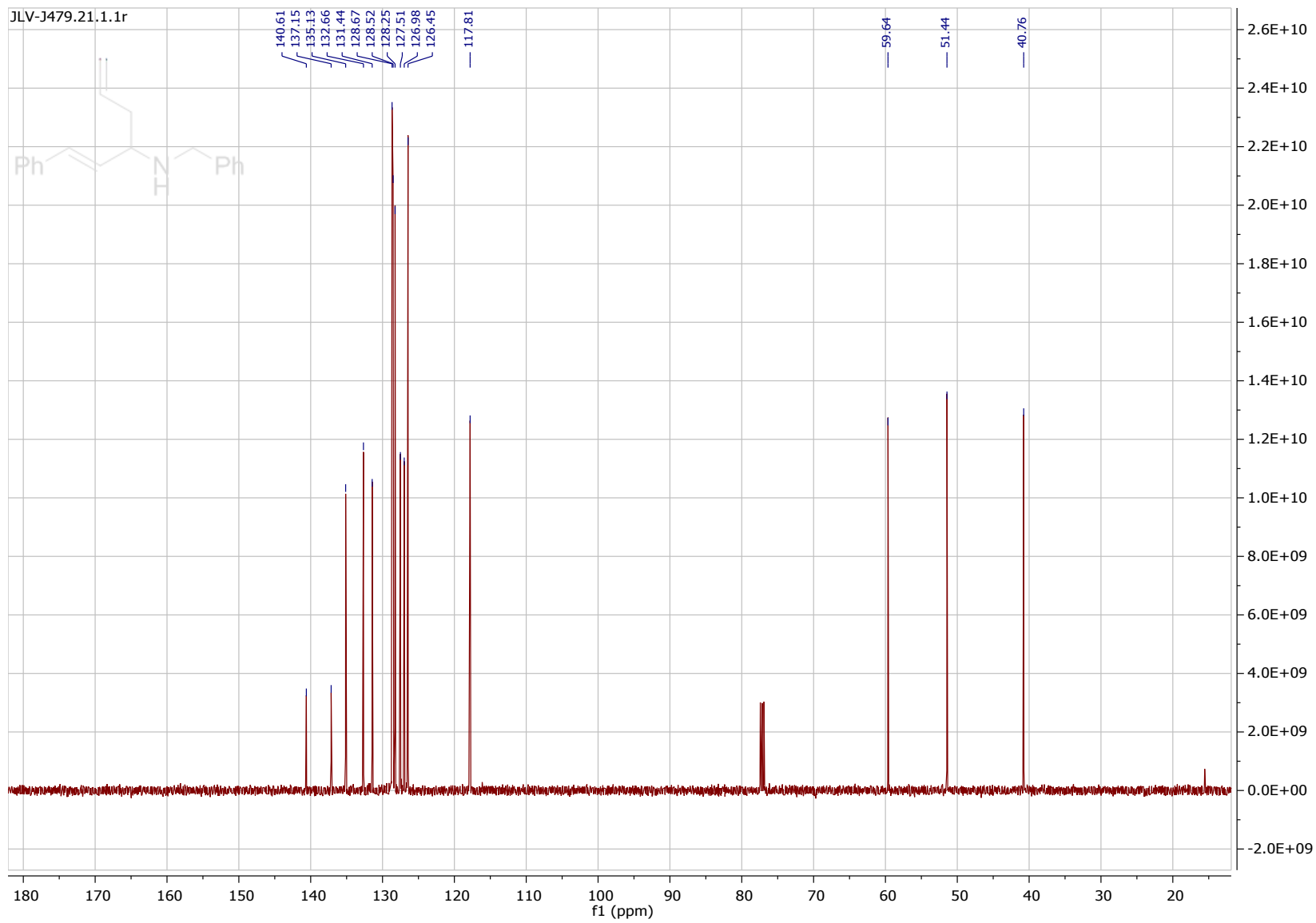
Chemical structure: CN(C)C(C=Cc1ccccc1)Cc2ccccc2

¹H NMR spectrum (400 MHz, CDCl₃) showing peaks from 1.6 to 7.6 ppm. The spectrum includes integration values and a chemical structure of (E)-1-((E)-3-phenyl-2-propenyl)-N,N-dimethylbenzylamine.

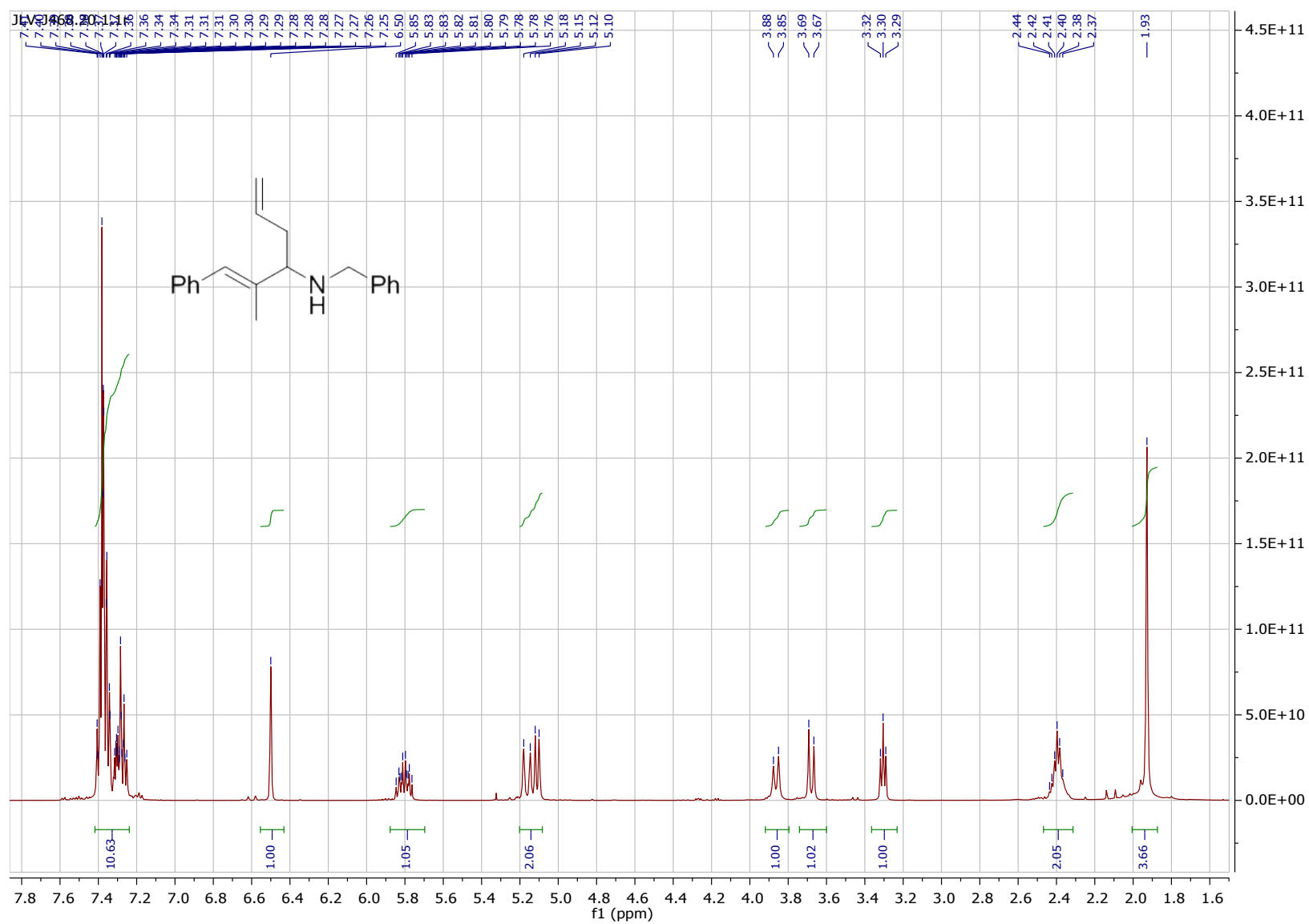
Chemical structure: CN(C)C(C=Cc1ccccc1)Cc2ccccc2

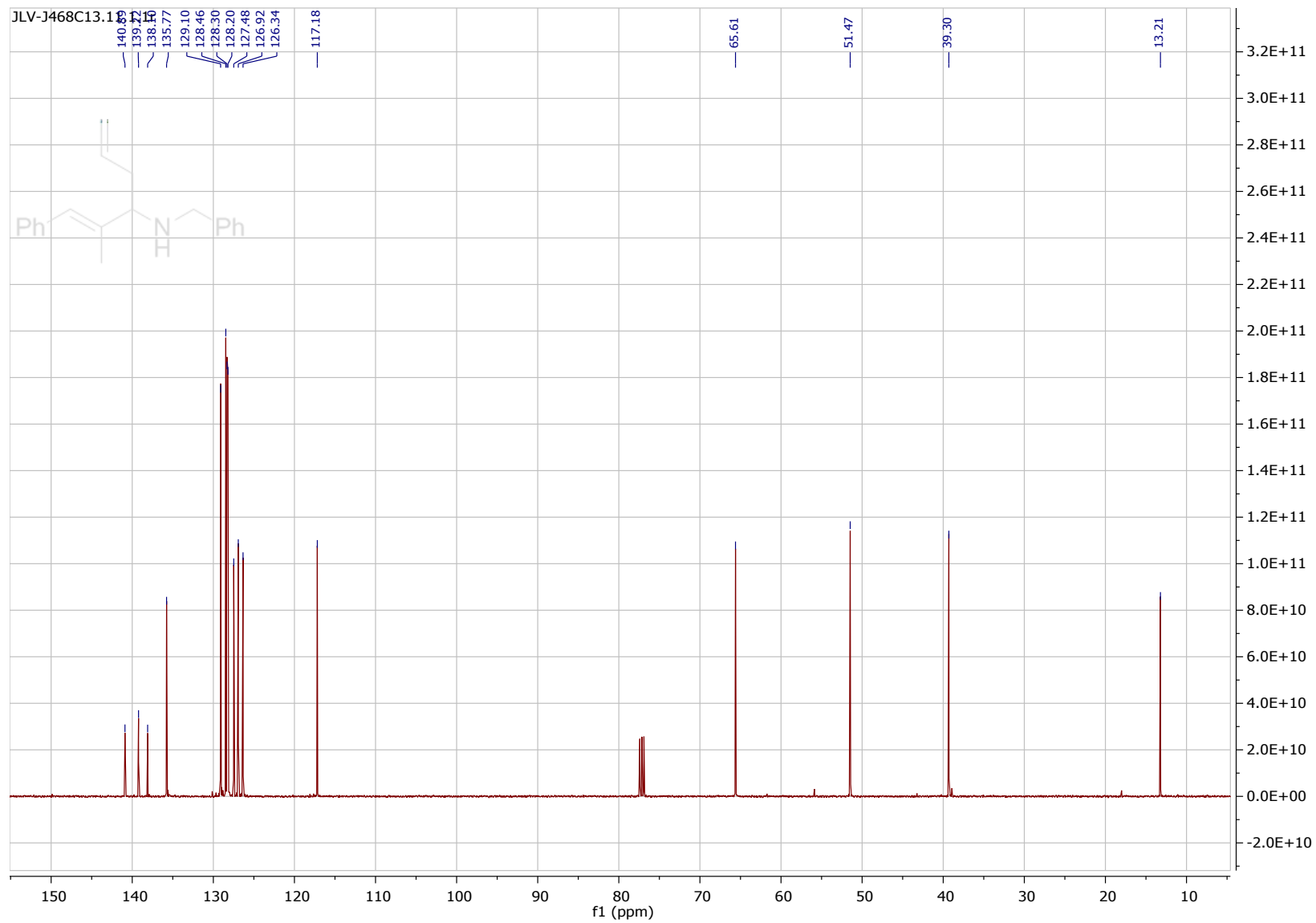
Integration values (from left to right): 2.06, 9.34, 1.00, 0.97, 1.00, 2.11, 1.01, 1.13, 0.98, 2.13.

Peak labels (from left to right): 7.44, 7.43, 7.42, 7.41, 7.40, 7.39, 7.38, 7.37, 7.36, 7.35, 7.34, 7.33, 7.32, 7.31, 7.30, 7.29, 7.28, 7.27, 6.57, 6.53, 6.17, 6.15, 6.13, 6.12, 5.88, 5.86, 5.85, 5.84, 5.83, 5.82, 5.81, 5.81, 5.81, 5.79, 3.94, 3.91, 3.76, 3.74, 3.37, 3.36, 3.34, 3.33, 2.47, 2.46, 2.45, 2.43, 2.42, 2.41, 2.39, 2.38, 2.37, 2.35, 1.81.

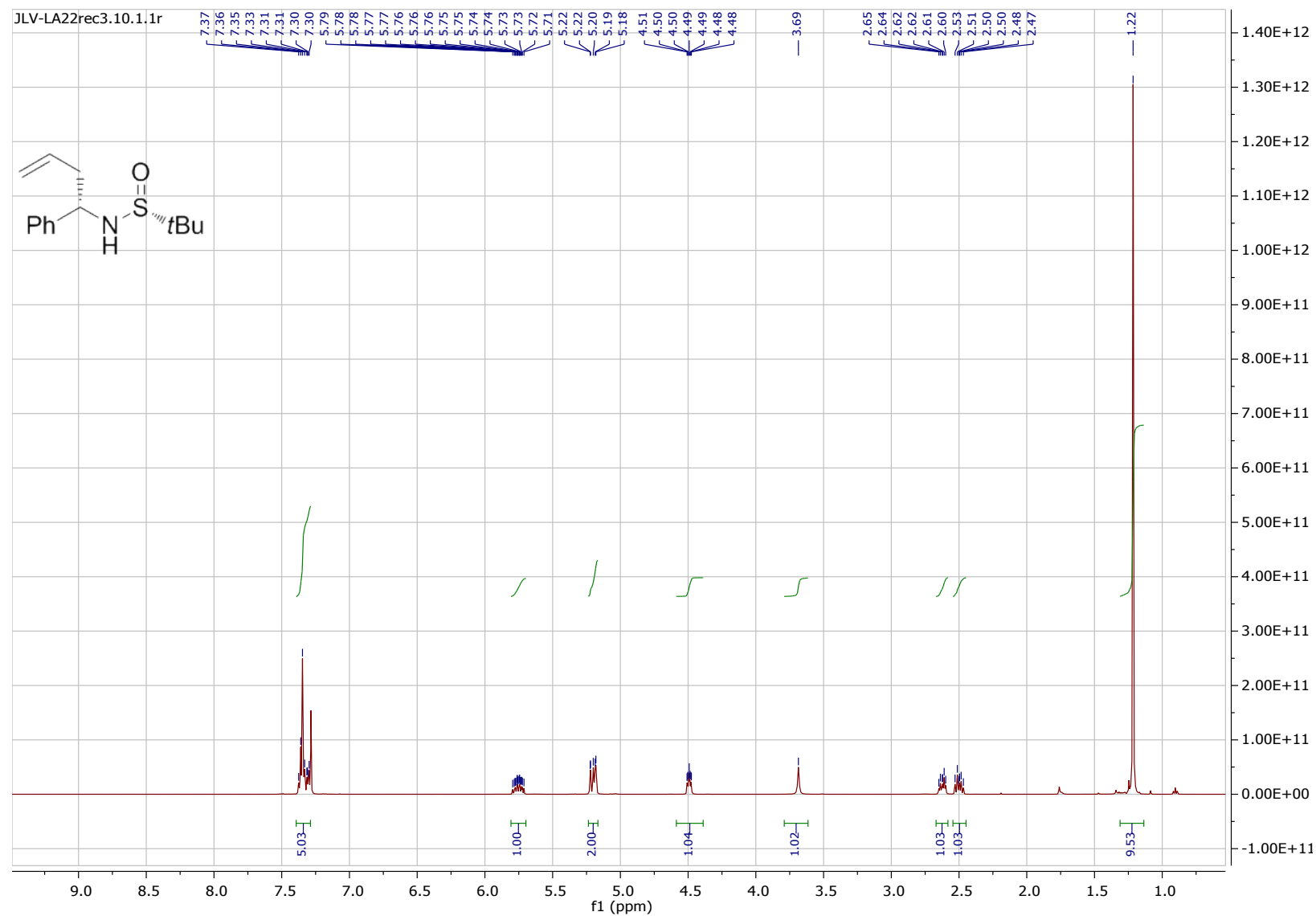


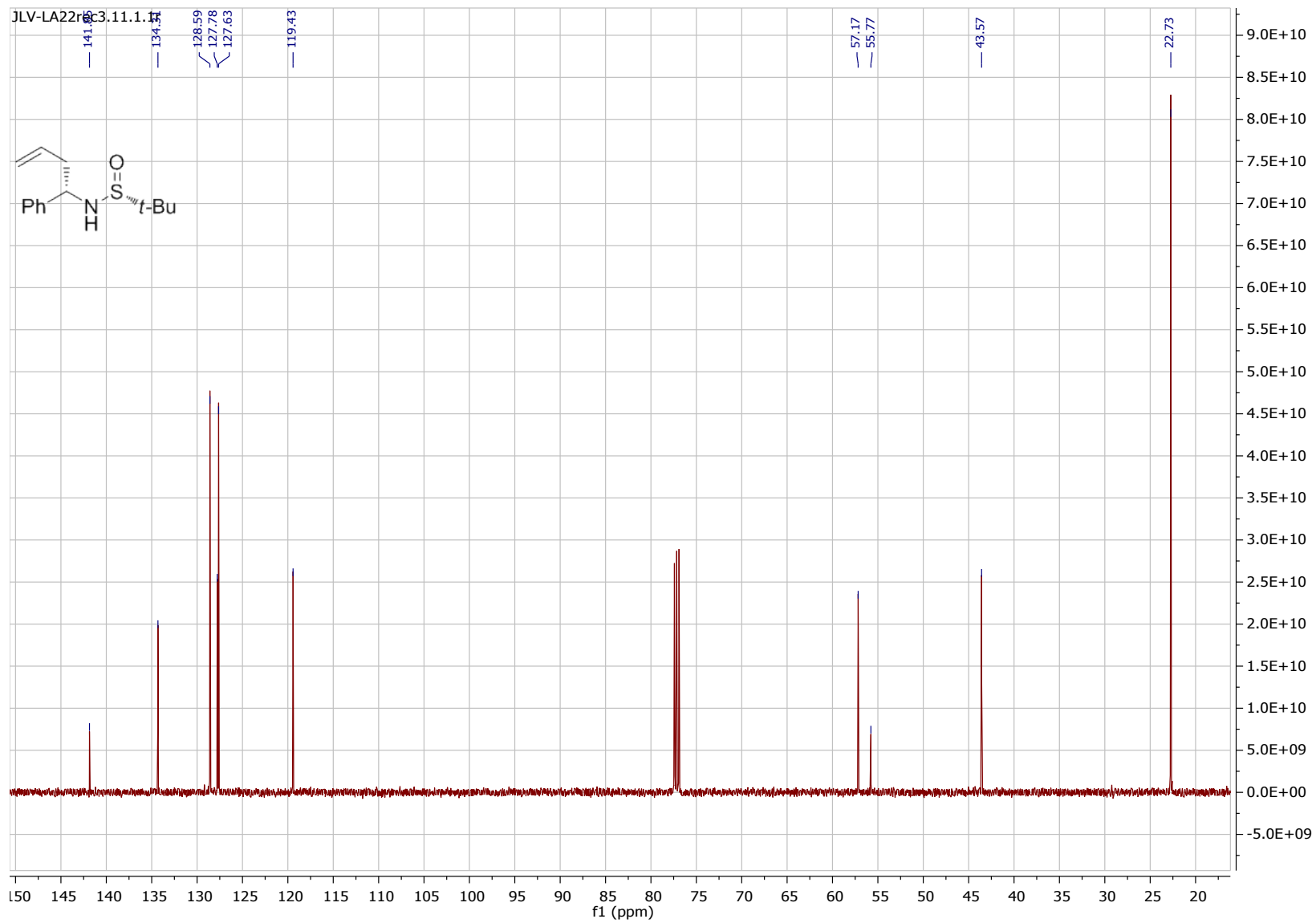
(E)-N-(Benzyl)-2-methyl-1-phenylhexa-1,5-dien-3-amine 1j



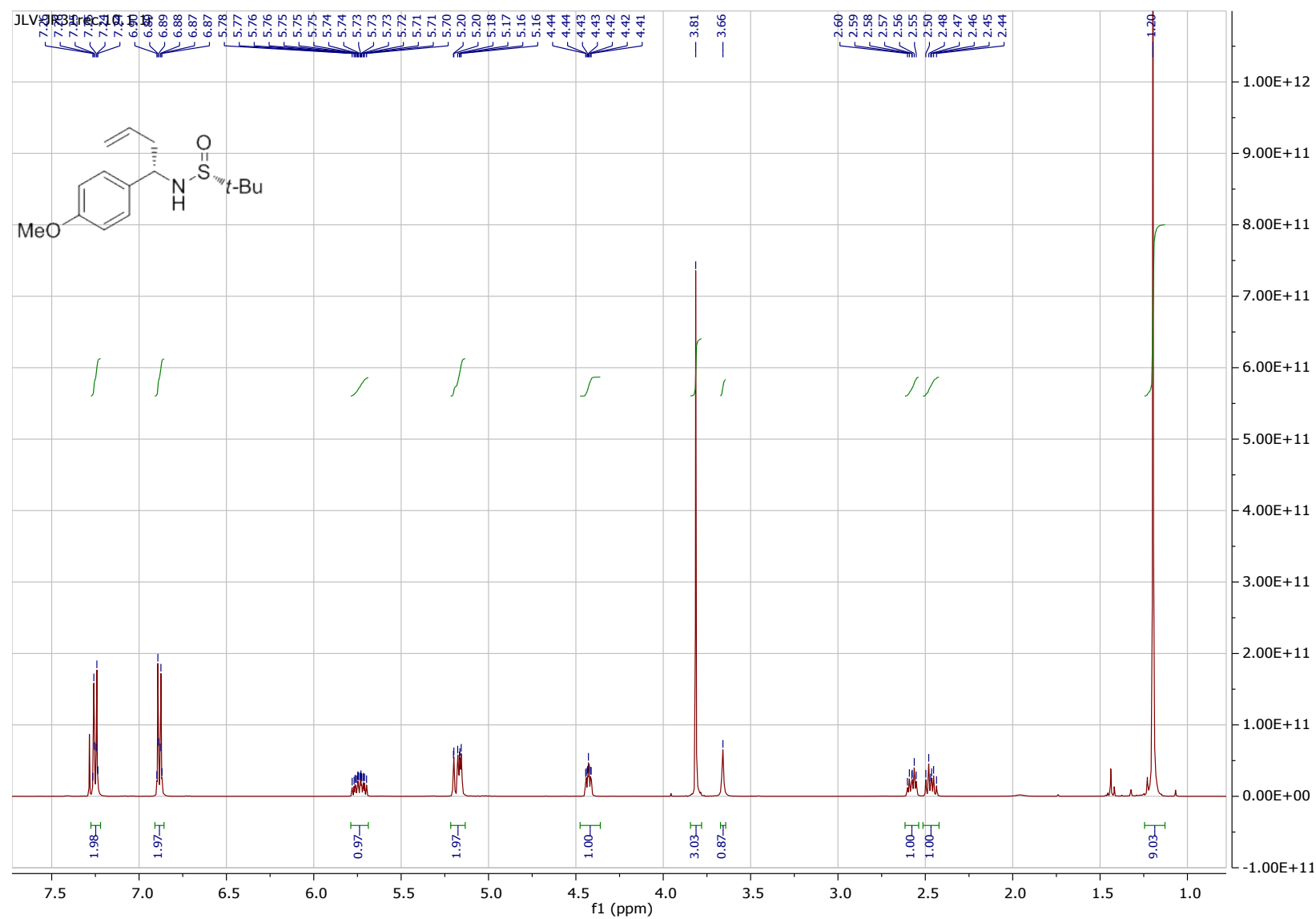


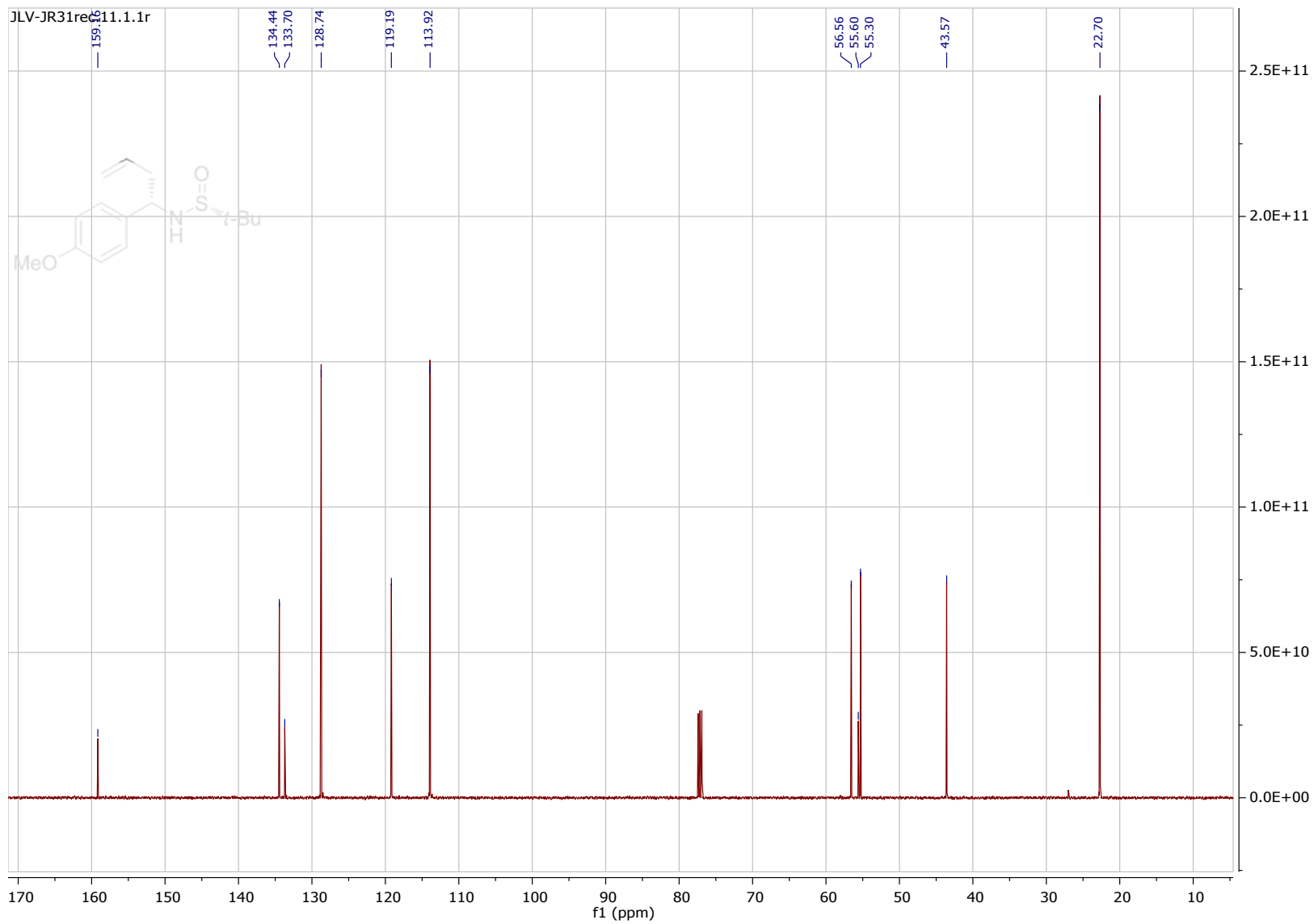
(R)-2-Methyl-N-[(S)-1-phenylbut-3-en-1-yl]propane-2-sulfonamide 5a



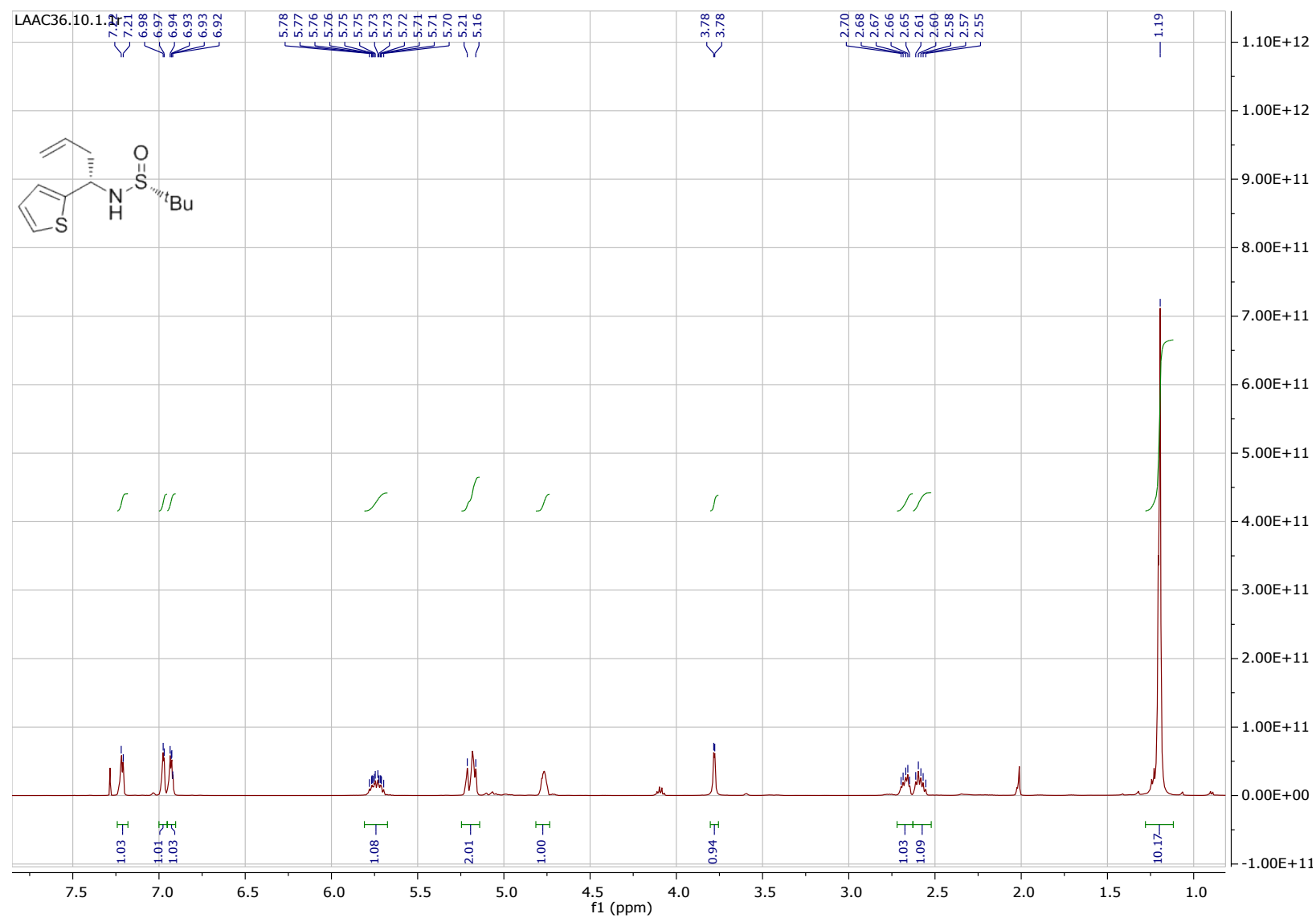


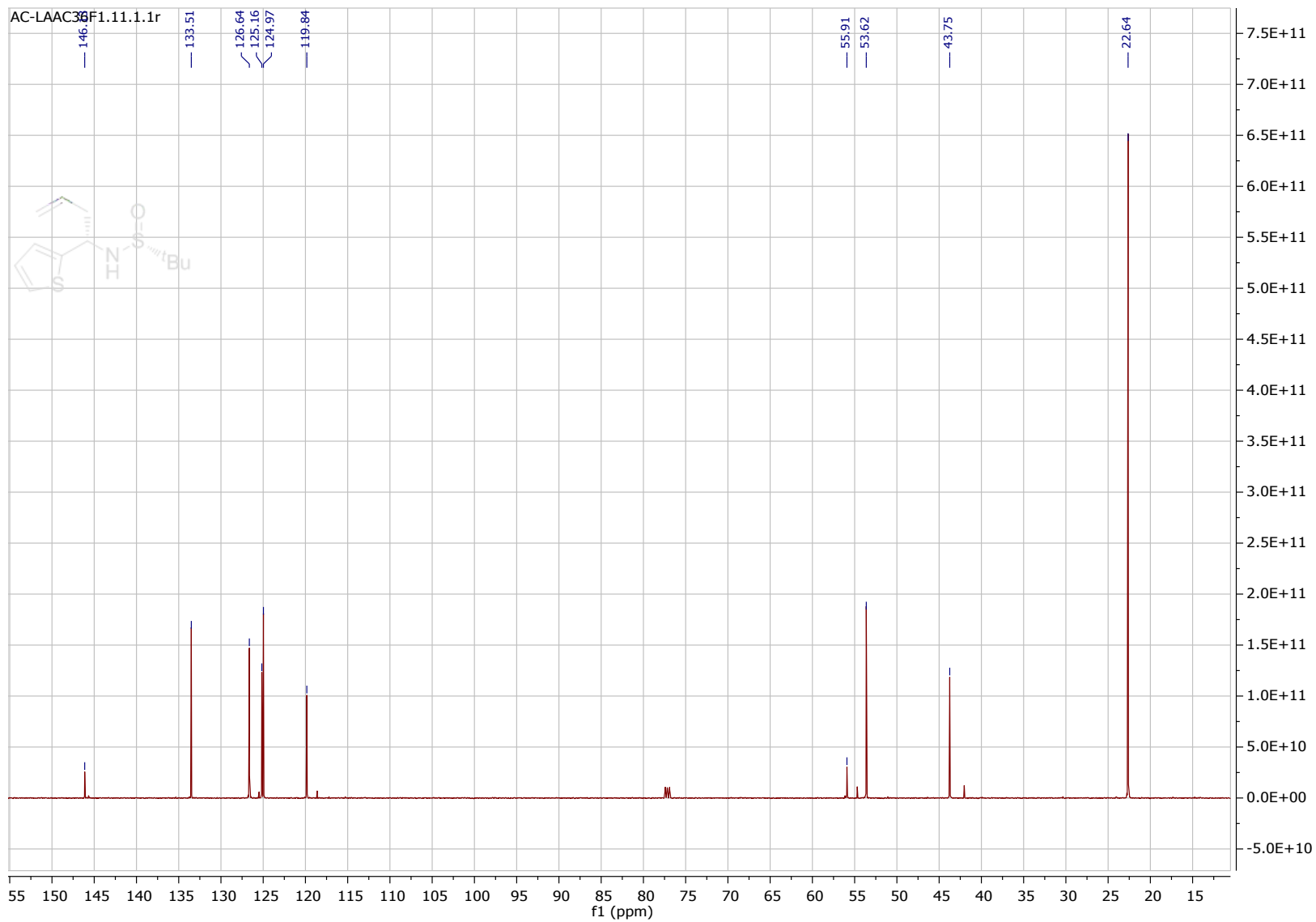
(R)-N-[(S)-1-(4-Methoxyphenyl)but-3-en-1-yl]-2-methylpropane-2-sulfonamide 5k



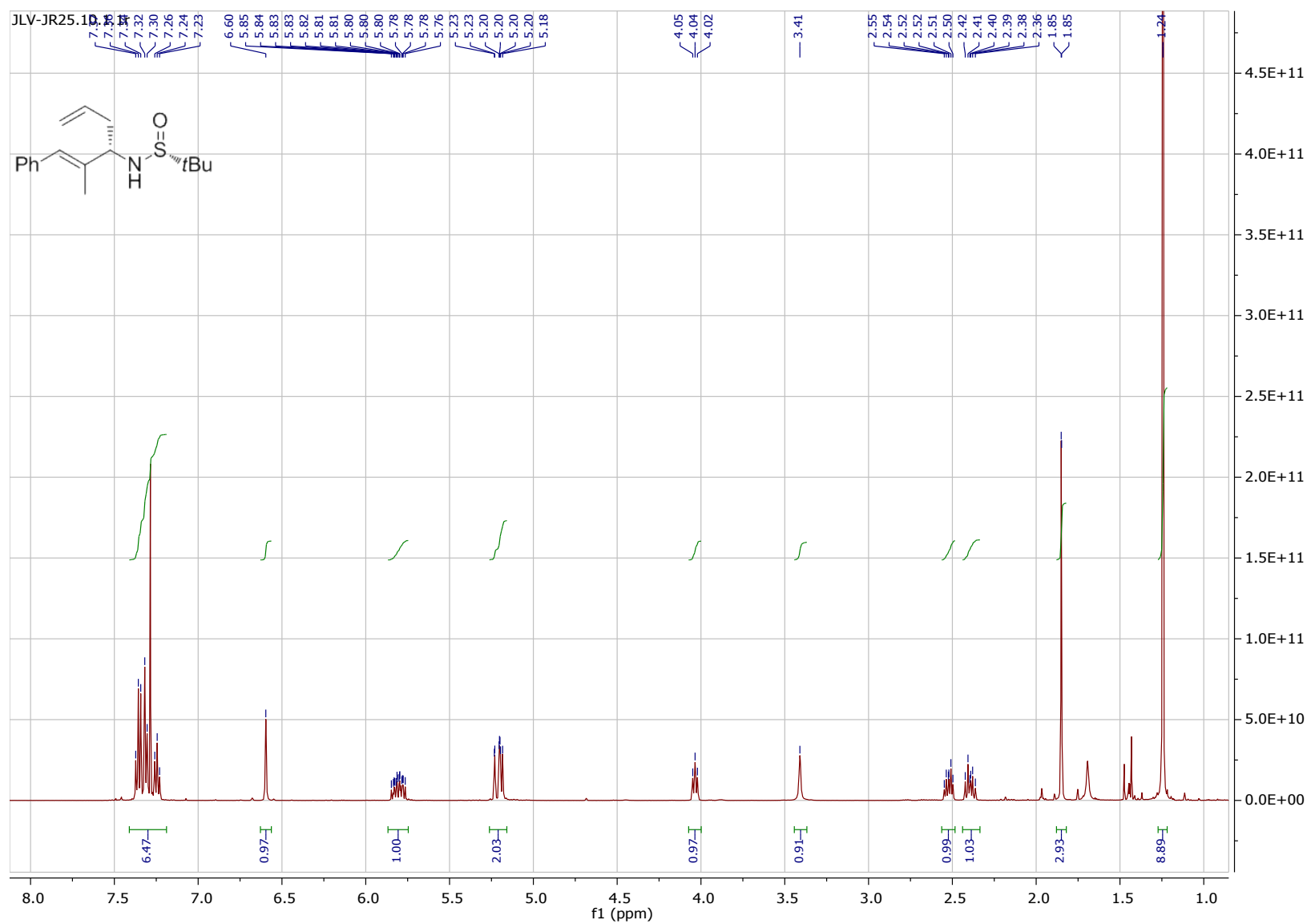


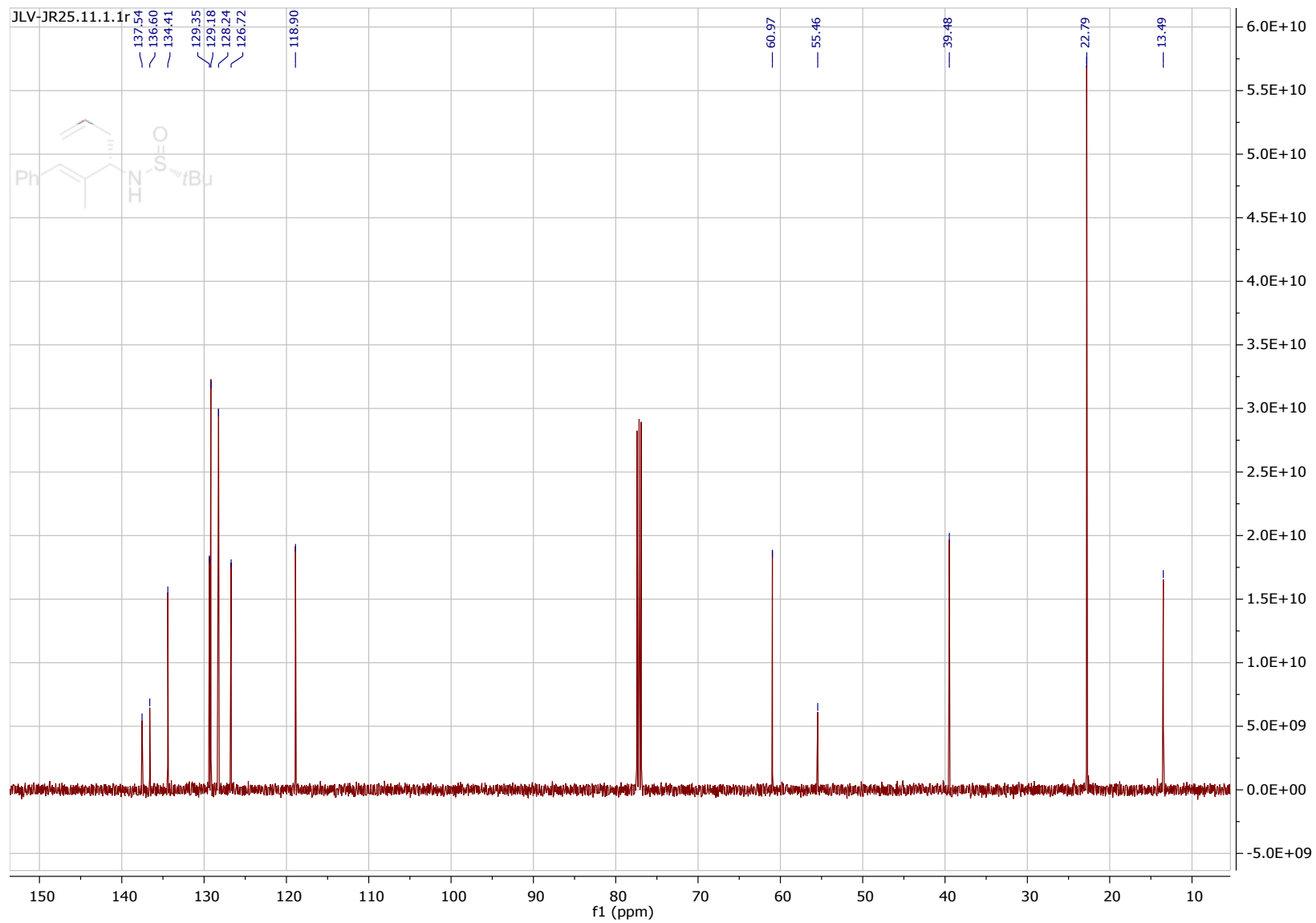
(R)-2-Methyl-N-[(S)-1-(thiophen-2-yl)but-3-en-1-yl]propane-2-sulfonamide 5e



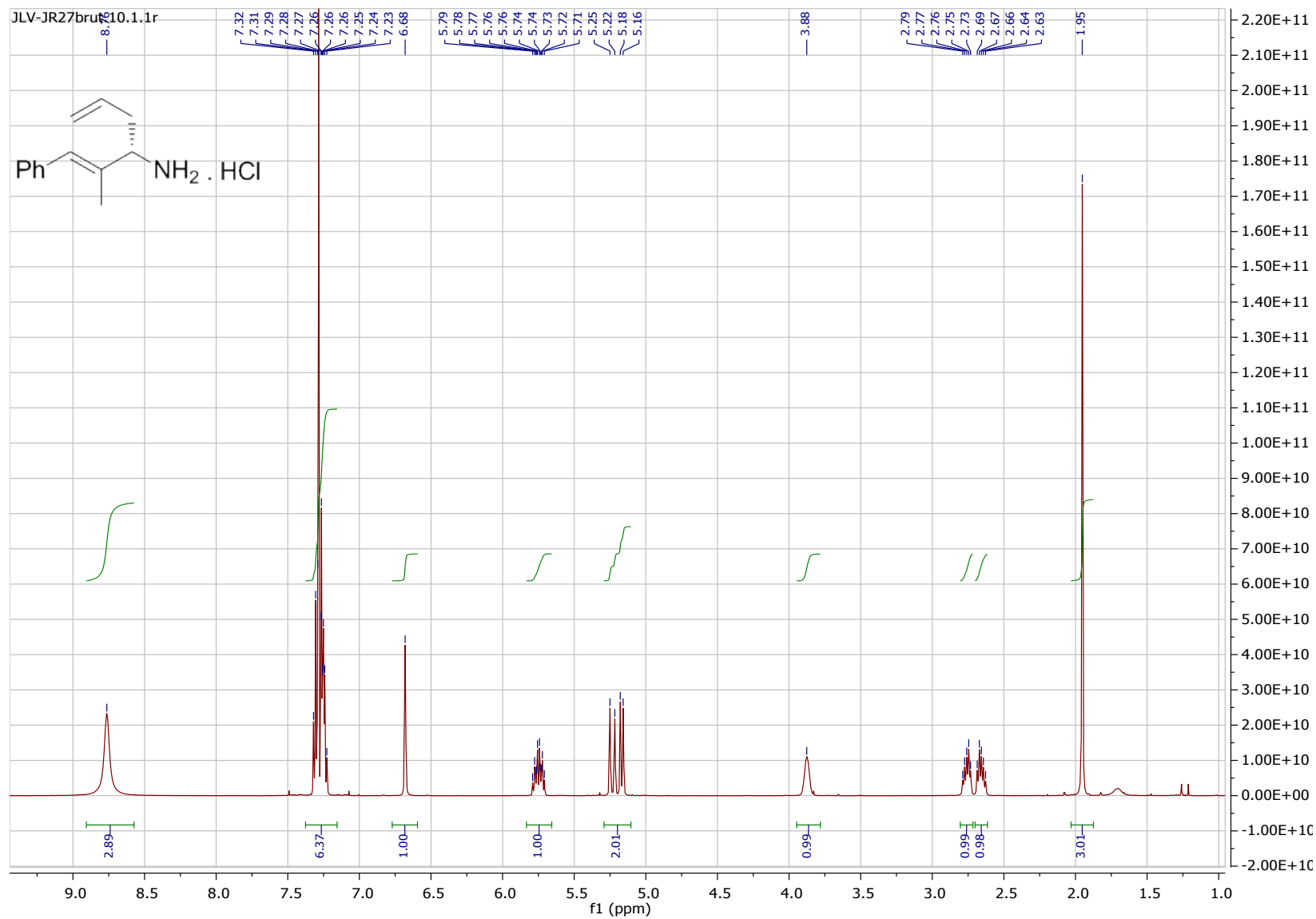


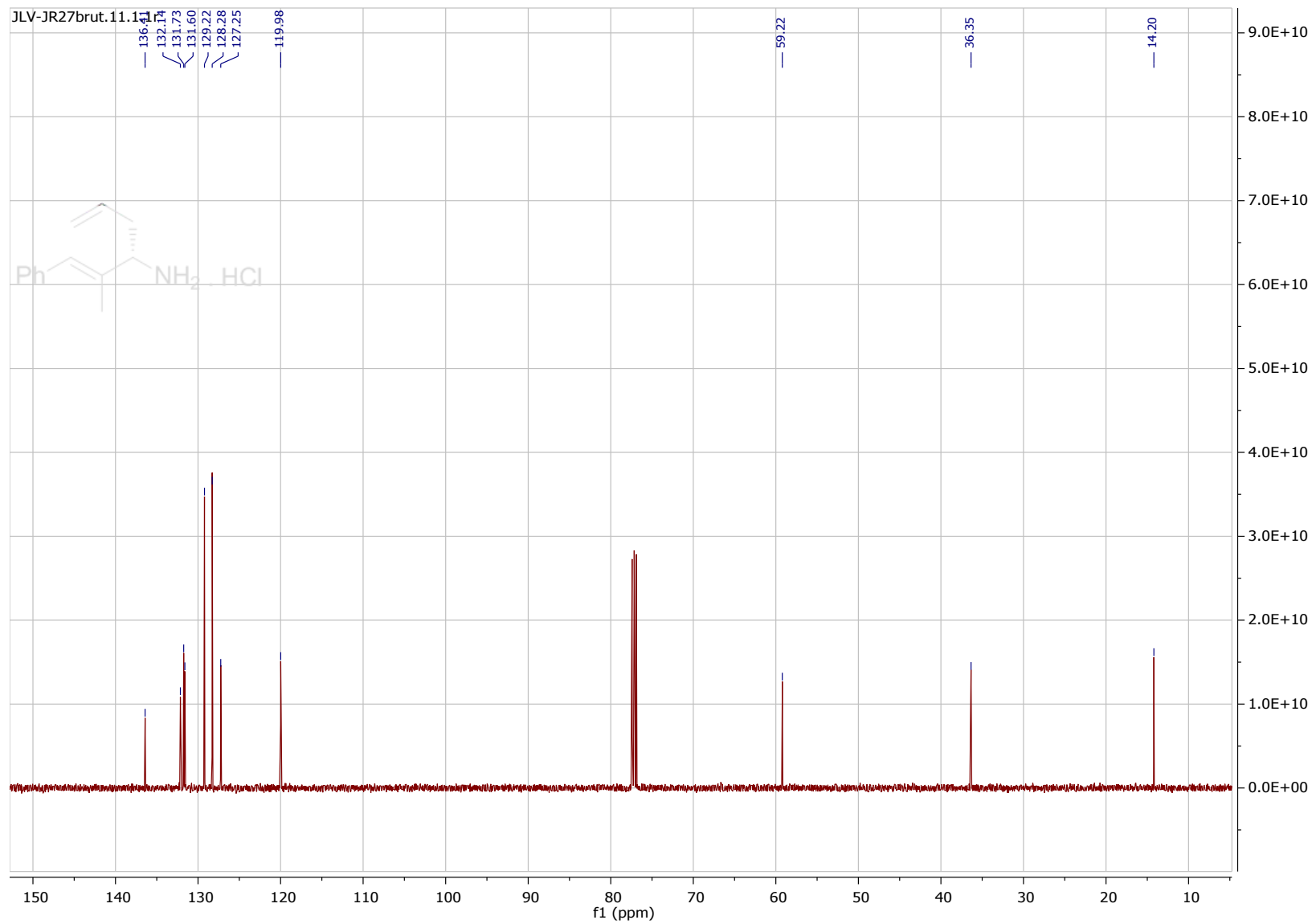
(R)-2-Methyl-N-[(S,E)-2-methyl-1-phenylhexa-1,5-dien-3-yl]propane-2-sulfonamide 5I



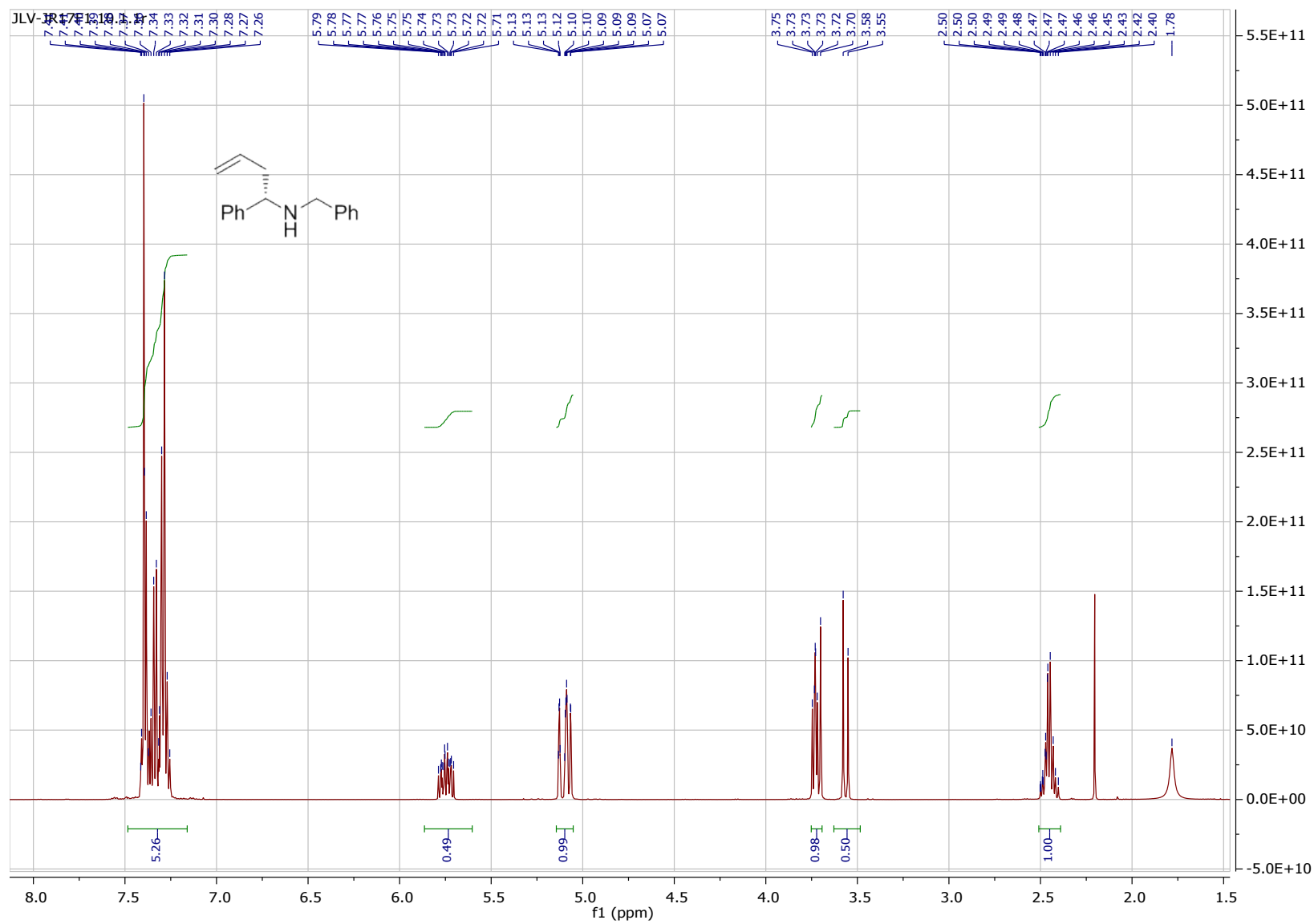


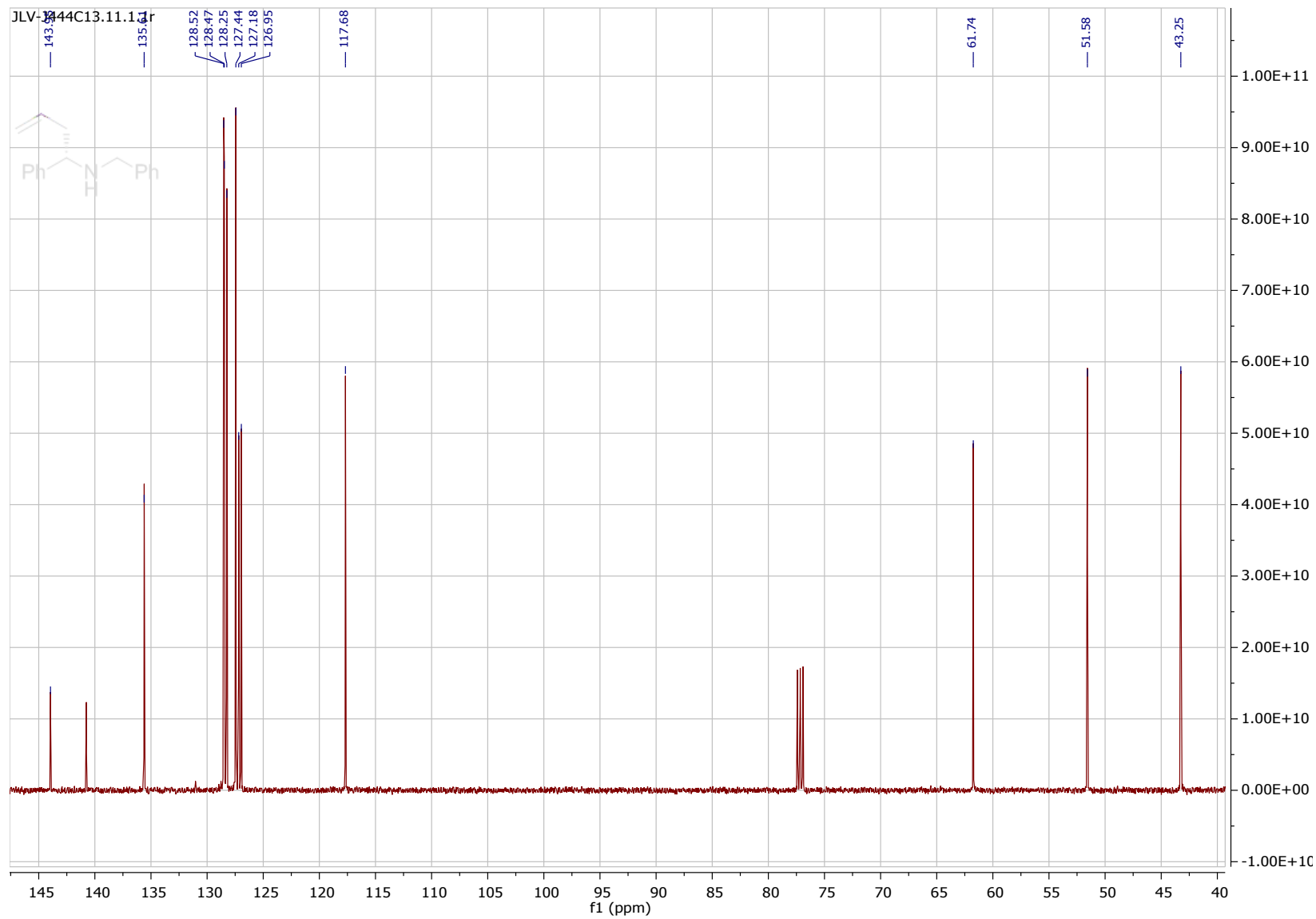
(S,E)-2-Methyl-1-phenylhexa-1,5-dien-3-amonium chloride



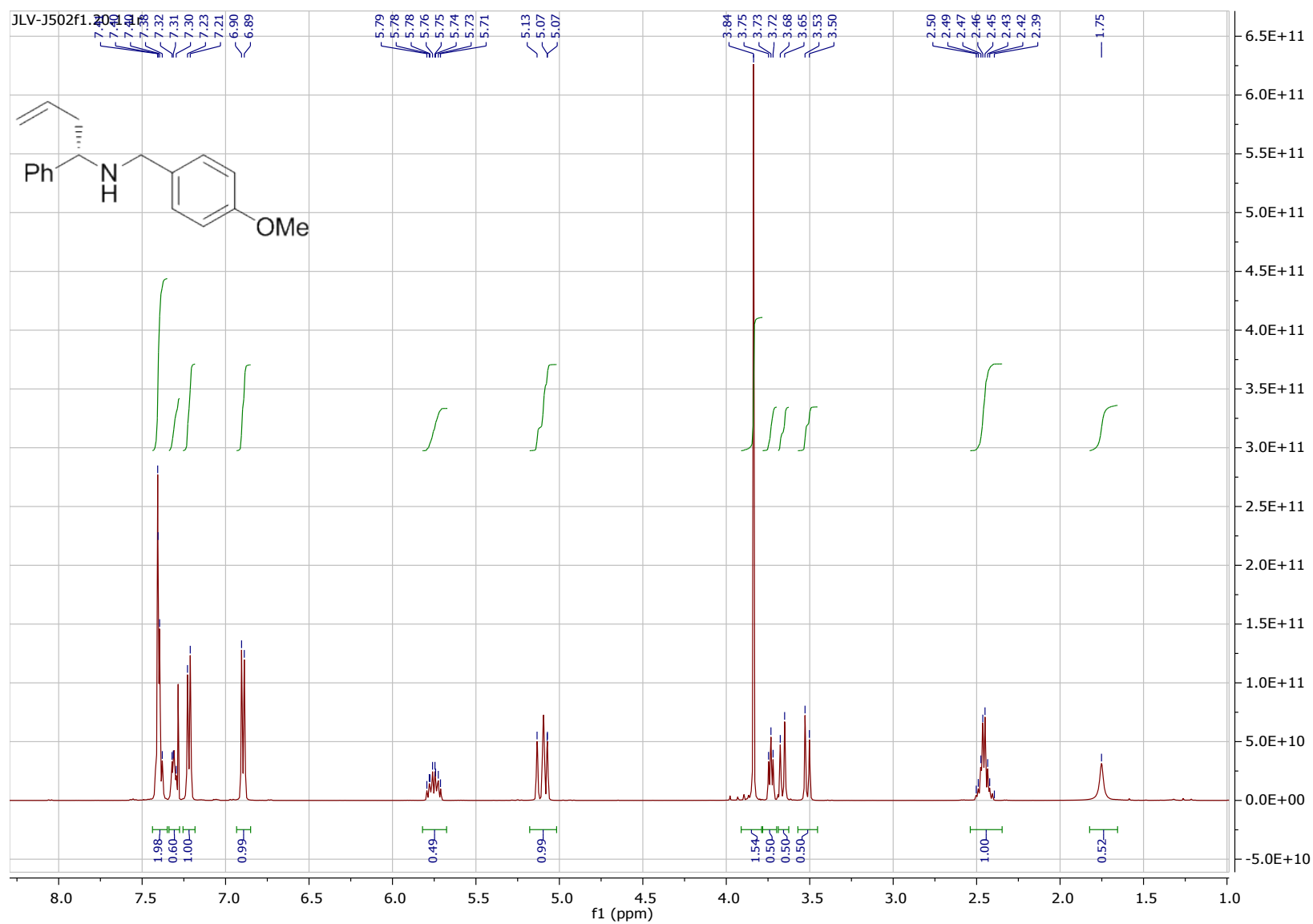


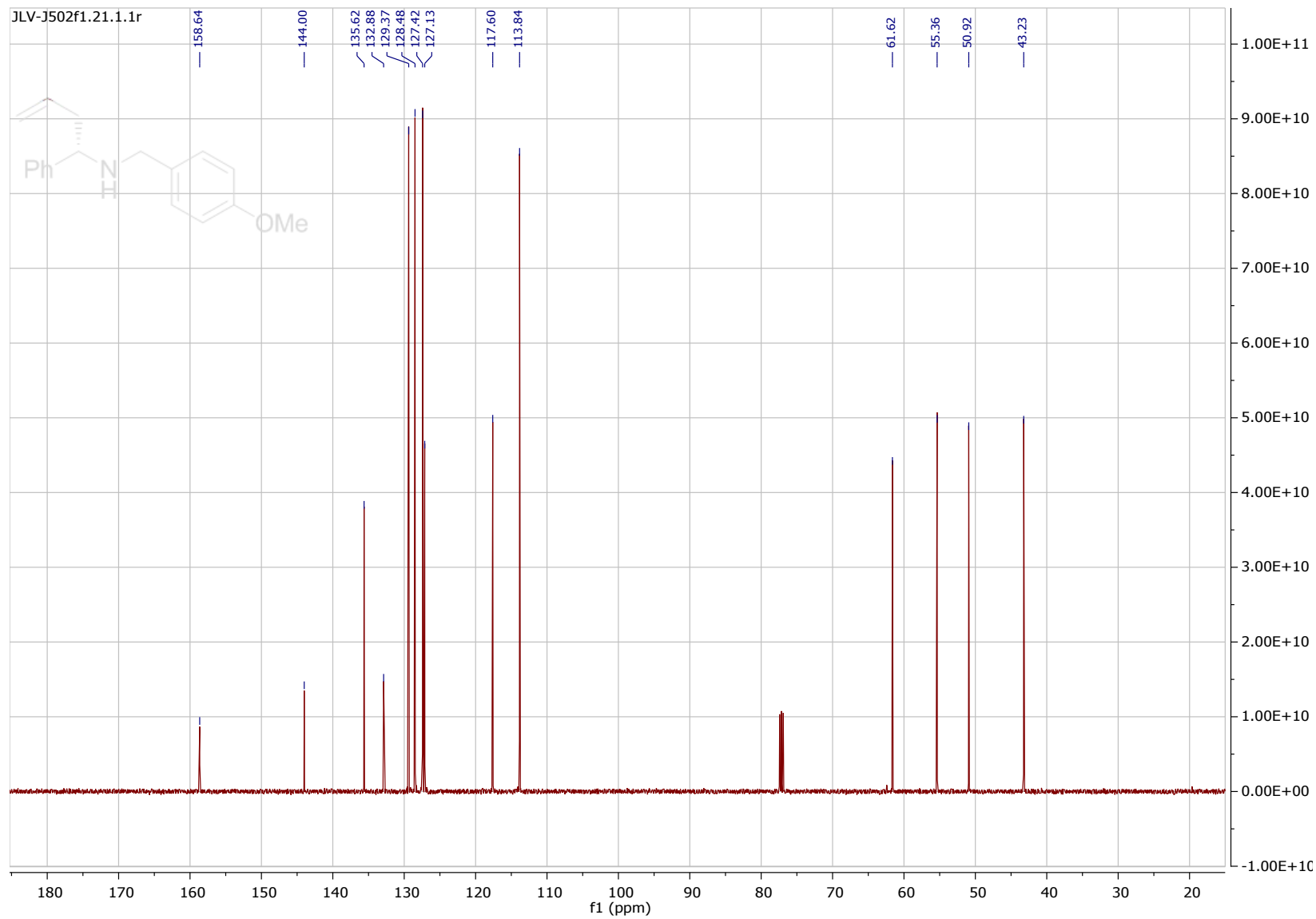
(S)-N-Benzyl-1-phenylbut-3-en-1-amine 1a



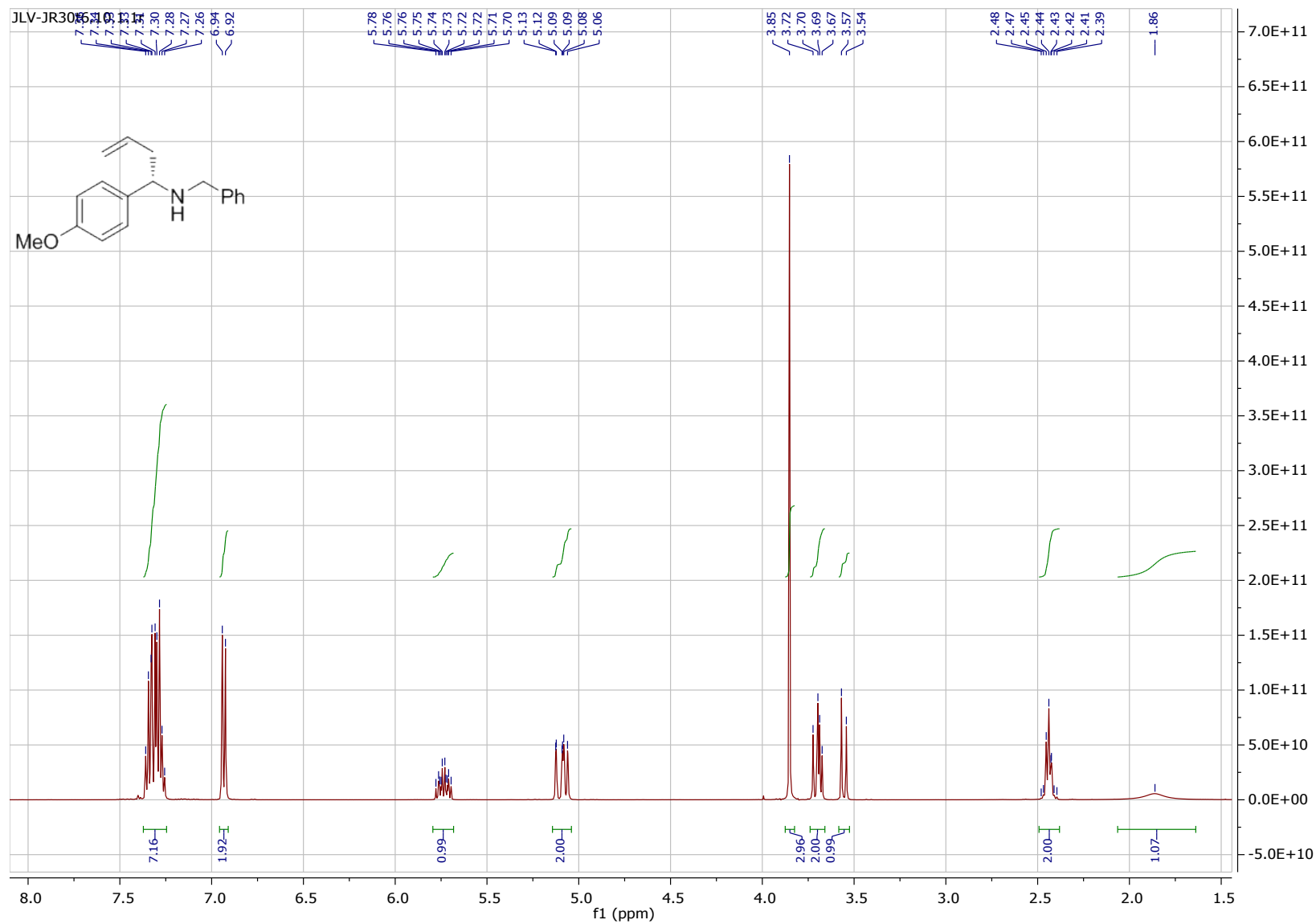


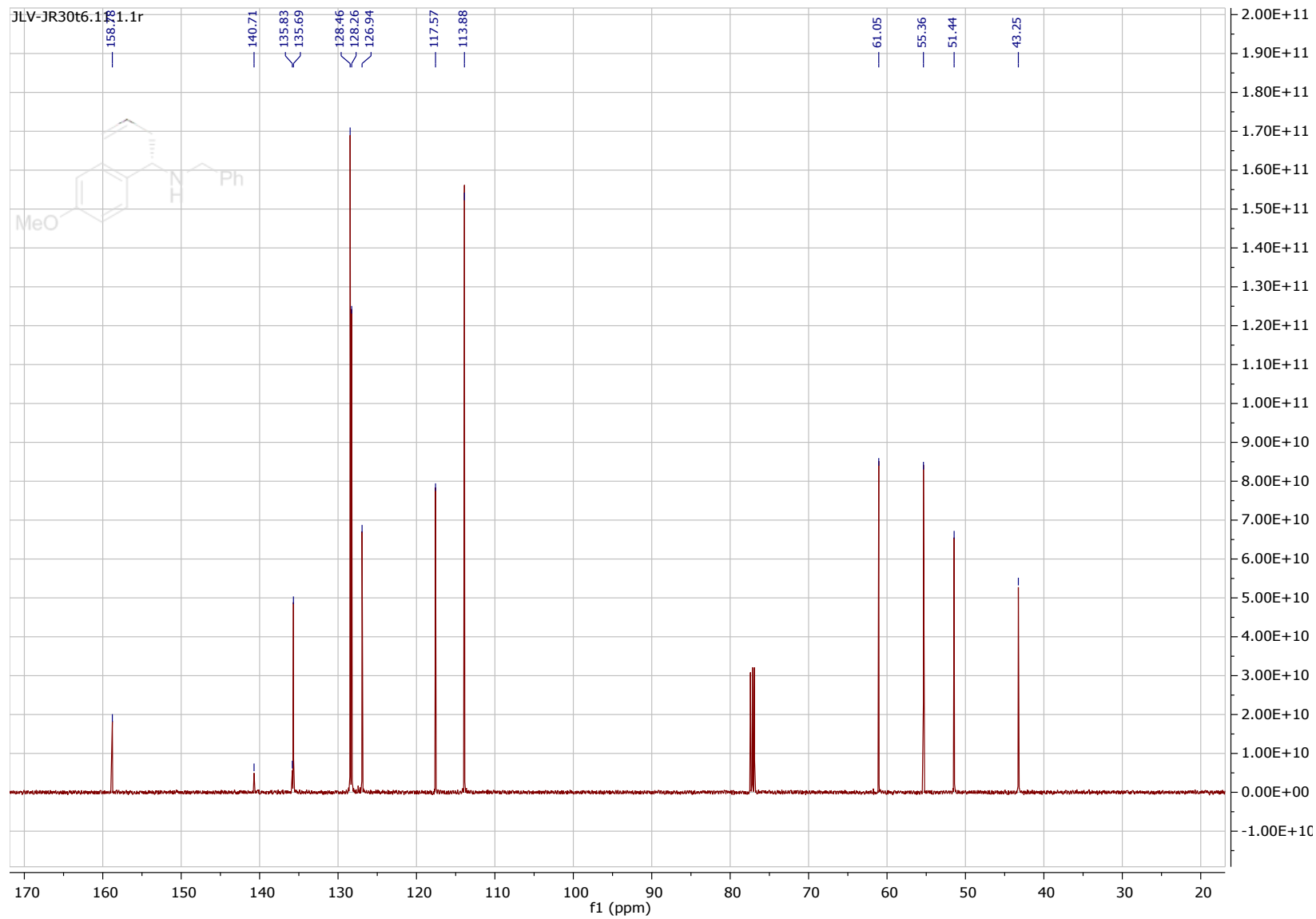
(S)-N-(4-Methoxybenzyl)-1-phenylbut-3-en-1-amine 1b



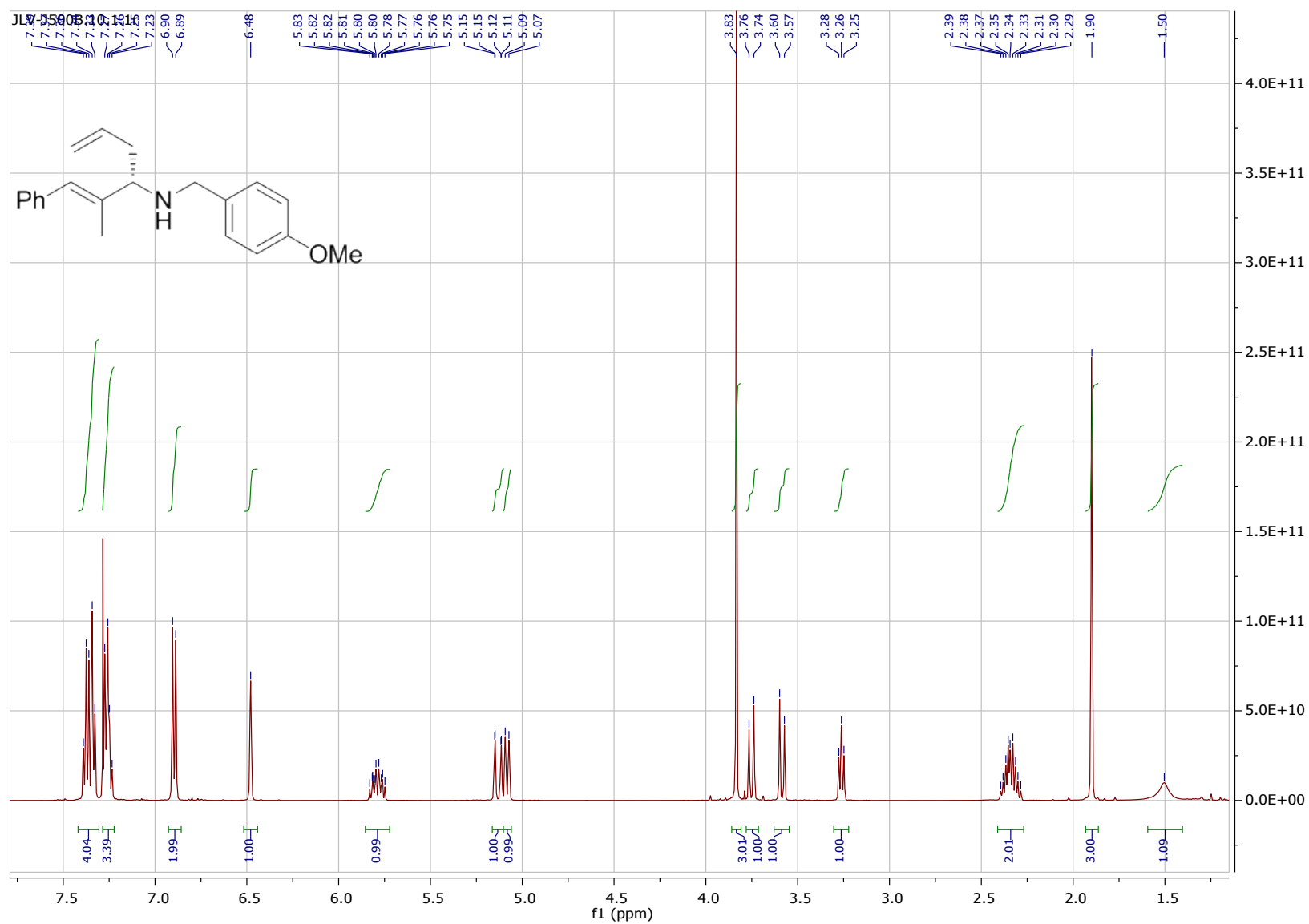


(S)-N-benzyl-1-(4-methoxyphenyl)but-3-en-1-amine 1k

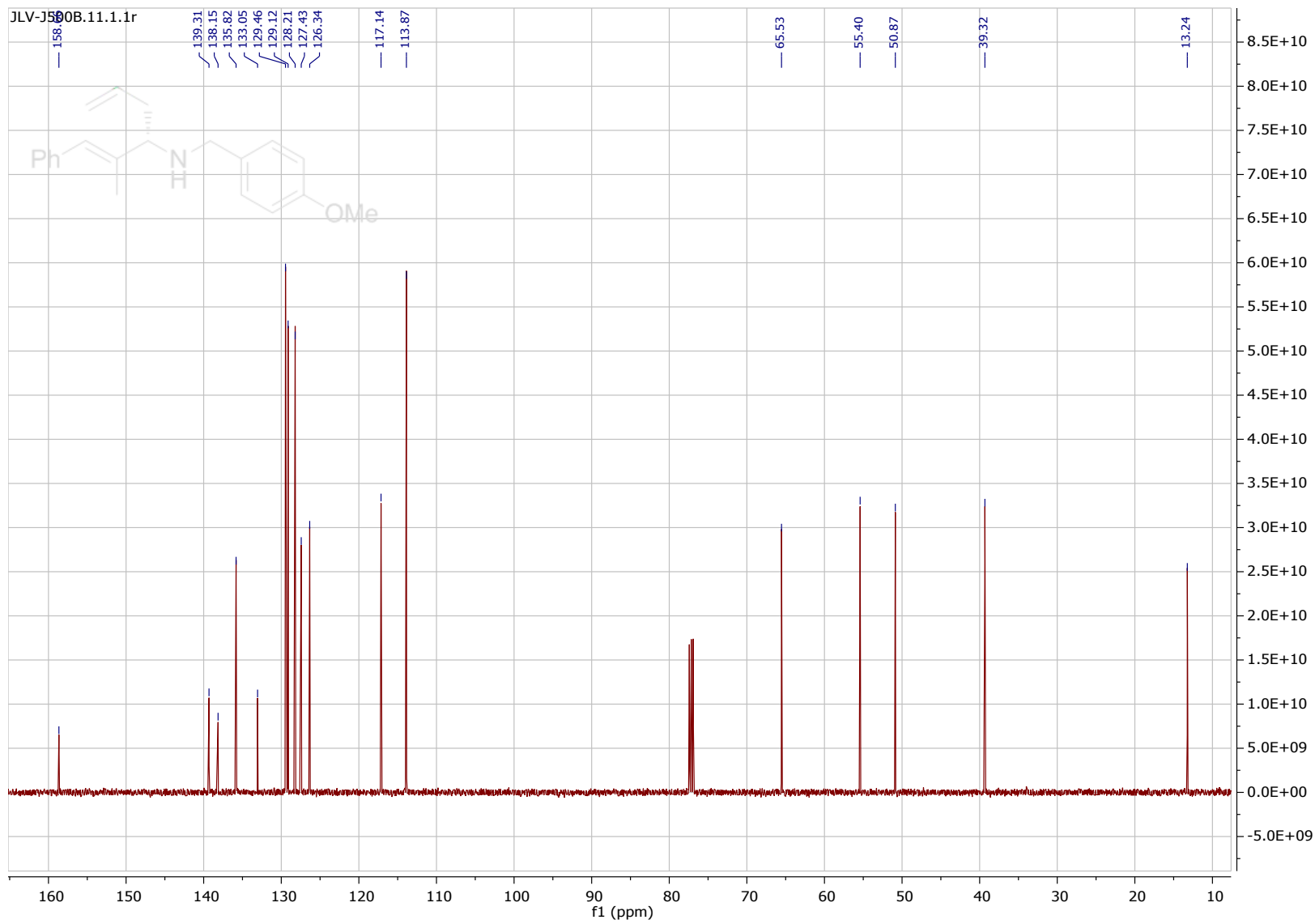




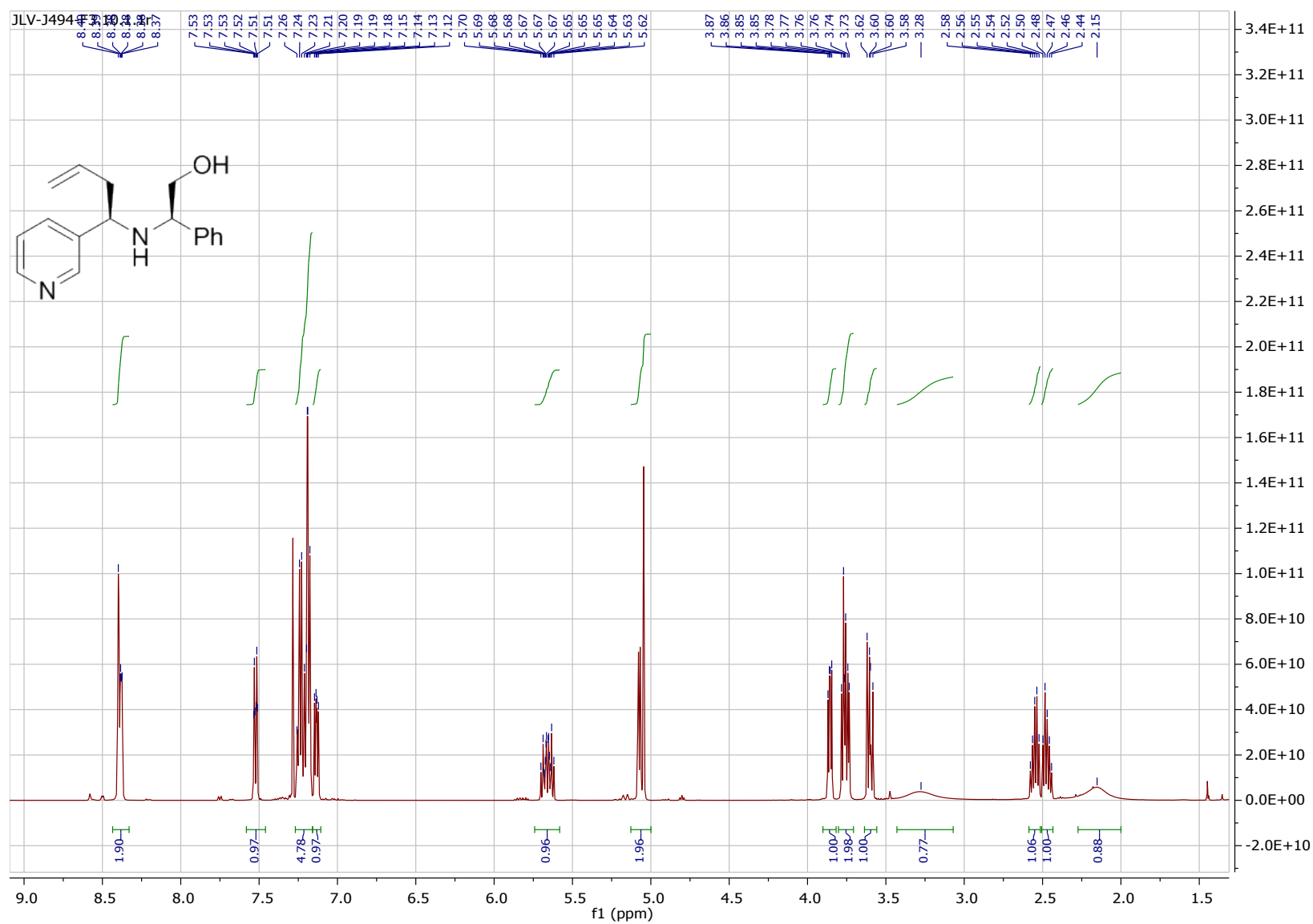
(S,E)-N-(4-Methoxybenzyl)-2-methyl-1-phenylhexa-1,5-dien-3-amine 11

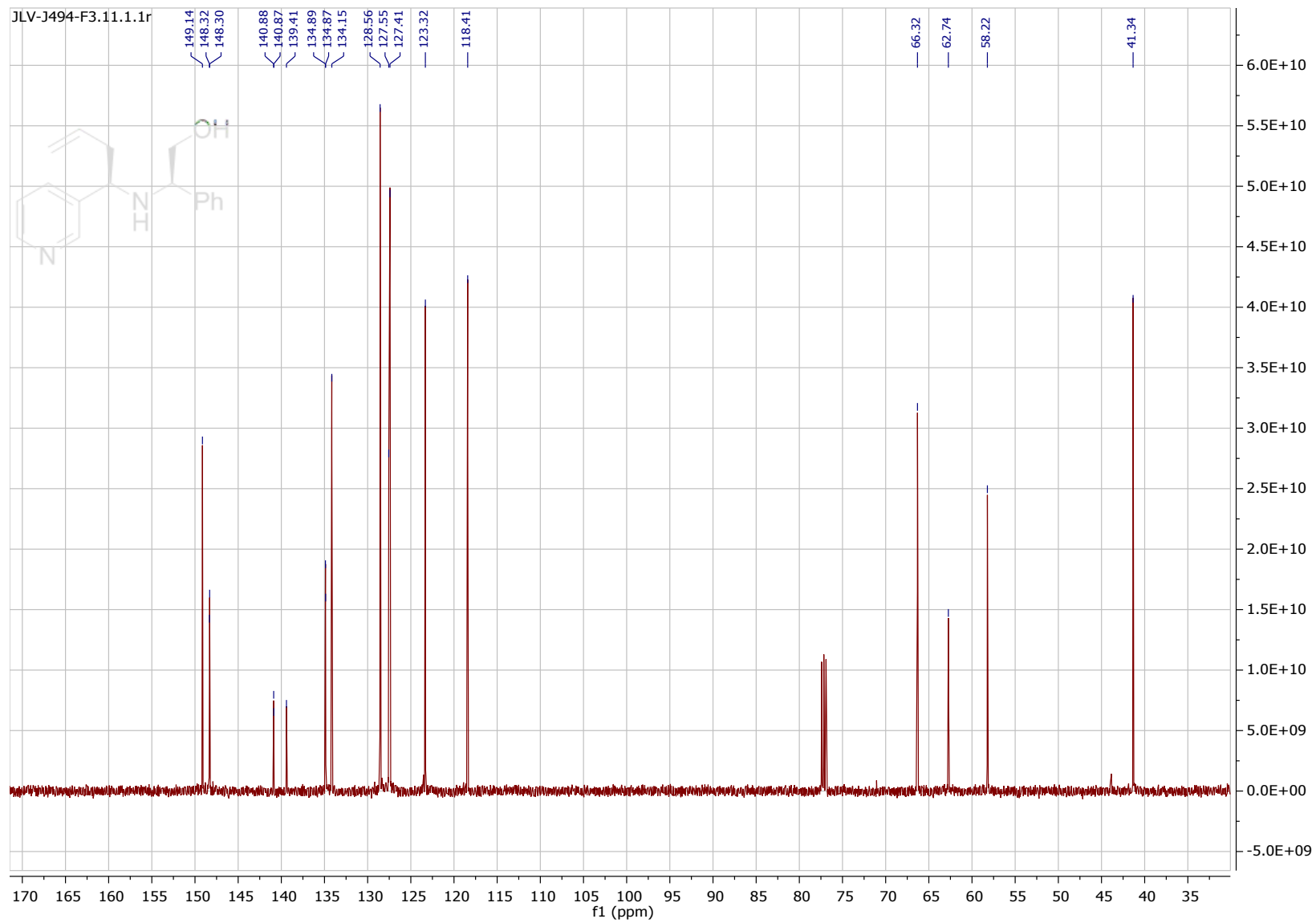


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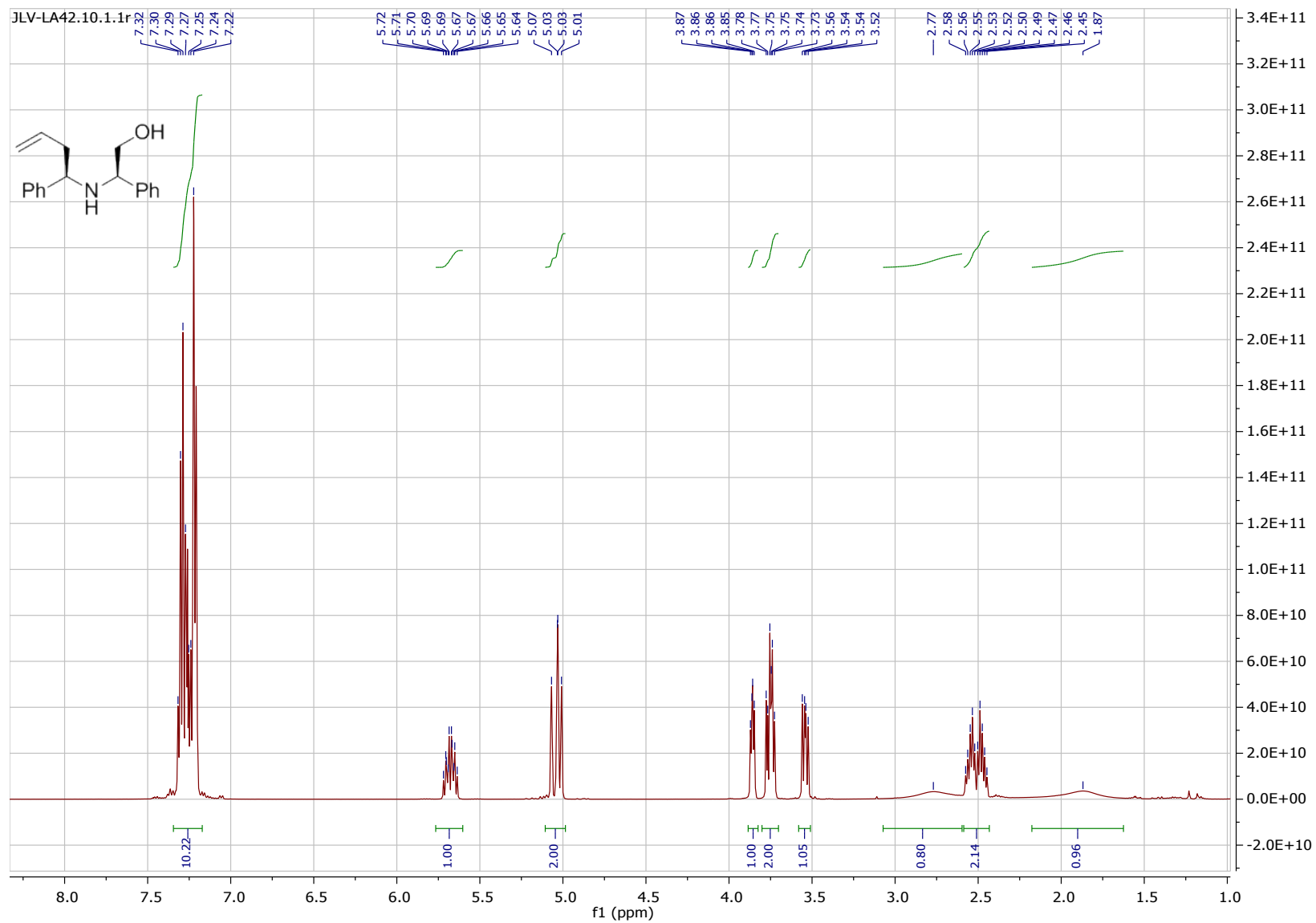


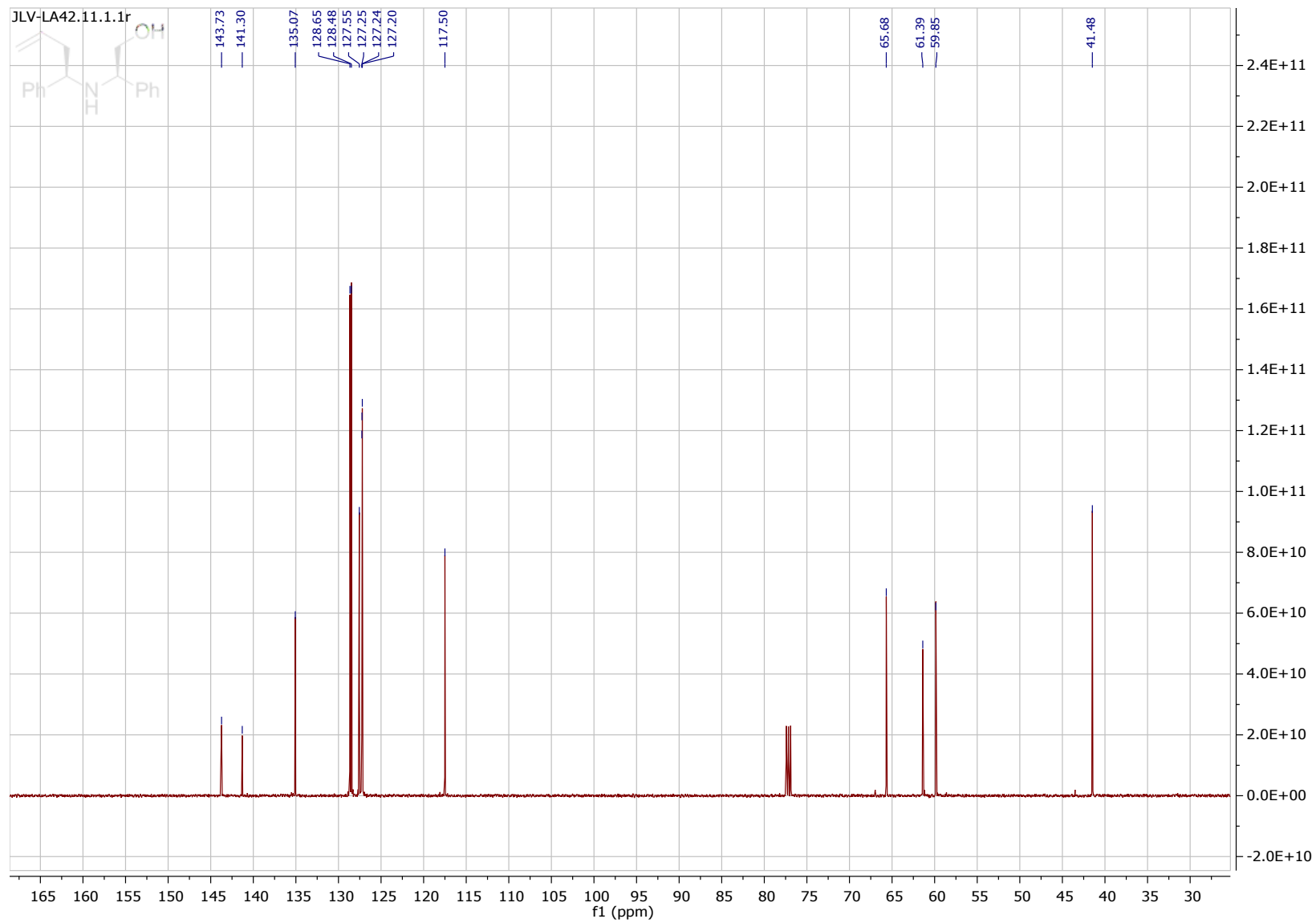
(R)-2-Phenyl-2-[(R)-1-(pyridin-3-yl)but-3-en-1-yl]amino}ethanol 5h



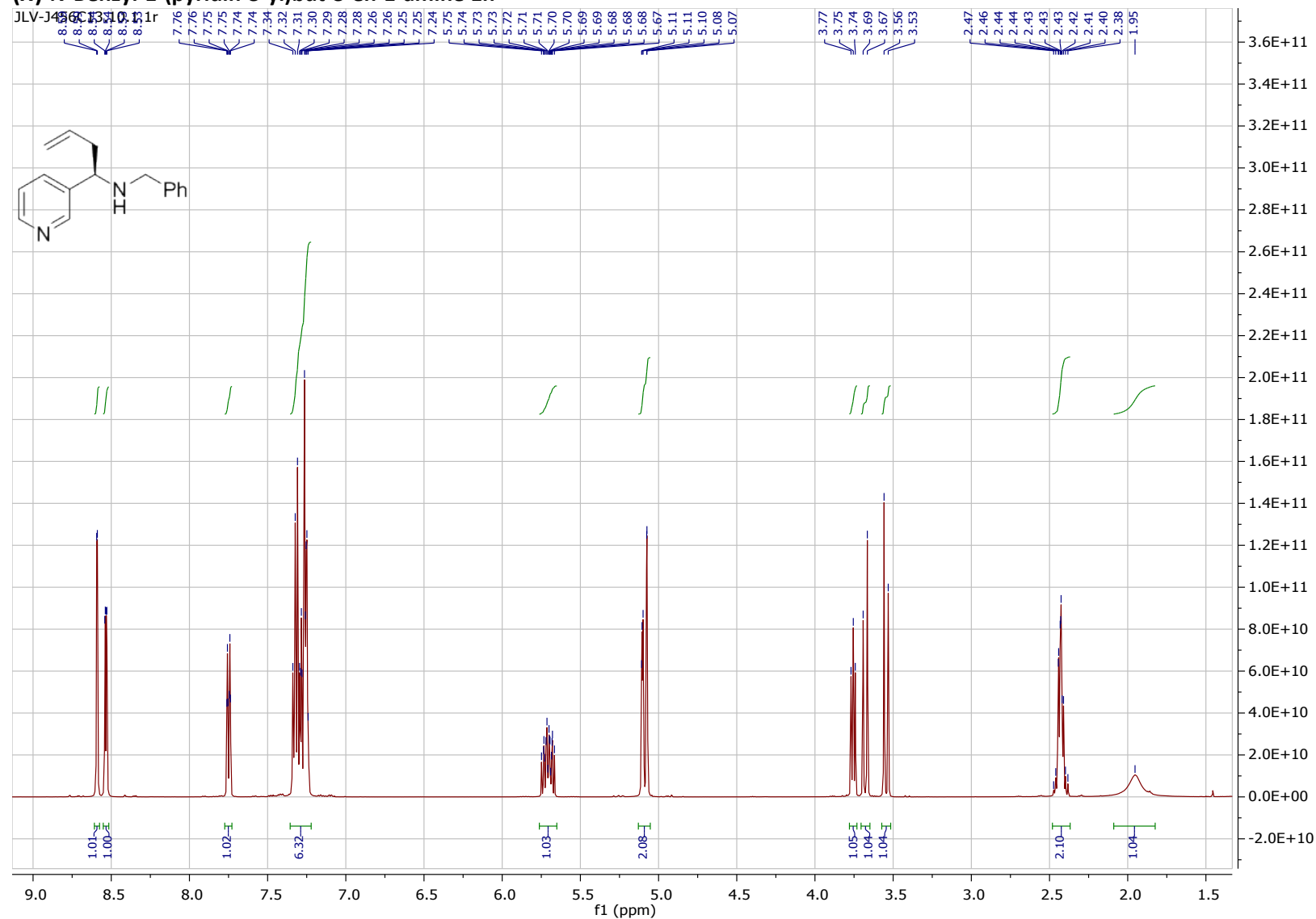


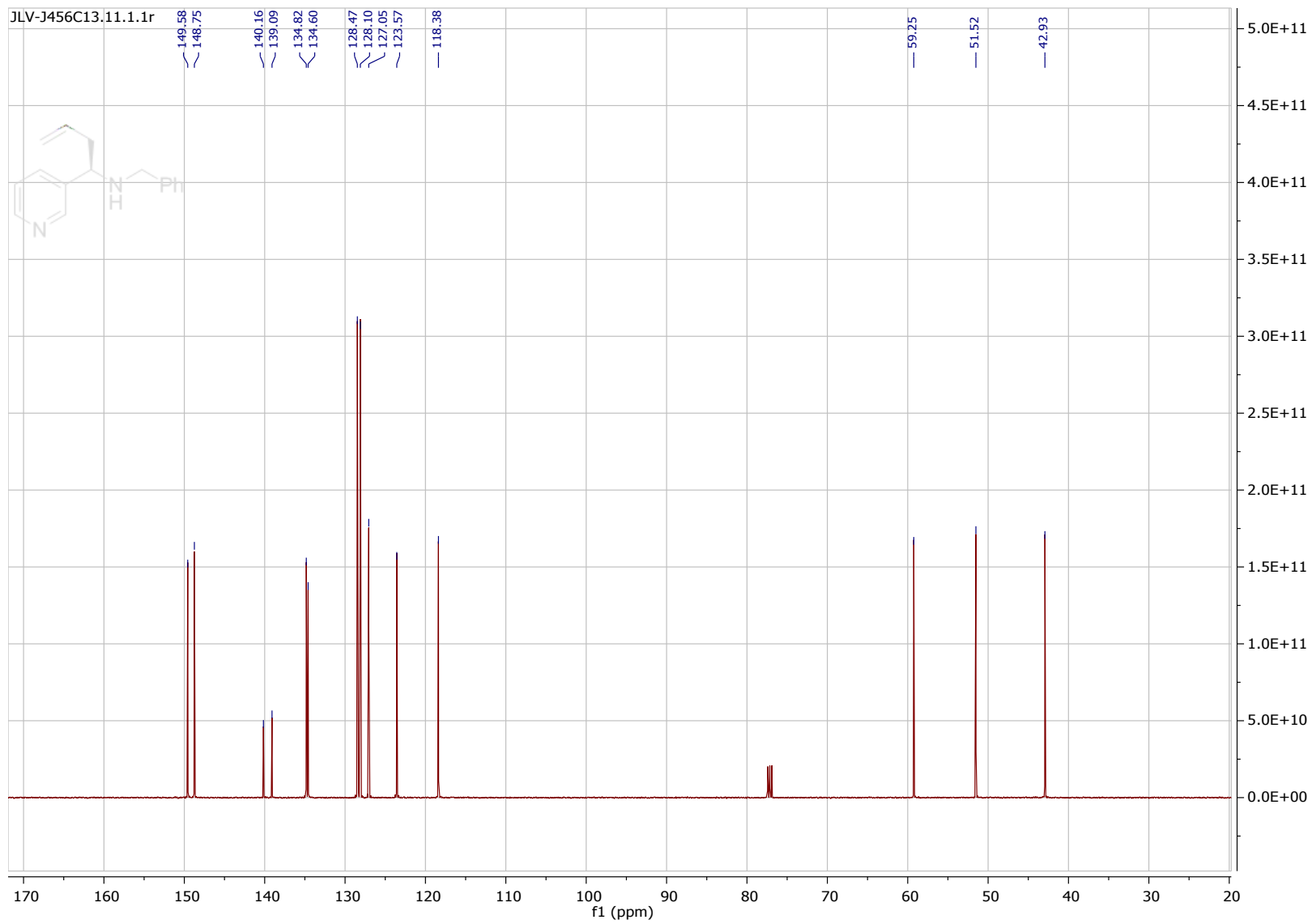
(R)-2-Phenyl-2-[(R)-1-phenylbut-3-en-1-yl]amino}ethanol



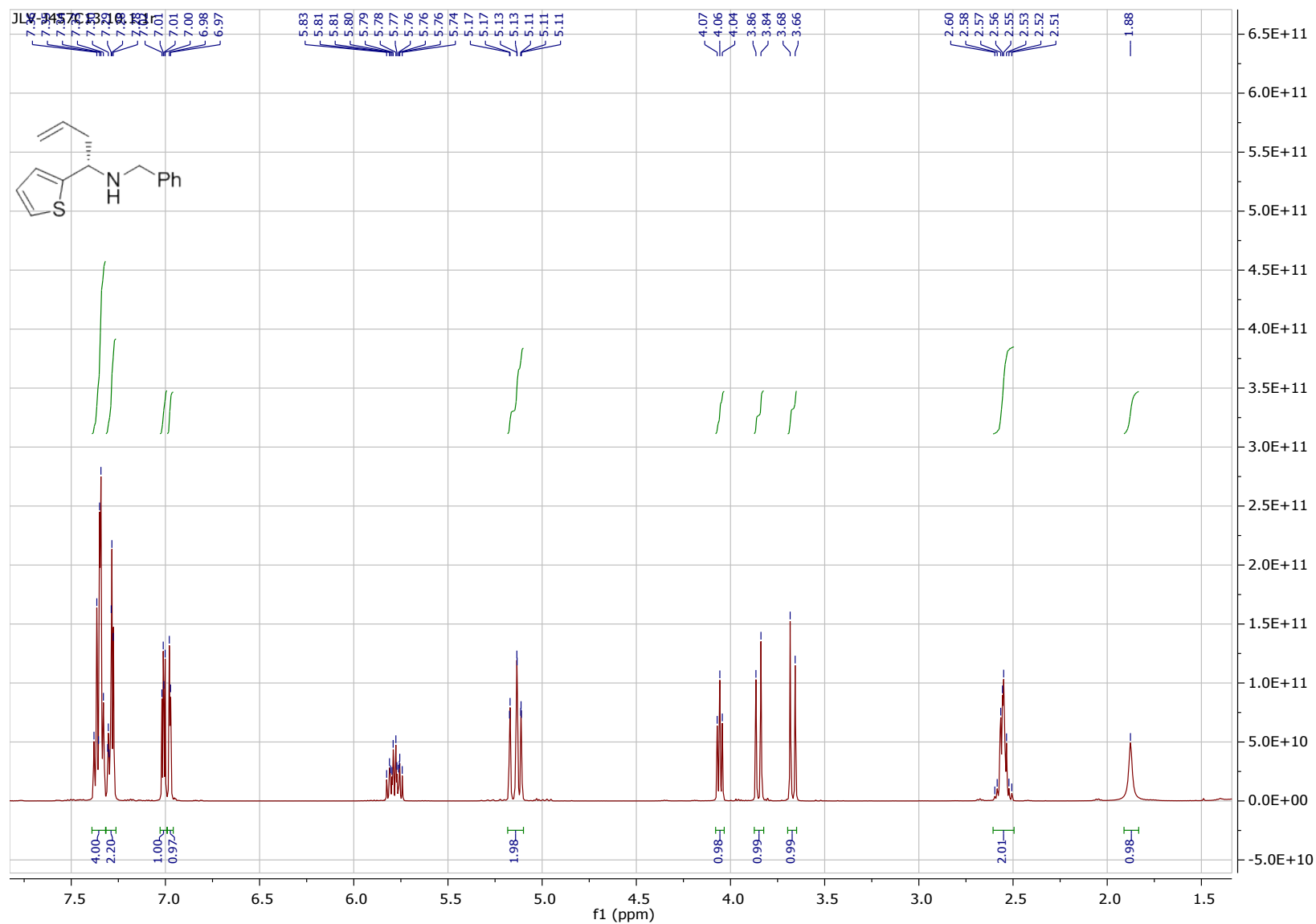


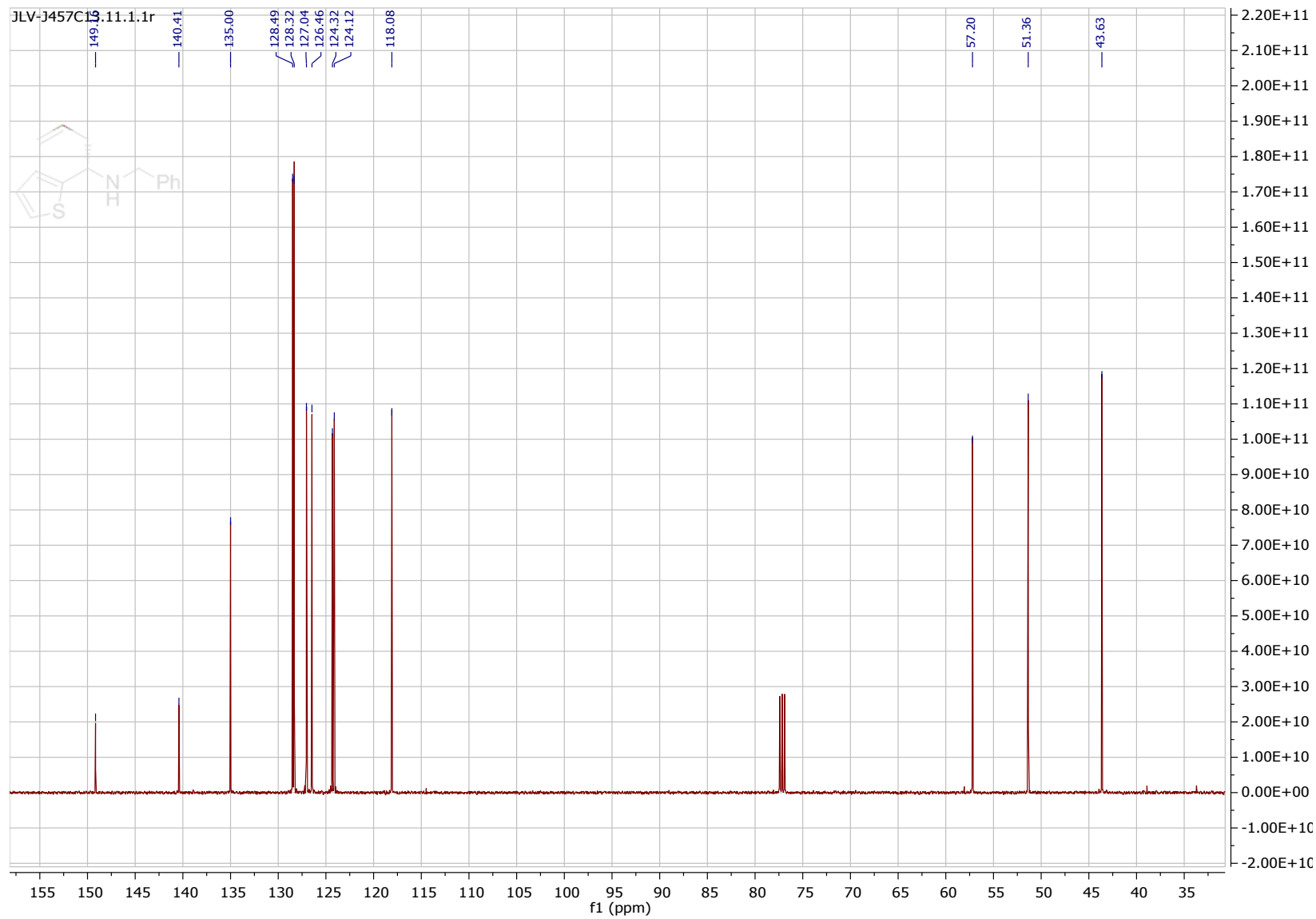
(R)-N-Benzyl-1-(pyridin-3-yl)but-3-en-1-amine 1h



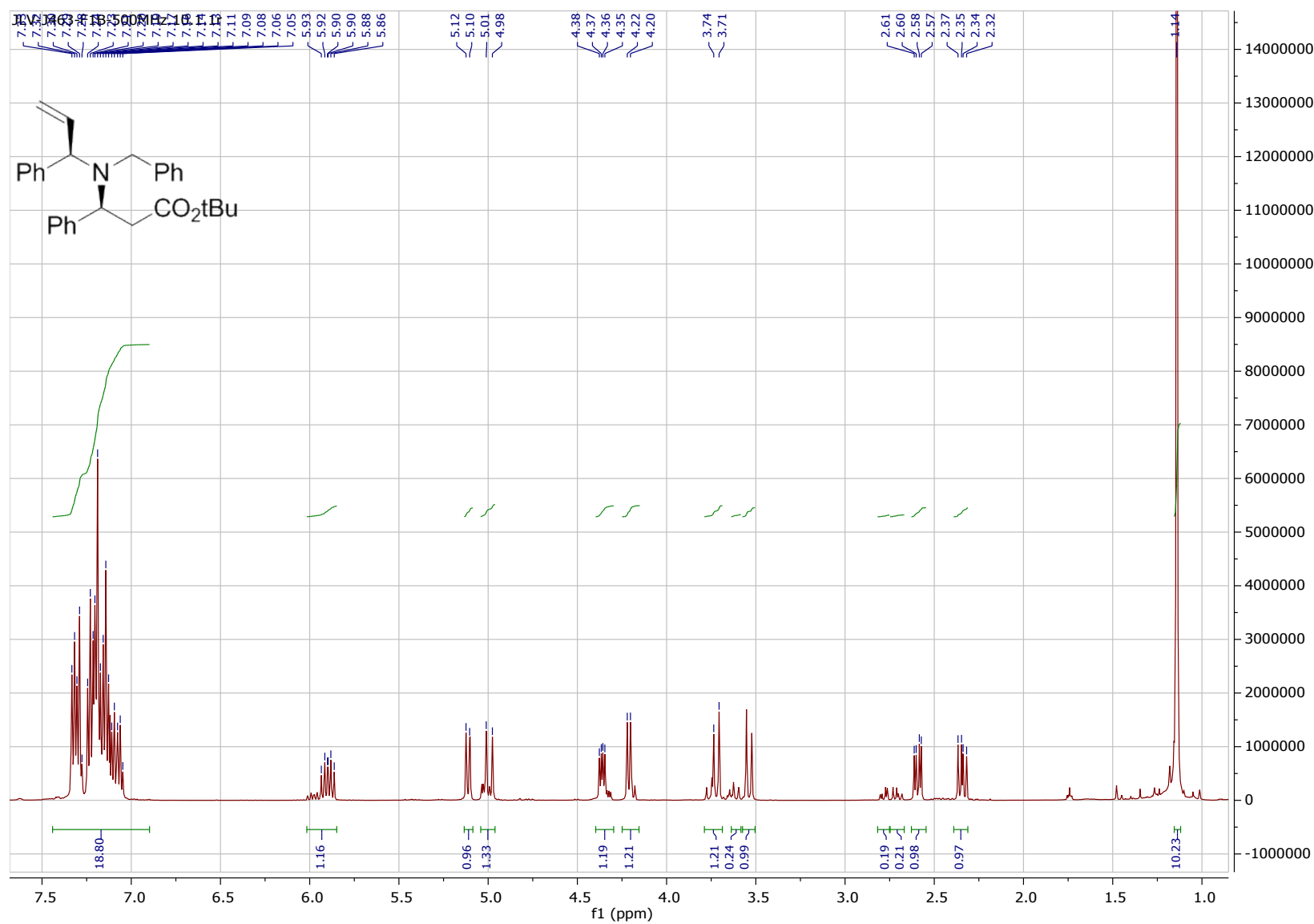


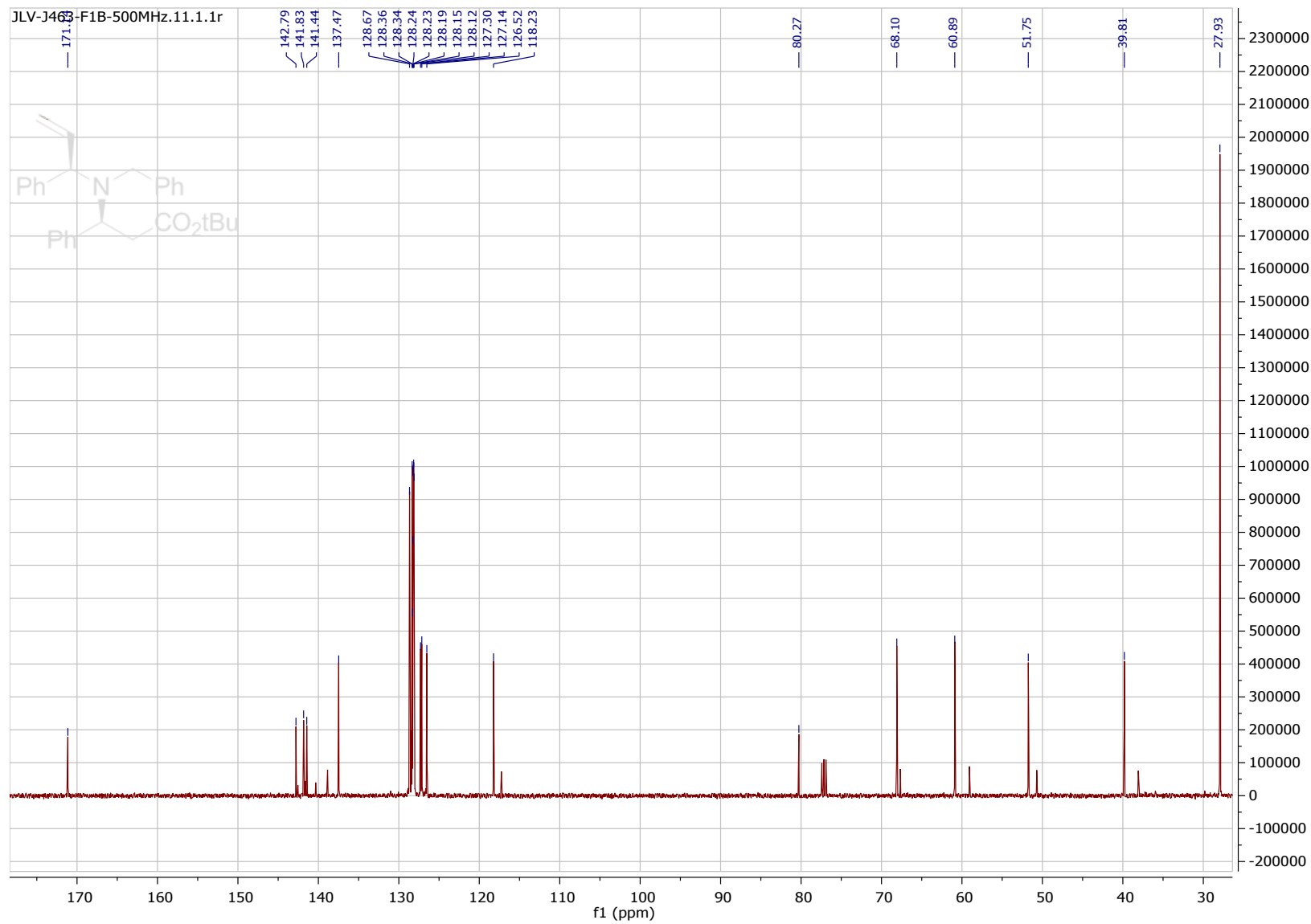
(S)-N-Benzyl-1-(thiophen-2-yl)but-3-en-1-amine 1e



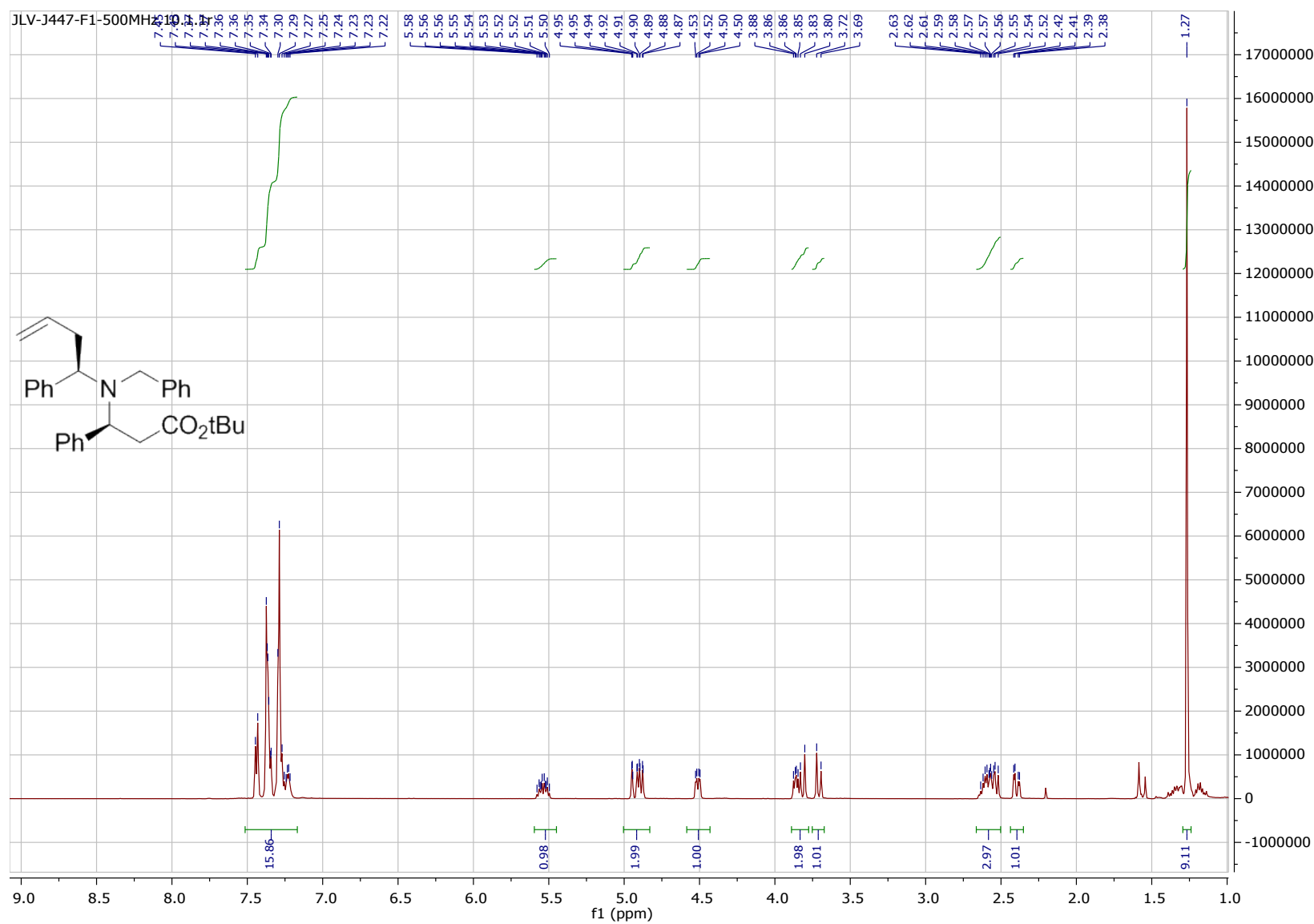


(S)-Tert-Butyl 3-(benzyl((R)-1-phenylallyl)amino)-3-phenylpropanoate 4

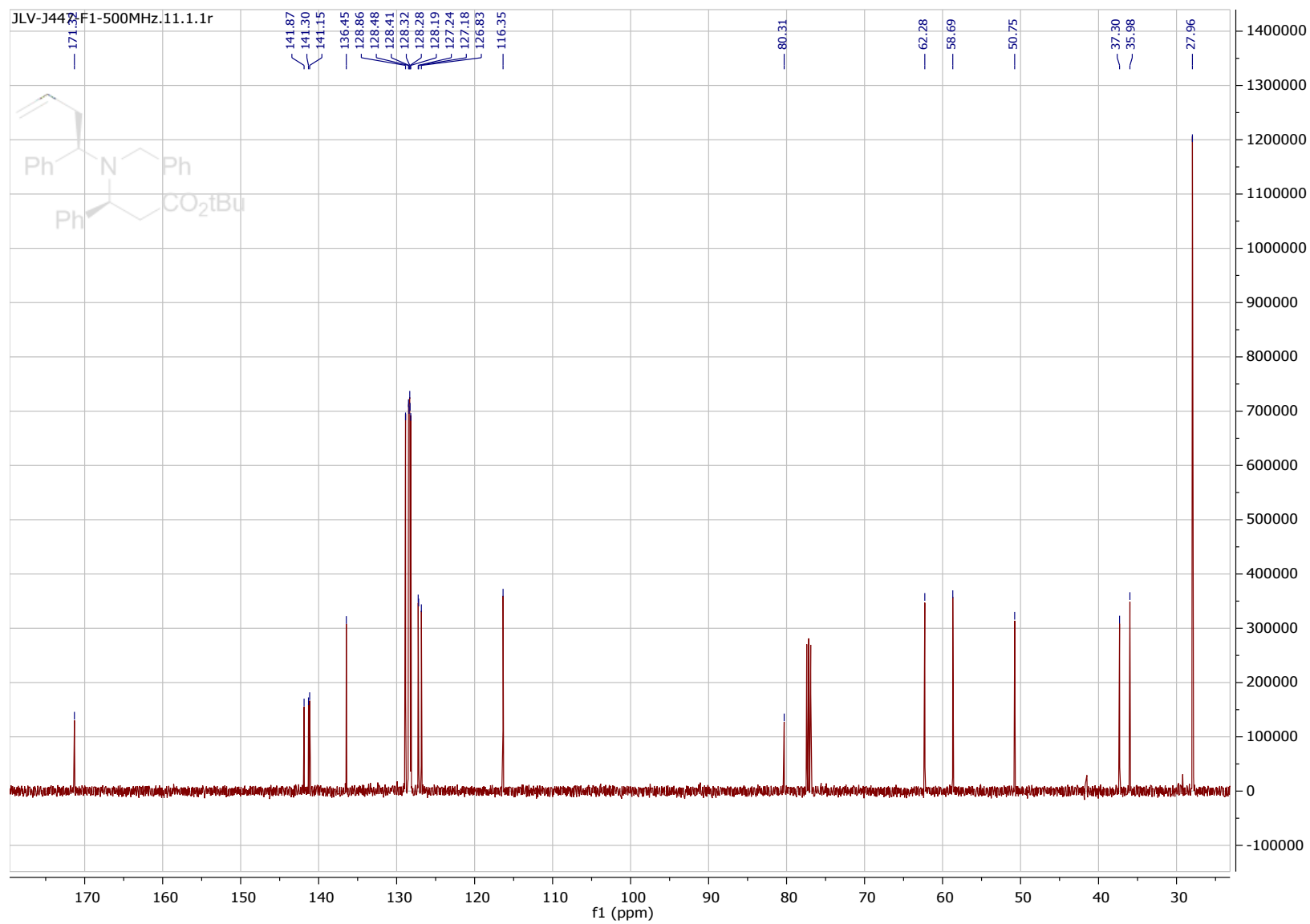




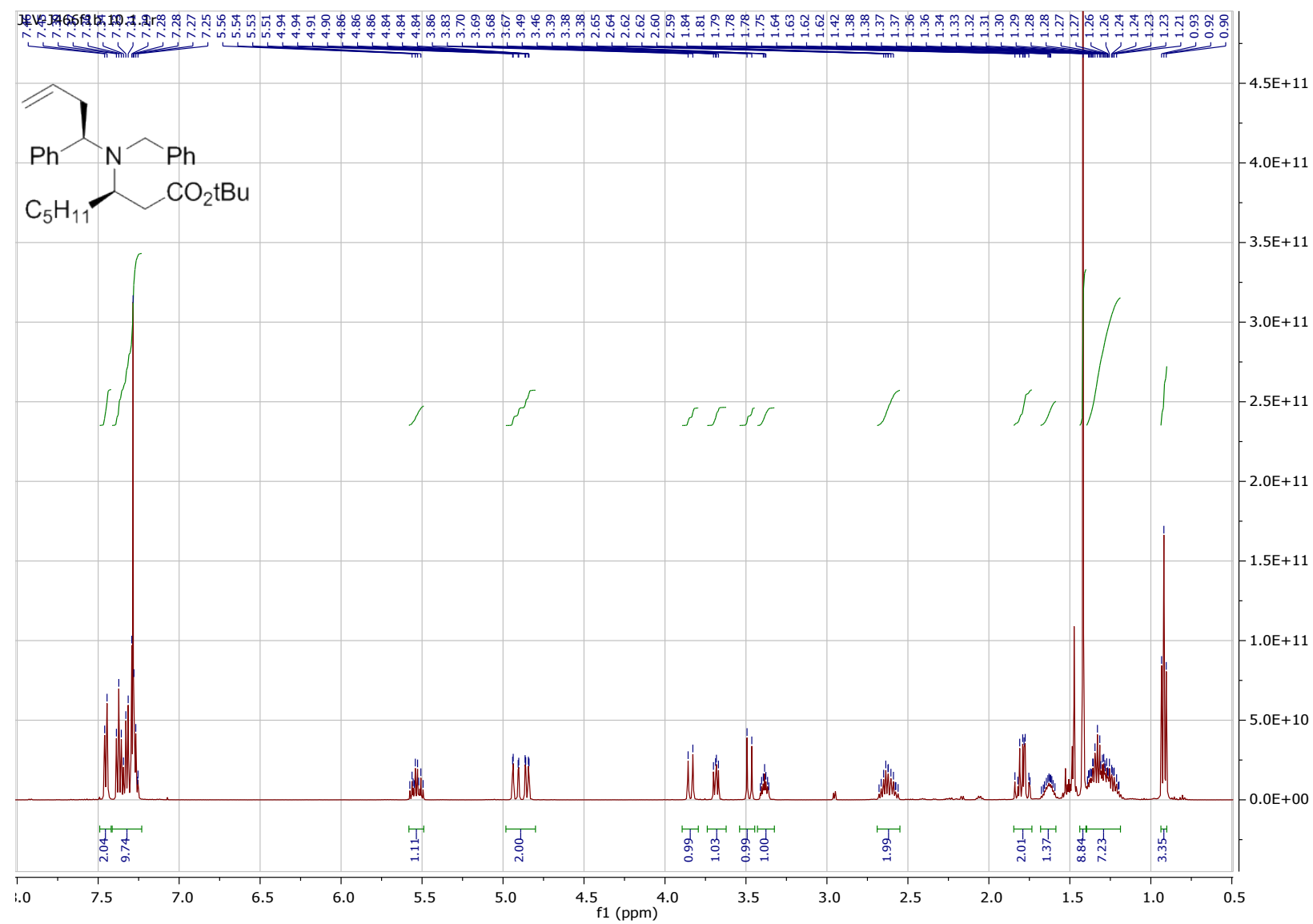
(S)-Tert-Butyl 3-{benzyl[(R)-1-phenylbut-3-en-1-yl]amino}-3-phenylpropanoate 3a

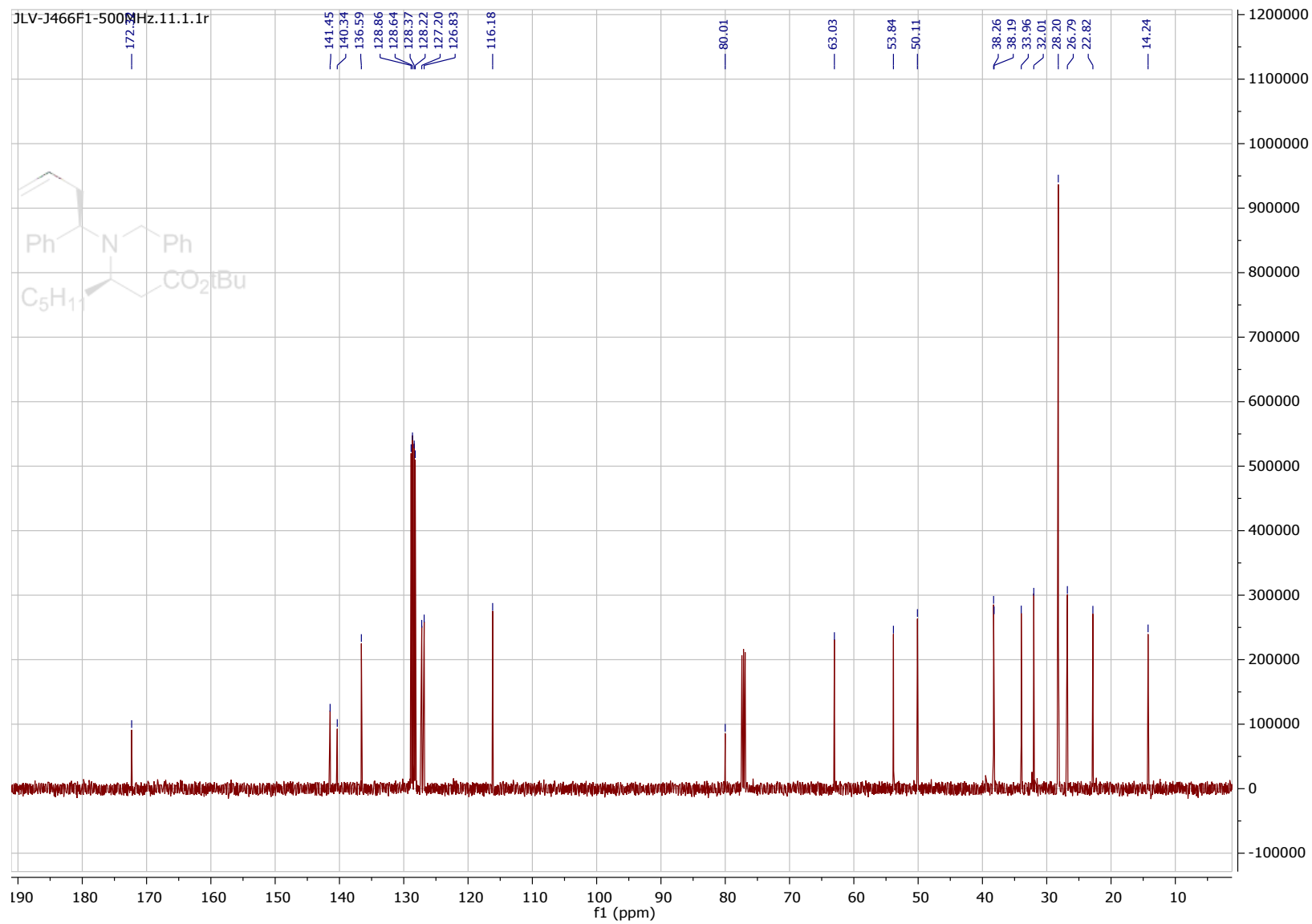


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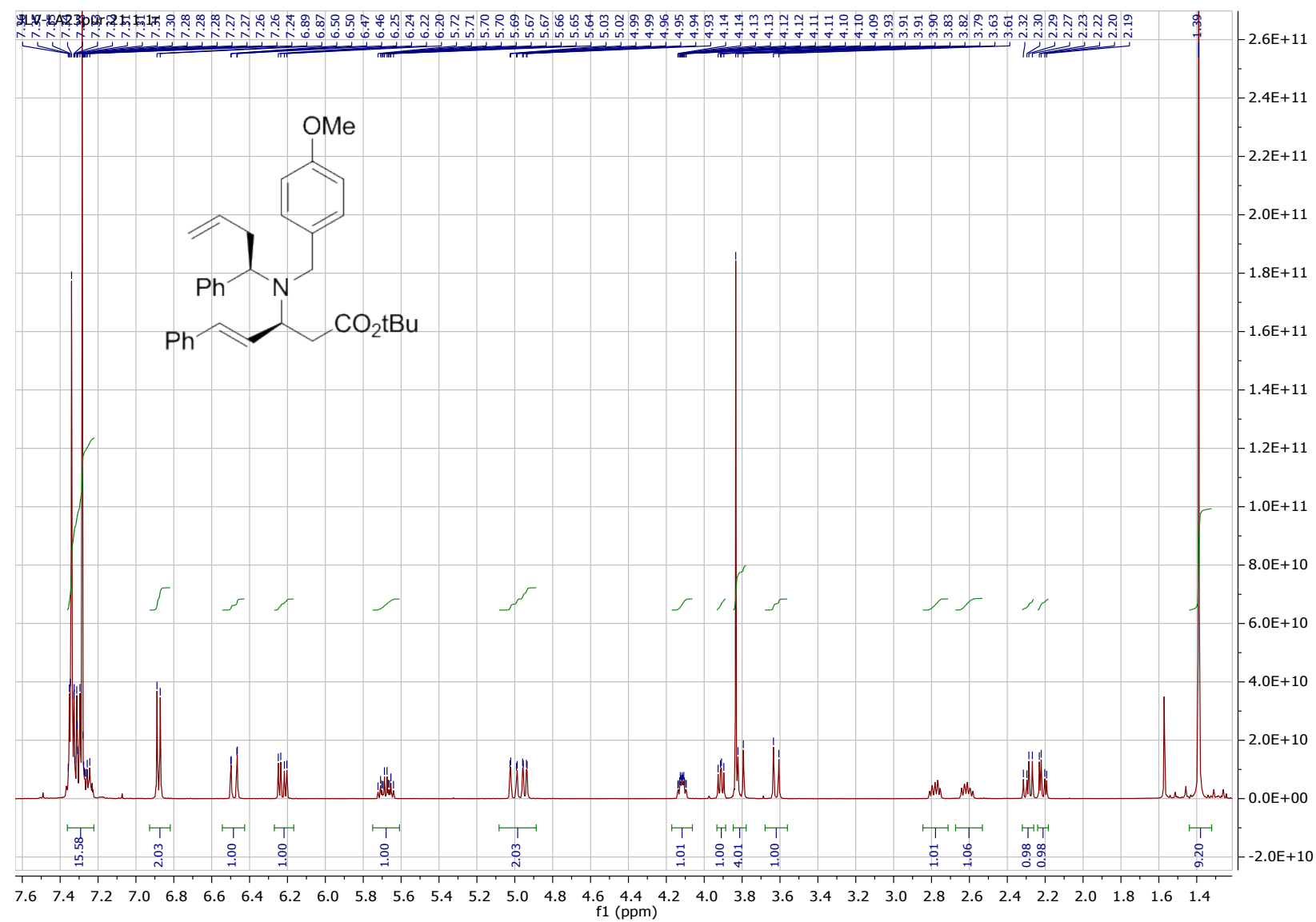


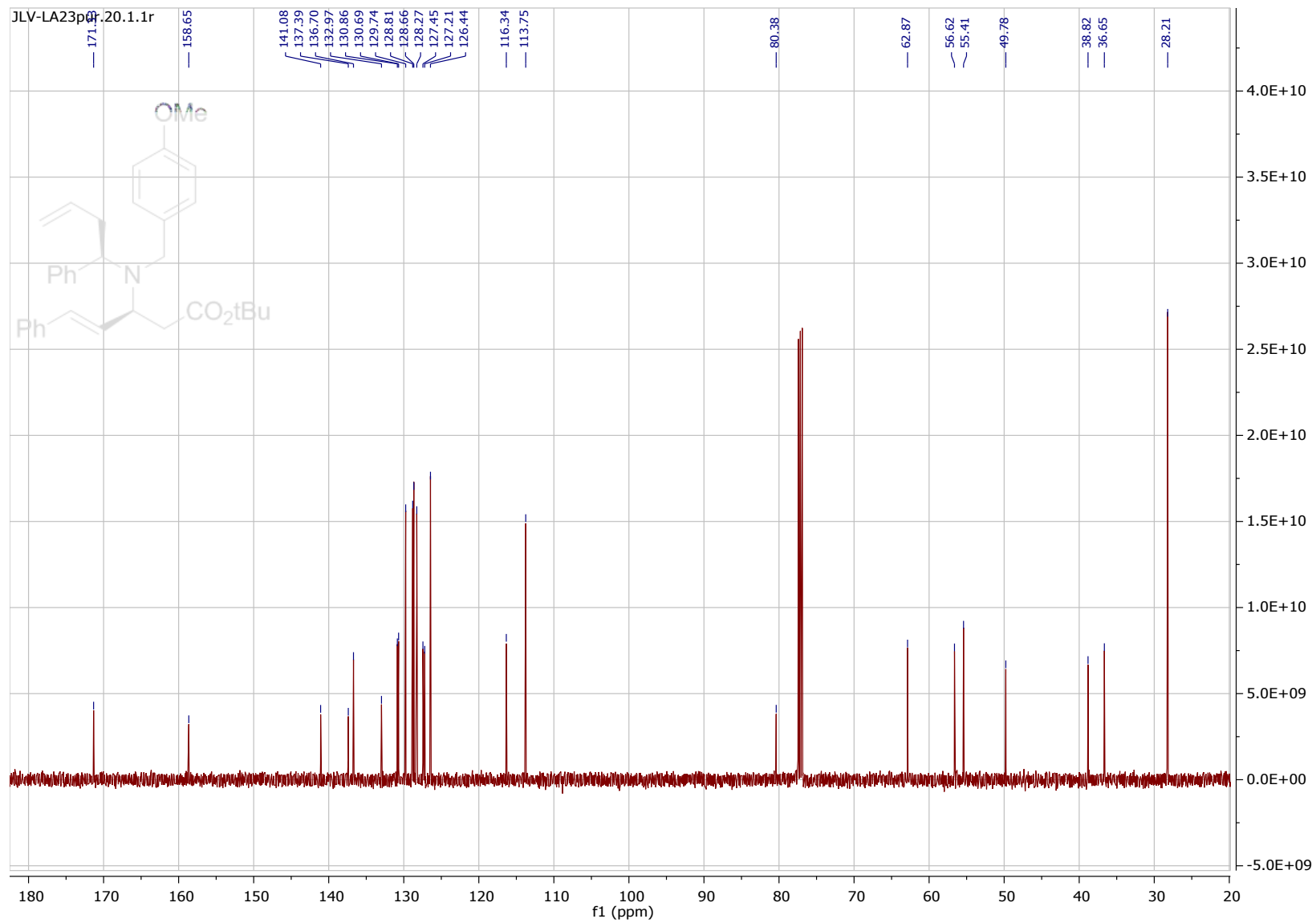
(±)-(R)-Tert-Butyl 3-{benzyl[(R)-1-phenylbut-3-en-1-yl]amino}octanoate 3b



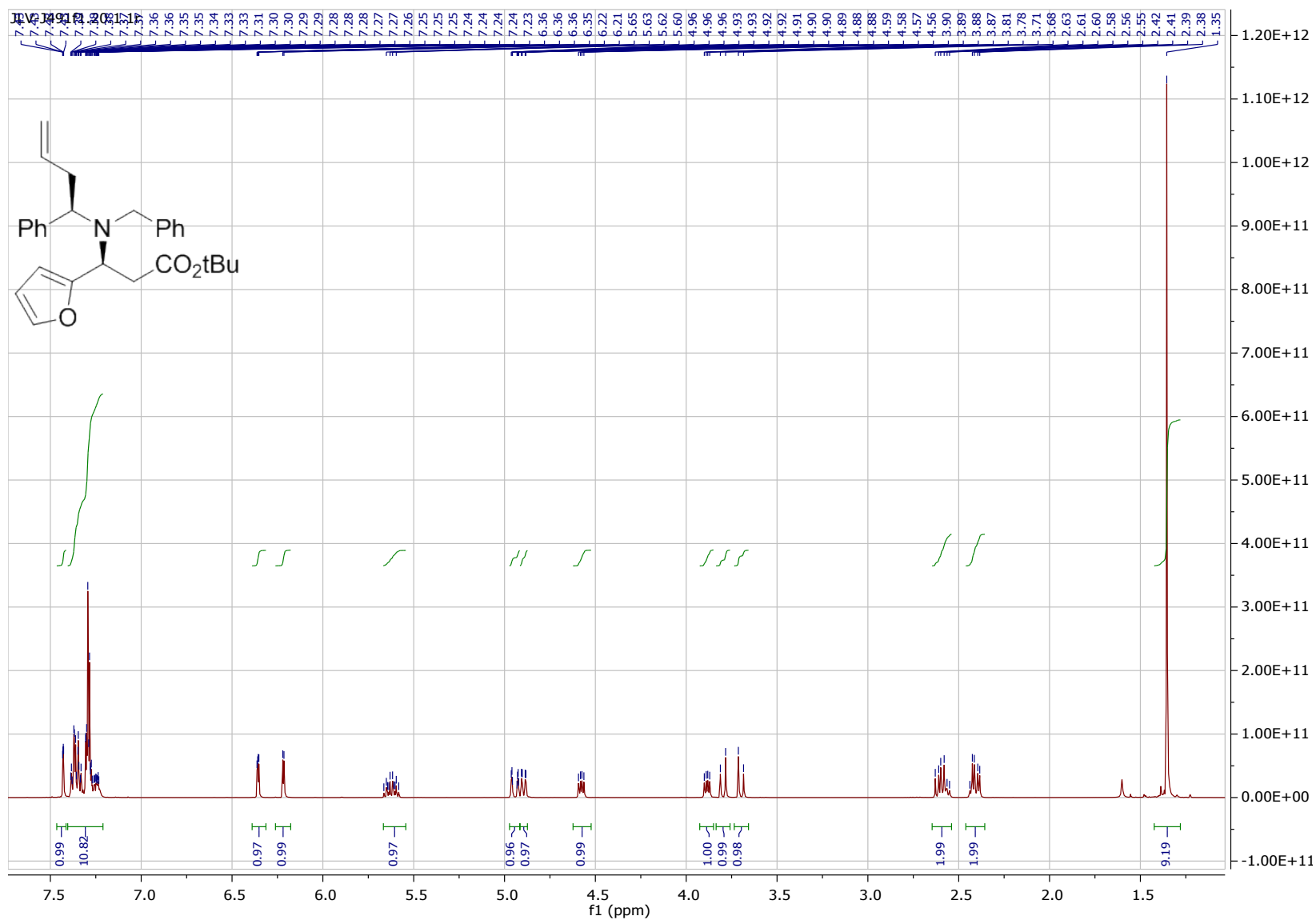


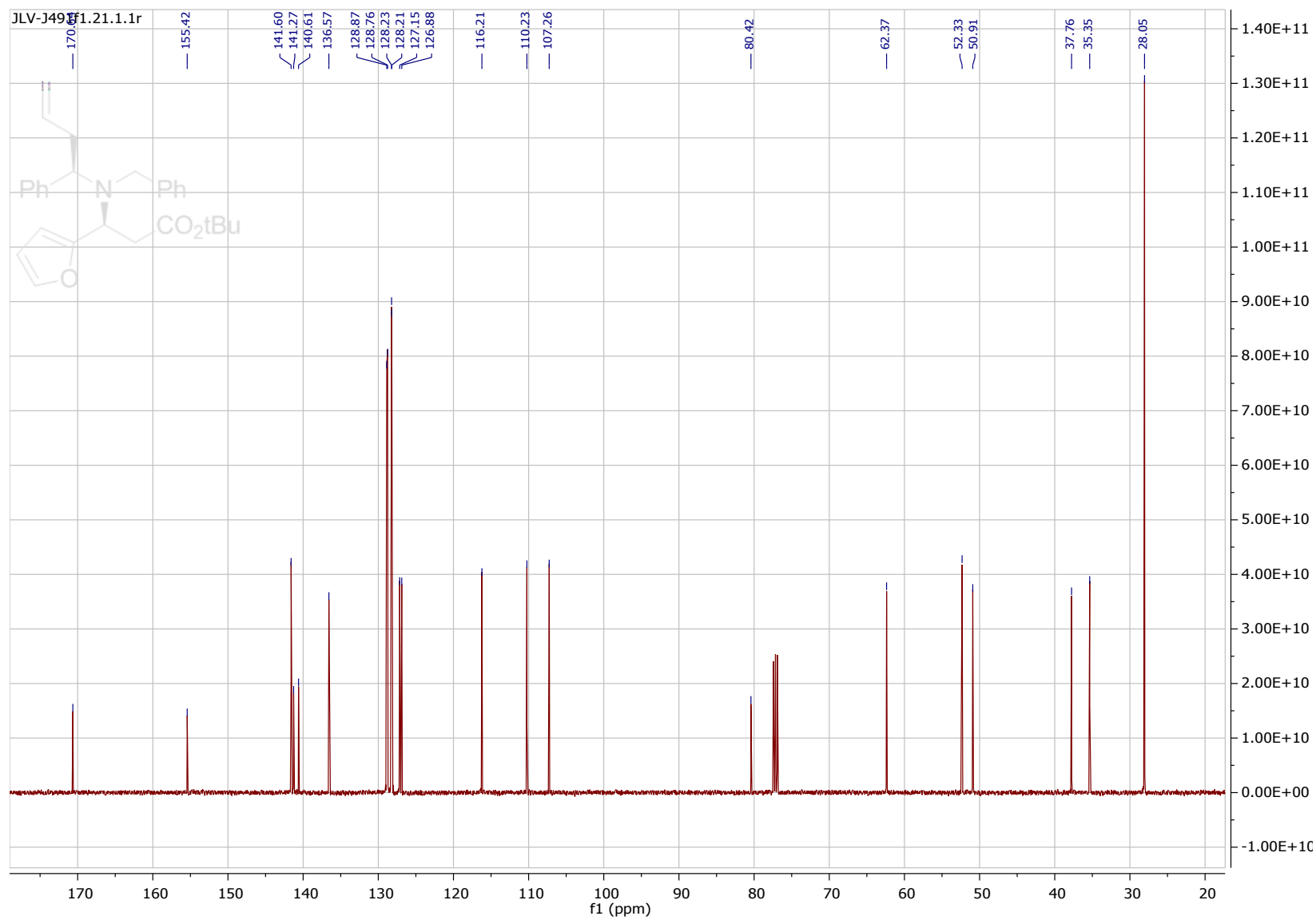
(±)-(S,E)-Tert-Butyl 3-((4-methoxybenzyl)((R)-1-phenylbut-3-en-1-yl)amino)-5-phenylpent-4-enoate 3c



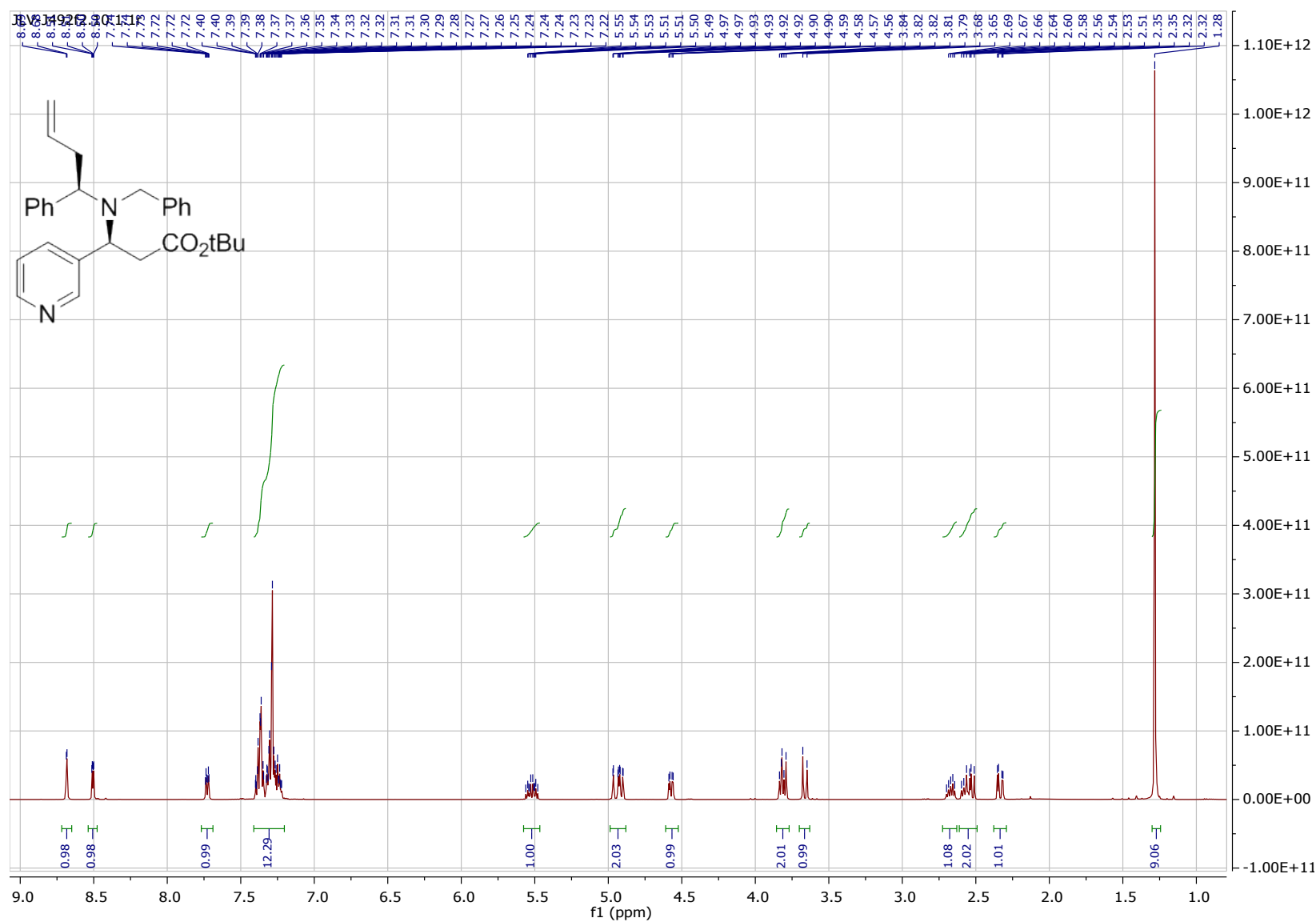


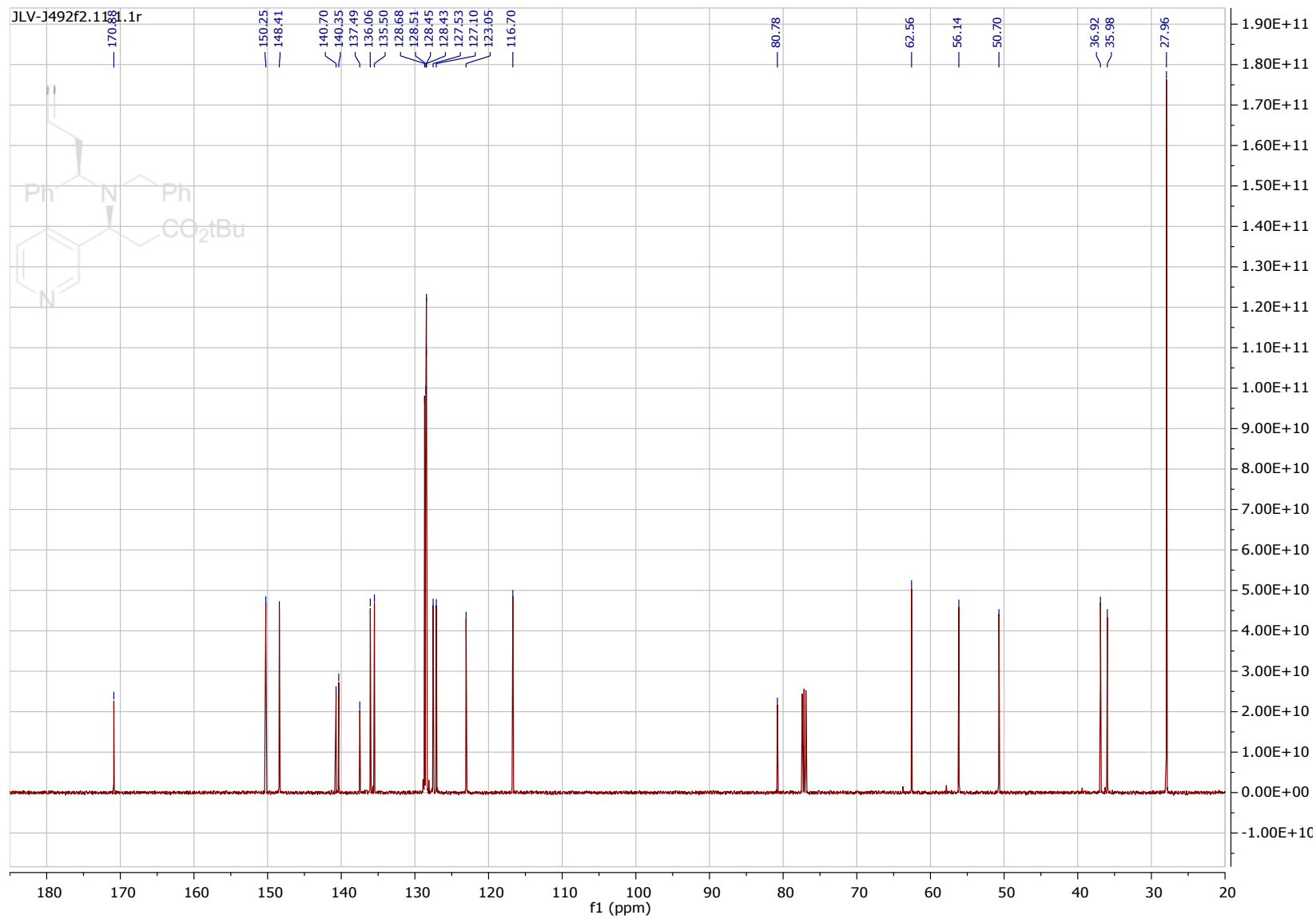
(±)-(S)-Tert-Butyl 3-{benzyl[(R)-1-phenylbut-3-en-1-yl]amino}-3-(furan-2-yl)propanoate 3d



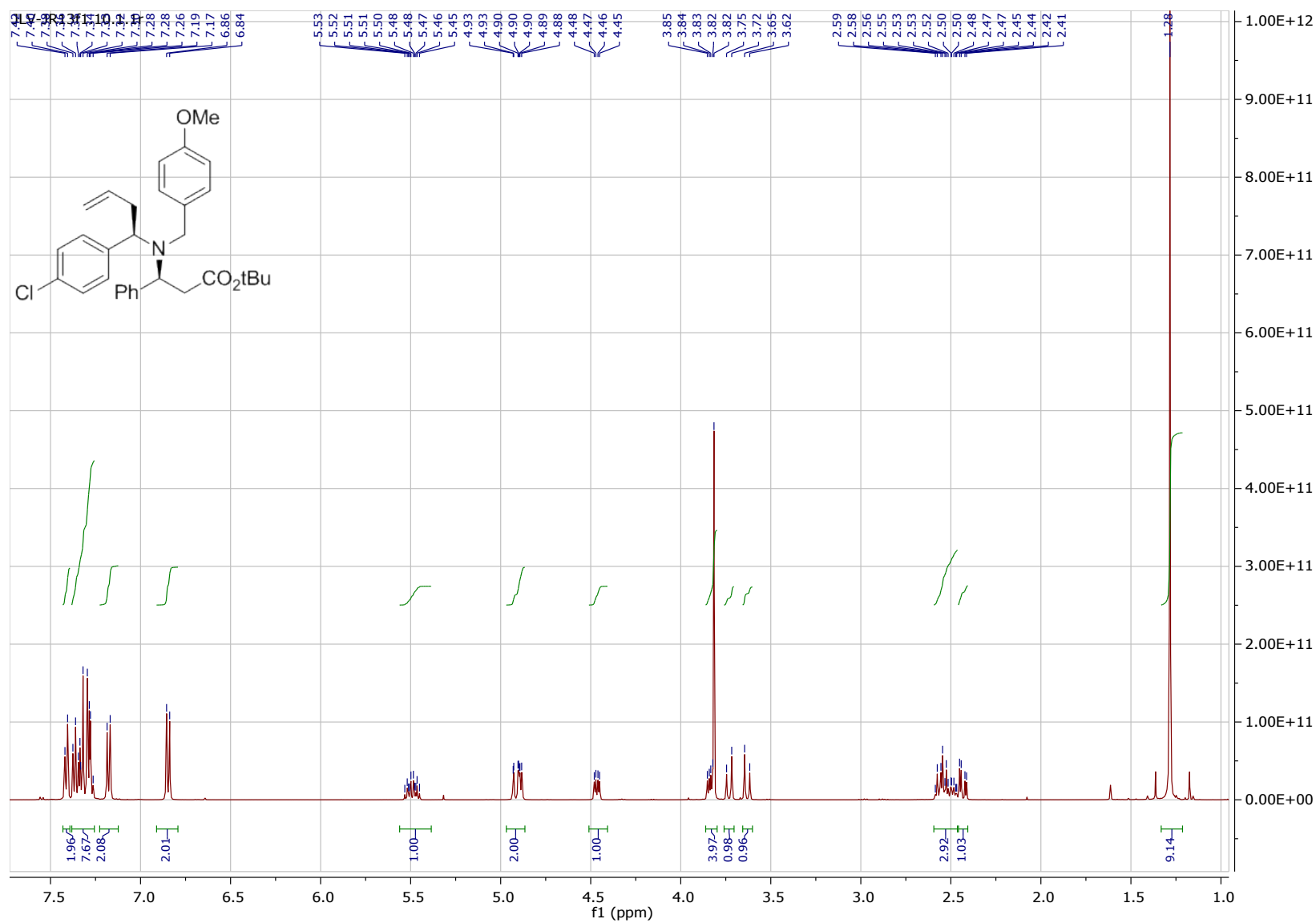


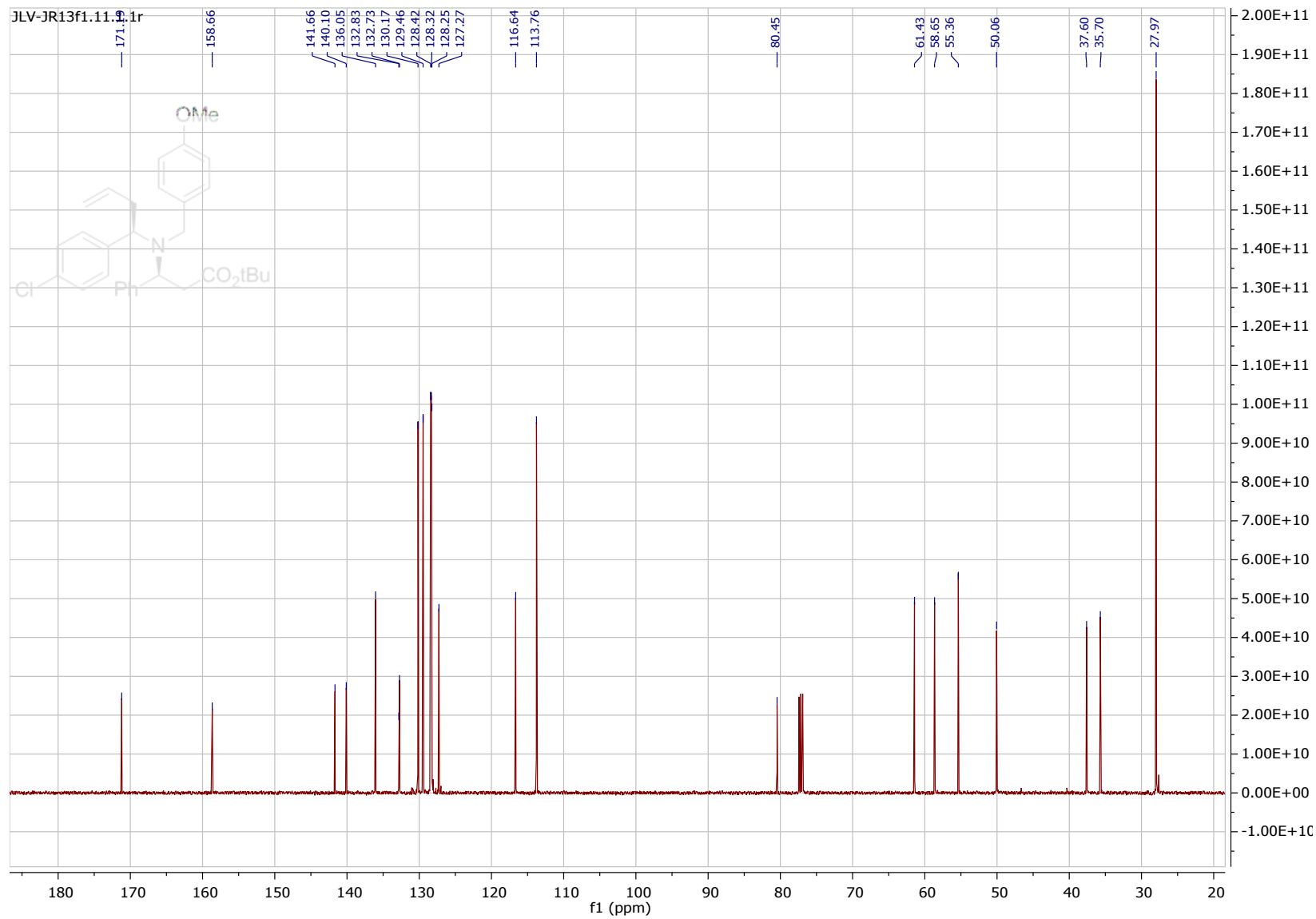
(±)-(*S*)-*Tert*-Butyl 3-{benzyl[*R*]-1-phenylbut-3-en-1-yl}amino-3-(pyridin-3-yl)propanoate **3e**



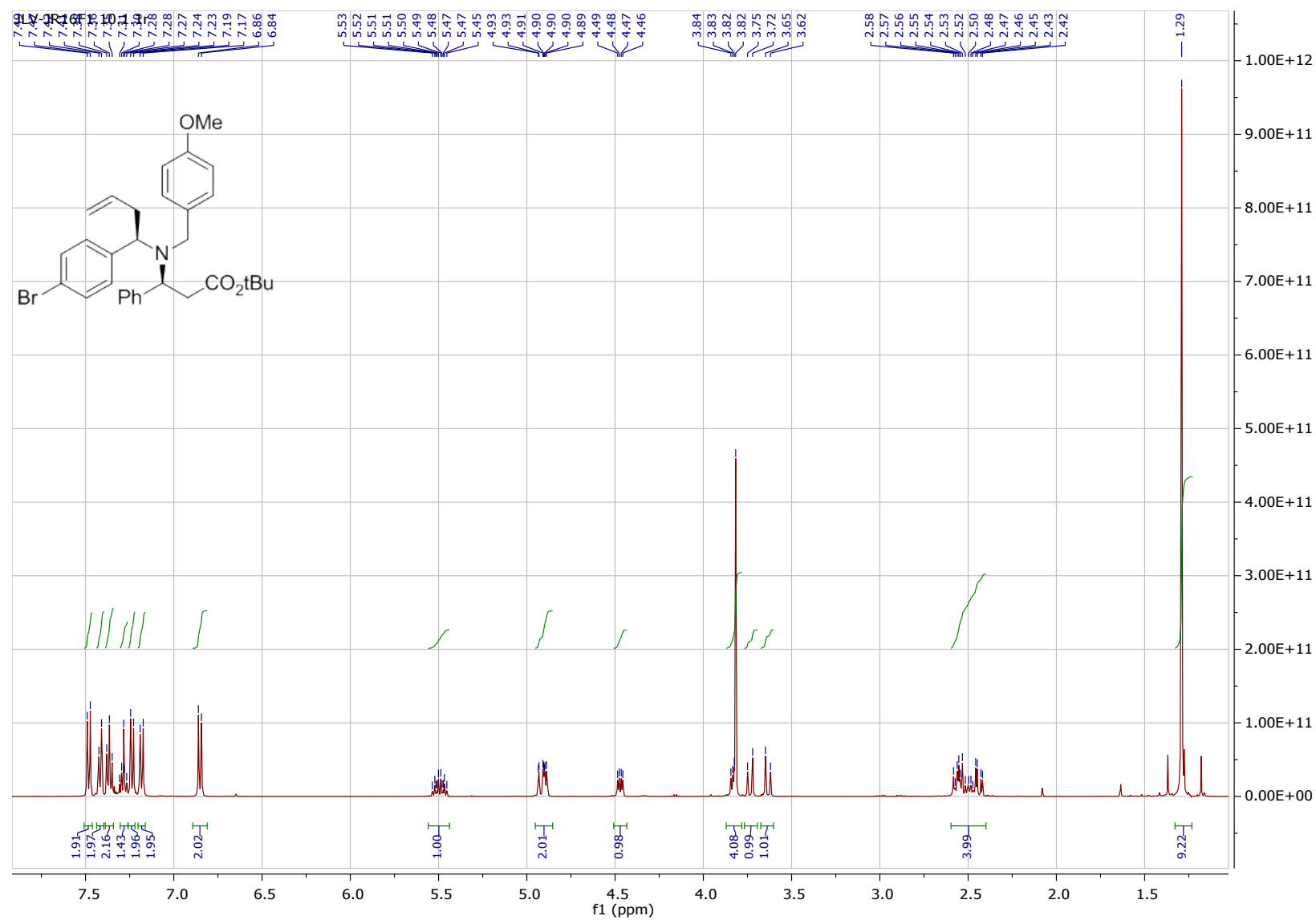


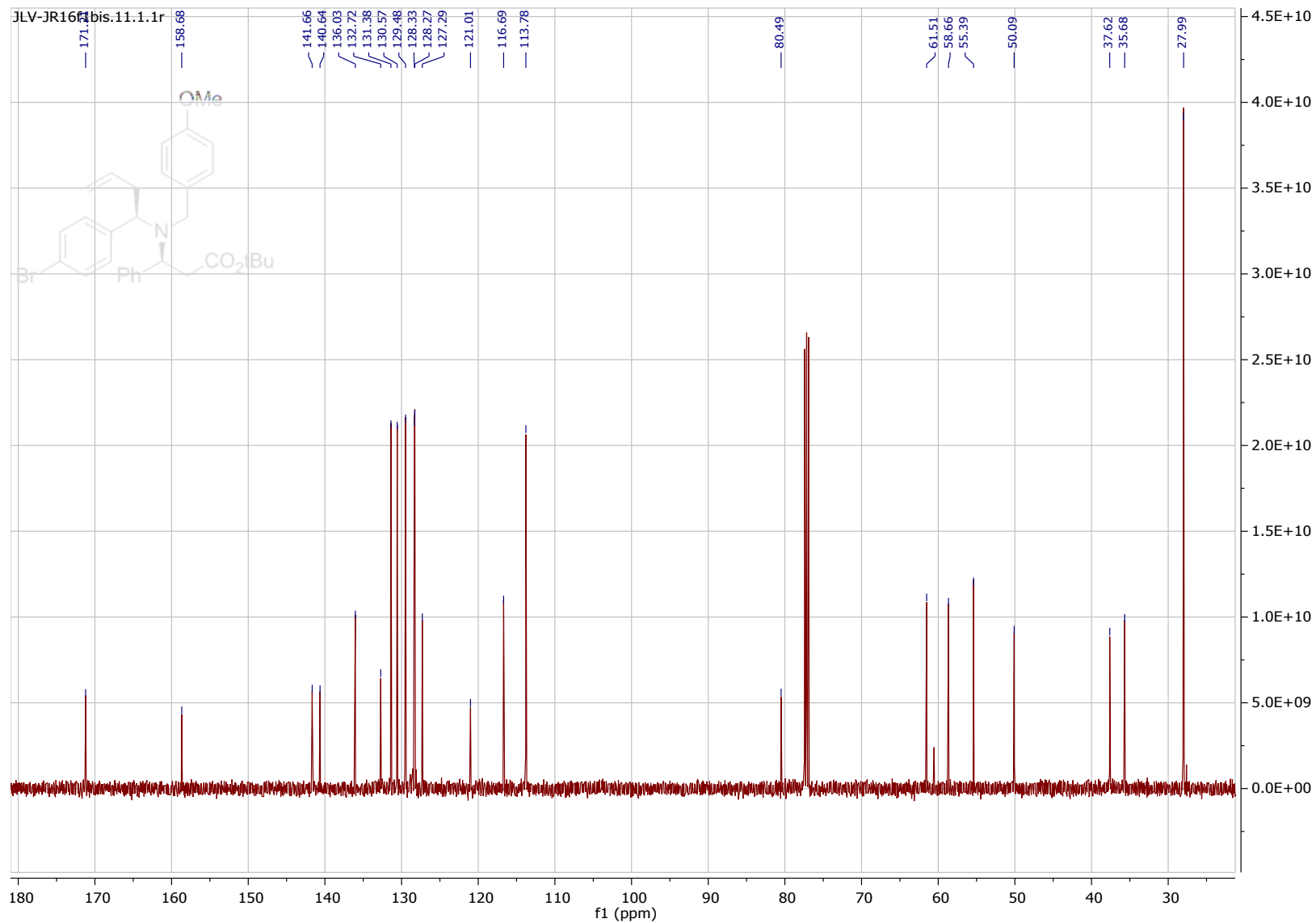
(±)-(*S*)-*Tert*-Butyl 3-[[*(R)*-1-(4-chlorophenyl)but-3-en-1-yl](4-methoxybenzyl)amino]-3-phenylpropanoate **3f**



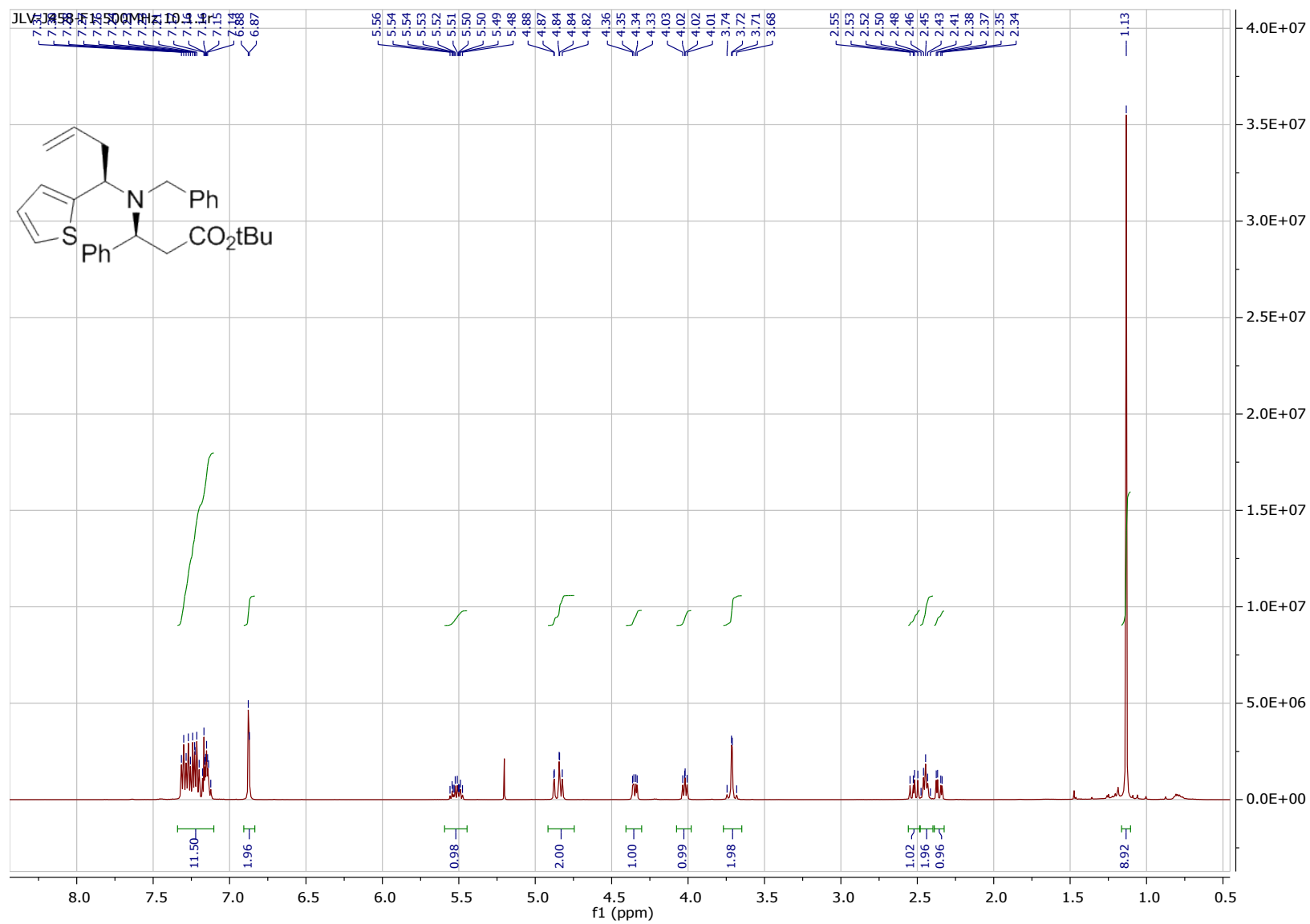


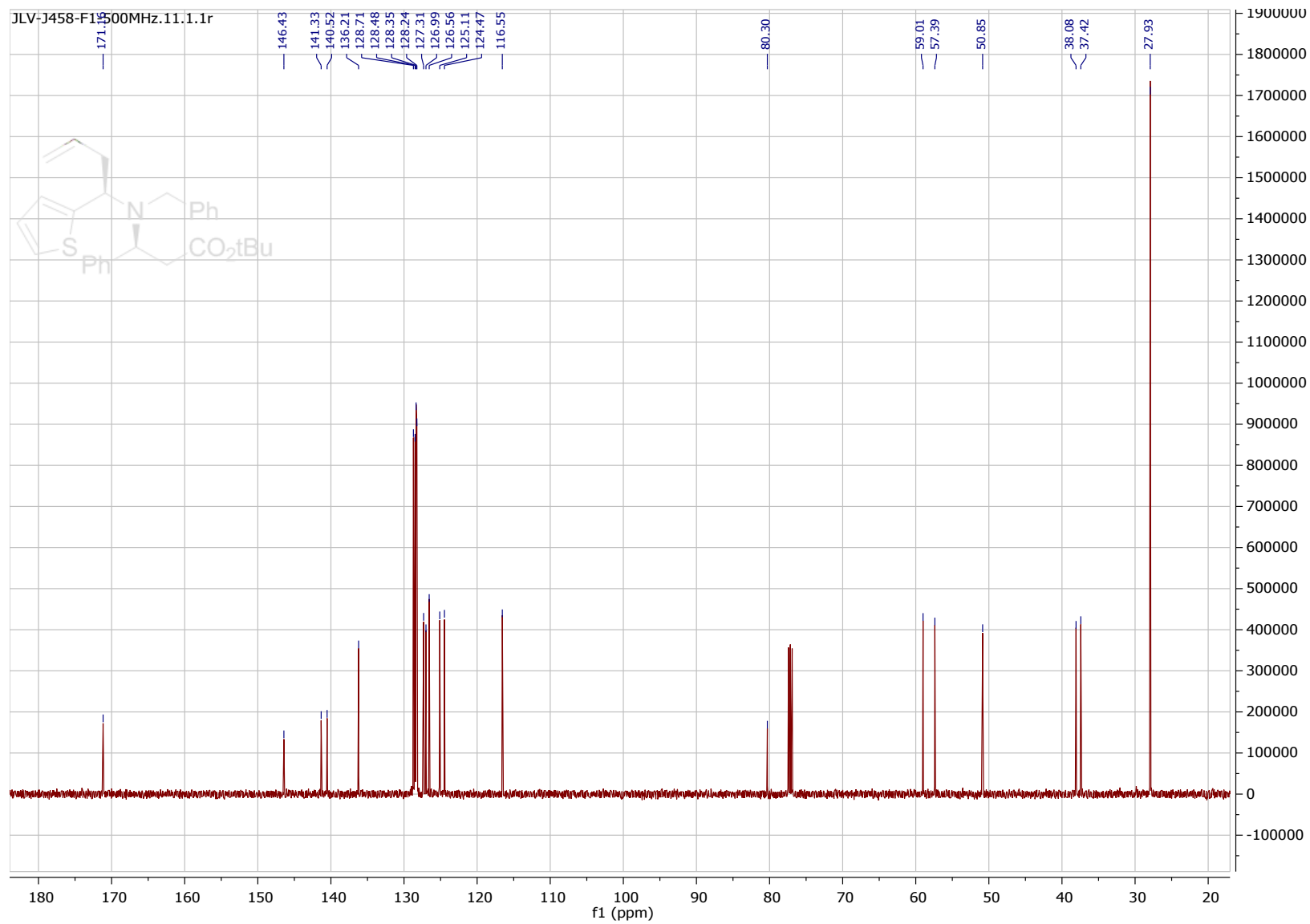
(±)-(*S*)-*Tert*-Butyl 3-[[(*R*)-1-(4-bromophenyl)but-3-en-1-yl](4-methoxybenzyl)amino}-3-phenylpropanoate **3g**



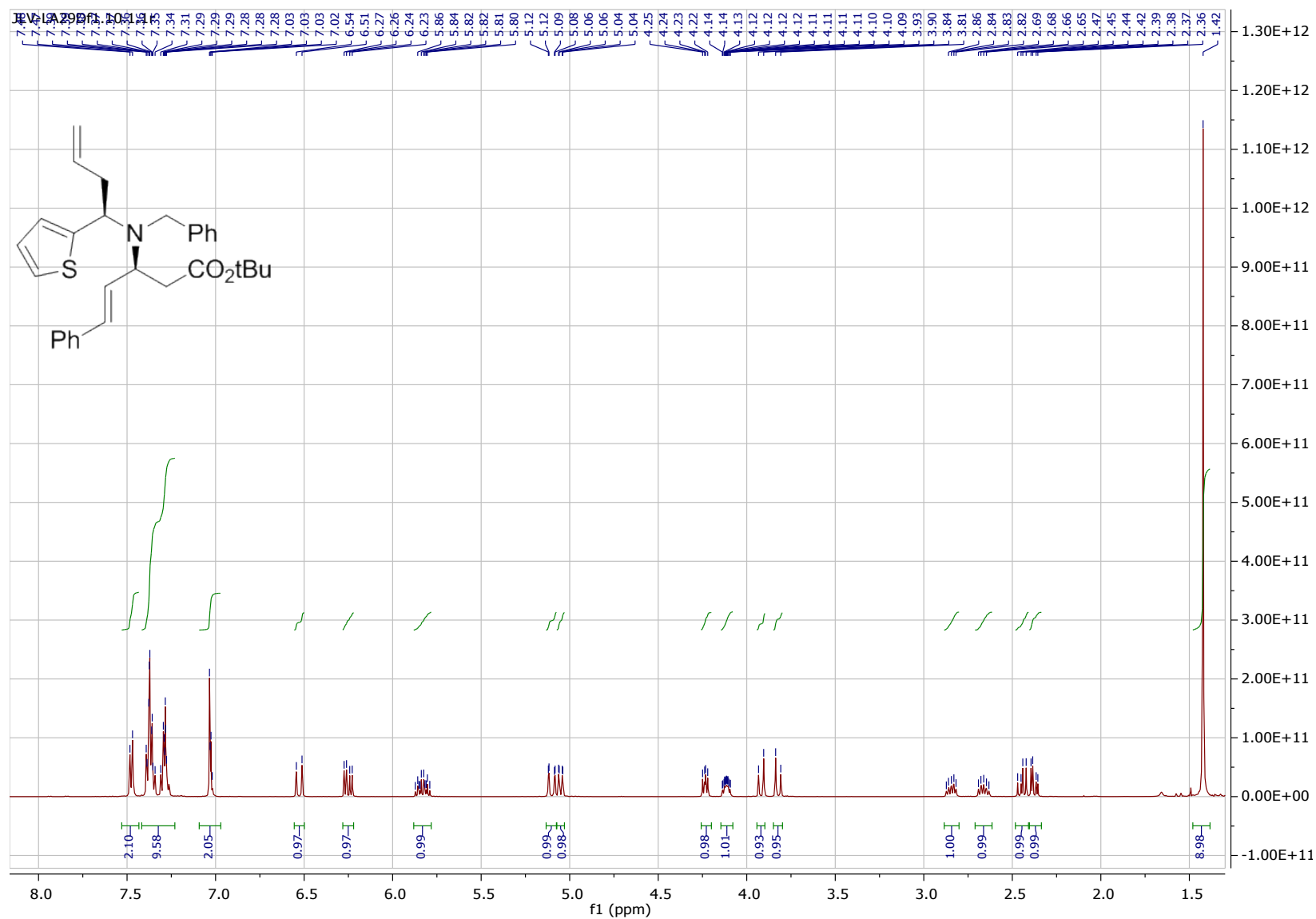


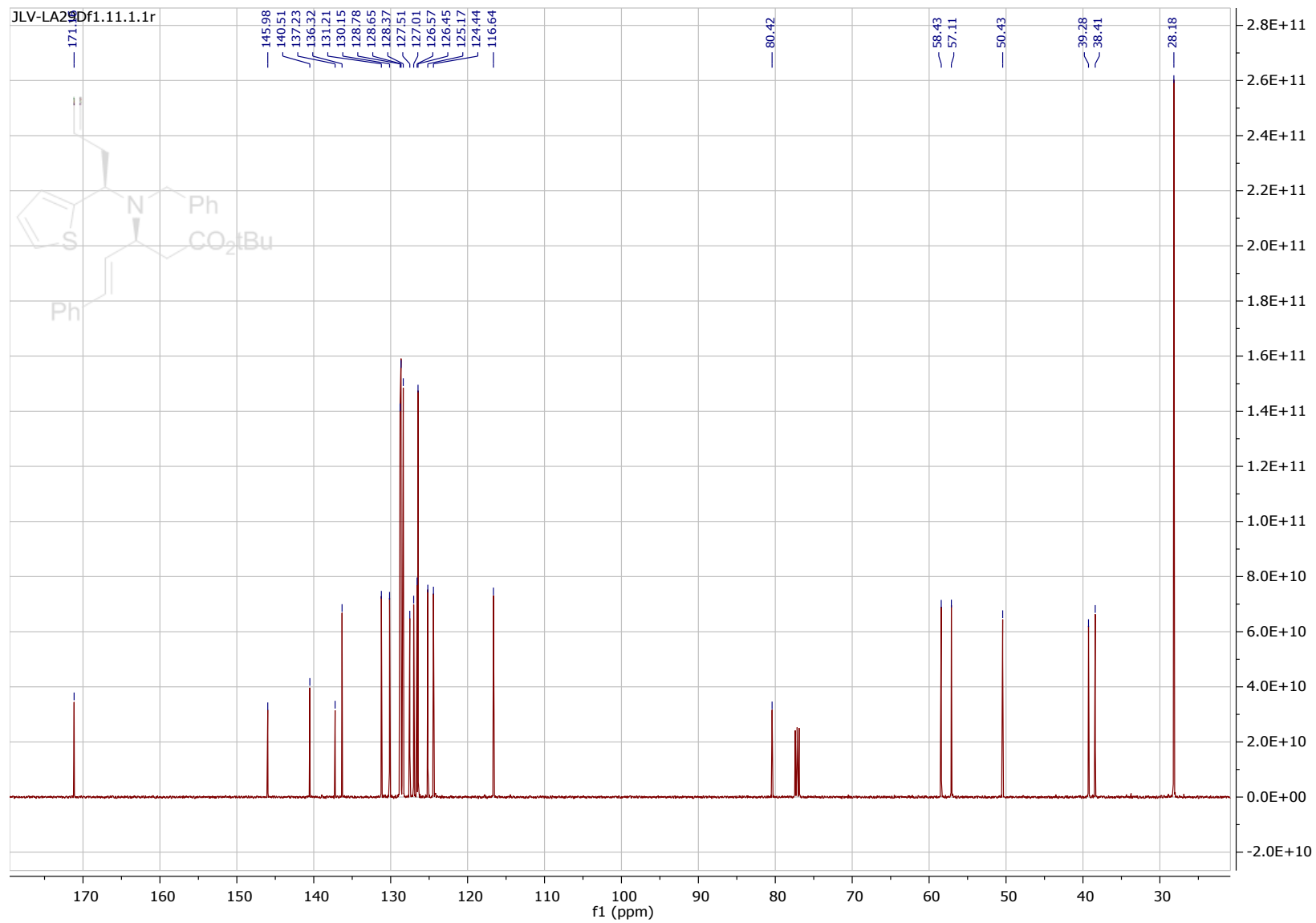
(±)-(*S*)-*Tert*-Butyl 3-{benzyl[(*R*)-1-(thiophen-2-yl)but-3-en-1-yl]amino}-3-phenylpropanoate **3h**





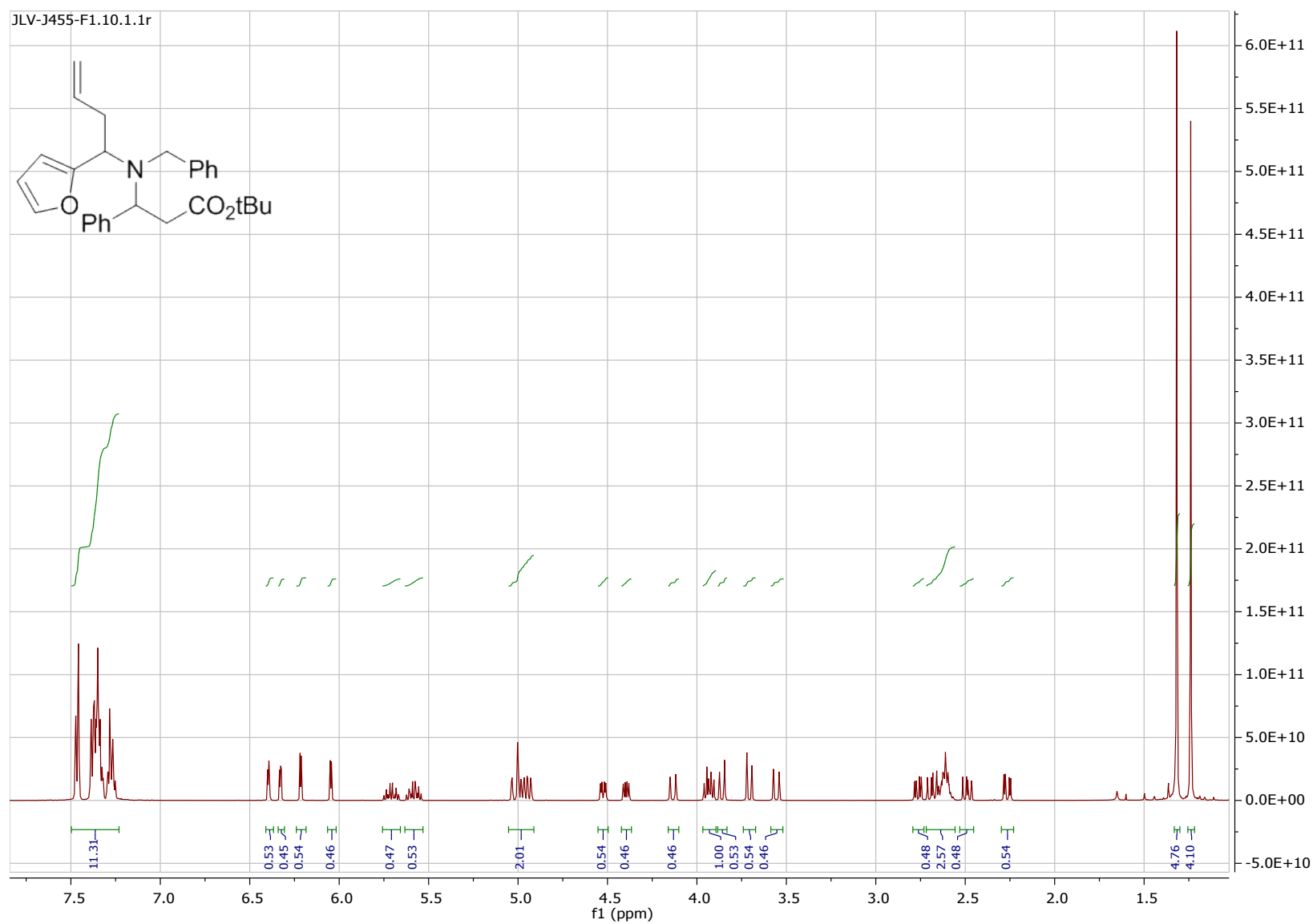
(±)-(S,E)-Tert-Butyl 3-{benzyl[(R)-1-(thiophen-2-yl)but-3-en-1-yl]amino}-5-phenylpent-4-enoate 3i

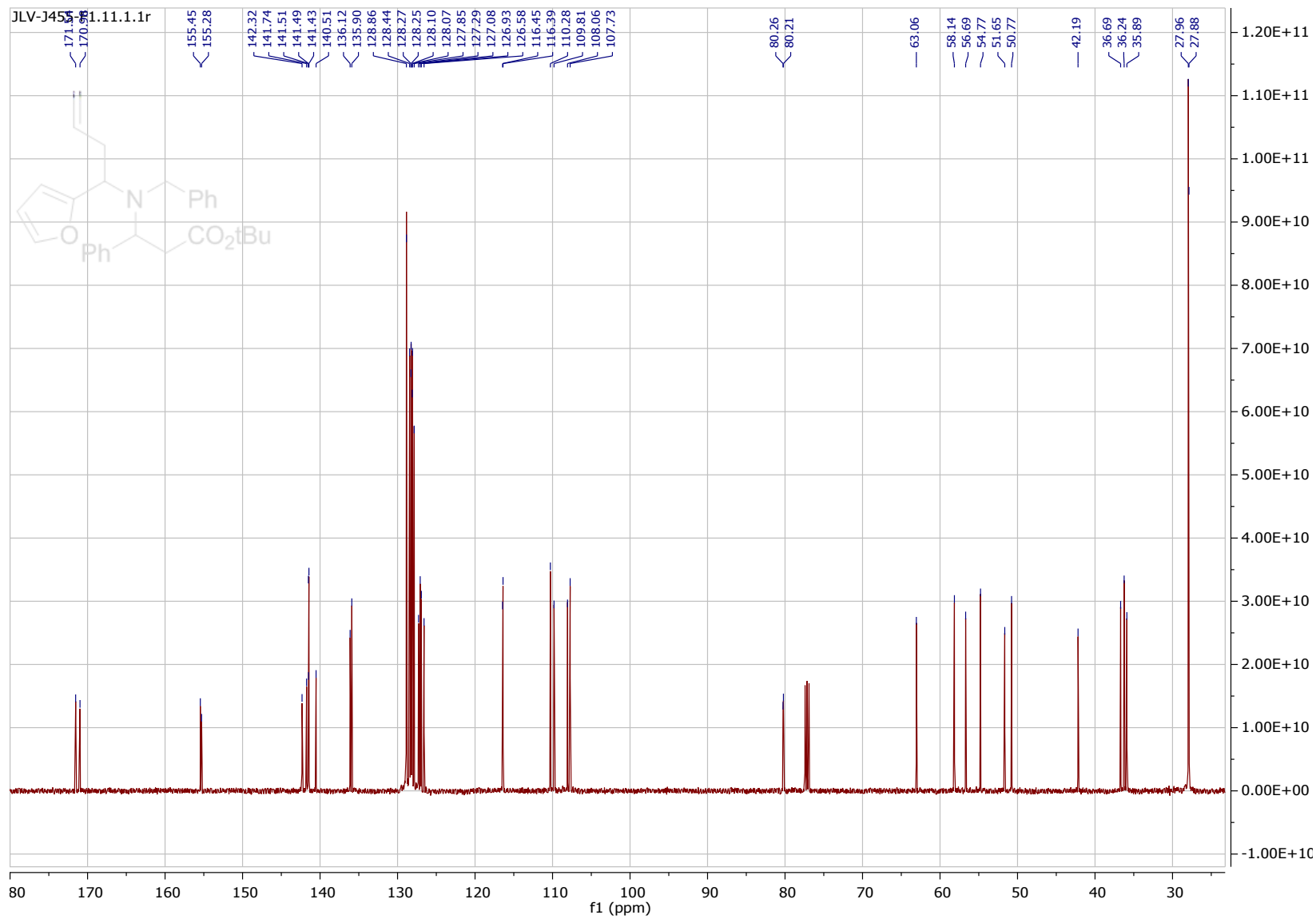




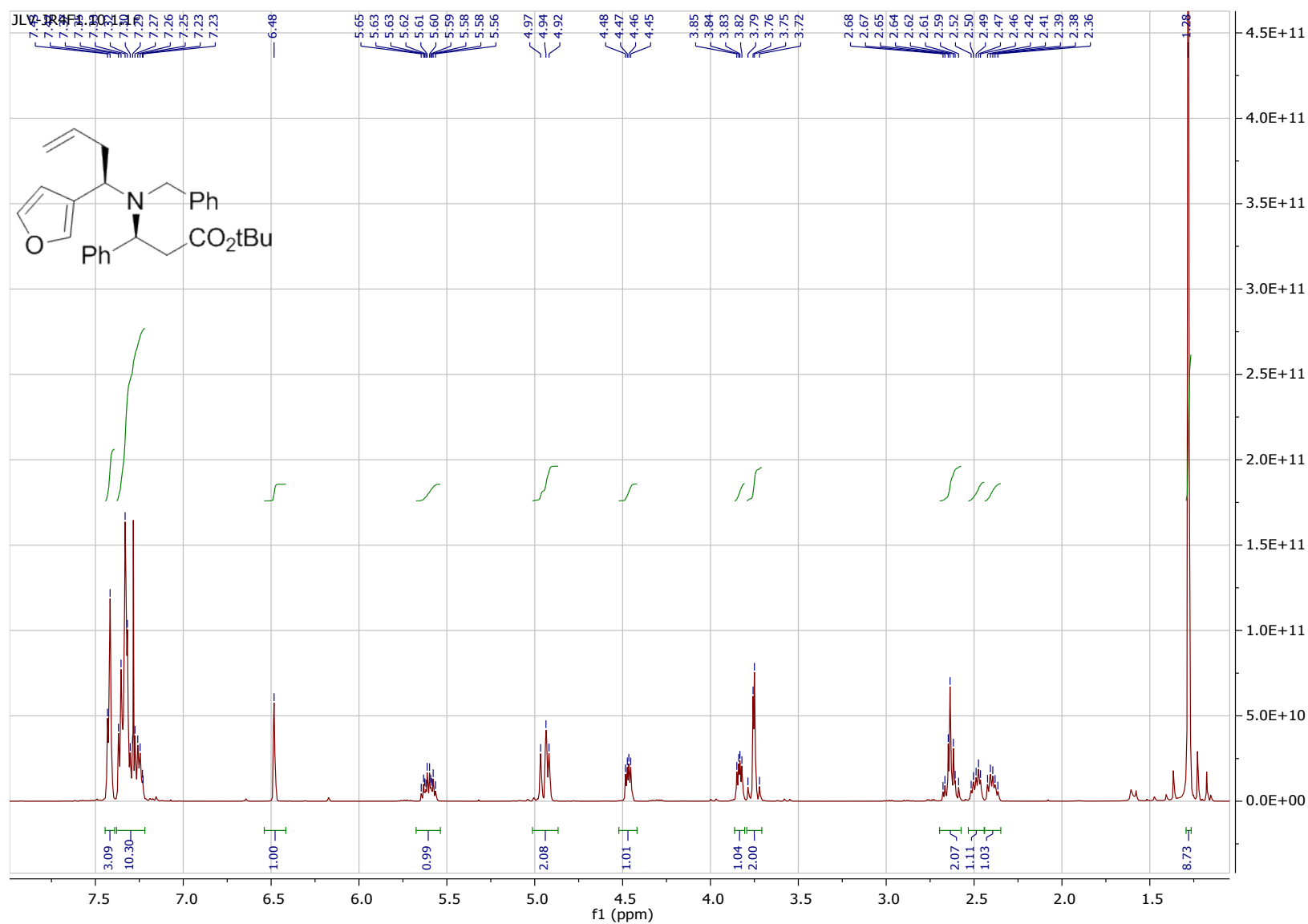
Tert-Butyl 3-{benzyl[-1-(furan-2-yl)but-3-en-1-yl]amino}-3-phenylpropanoate 3j

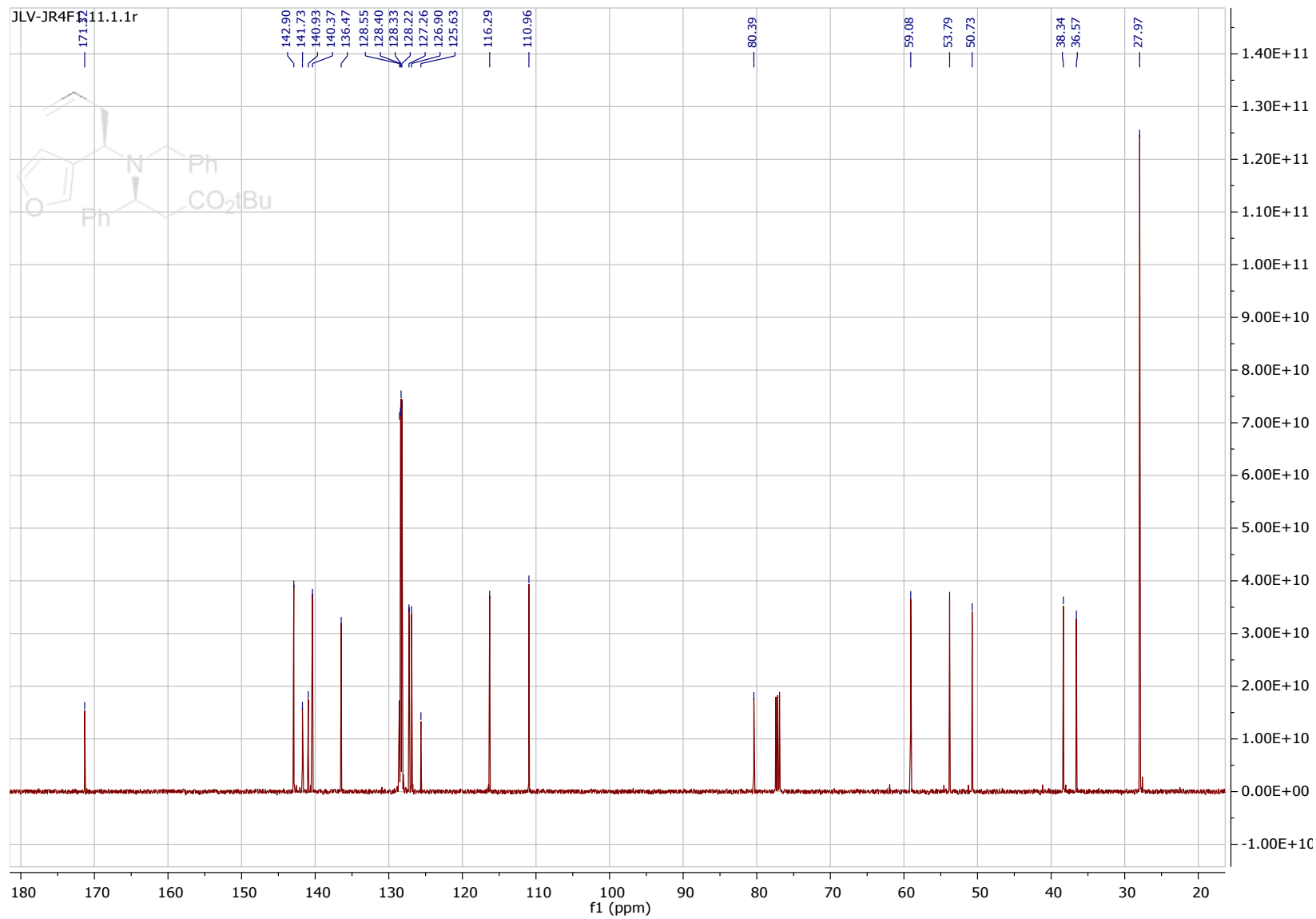
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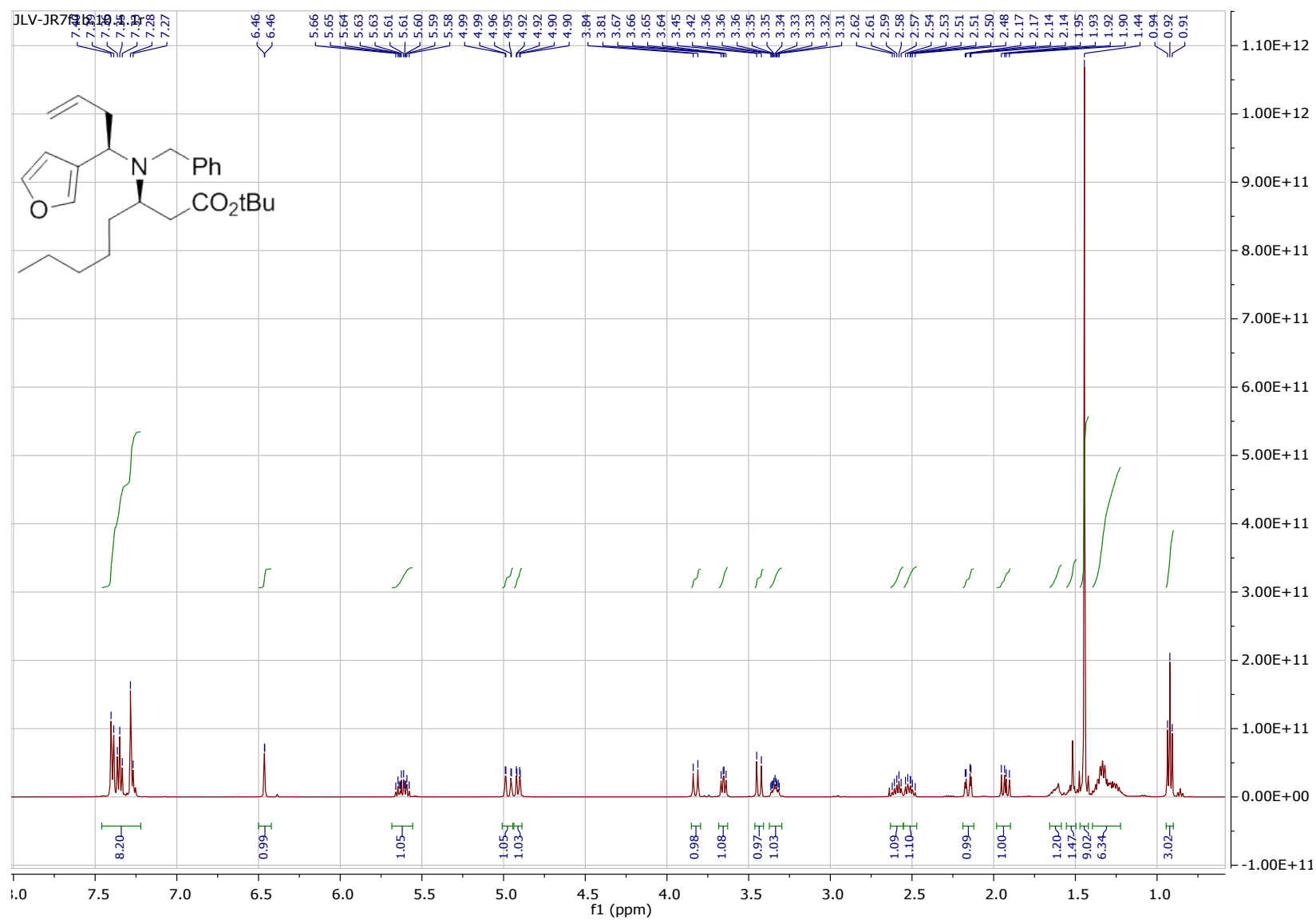


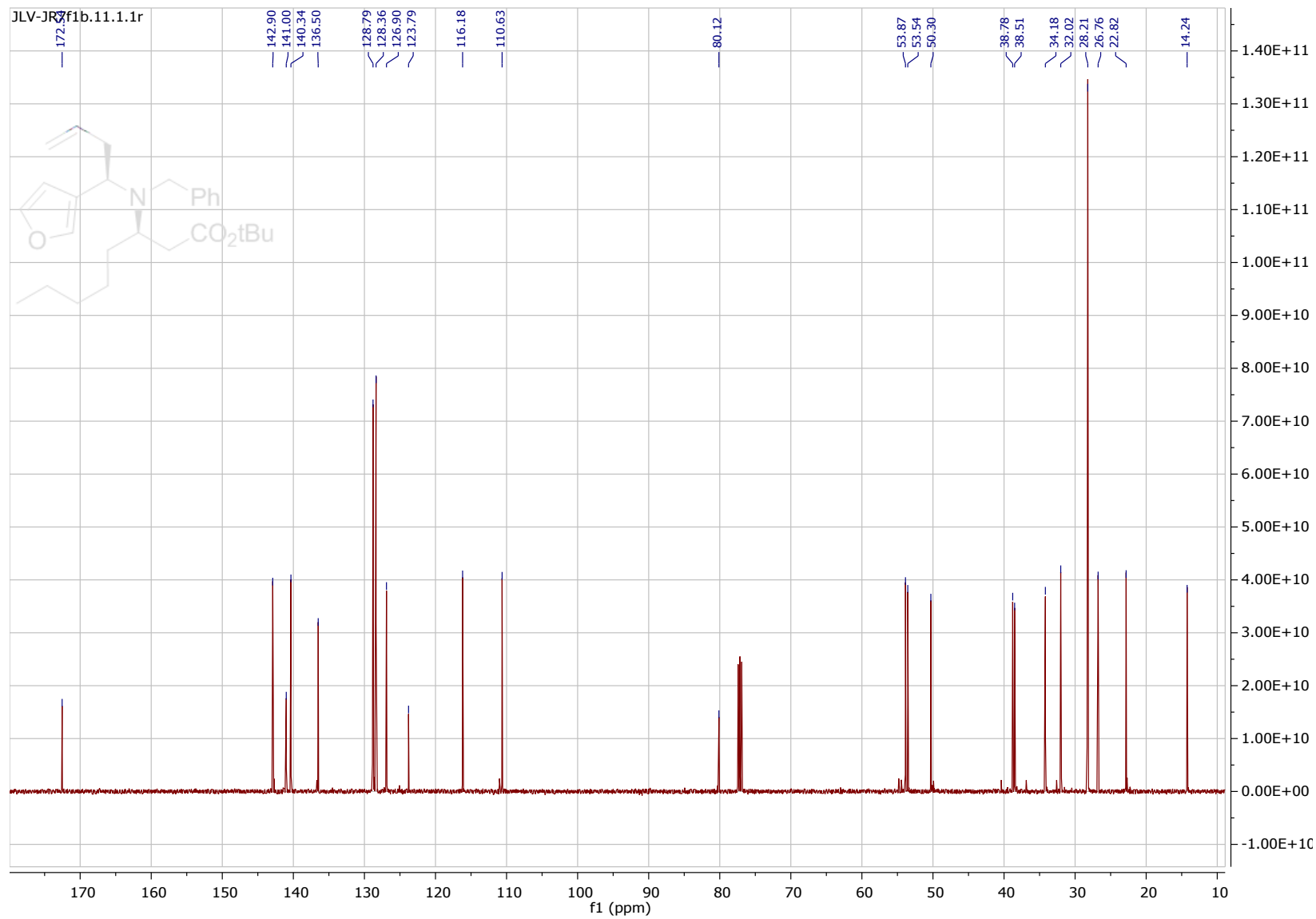
(±)-(S)-Tert-Butyl 3-{benzyl[(R)-1-(furan-3-yl)but-3-en-1-yl]amino}-3-phenylpropanoate 3k





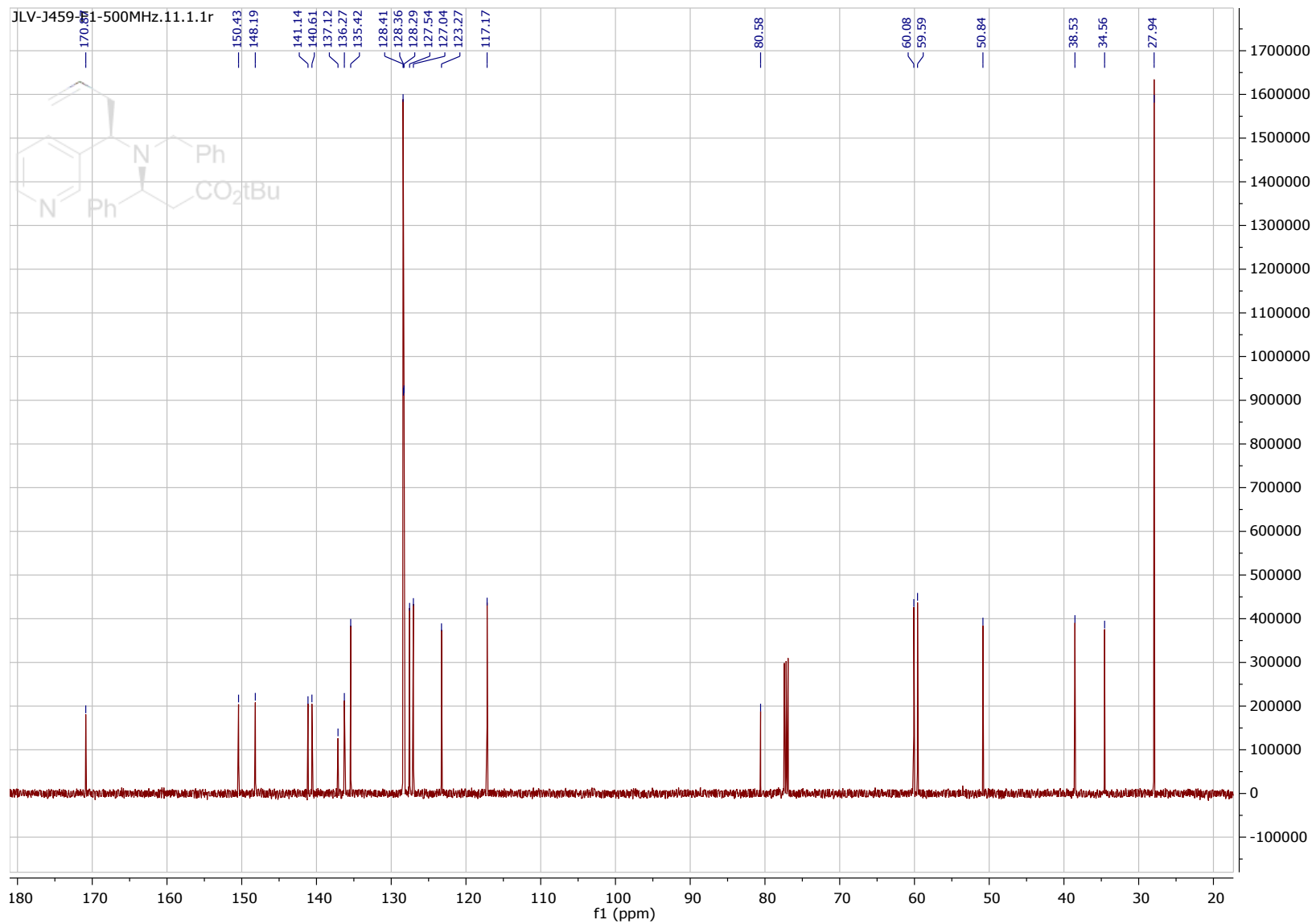
(±)-(*R*)-Tert-Butyl 3-{benzyl[(*R*)-1-(furan-3-yl)but-3-en-1-yl]amino}octanoate 3l



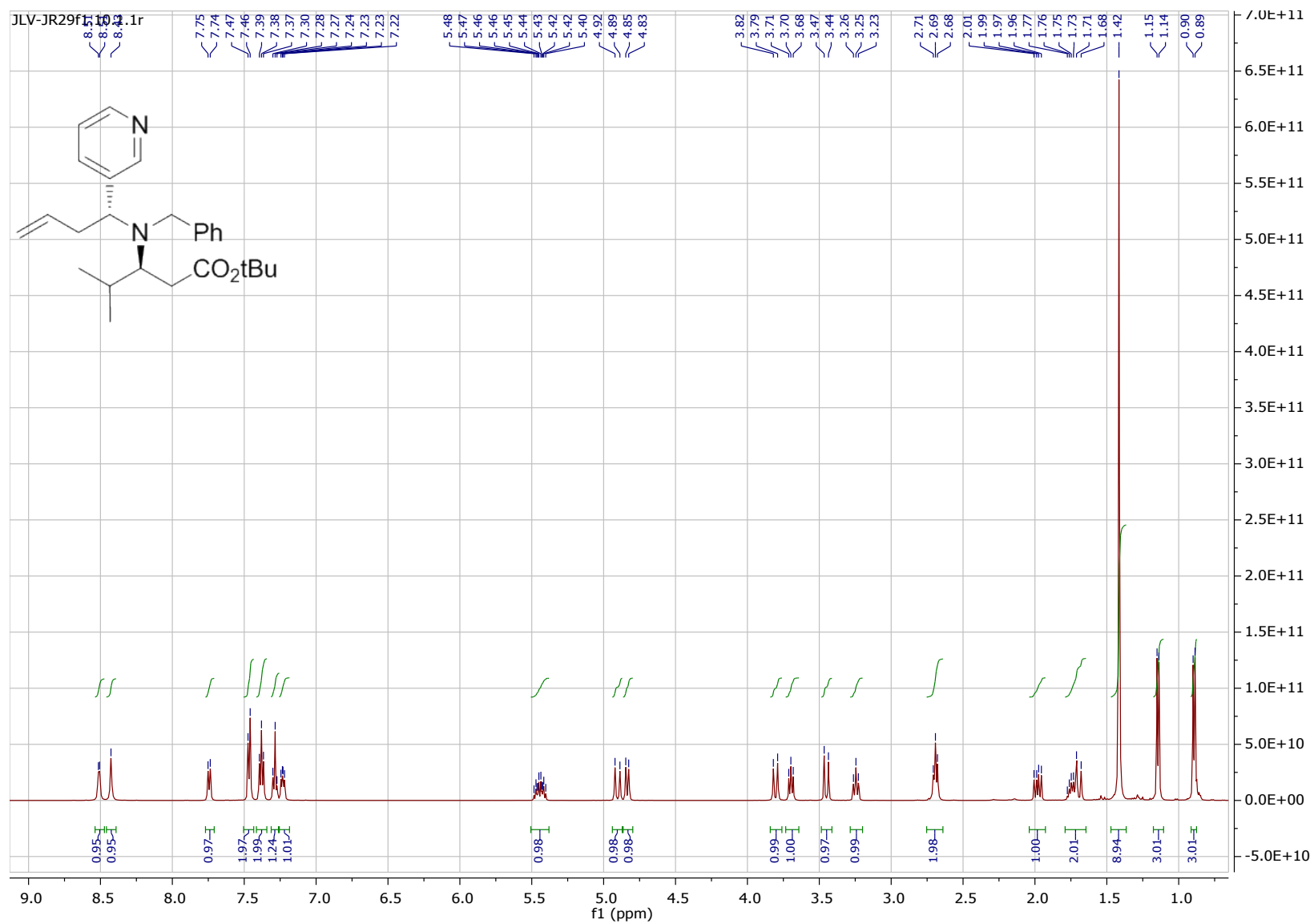


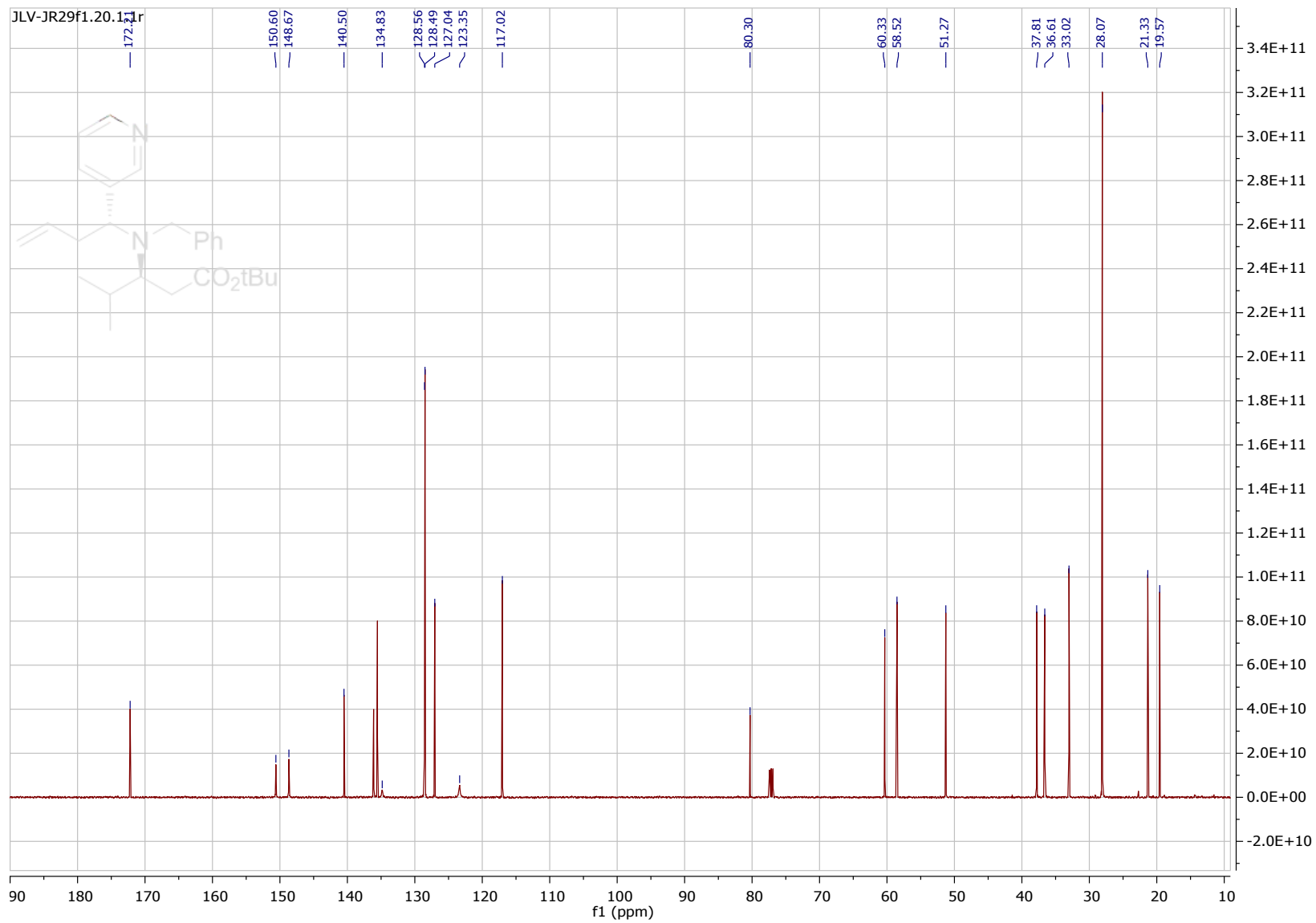
Chemical structure: C=CC[C@H]1CC[C@@H](C1Cc2ccccc2)C(=O)OCC(C)(C)C

¹H NMR spectrum (400 MHz, CDCl₃) showing peaks from 0.5 to 8.5 ppm. The spectrum includes a broad peak at ~7.2 ppm (NH), aromatic signals between 7.0-7.5 ppm, a multiplet at ~4.8 ppm, a doublet at ~3.8 ppm, a multiplet at ~2.5 ppm, and a large peak at ~1.2 ppm (tBu). Integration values are shown below the baseline.

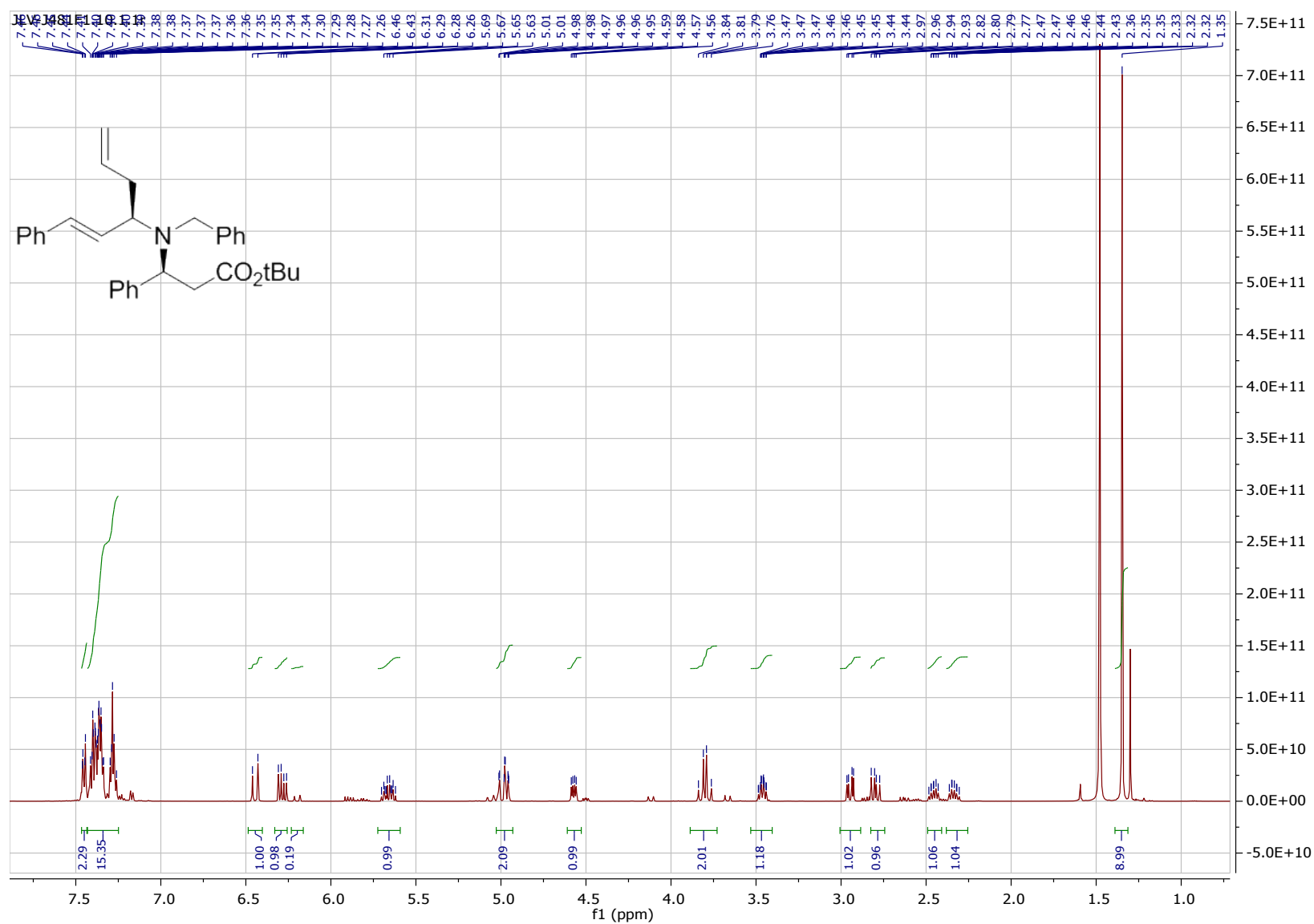


(±)-(S)-Tert-Butyl 3-{benzyl[(R)-1-(pyridin-3-yl)but-3-en-1-yl]amino}-4-methylpentanoate 3n

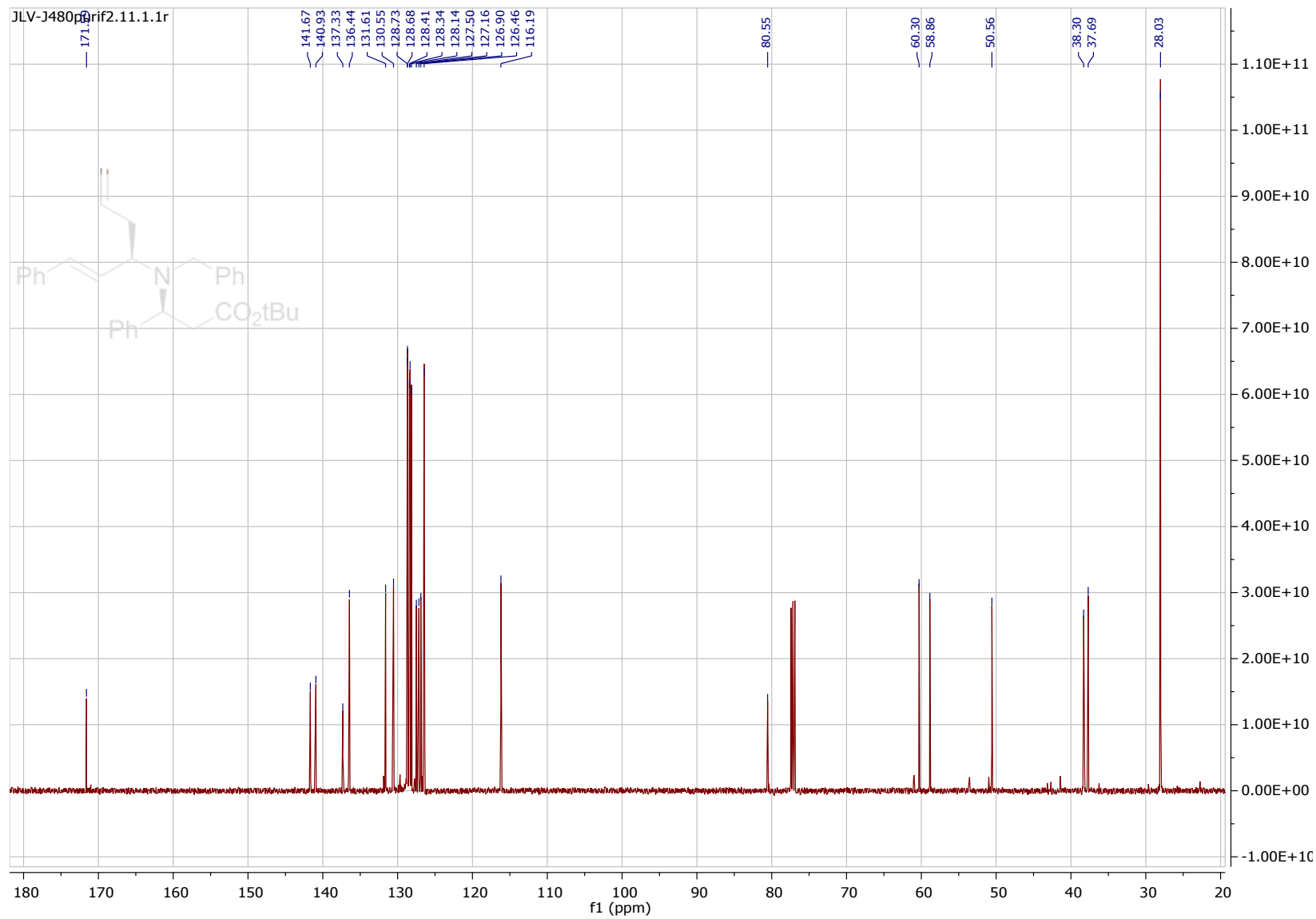




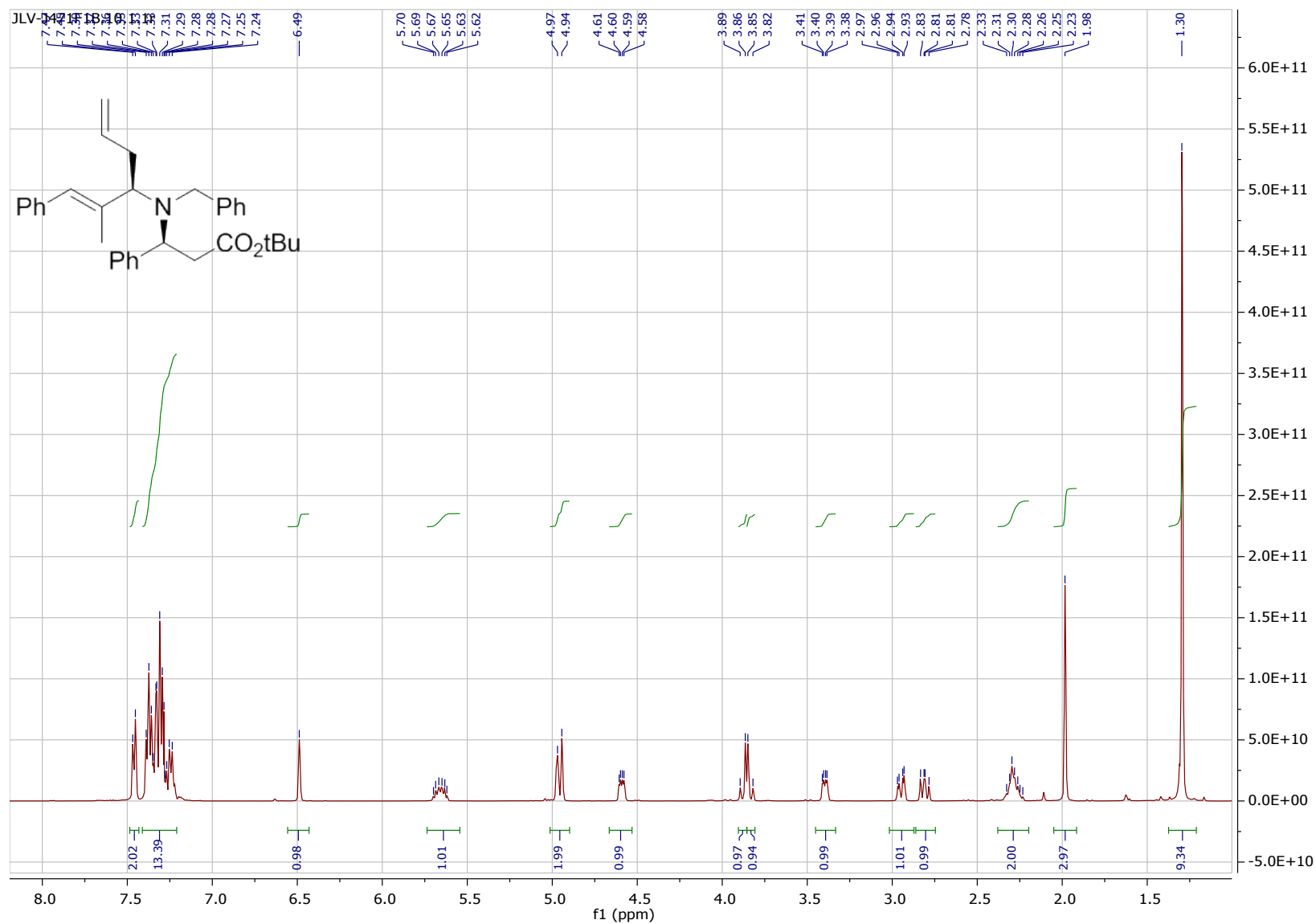
(±)-(*S*)-*Tert*-Butyl 3-{benzyl[(*R,E*)-1-phenylhexa-1,5-dien-3-yl]amino}-3-phenylpropanoate **3o**

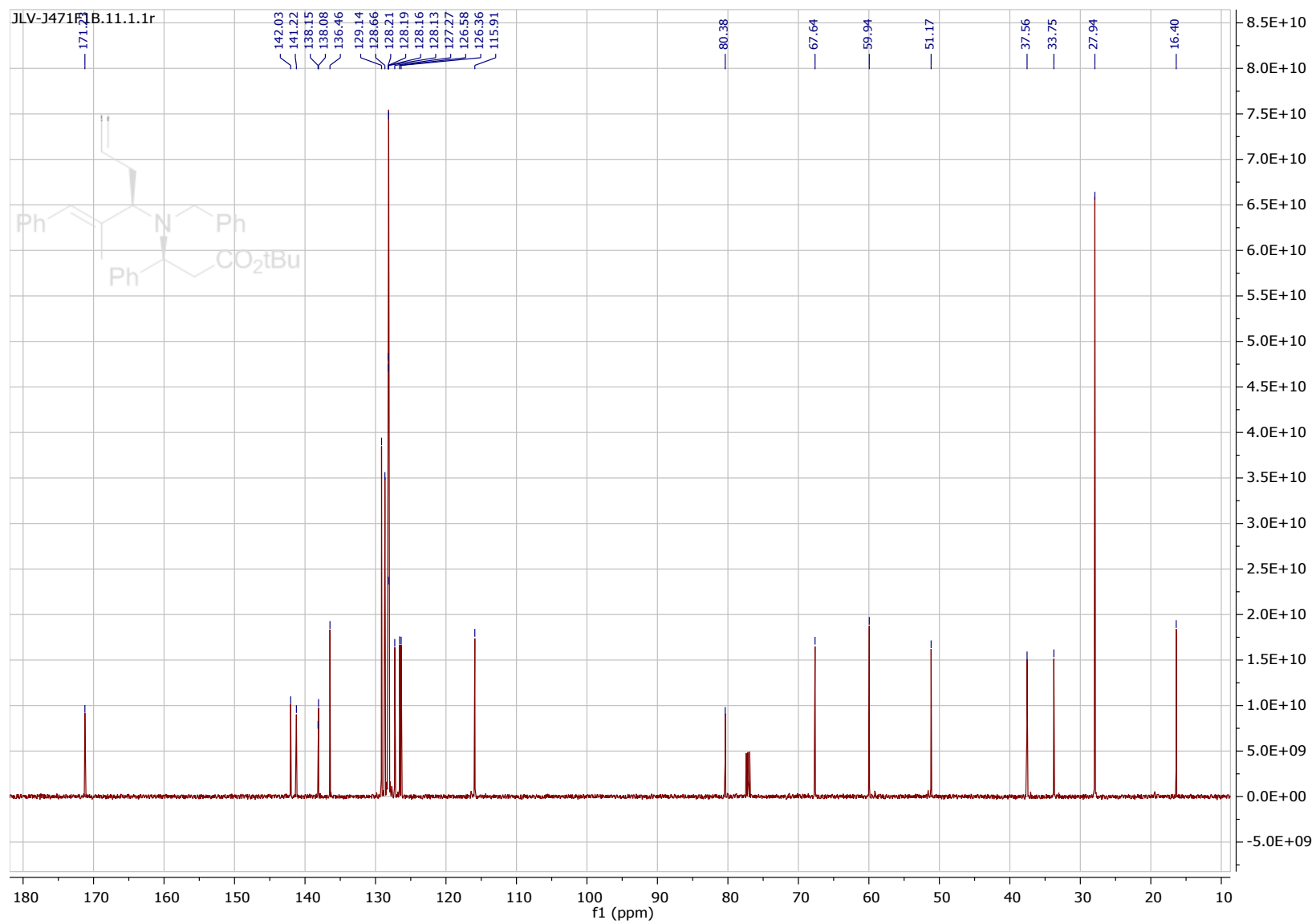


JLV-J480p
Prif2.11.1.1r

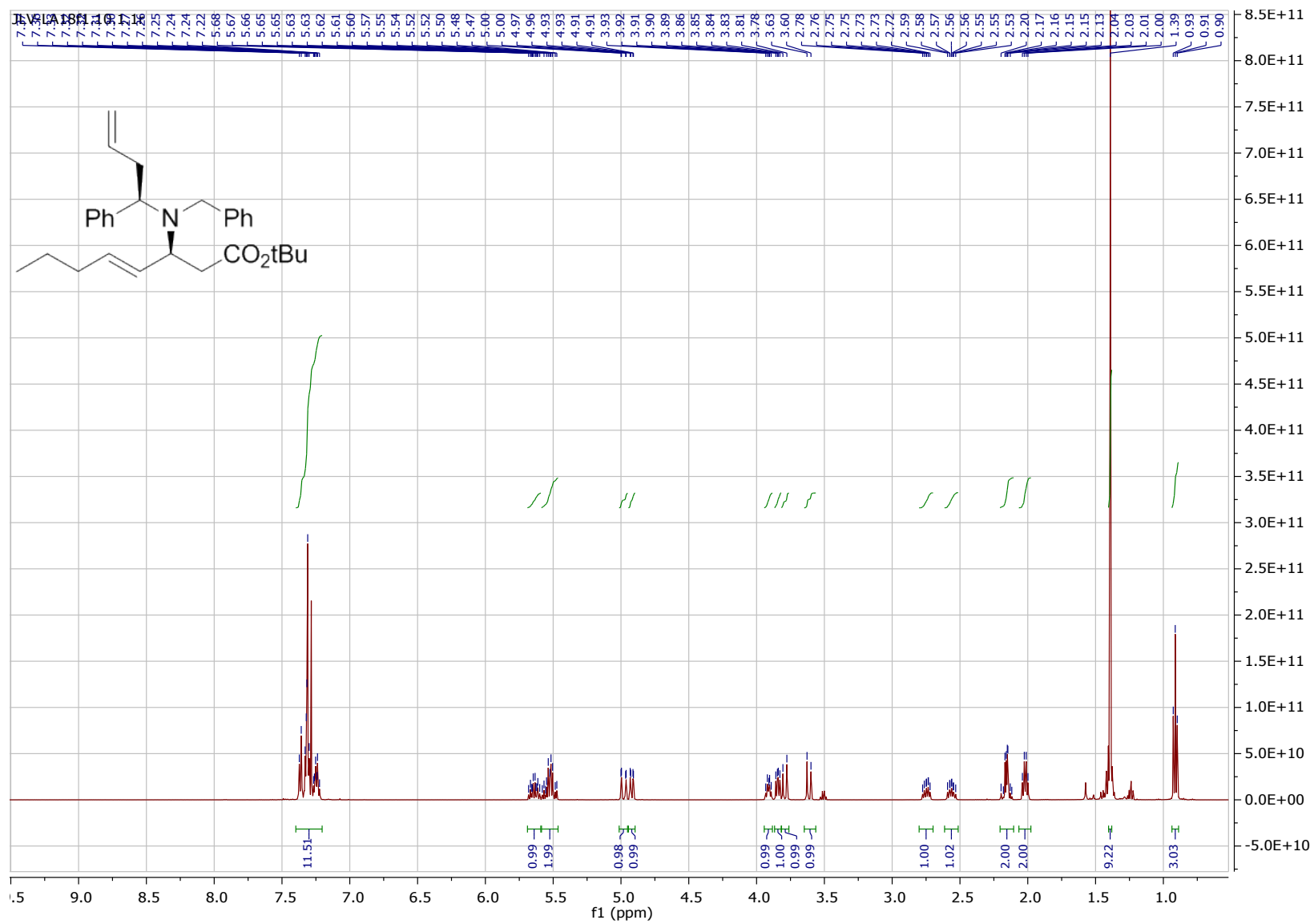


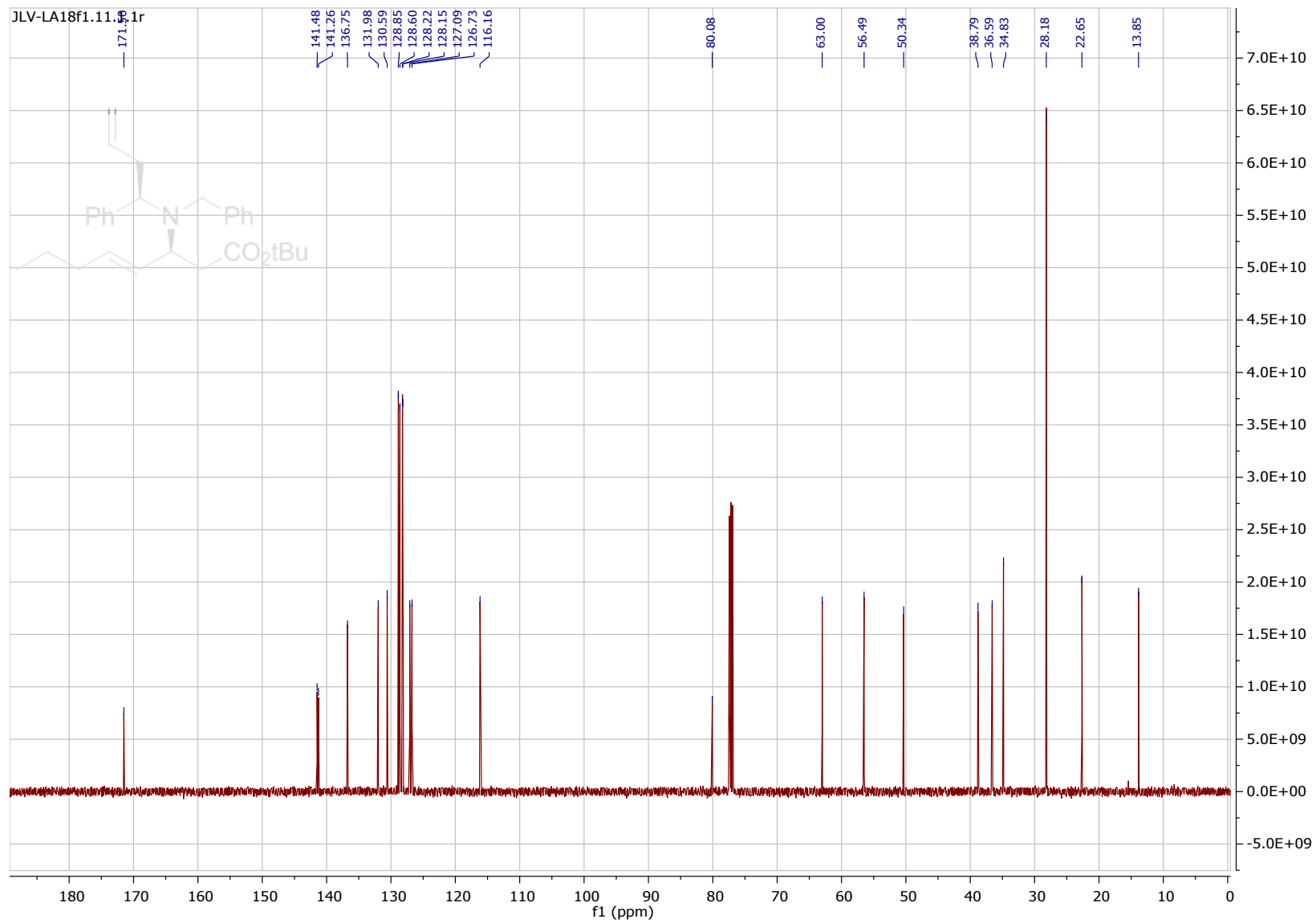
(±)-(*S*)-*Tert*-Butyl 3-(benzyl(*R,E*)-2-methyl-1-phenylhexa-1,5-dien-3-yl)amino)-3-phenylpropanoate 3p



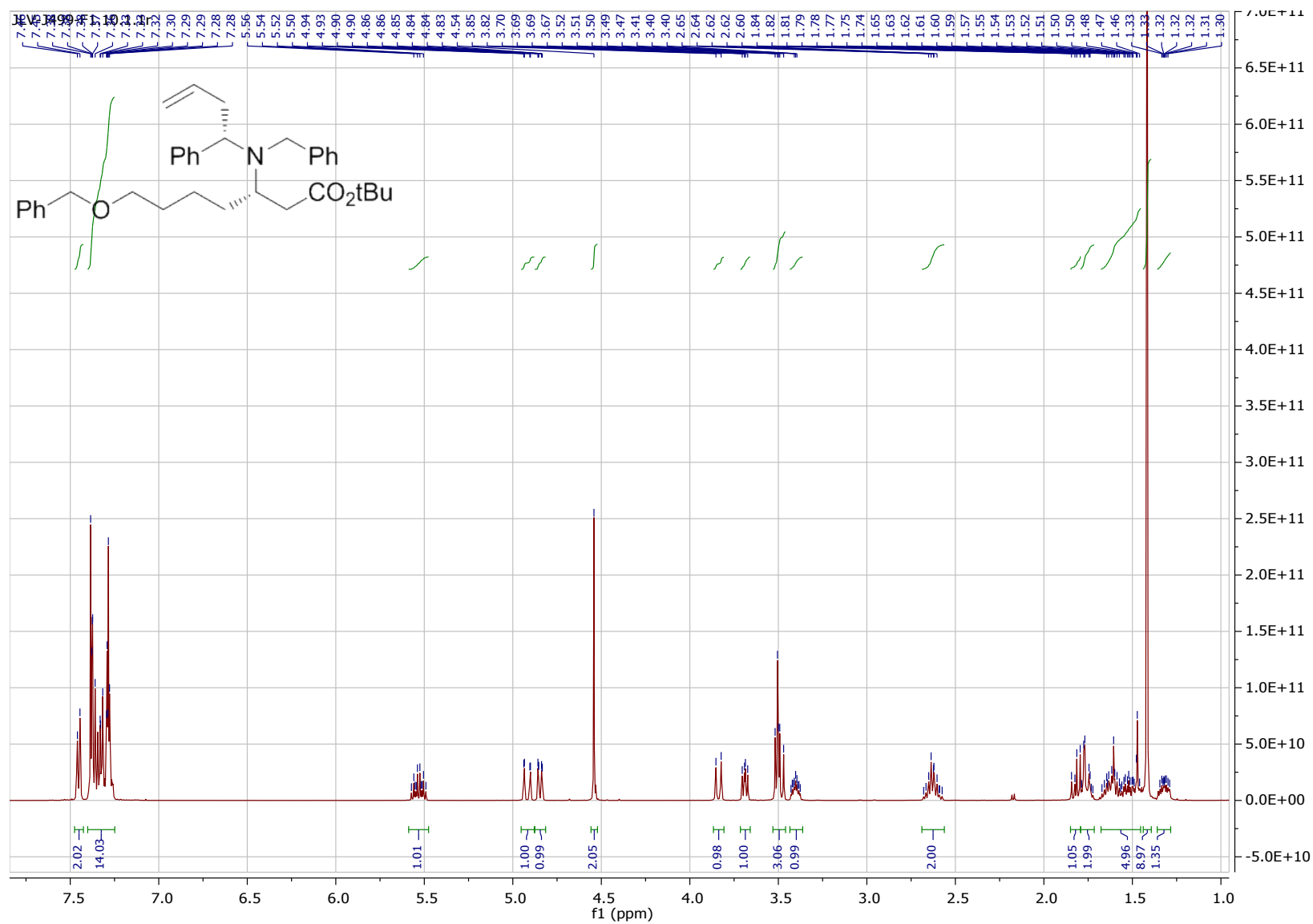


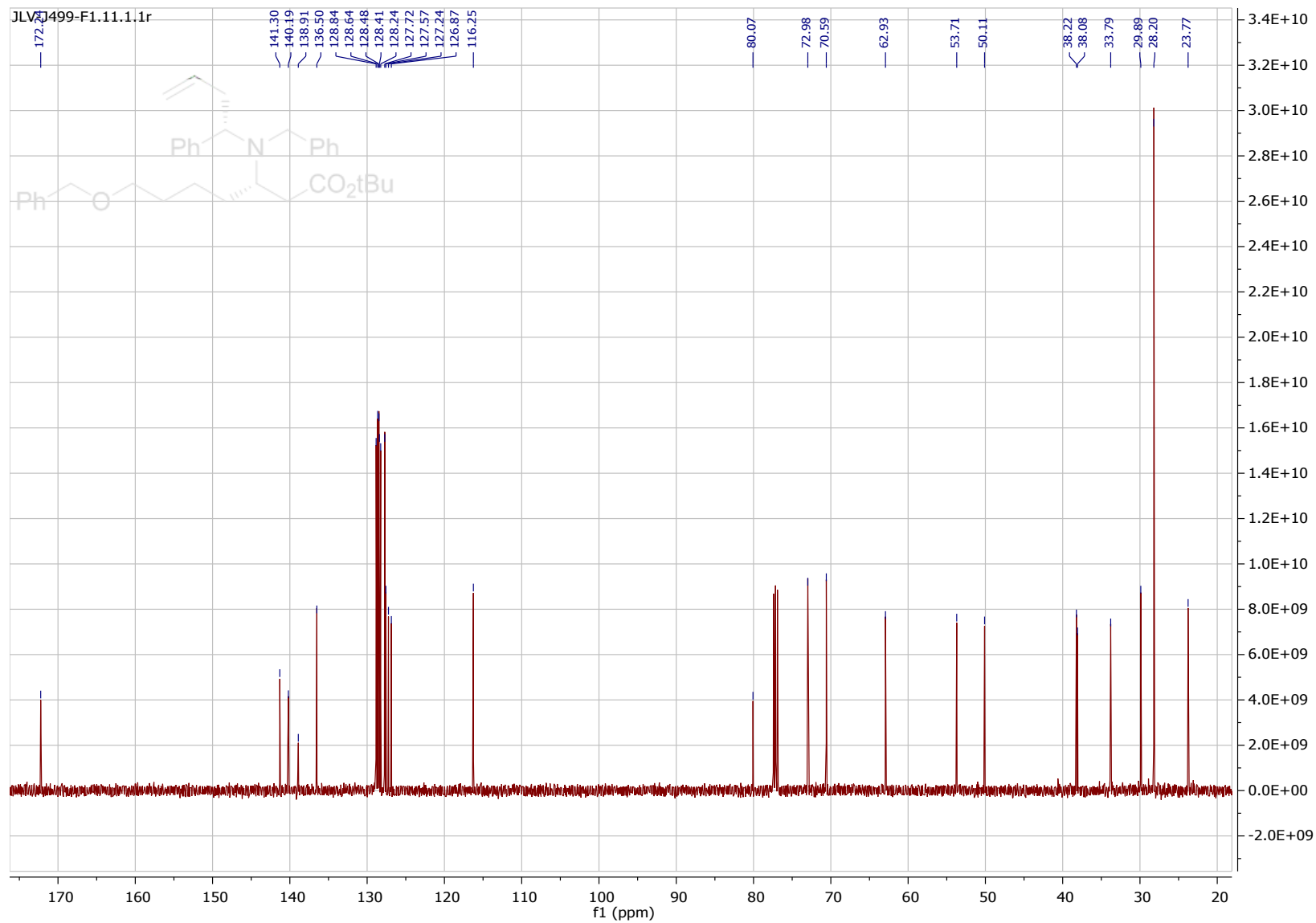
(S,E)-Tert-Butyl 3-{benzyl[(R)-1-phenylbut-3-en-1-yl]amino}oct-4-enoate 3q



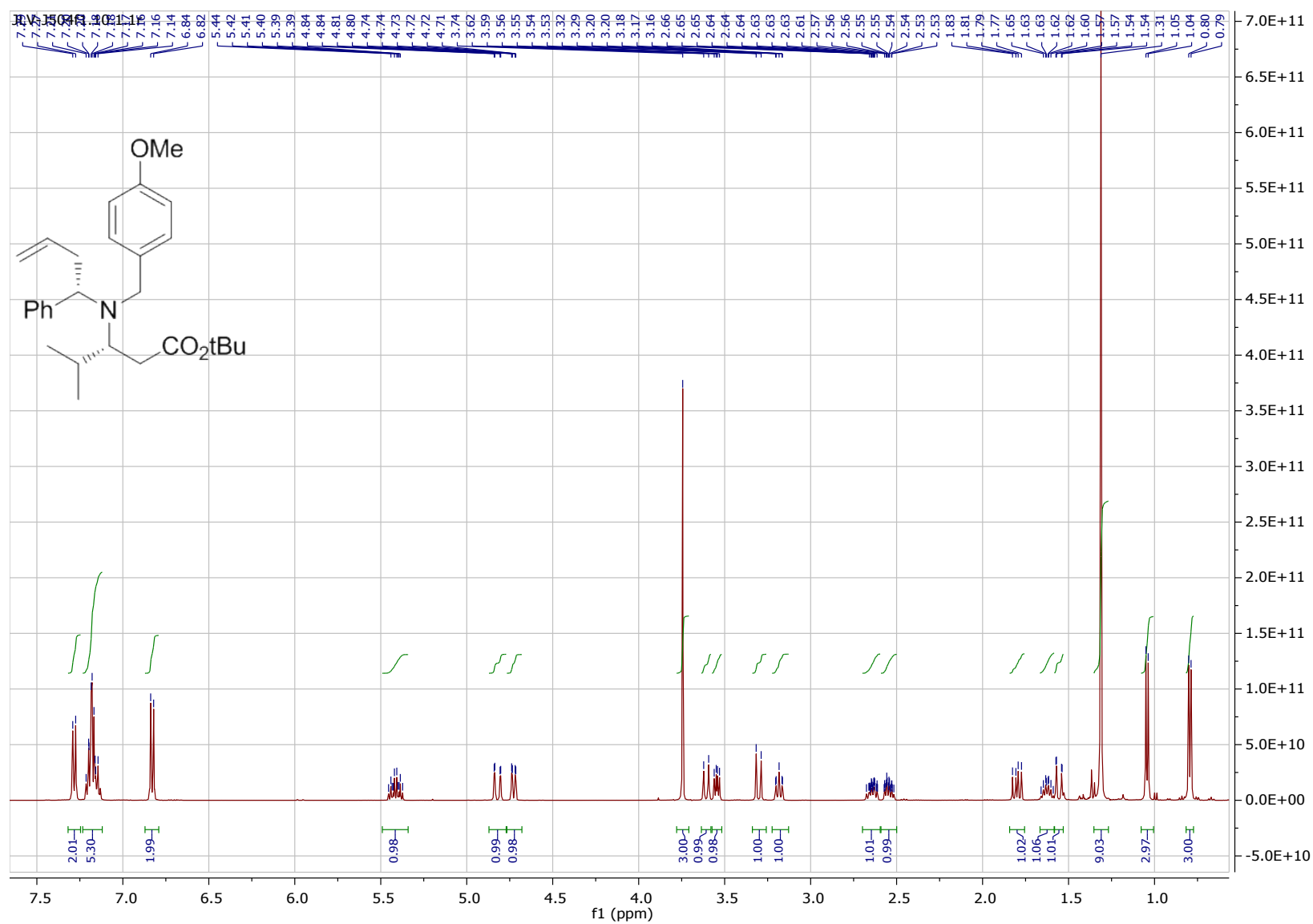


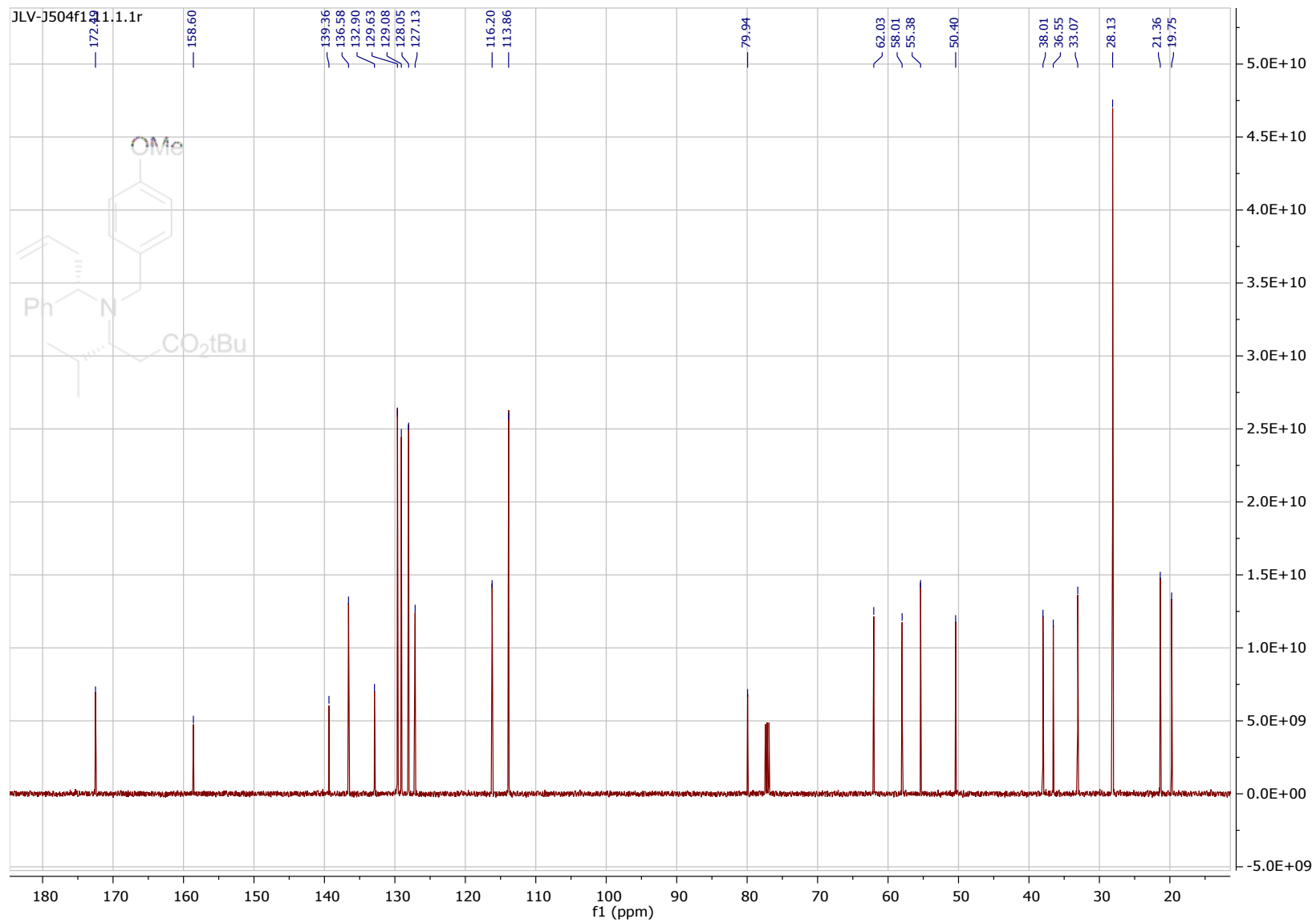
(S)-Tert-Butyl 3-{benzyl[(S)-1-phenylbut-3-en-1-yl]amino}-7-(benzyloxy)heptanoate 3r



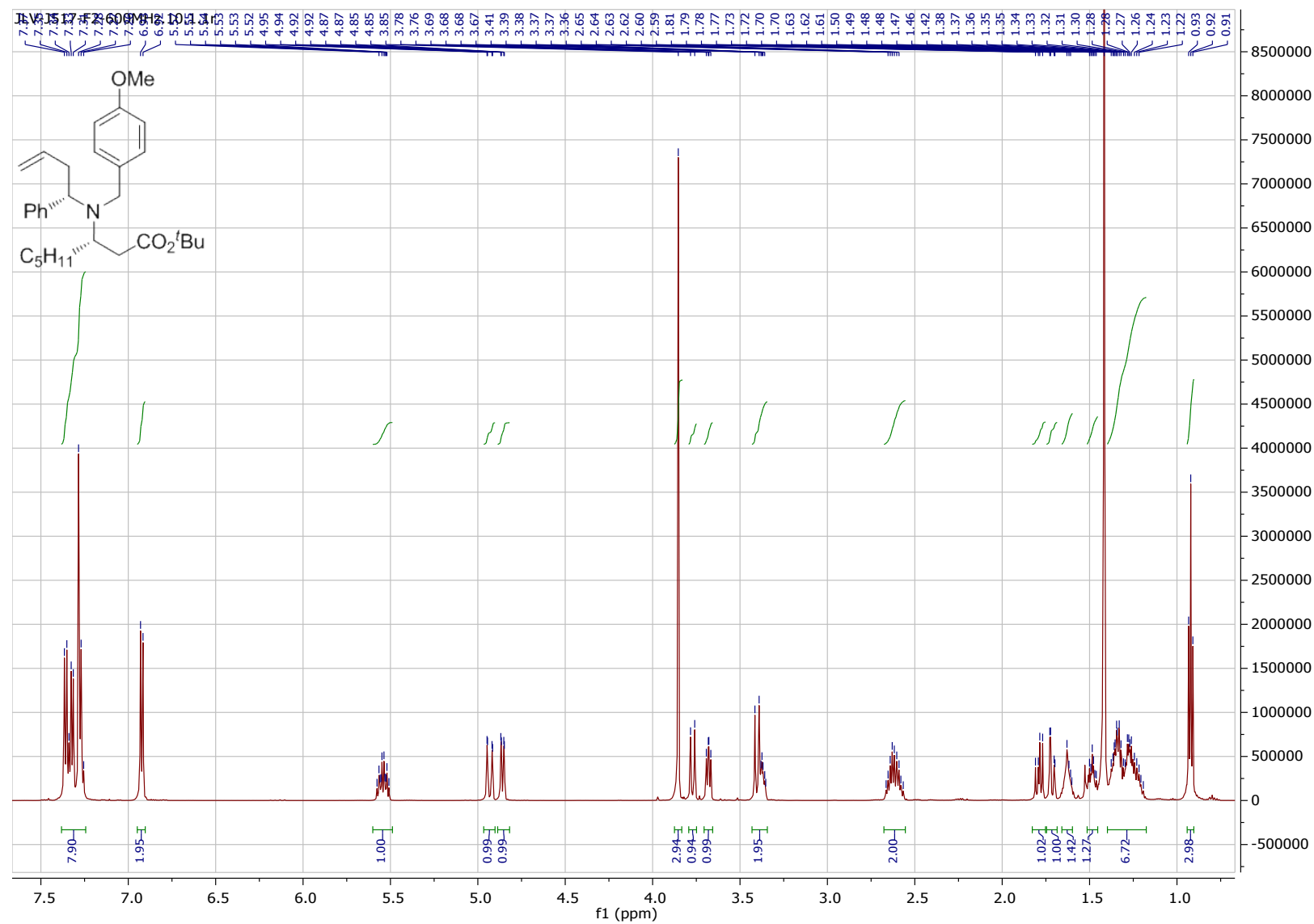


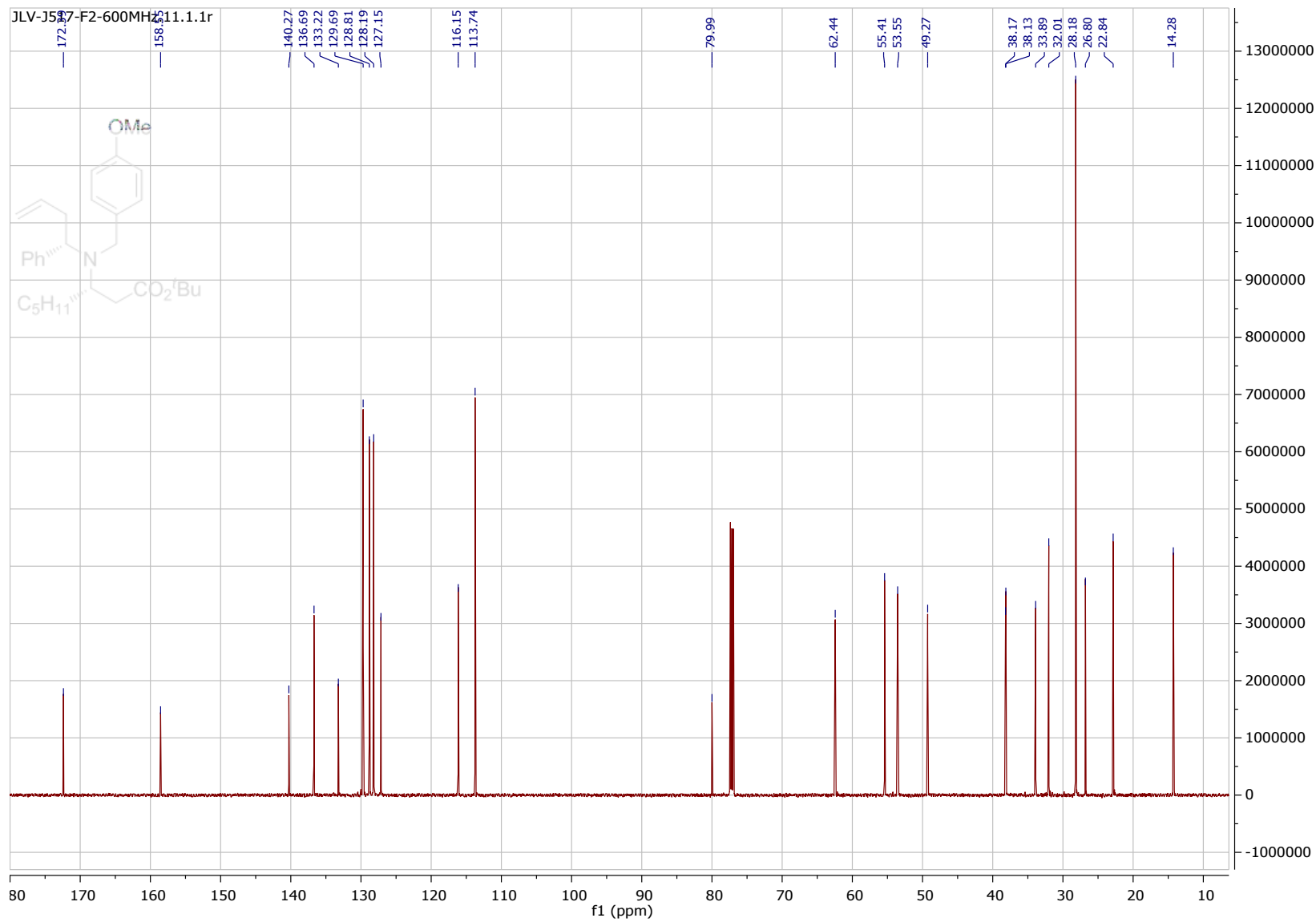
(R)-Tert-Butyl 3-((4-methoxybenzyl)[(S)-1-phenylbut-3-en-1-yl]amino)-4-methylpentanoate 3s



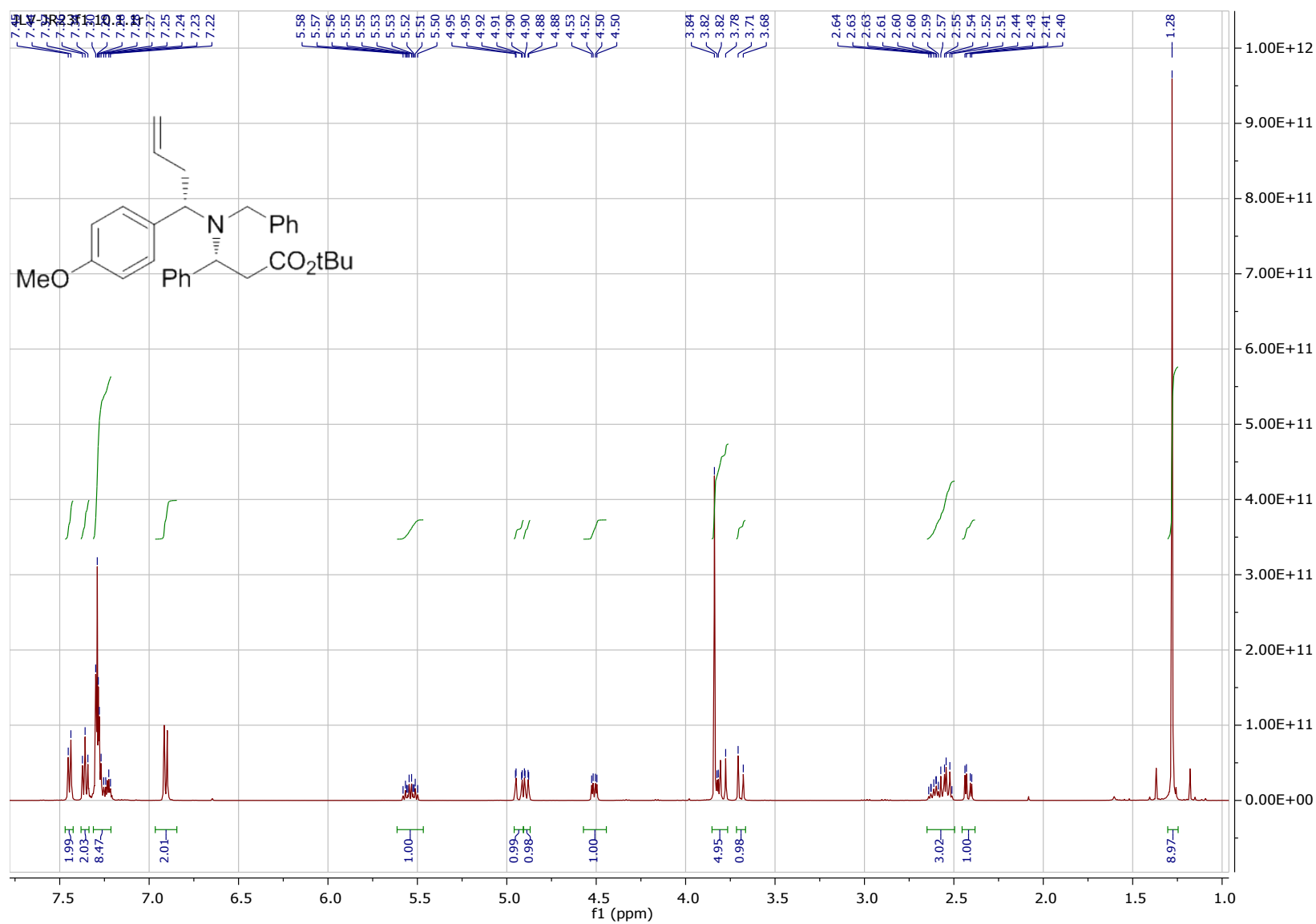


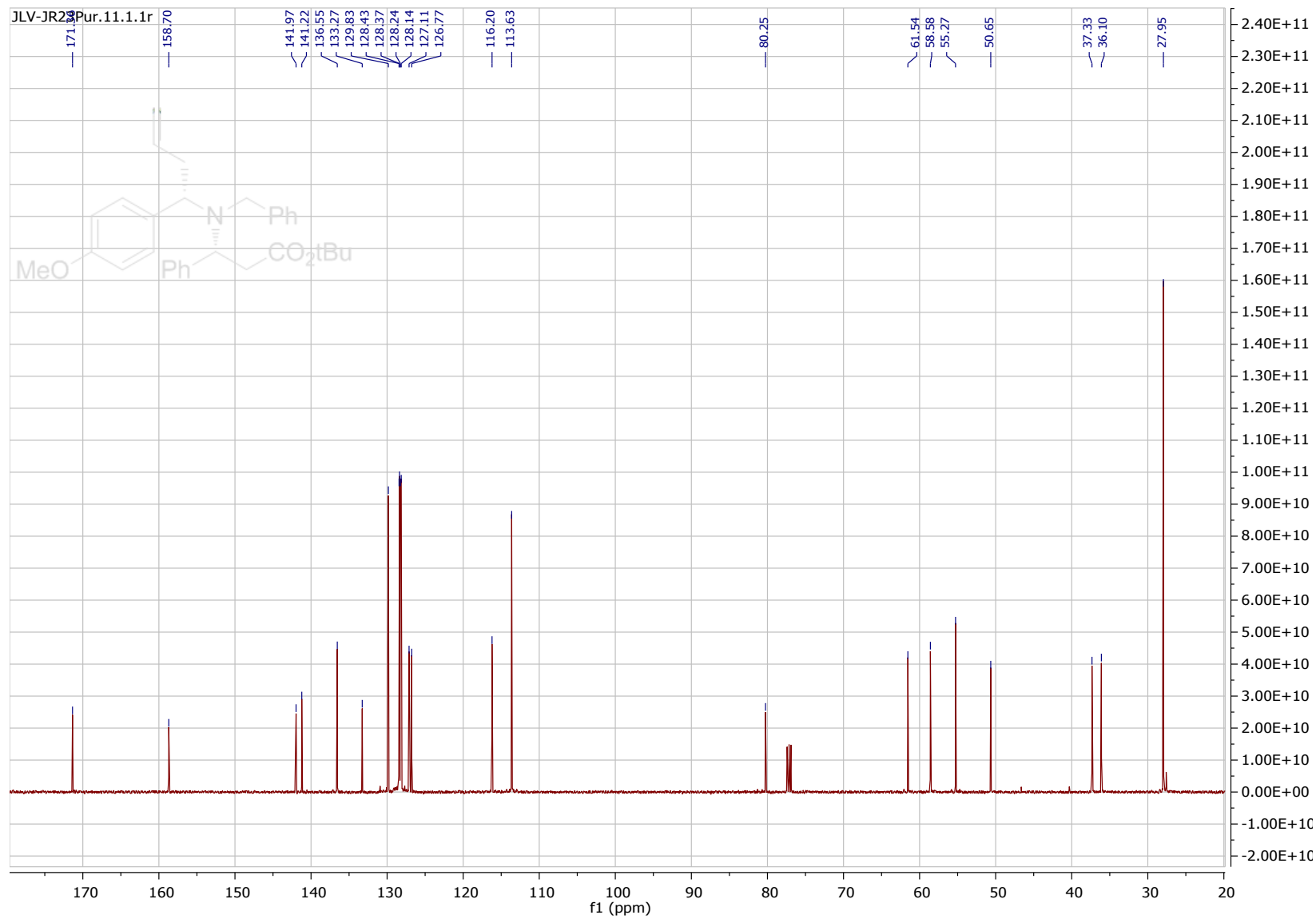
(S)-Tert-Butyl 3-((4-methoxybenzyl)[(S)-1-phenylbut-3-en-1-yl]amino)octanoate 3t



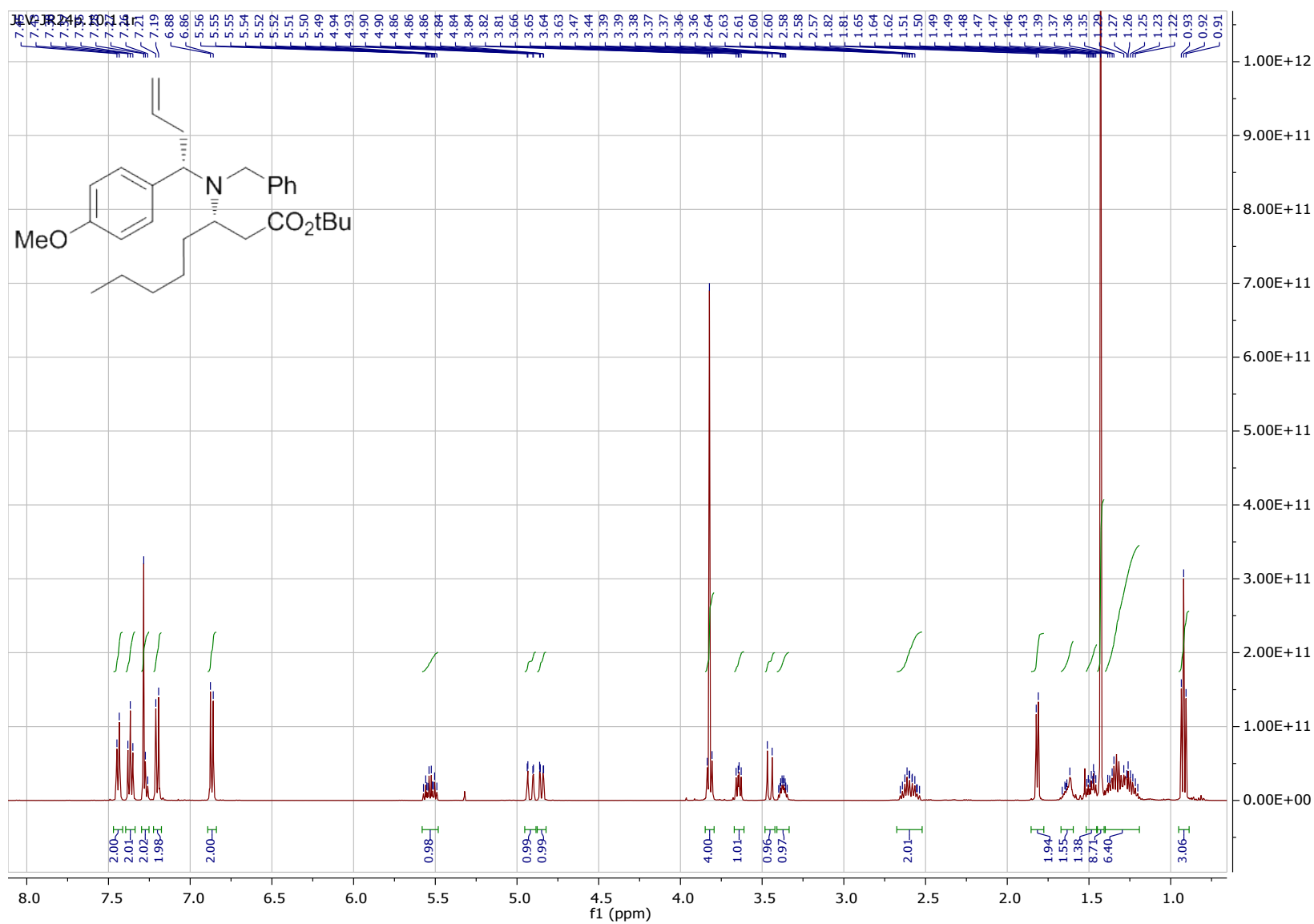


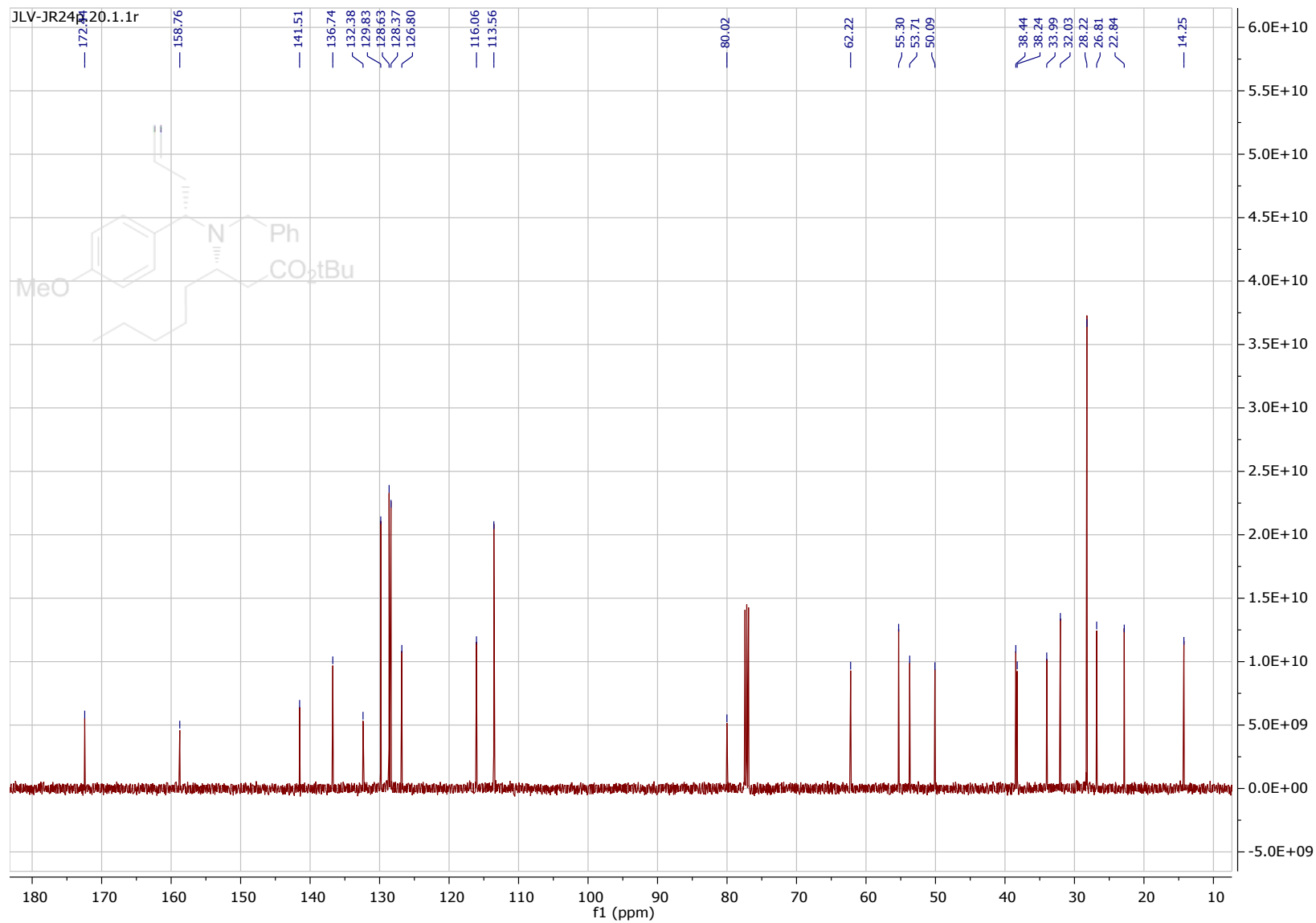
(R)-Tert-Butyl 3-{benzyl[(S)-1-(4-methoxyphenyl)but-3-en-1-yl]amino}-3-phenylpropanoate 3u



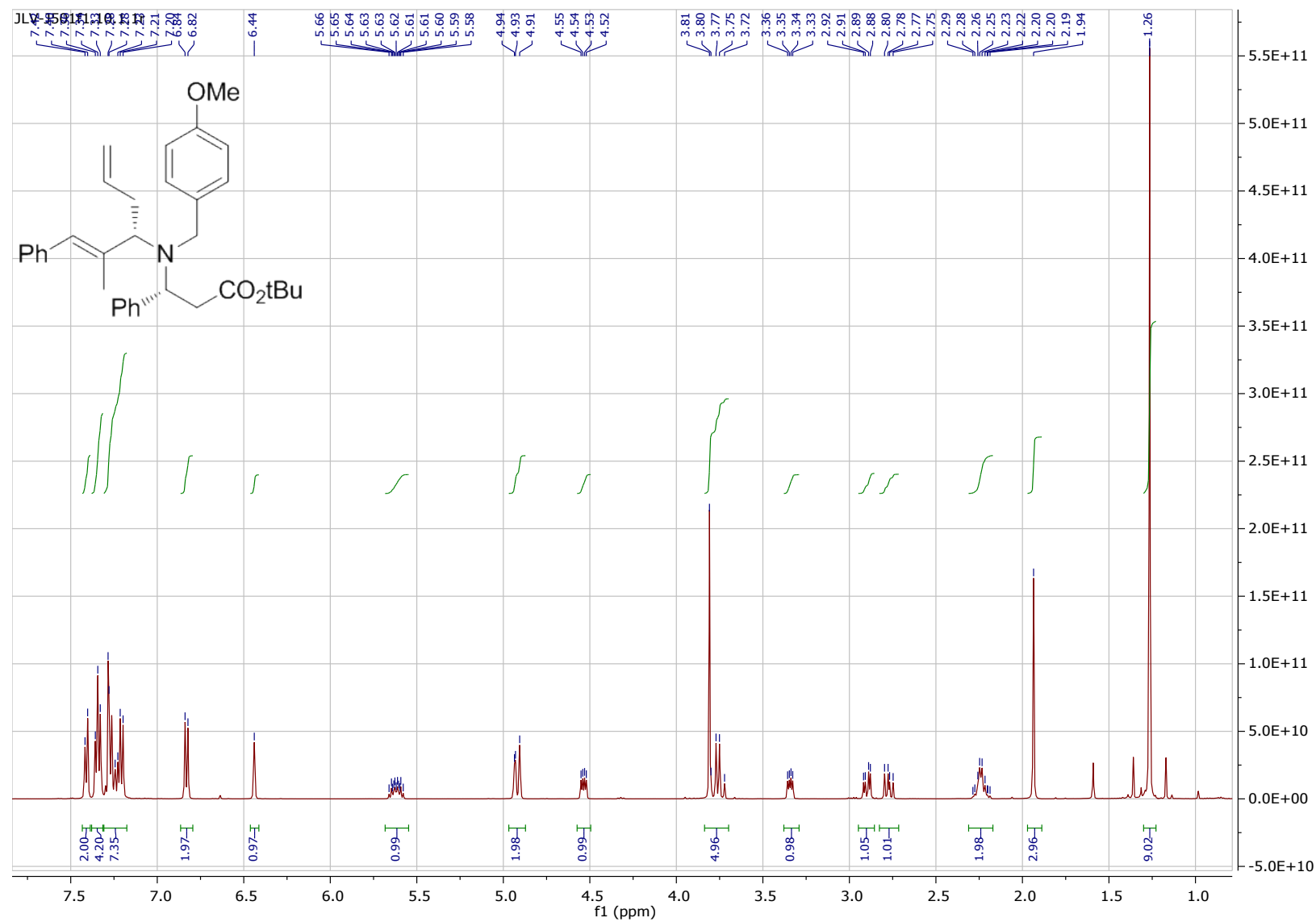


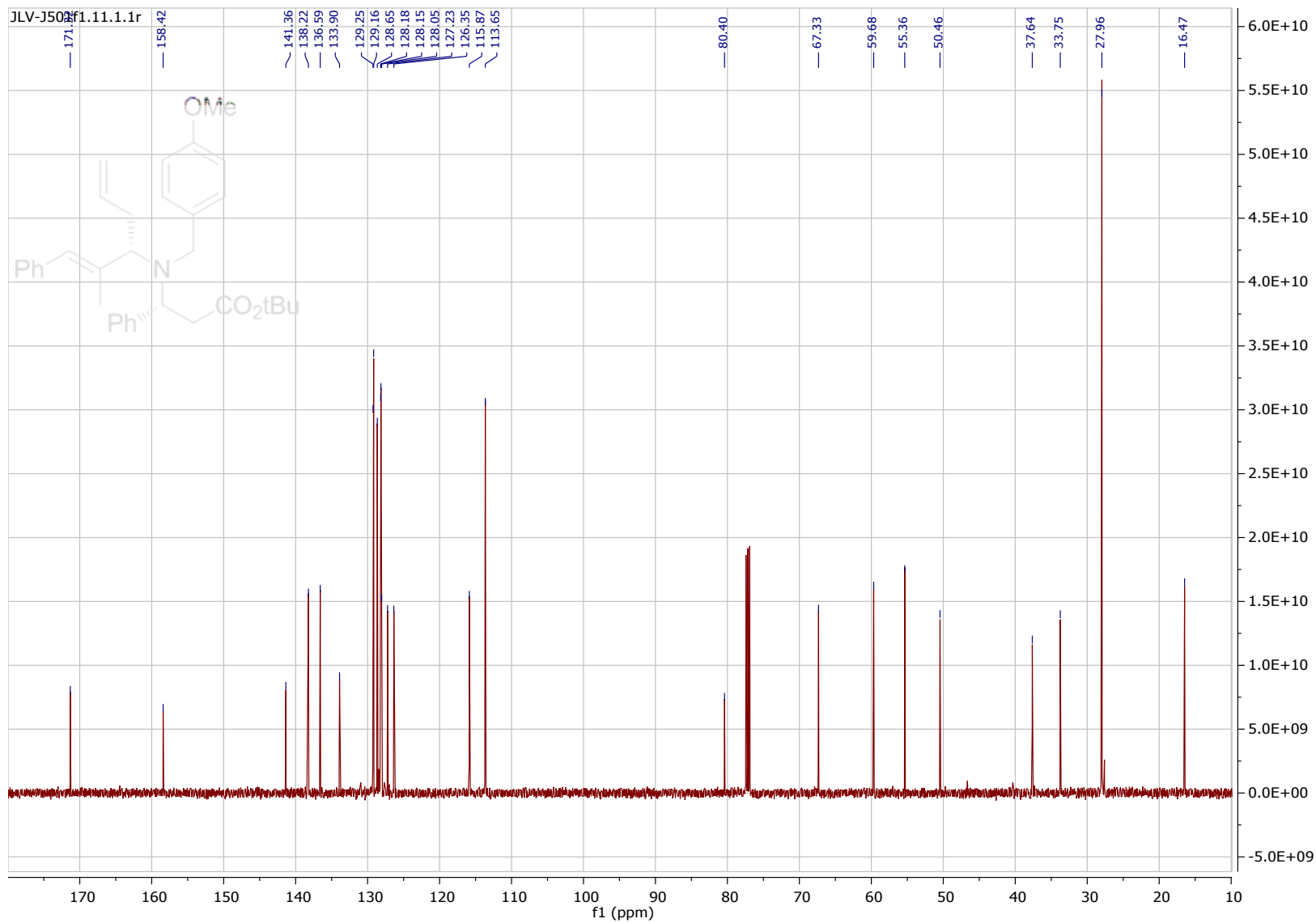
(S)-Tert-Butyl 3-{benzyl[(S)-1-(4-methoxyphenyl)but-3-en-1-yl]amino}octanoate 3v



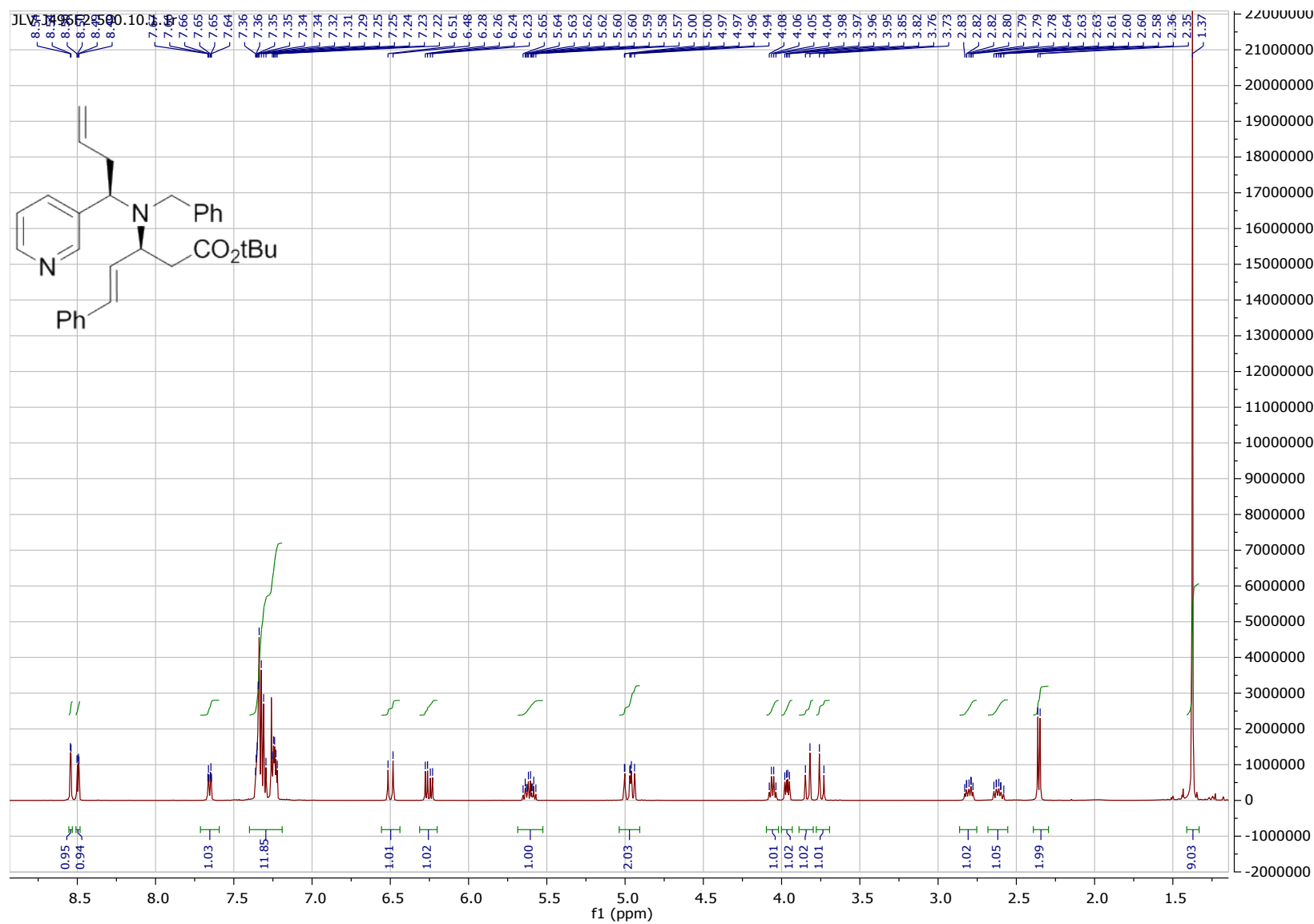


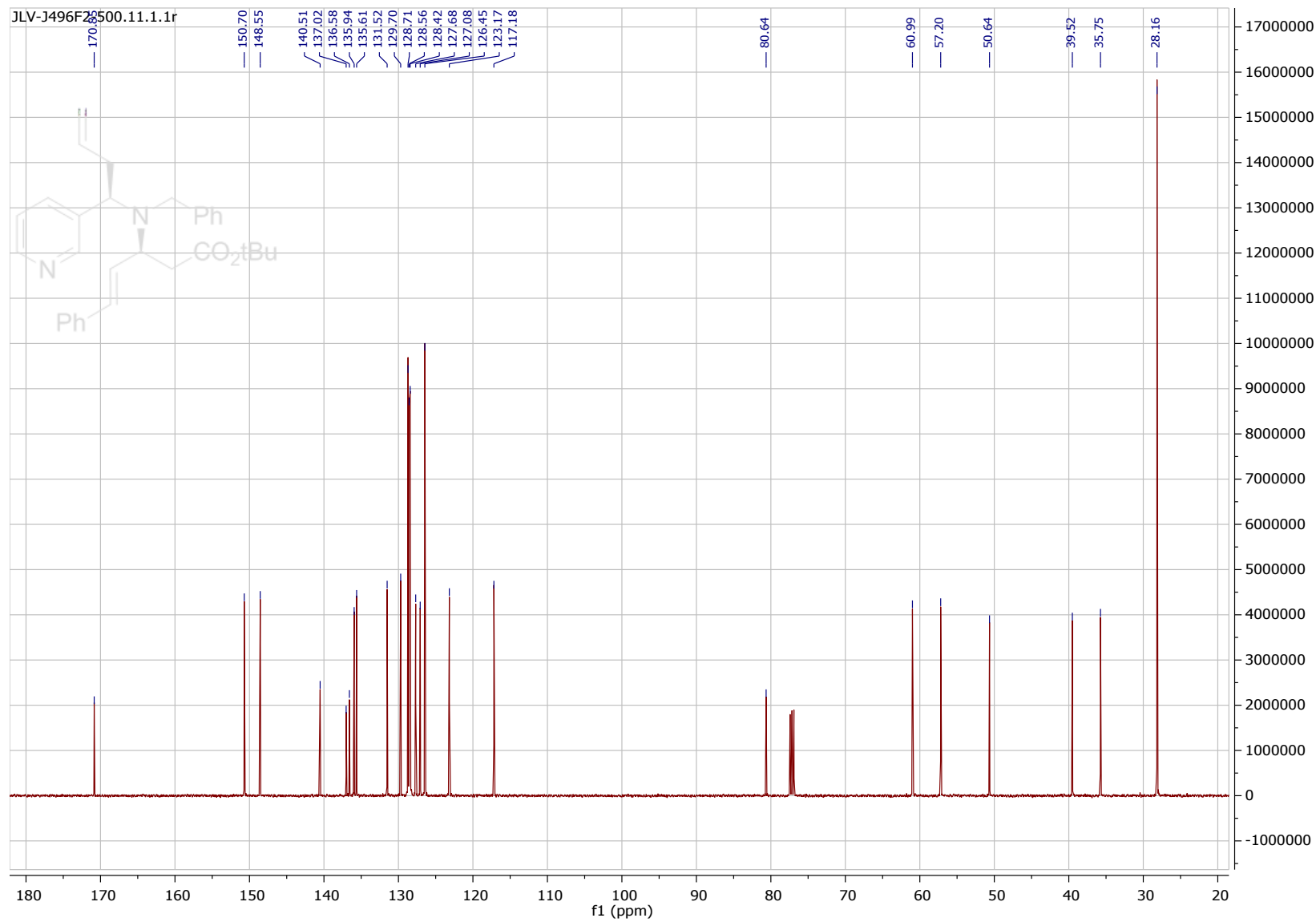
(R)-Tert-Butyl 3-((4-methoxybenzyl)[(S,E)-2-methyl-1-phenylhexa-1,5-dien-3-yl]amino)-3-phenylpropanoate 3w



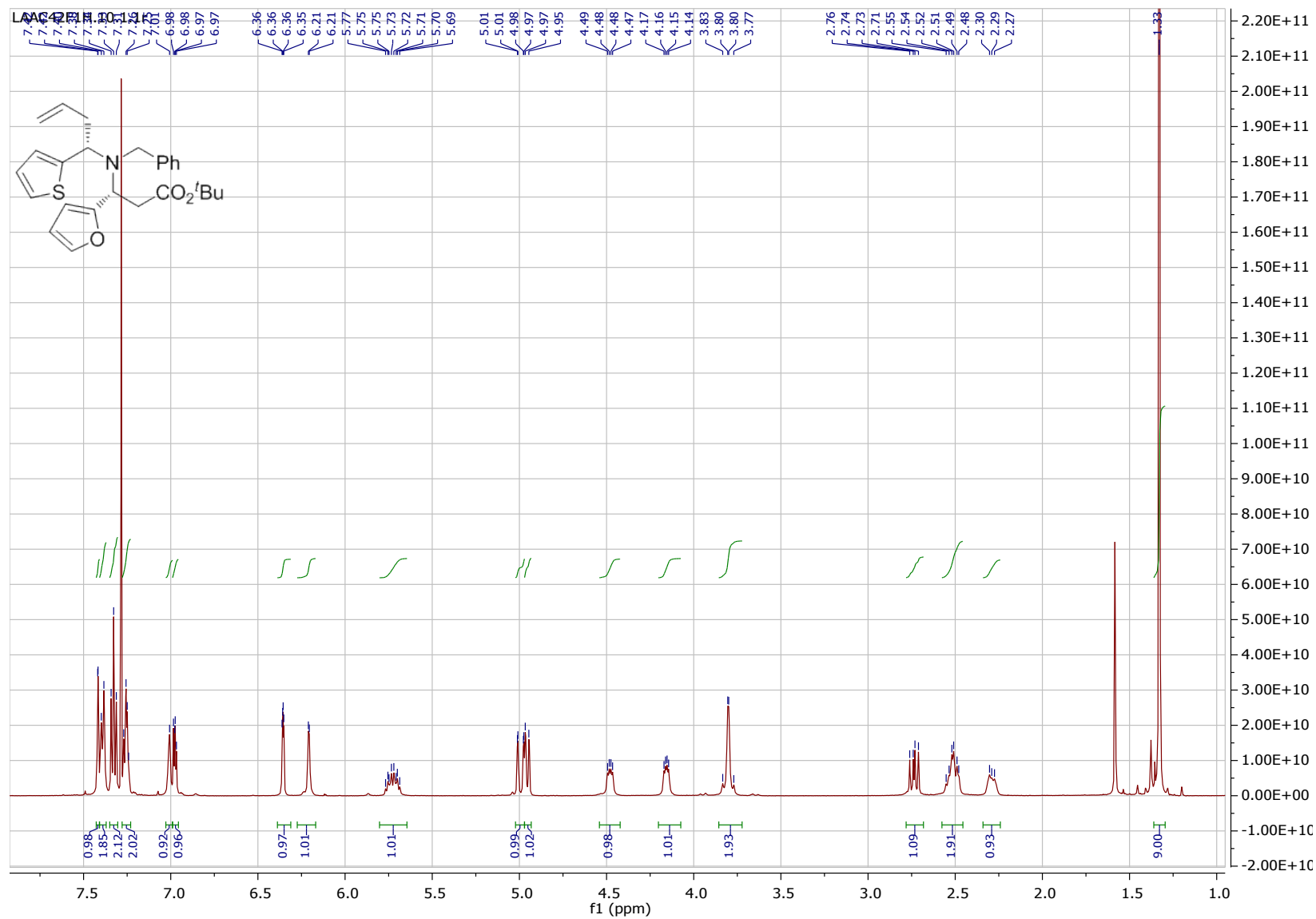


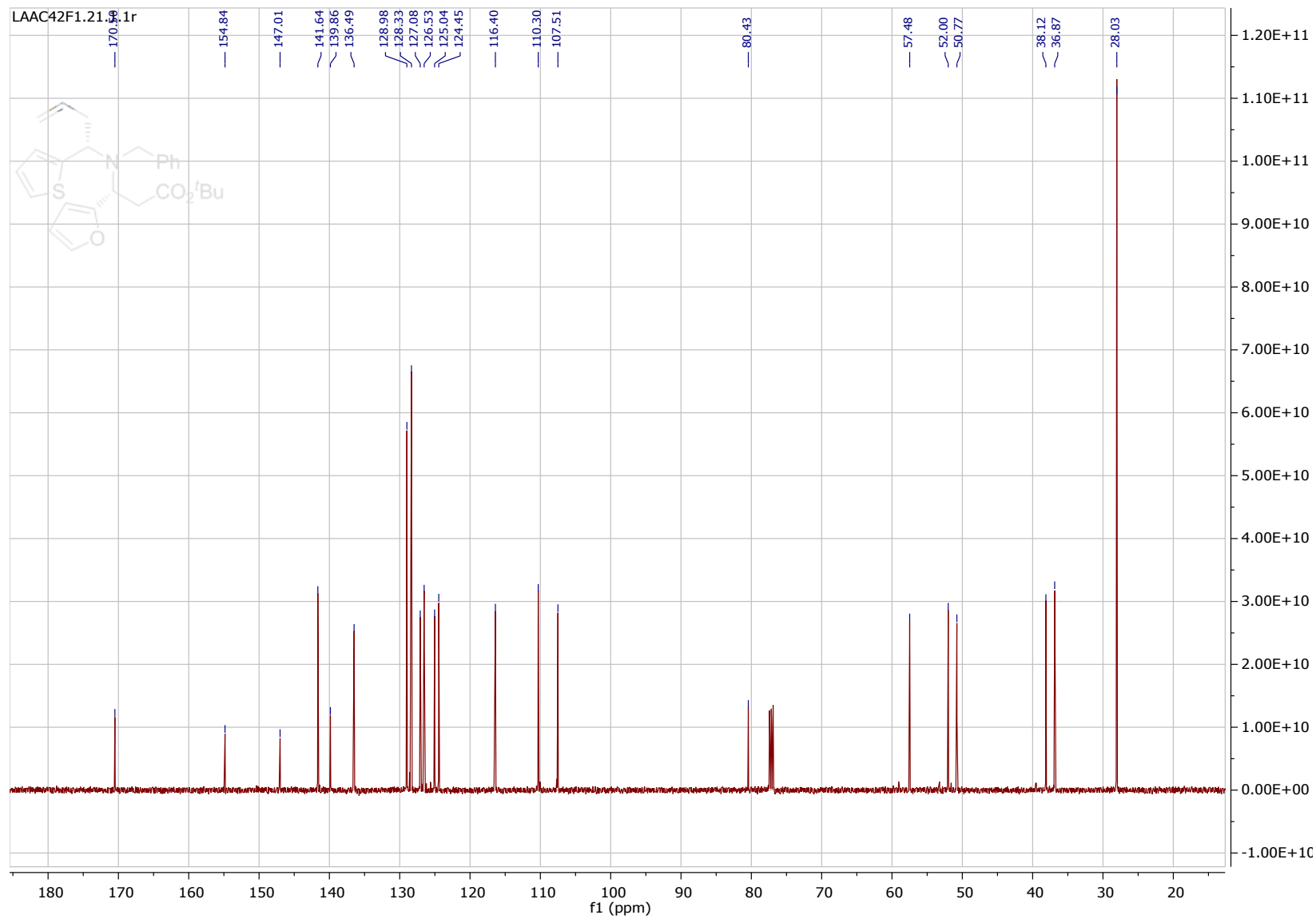
(S,E)-Tert-Butyl 3-{benzyl[(R)-1-(pyridin-3-yl)but-3-en-1-yl]amino}-5-phenylpent-4-enoate 3x



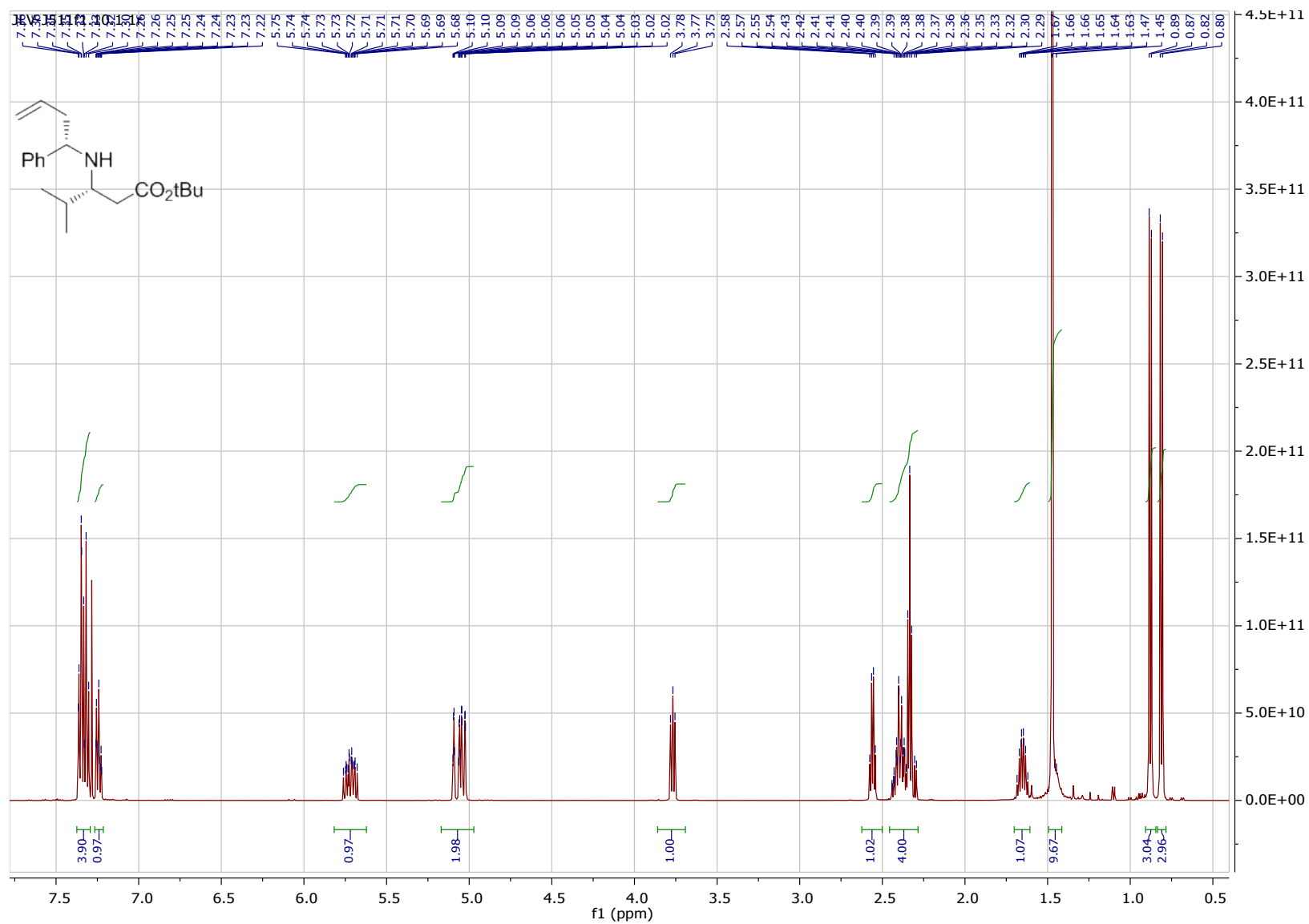


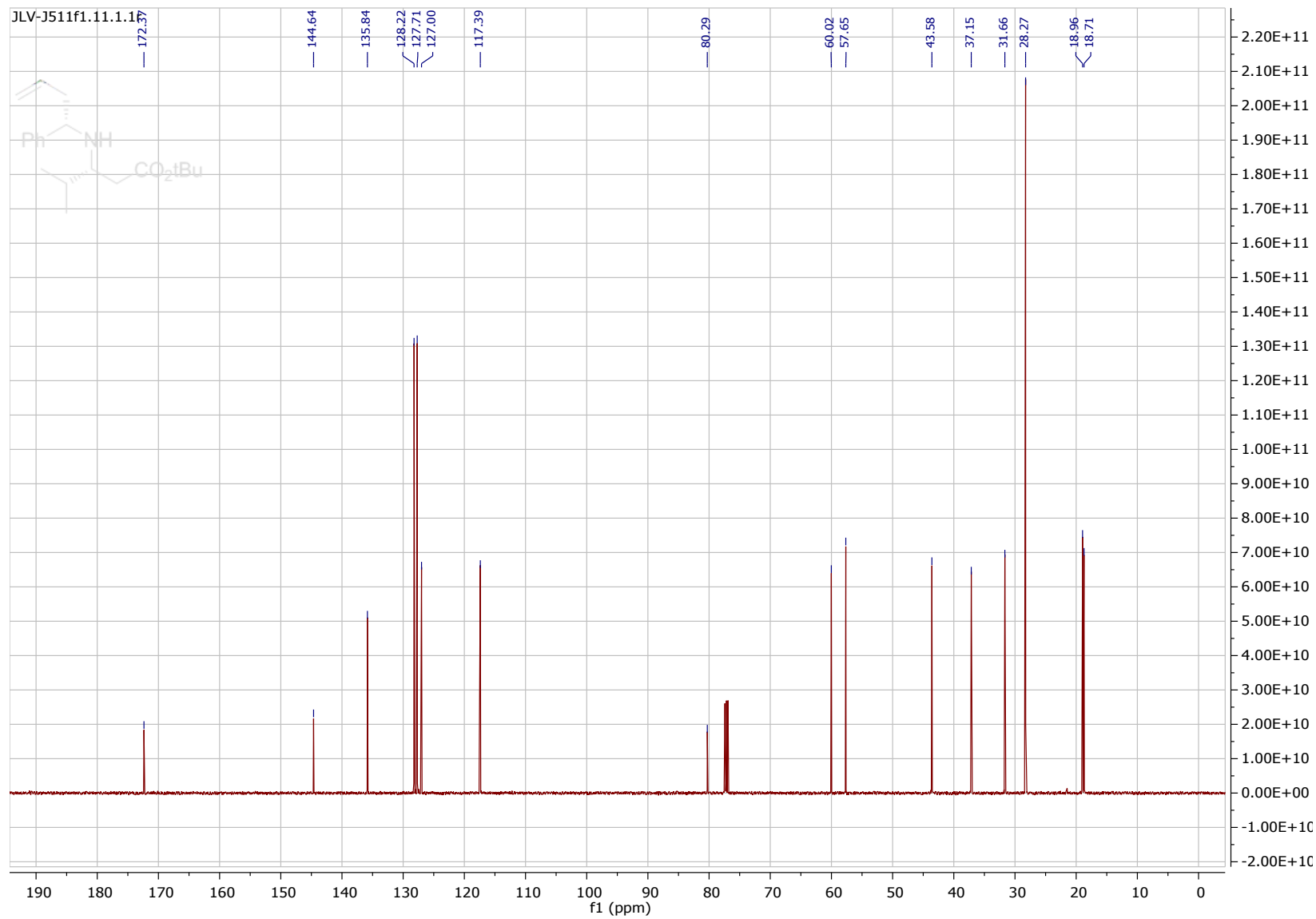
(R)-Tert-Butyl 3-{Benzyl[(S)-1-(thiophen-2-yl)but-3-en-1-yl]amino}-3-(furan-2-yl)propanoate 3y



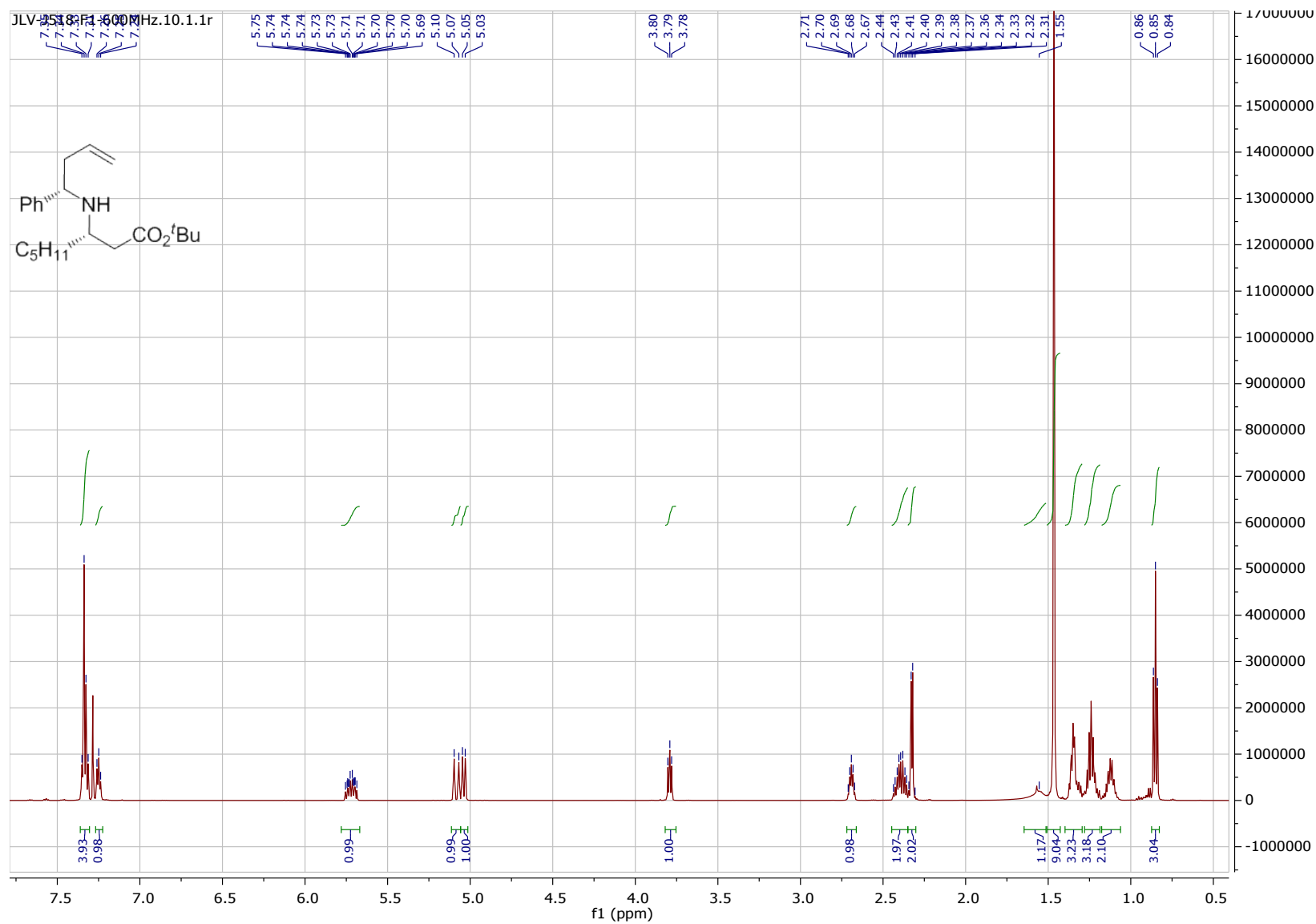


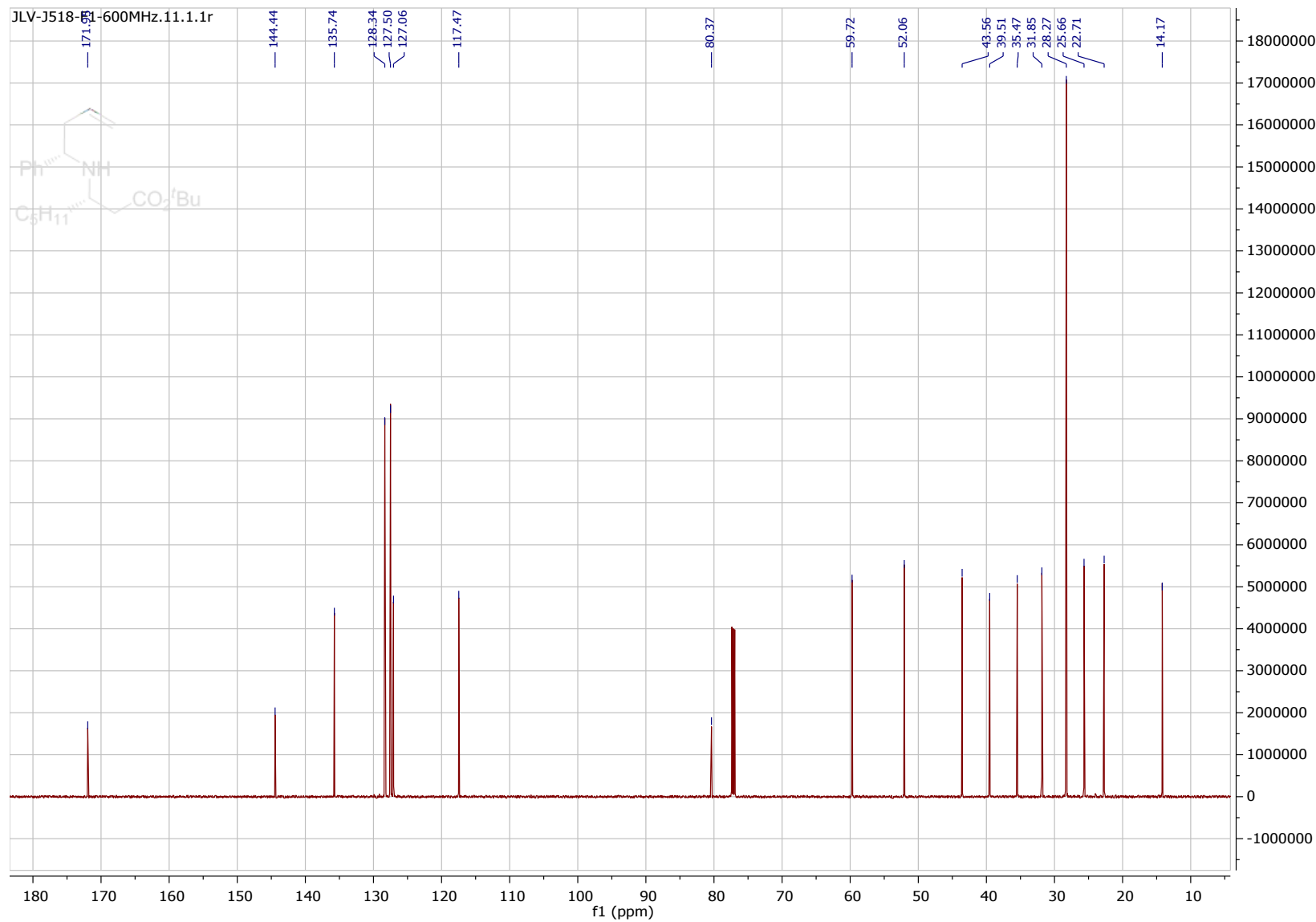
(R)-Tert-Butyl 4-methyl-3-[[[(S)-1-phenylbut-3-en-1-yl]amino]pentanoate 6s



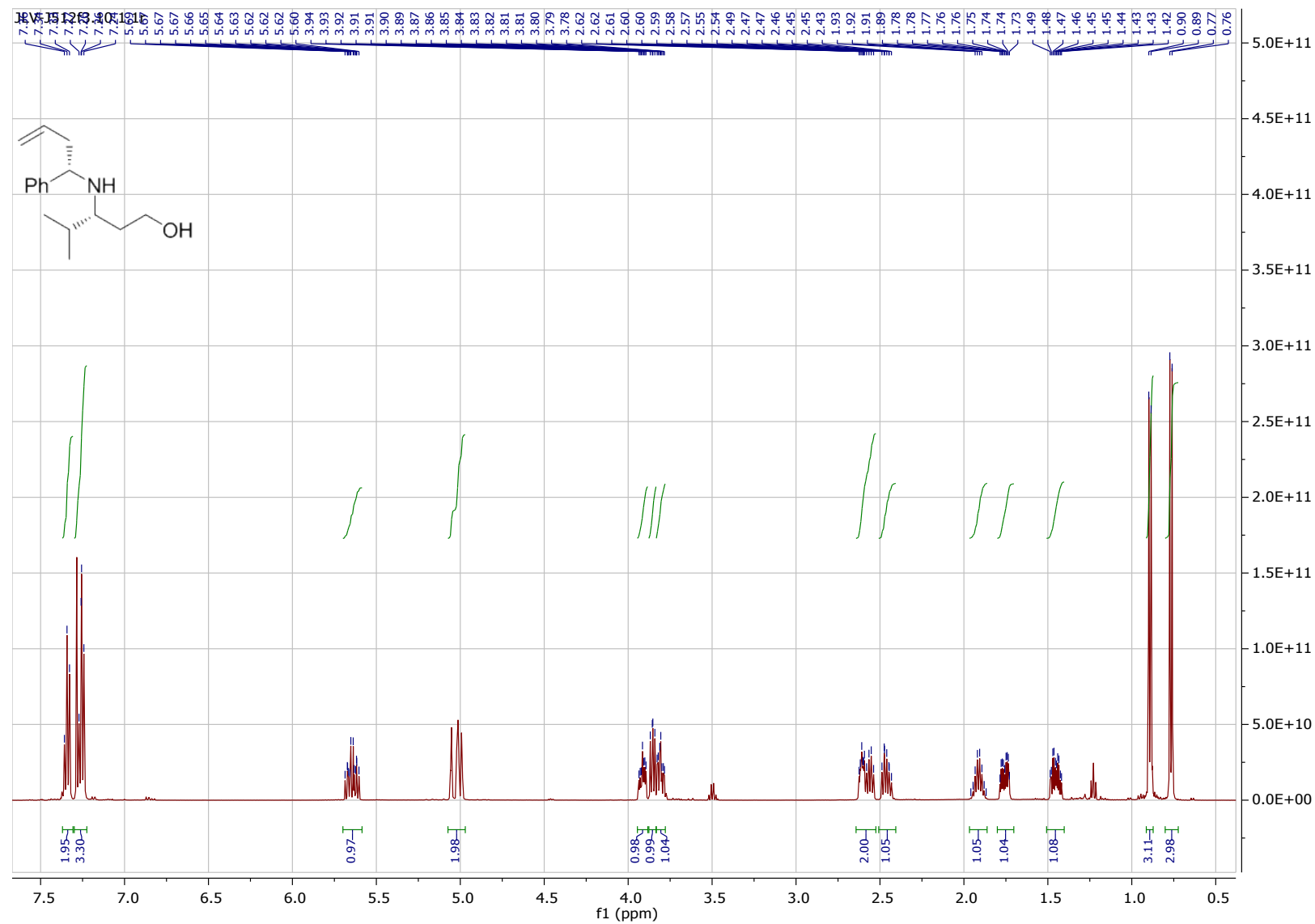


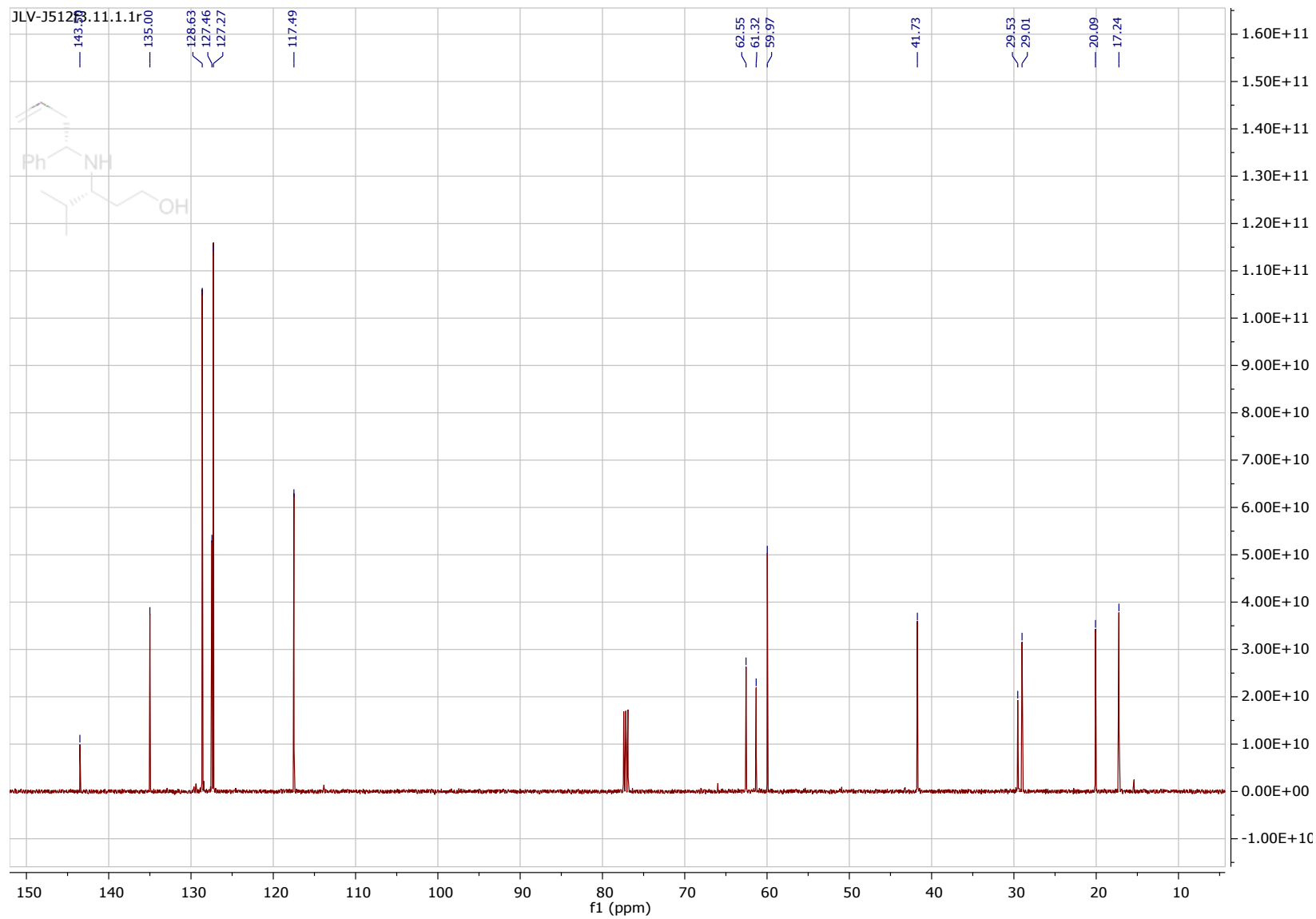
(S)-Tert-Butyl 3-[[*(S)*-1-phenylbut-3-en-1-yl]amino]octanoate 6t



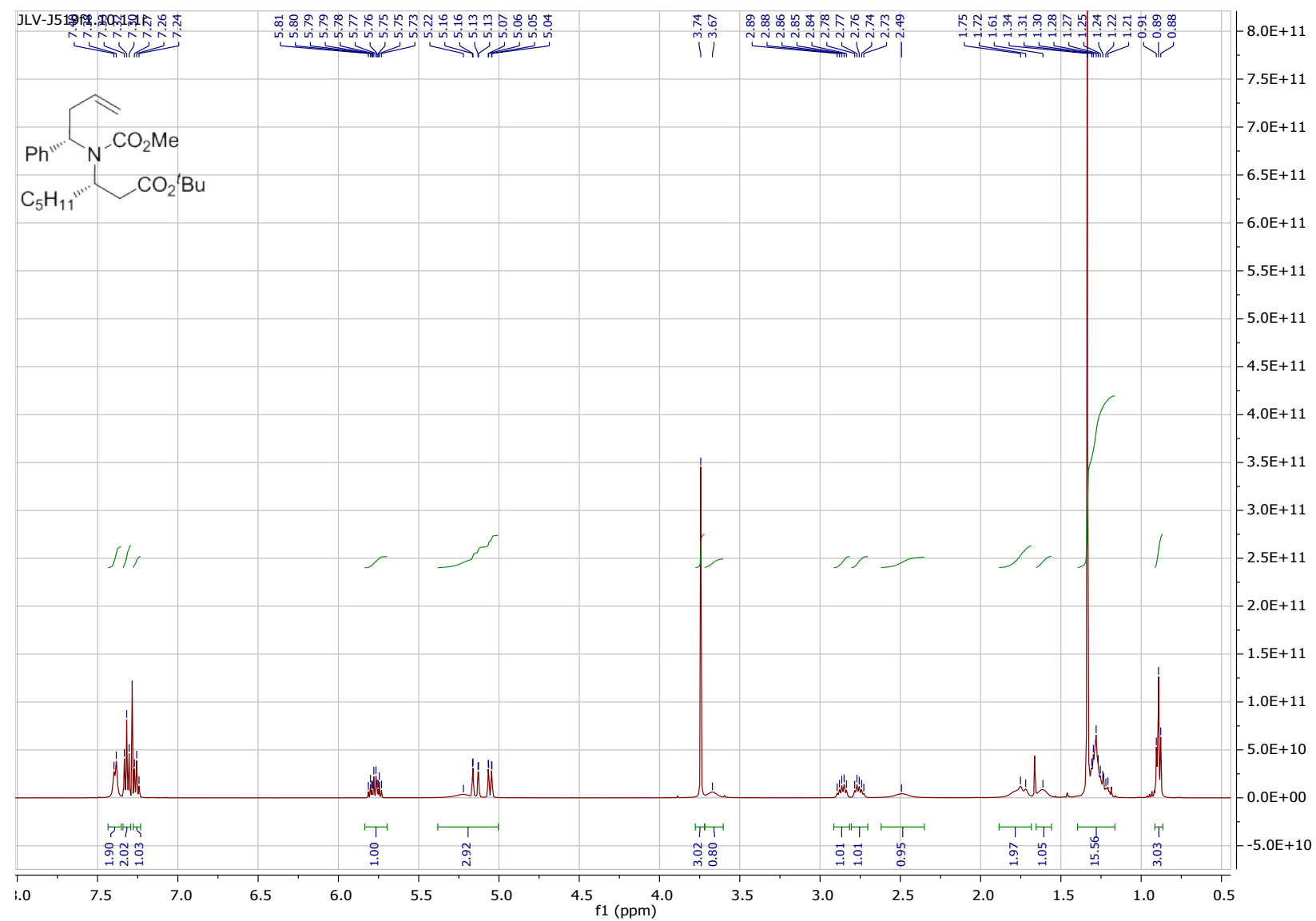


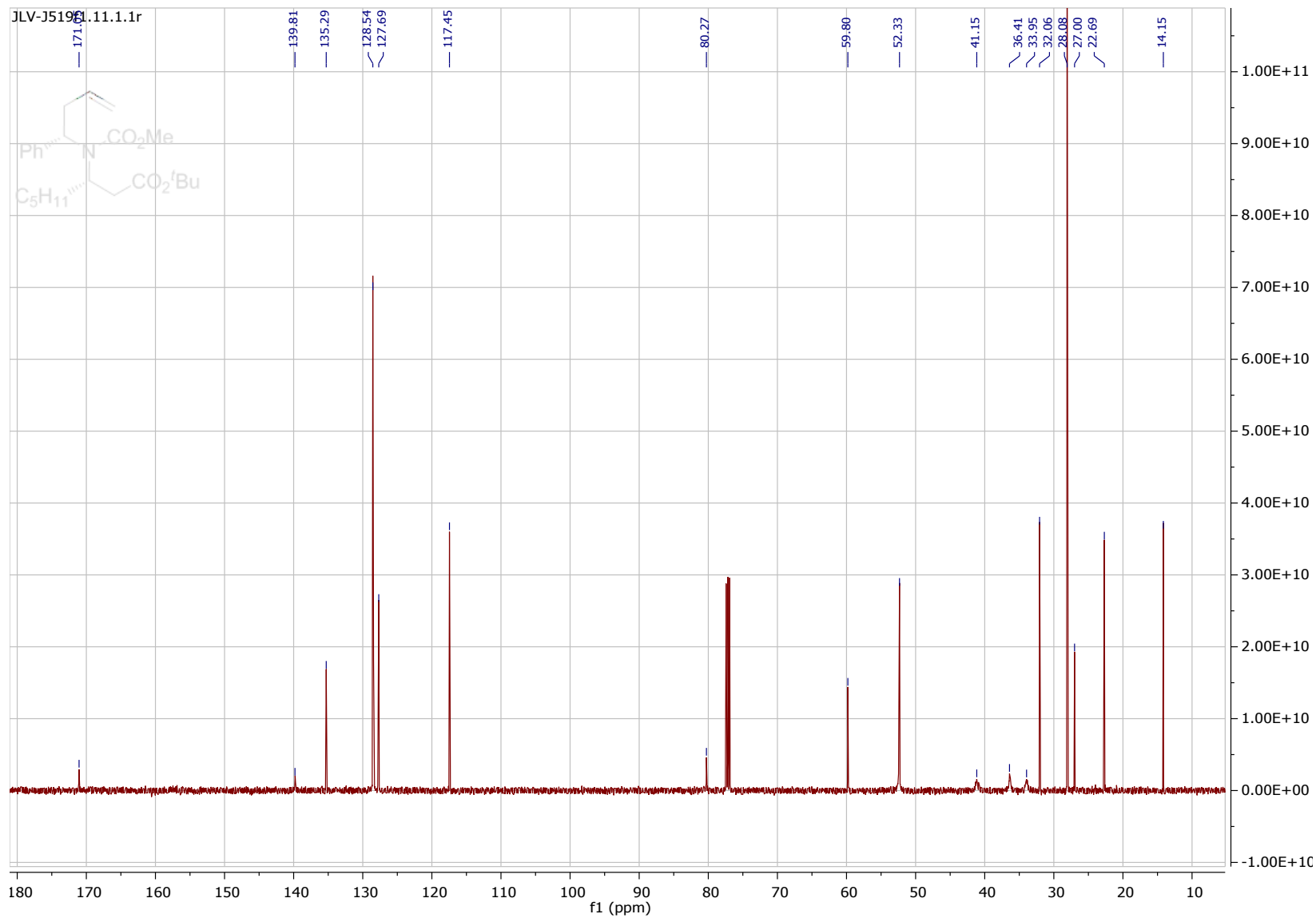
(R)-4-Methyl-3-[[[(S)-1-phenylbut-3-en-1-yl]amino]pentan-1-ol 7



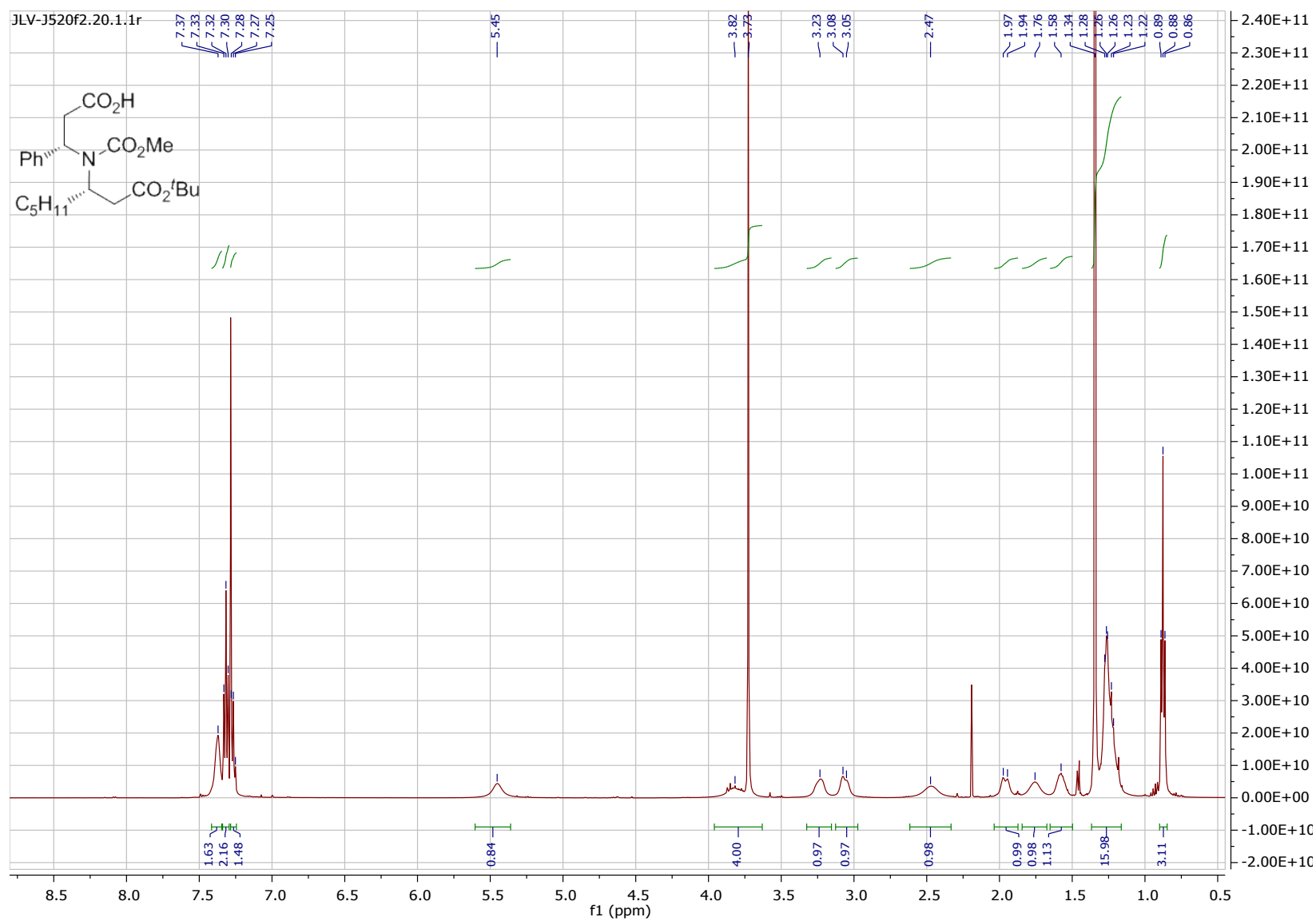


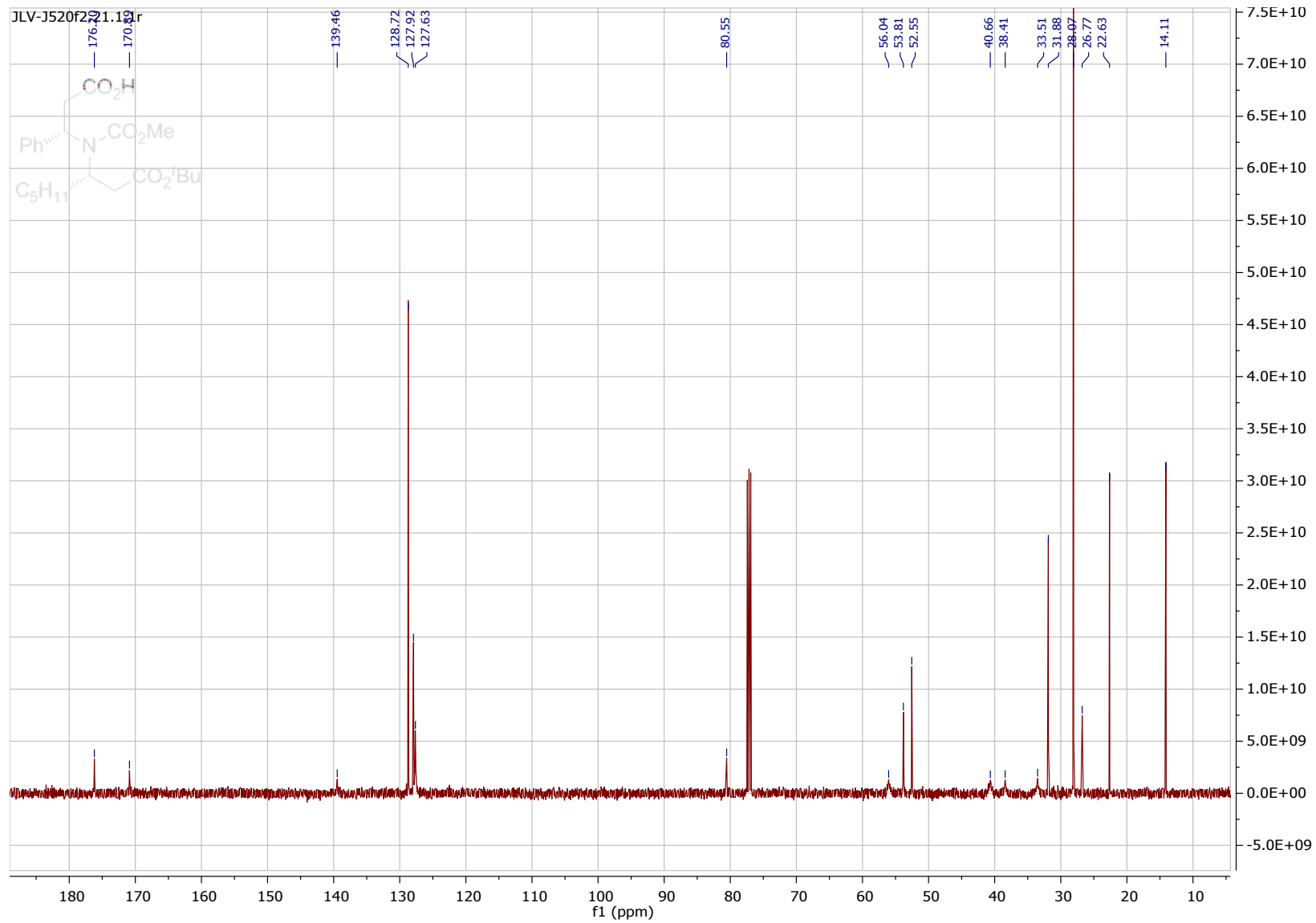
(S)-Tert-Butyl 3-((methoxycarbonyl)[(S)-1-phenylbut-3-en-1-yl]amino)octanoate 8



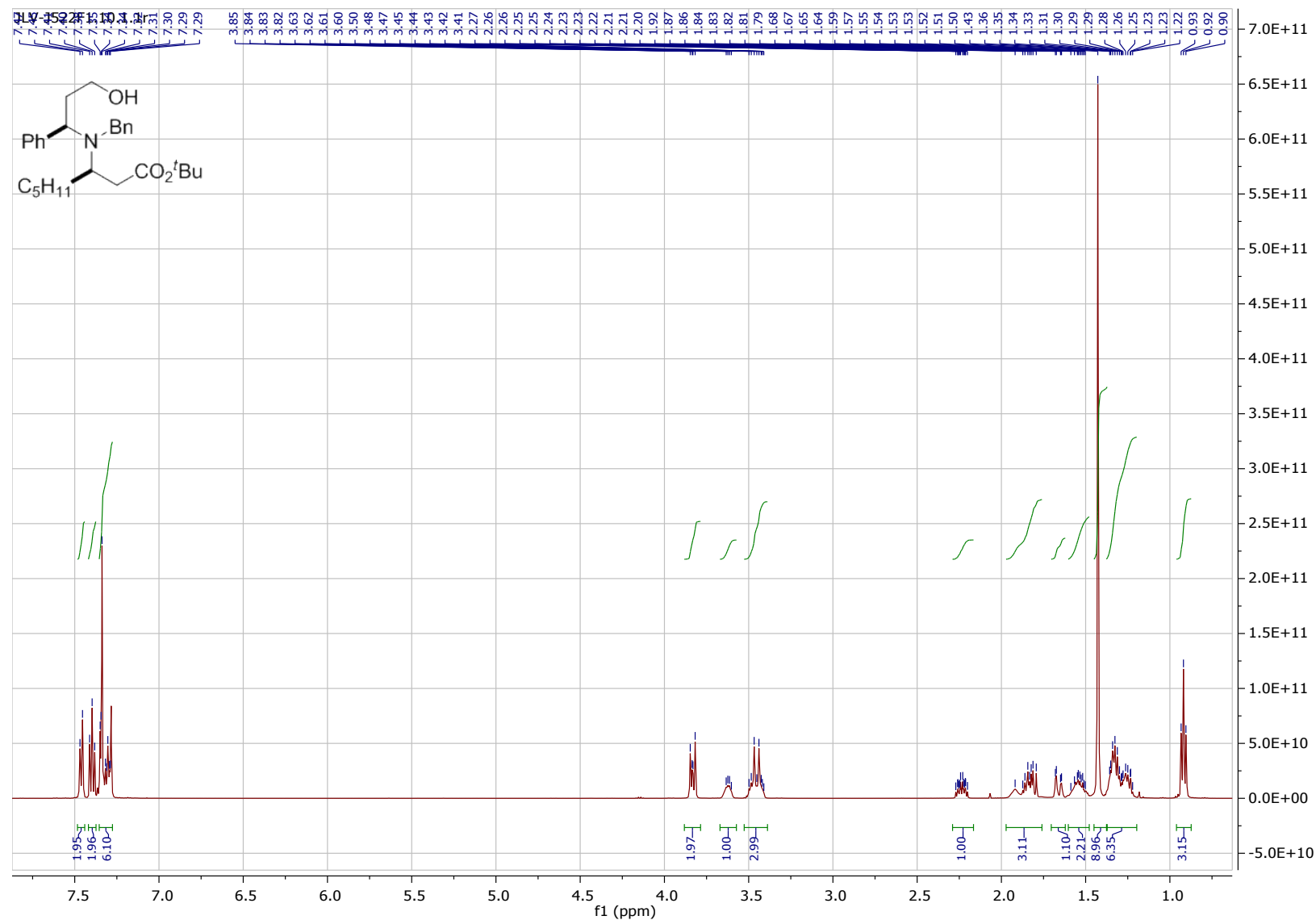


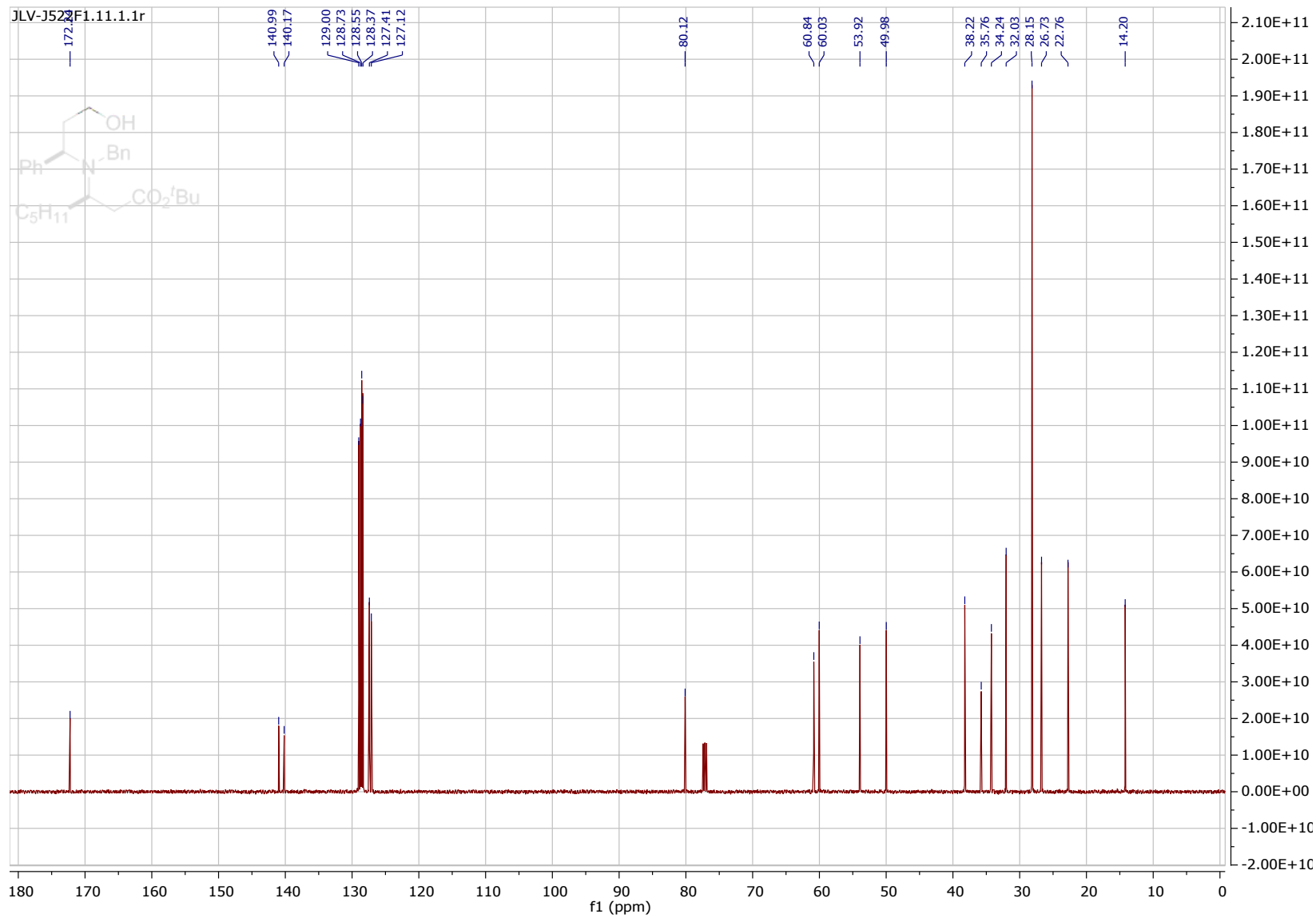
(S)-3-[[[(S)-1-(*Tert*-Butoxy)-1-oxooctan-3-yl](methoxycarbonyl)amino]-3-phenylpropanoic acid 9



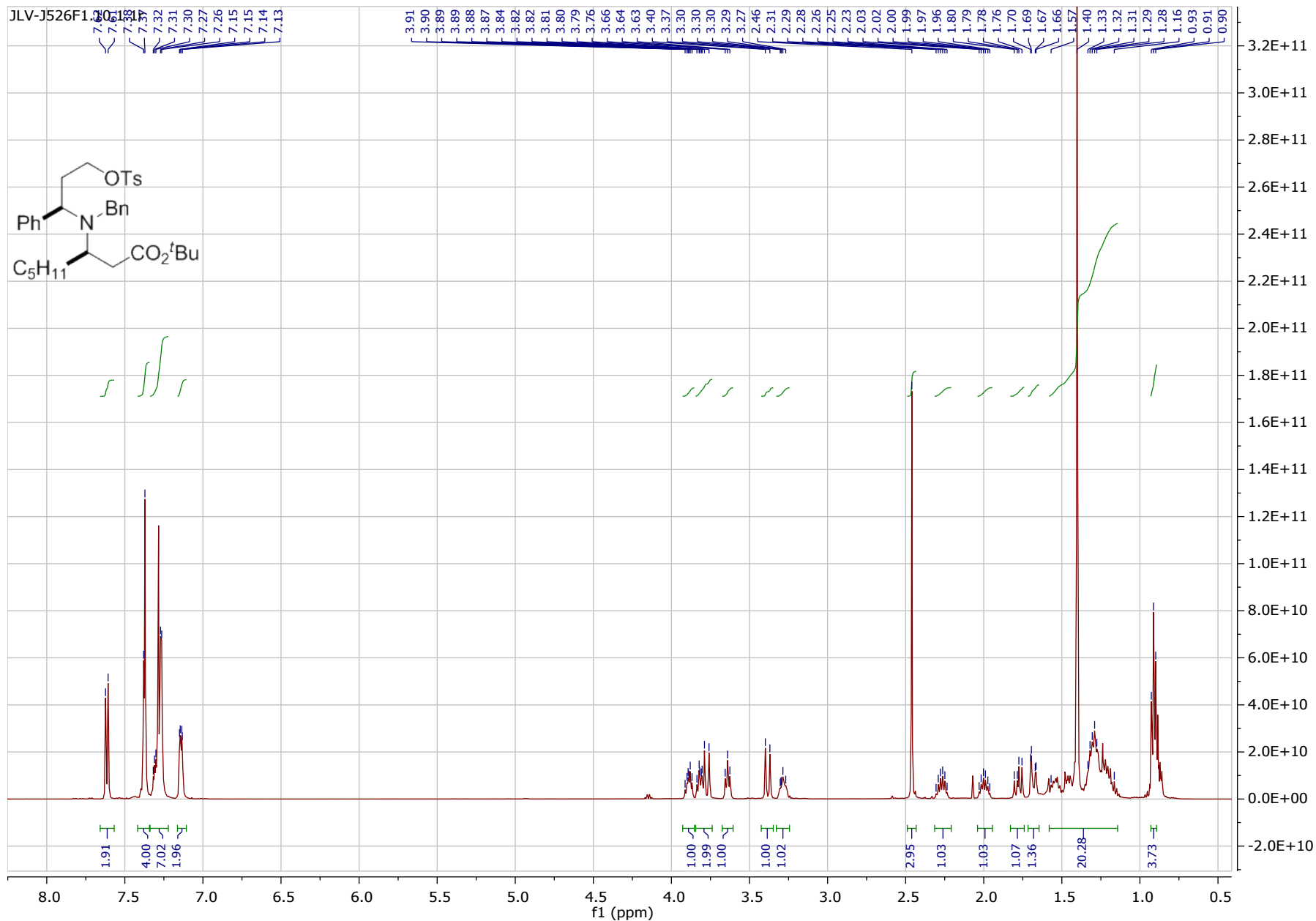


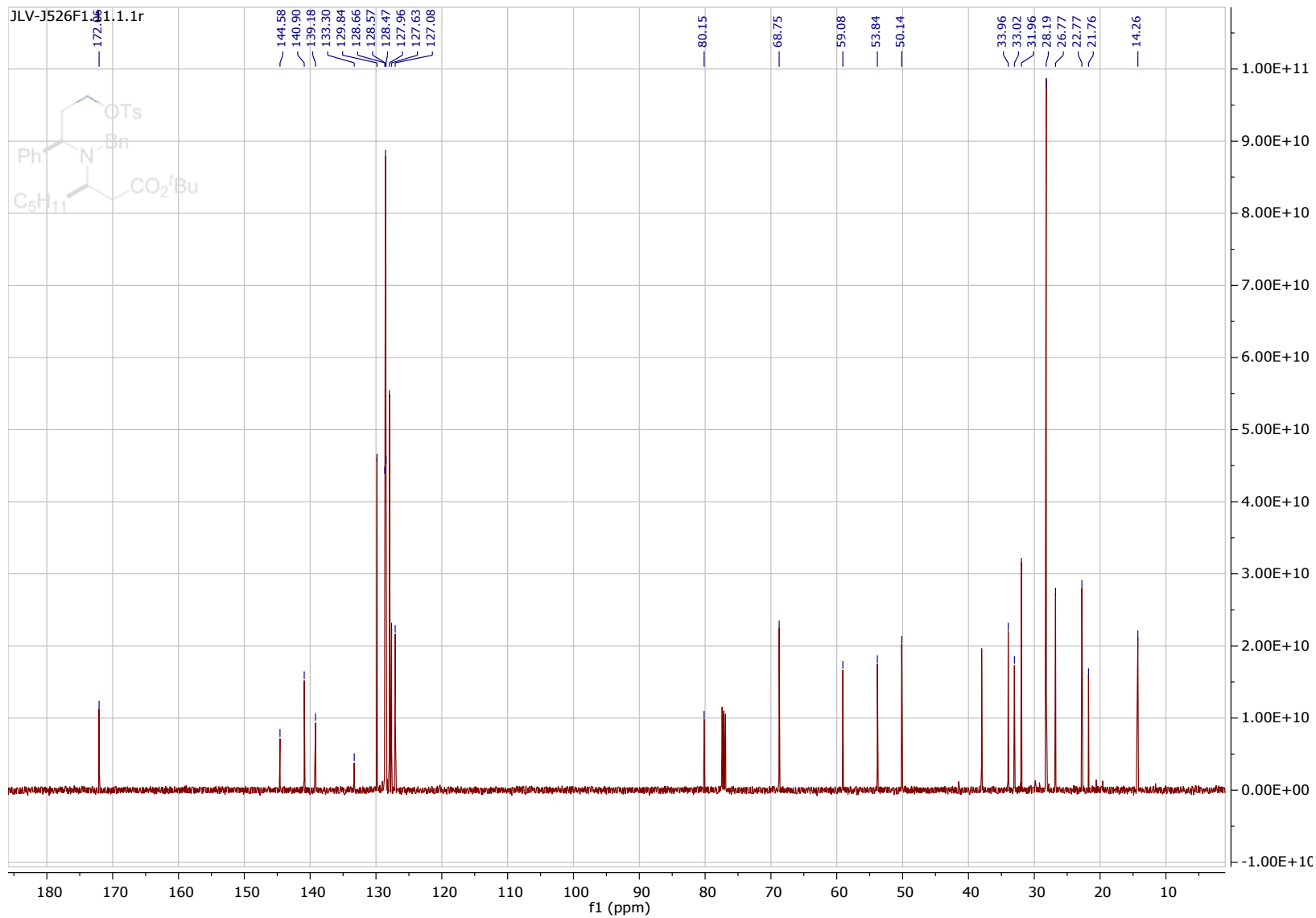
(±)-(R)-Tert-Butyl 3-{benzyl[(R)-3-hydroxy-1-phenylpropyl]amino}octanoate 10



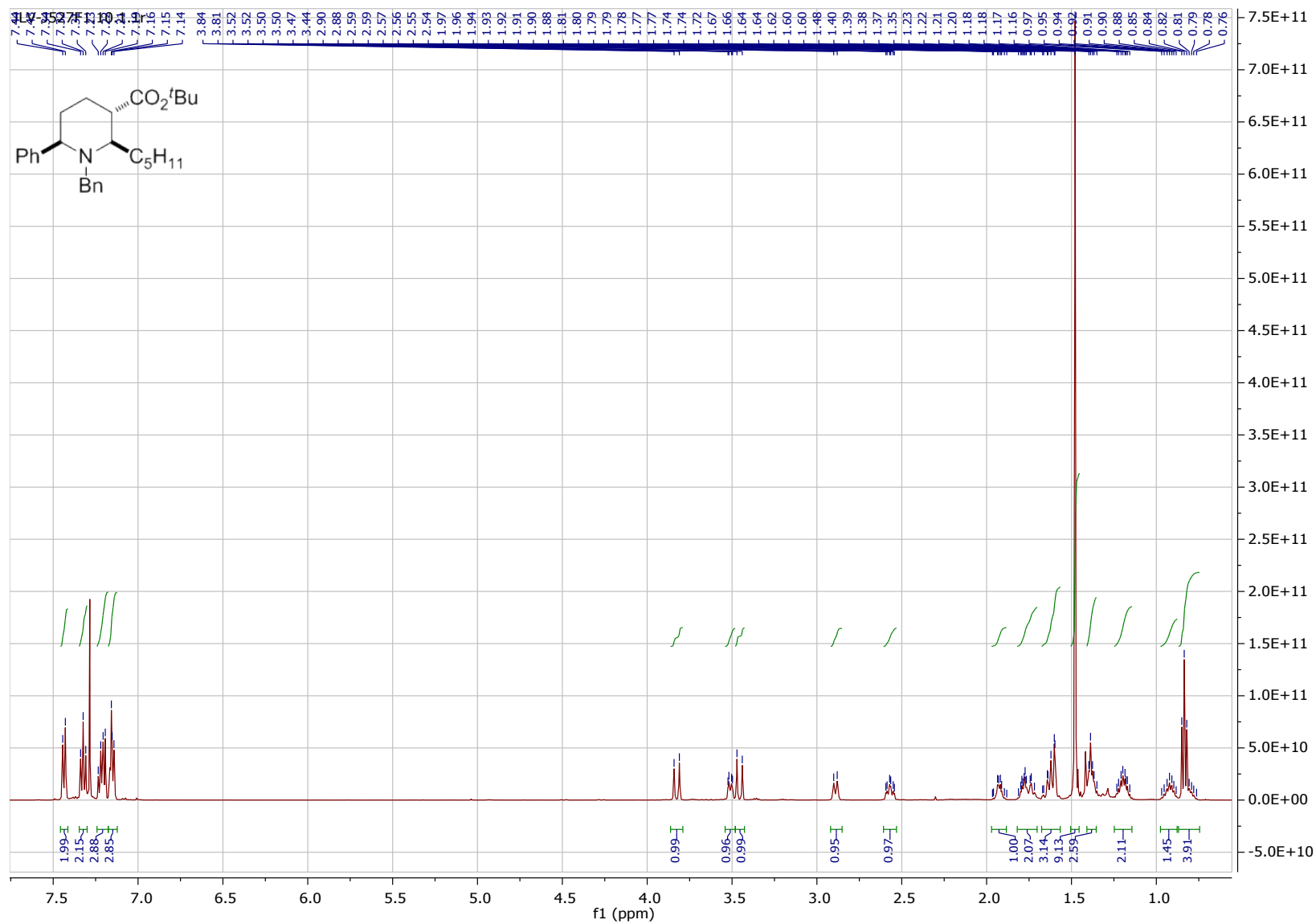


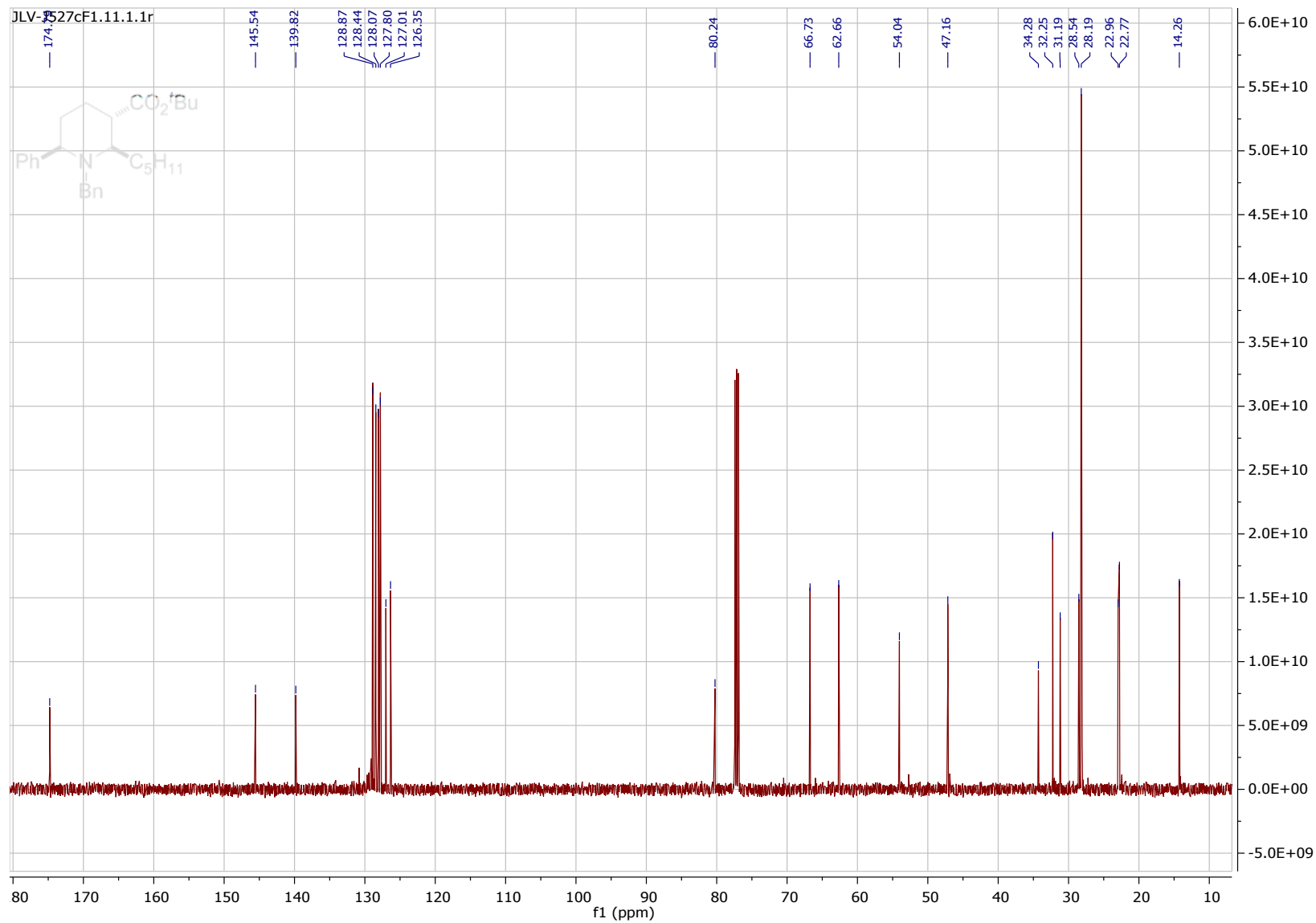
(±)-(*R*)-*Tert*-Butyl 3-{benzyl[(*R*)-1-phenyl-3-(tosyloxy)propyl]amino}octanoate



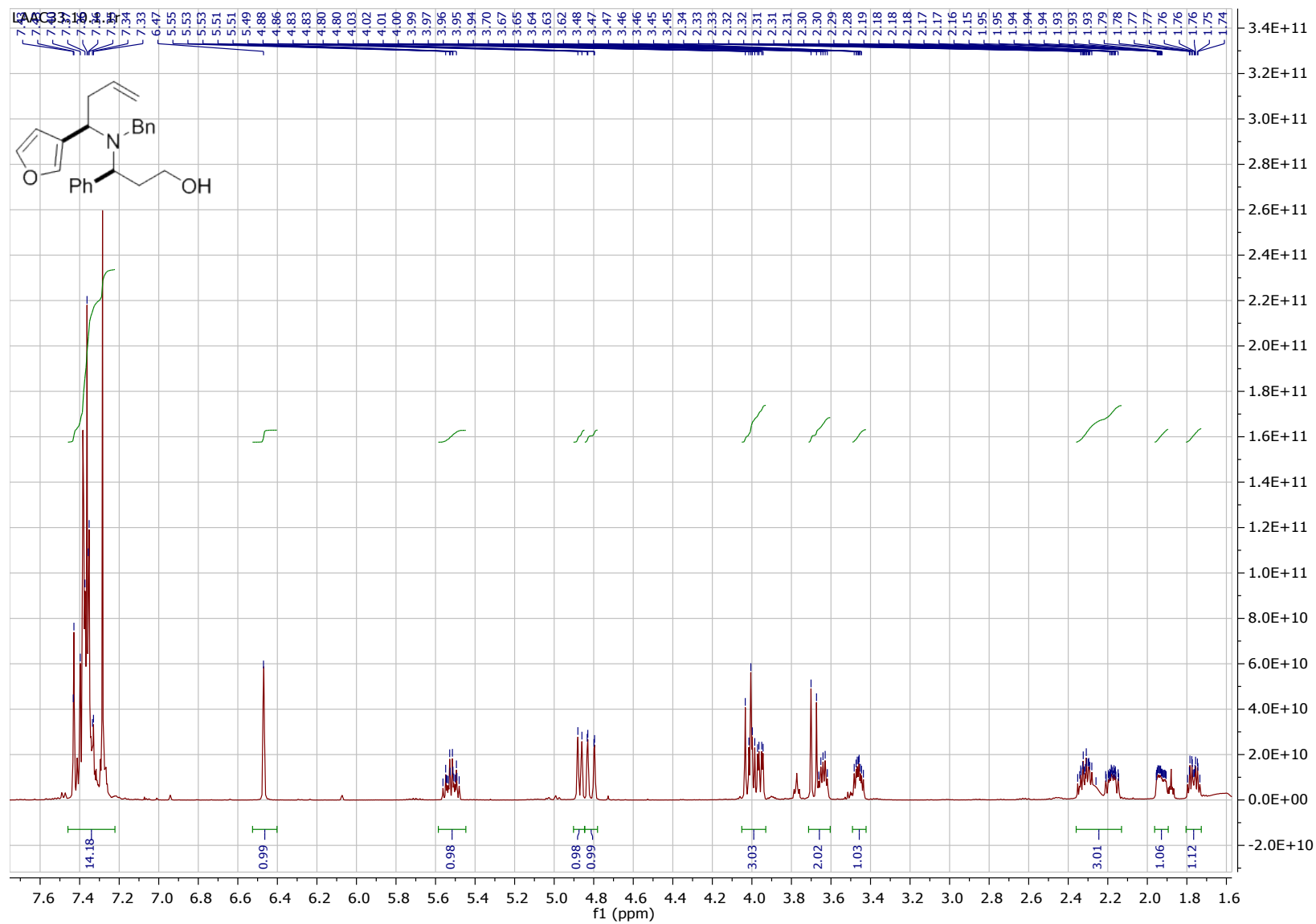


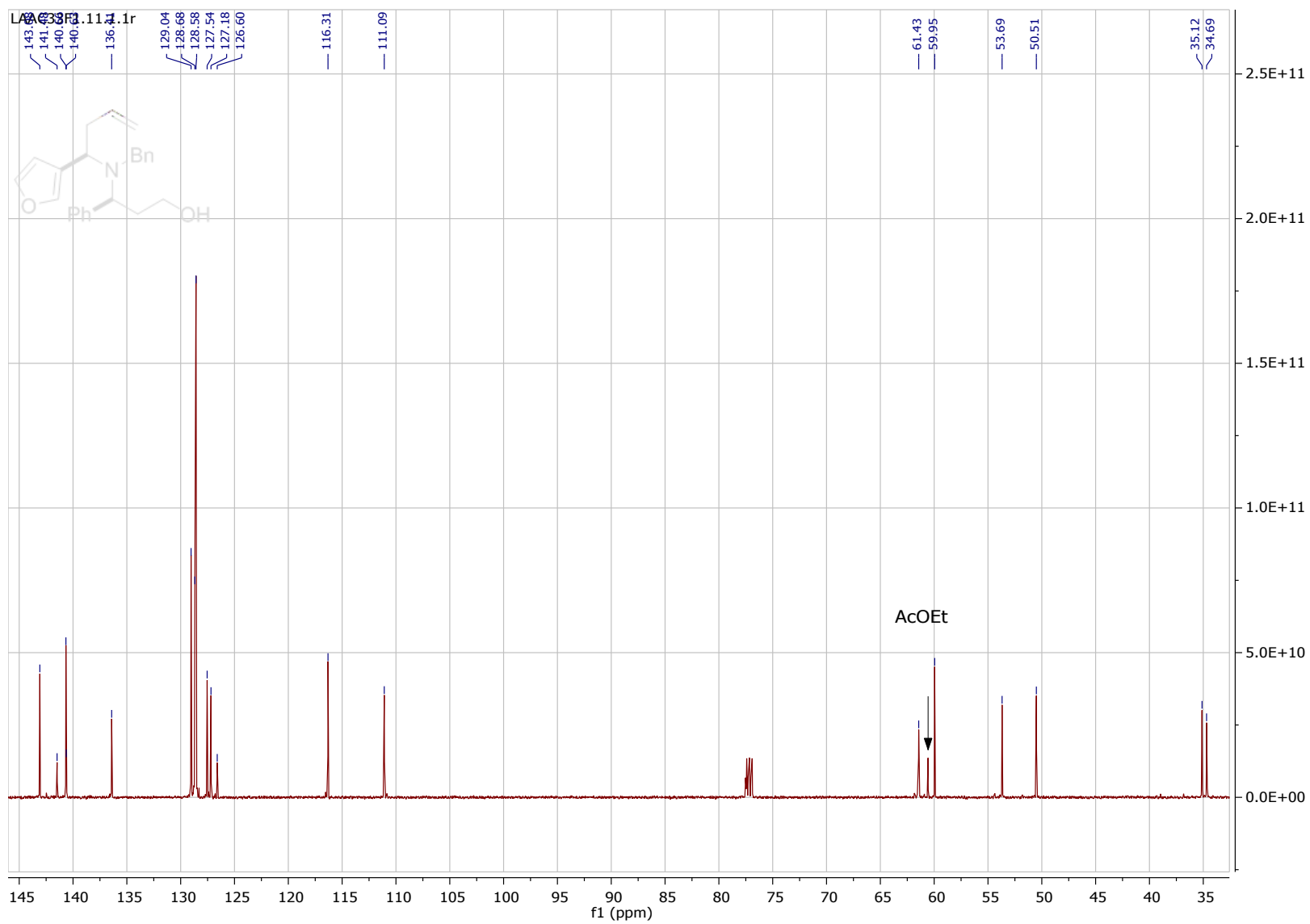
(±)-(2R,3S,6R)-Tert-Butyl 1-benzyl-2-pentyl-6-phenylpiperidine-3-carboxylate 11



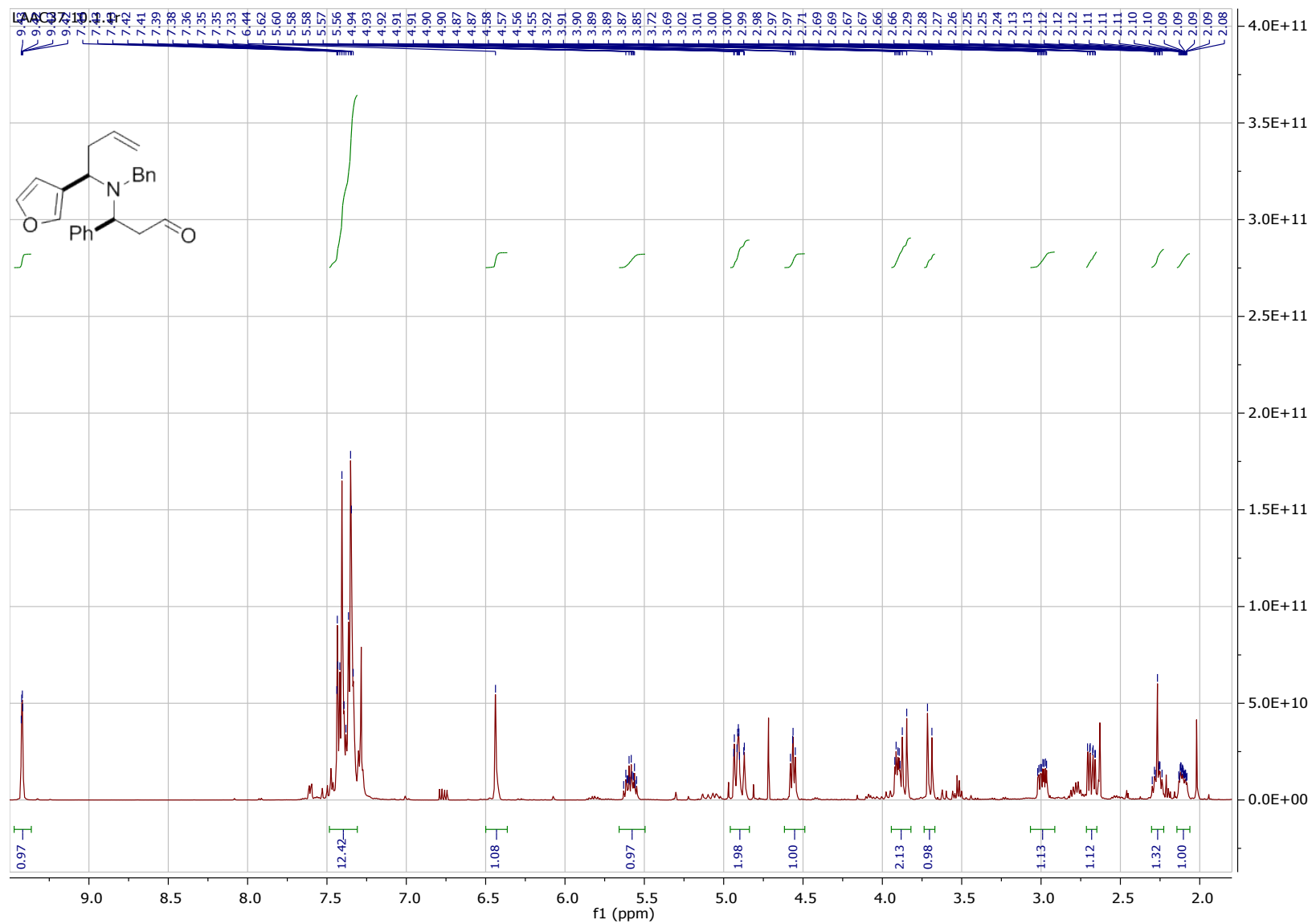


(±)-3-{Benzyl[(*R*)-1-(furan-3-yl)but-3-en-1-yl]amino}-3-phenylpropan-1-ol

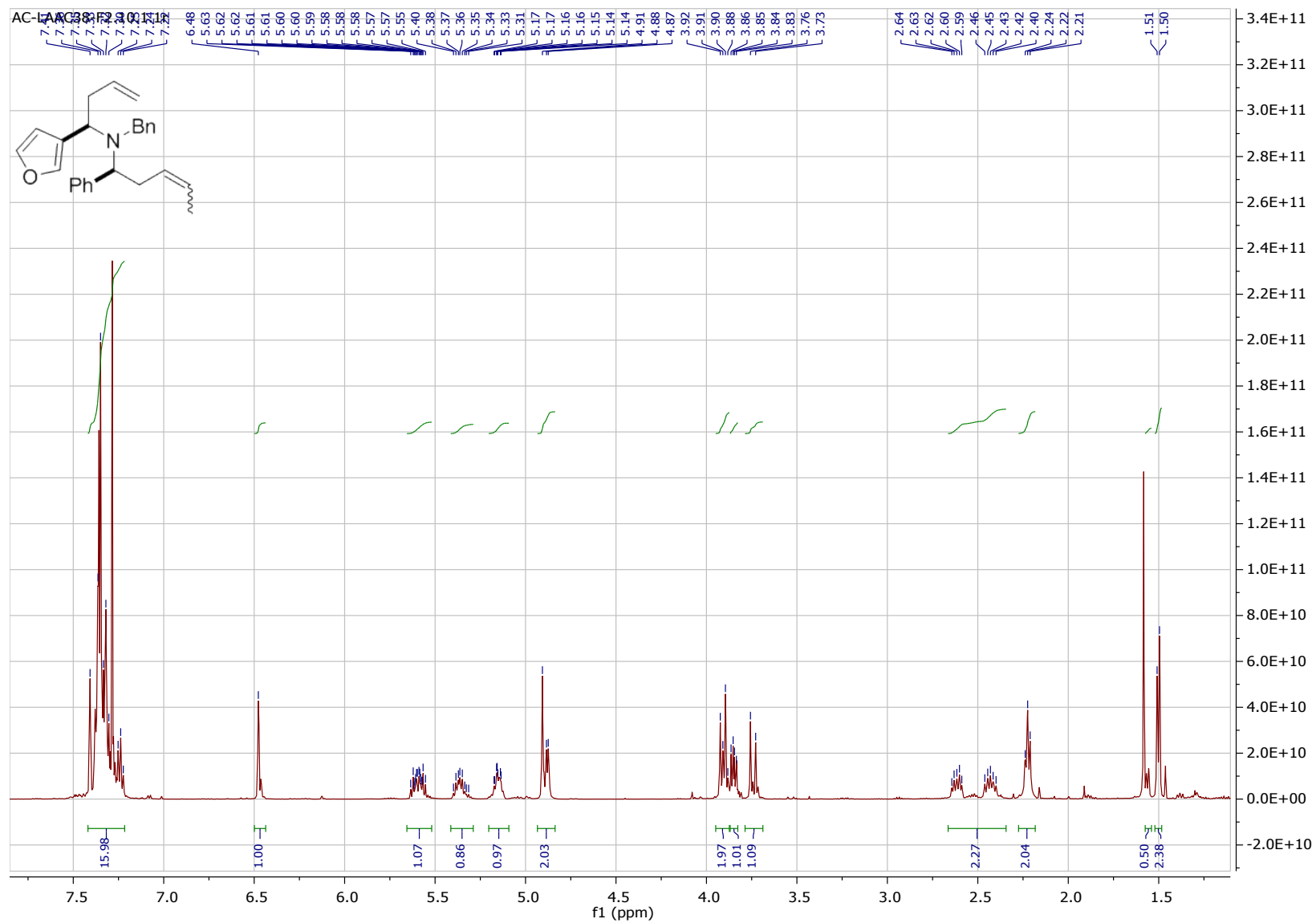


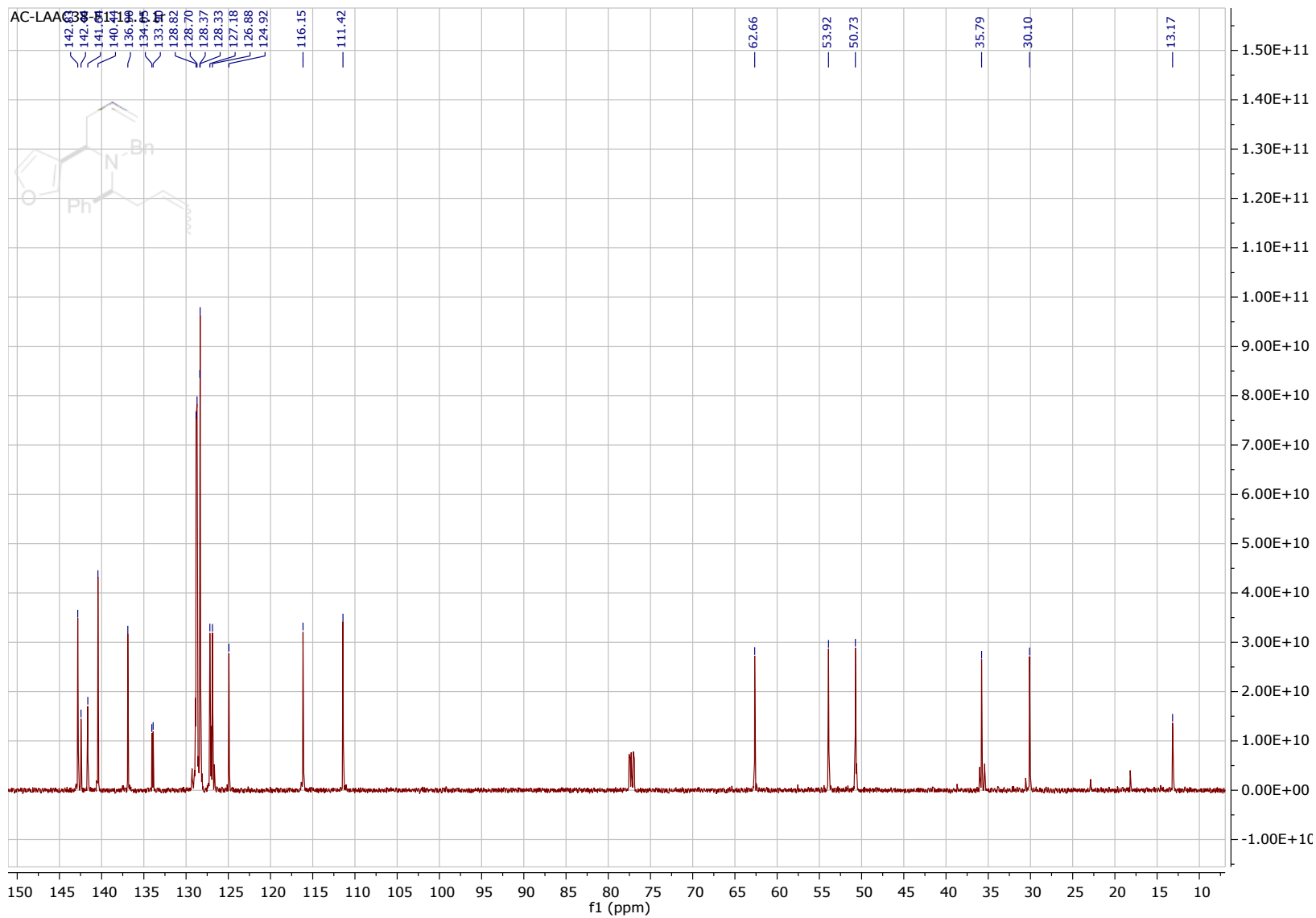


(±)-(S)-3-{Benzyl[(R)-1-(furan-3-yl)but-3-en-1-yl]amino}-3-phenylpropanal

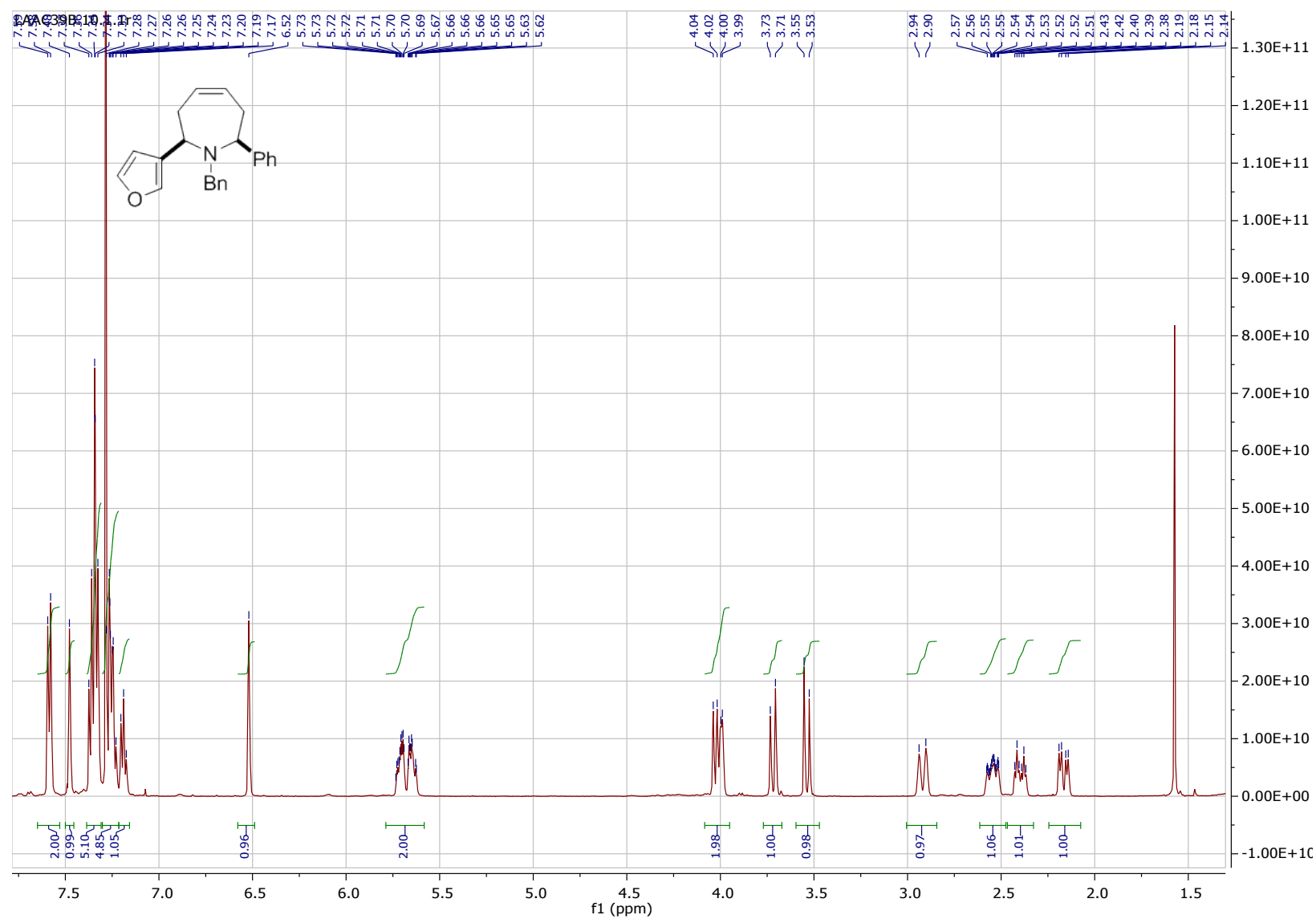


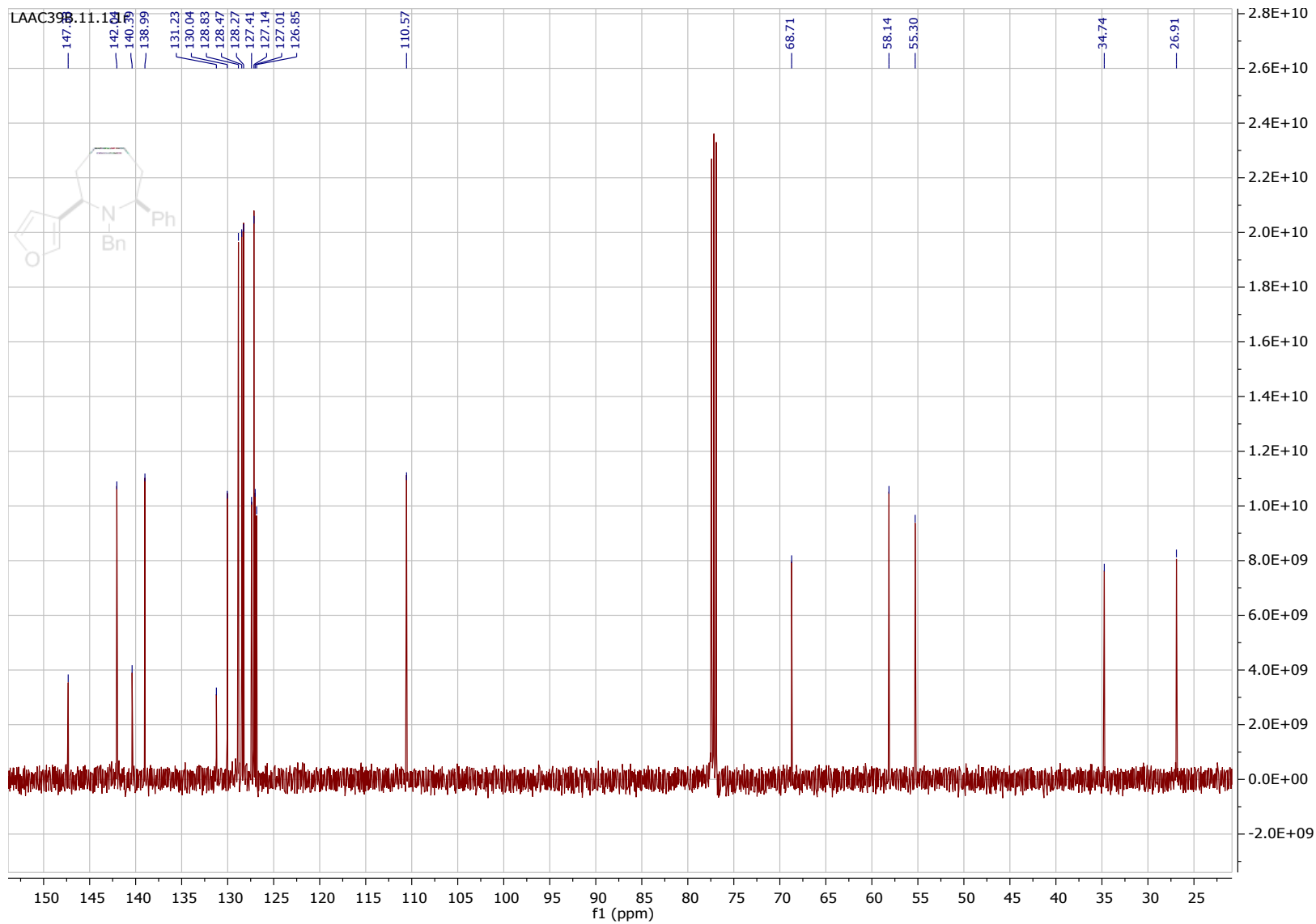
(±)-(R,Z)-N-Benzyl-N-[(R)-1-phenylbut-3-en-1-yl]dec-2-en-5-amine 12



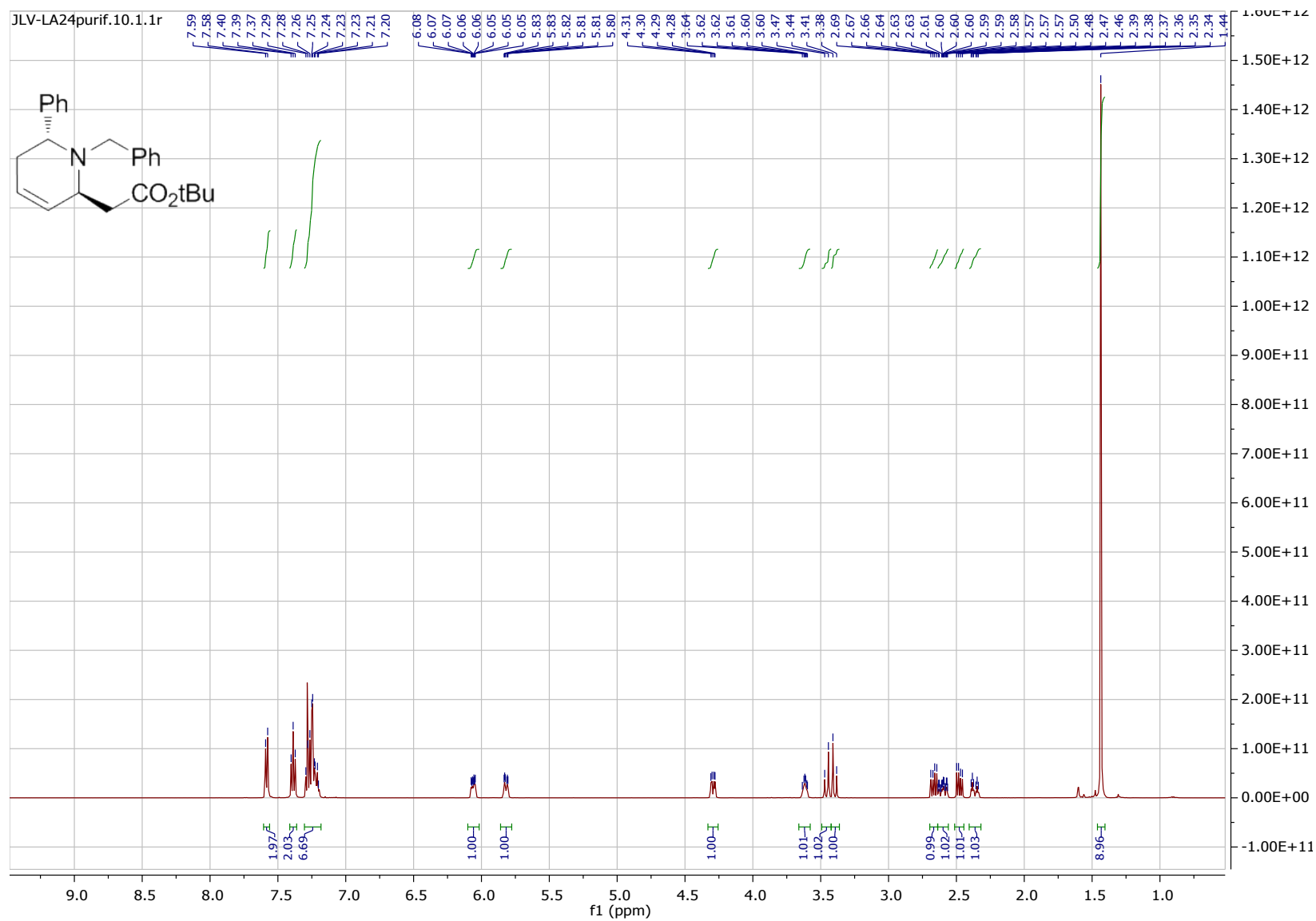


(2*R*,7*S*)-1-Benzyl-2-(furan-3-yl)-7-phenyl-2,3,6,7-tetrahydro-1*H*-azepine 13

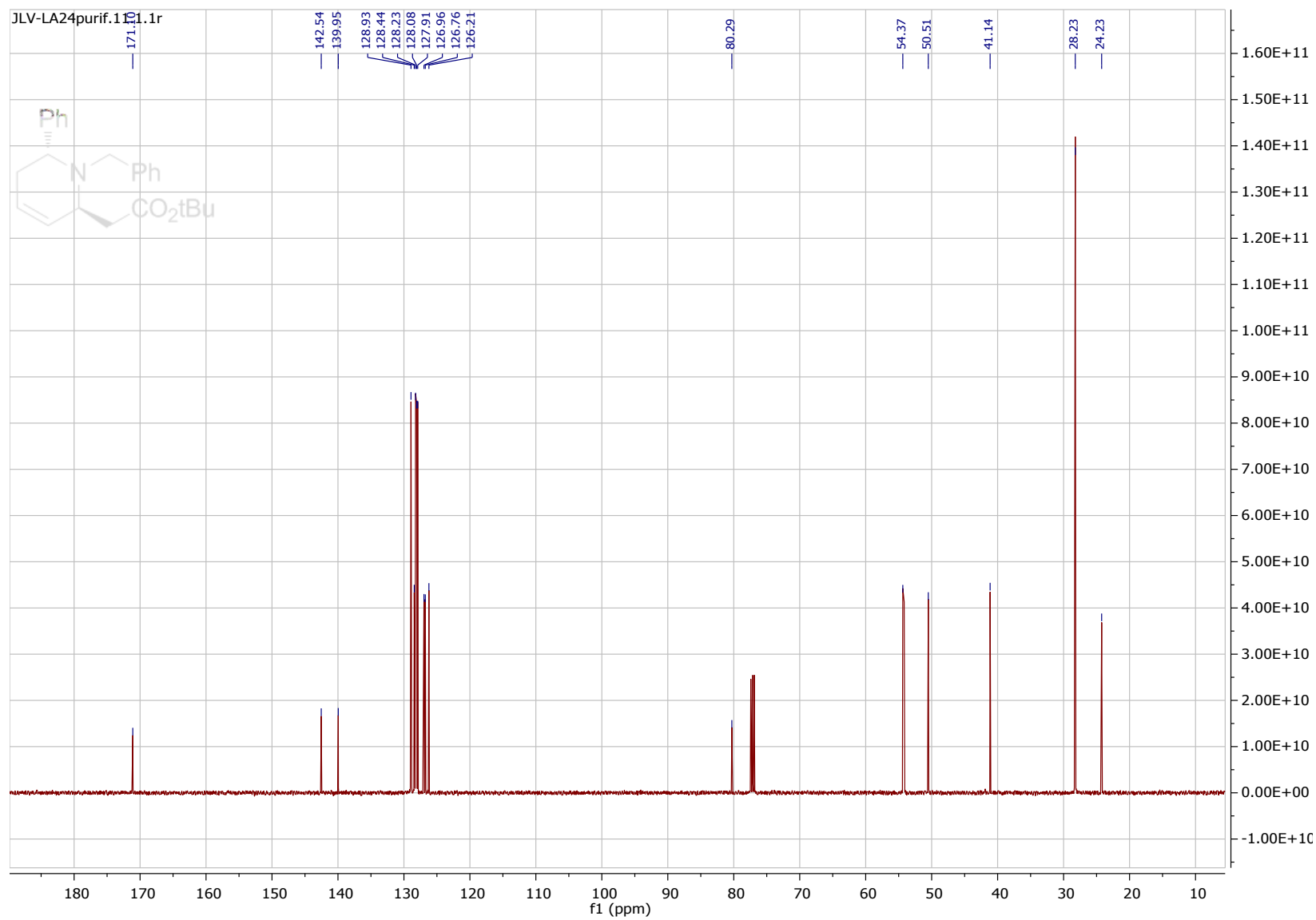




Tert-Butyl 2-[(2*S*,6*R*)-1-Benzyl-6-phenyl-1,2,5,6-tetrahydropyridin-2-yl]acetate 14



JLV-LA24purif.110.1r



Tert-Butyl 2-[(2R,6R)-6-phenylpiperidin-2-yl]acetate

