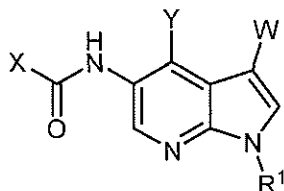


Claims

1. A compound of Formula I:



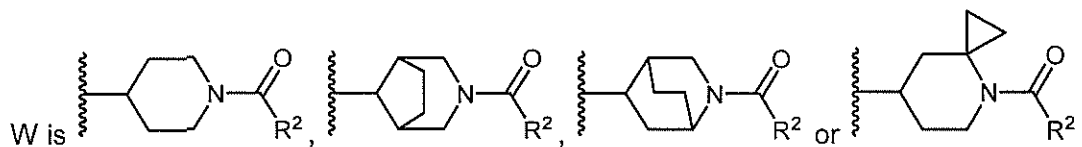
I

or a pharmaceutically acceptable salt, pharmaceutically active metabolite, pharmaceutically acceptable prodrug, or pharmaceutically acceptable solvate thereof, wherein,

Y is -CF₃;

X is phenyl optionally substituted with one, two, three, four or five substituents independently selected from the group consisting of -CH₃, -CH₂CH₃, -CH₂OH, -OH, -OCH₃, -SCH₃, -OCH₂CH₃, -OCH₂CH₂OH, -OCH₂CH₂OCH₃, -F, -Cl, -Br and -CN;

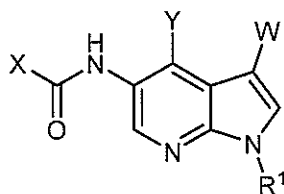
R¹ is -CH₃ or -CH₂CH₃;



each optionally substituted with one, two, three, four or five -CH₃; and

R² is (C₁-C₆)alkyl, (C₃-C₁₀)cycloalkyl, phenyl, tetrahydrothiophenyl, thietanyl or indanyl, optionally substituted with one, two, three, four or five substituents independently selected for each occurrence from the group consisting of -F, -Cl, -Br, -OH, (C₁-C₃)alkyl, (C₁-C₃)haloalkyl and (C₃-C₁₀)cycloalkyl.

2. A compound of Formula I:



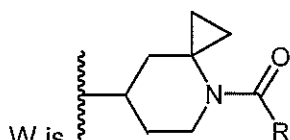
I

or a pharmaceutically acceptable salt, pharmaceutically active metabolite, pharmaceutically acceptable prodrug, or pharmaceutically acceptable solvate thereof, wherein,

Y is -CH₃ or -CF₃;

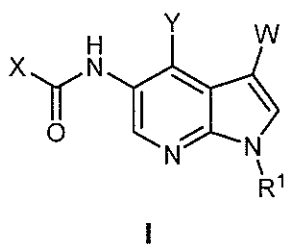
X is phenyl optionally substituted with one, two, three, four or five substituents independently selected from the group consisting of -CH₃, -CH₂CH₃, -CH₂OH, -OH, -OCH₃, -SCH₃, -OCH₂CH₃, -OCH₂CH₂OH, -OCH₂CH₂OCH₃, -F, -Cl, -Br and -CN;

R¹ is -CH₃ or -CH₂CH₃;



R² is (C₁-C₆)alkyl, (C₃-C₁₀)cycloalkyl, phenyl, tetrahydrothiophenyl, thietanyl or indanyl, optionally substituted with one, two, three, four or five substituents independently selected for each occurrence from the group consisting of -F, -Cl, -Br, -OH, (C₁-C₃)alkyl, (C₁-C₃)haloalkyl and (C₃-C₁₀)cycloalkyl.

3. A compound of Formula I:

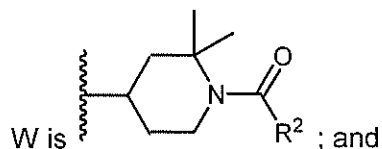


or a pharmaceutically acceptable salt, pharmaceutically active metabolite, pharmaceutically acceptable prodrug, or pharmaceutically acceptable solvate thereof, wherein,

Y is -CH₃ or -CF₃;

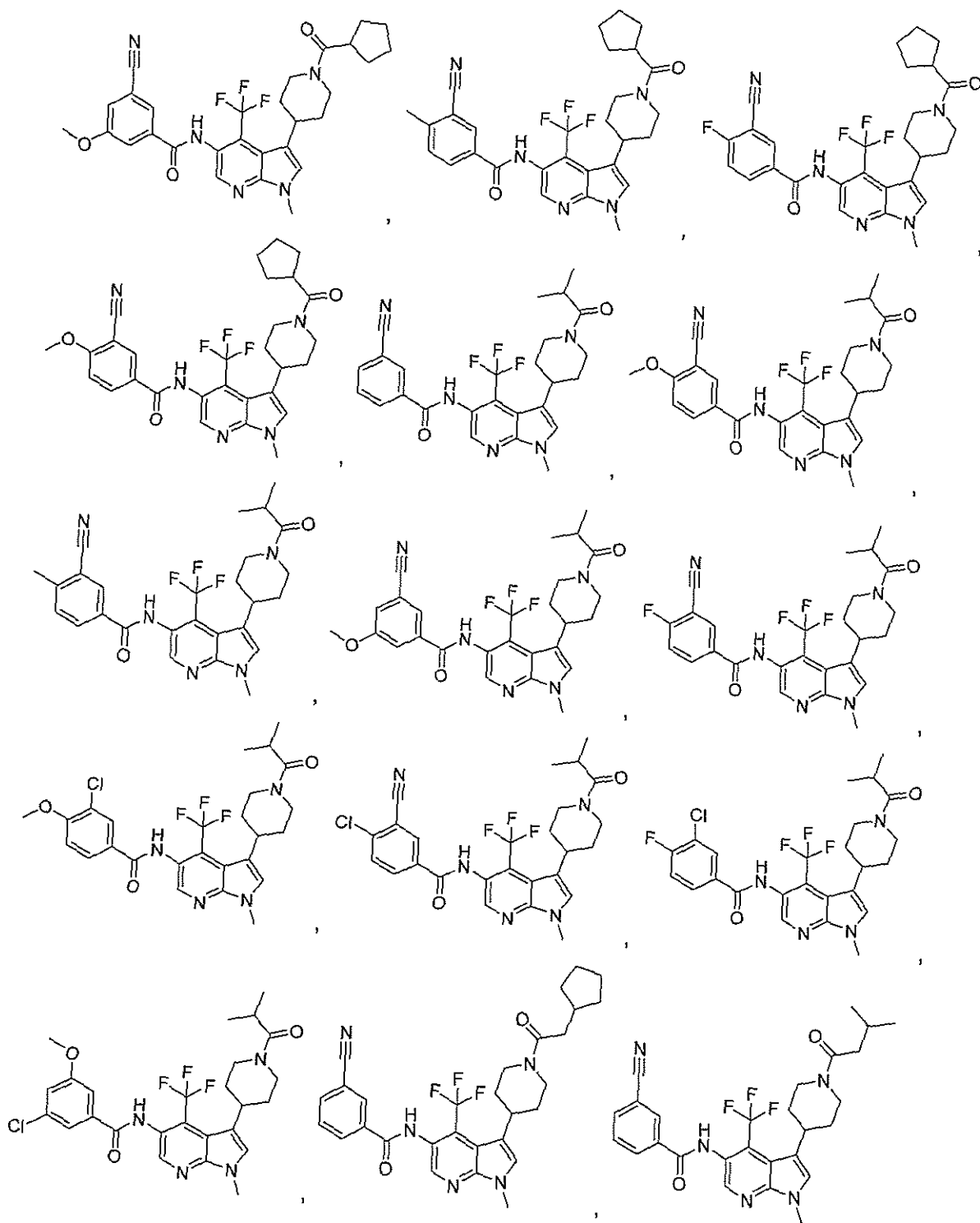
X is phenyl optionally substituted with one, two, three, four or five substituents independently selected from the group consisting of -CH₃, -CH₂CH₃, -CH₂OH, -OH, -OCH₃, -SCH₃, -OCH₂CH₃, -OCH₂CH₂OH, -OCH₂CH₂OCH₃, -F, -Cl, -Br and -CN;

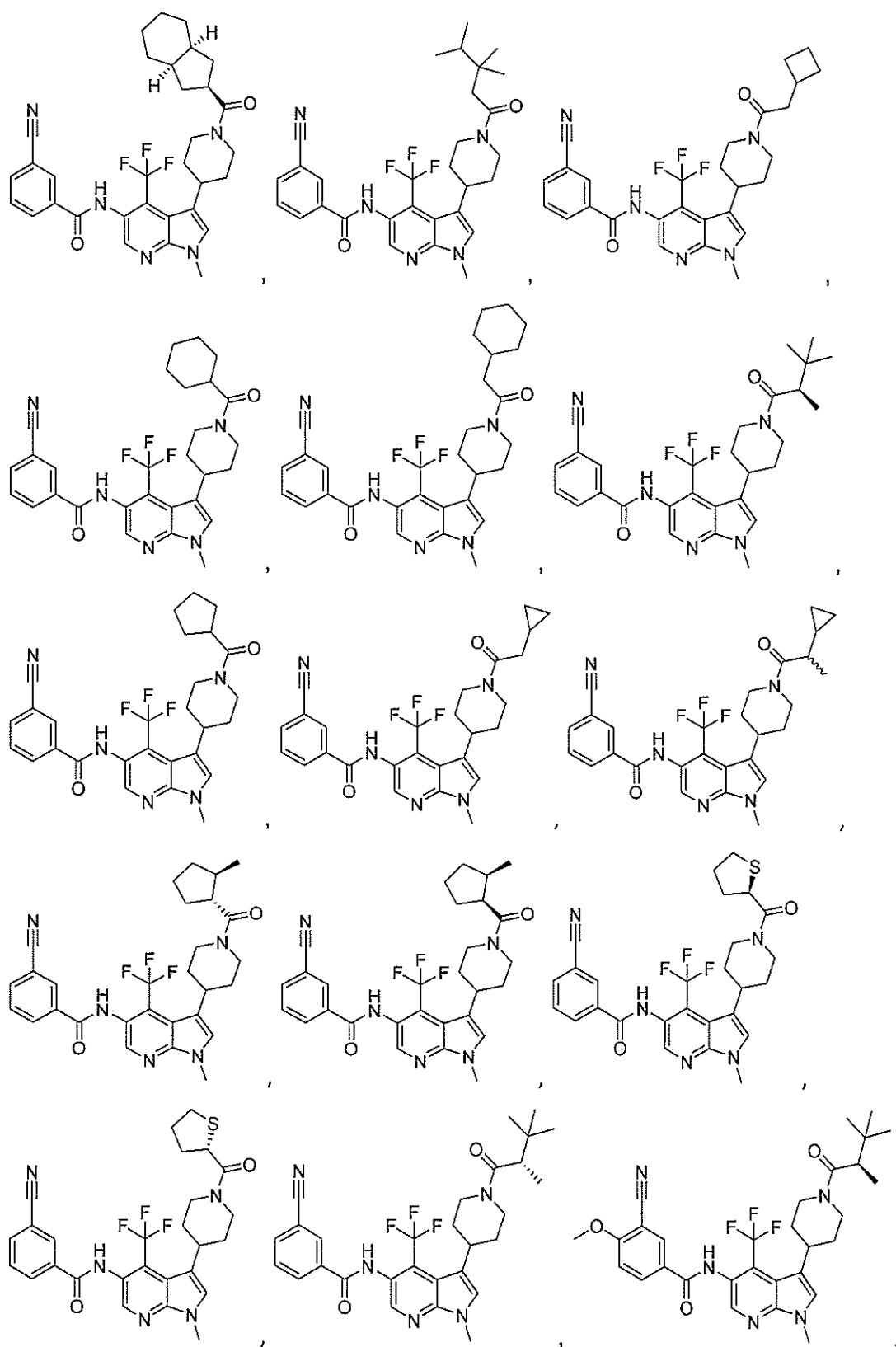
R¹ is -CH₃ or -CH₂CH₃;

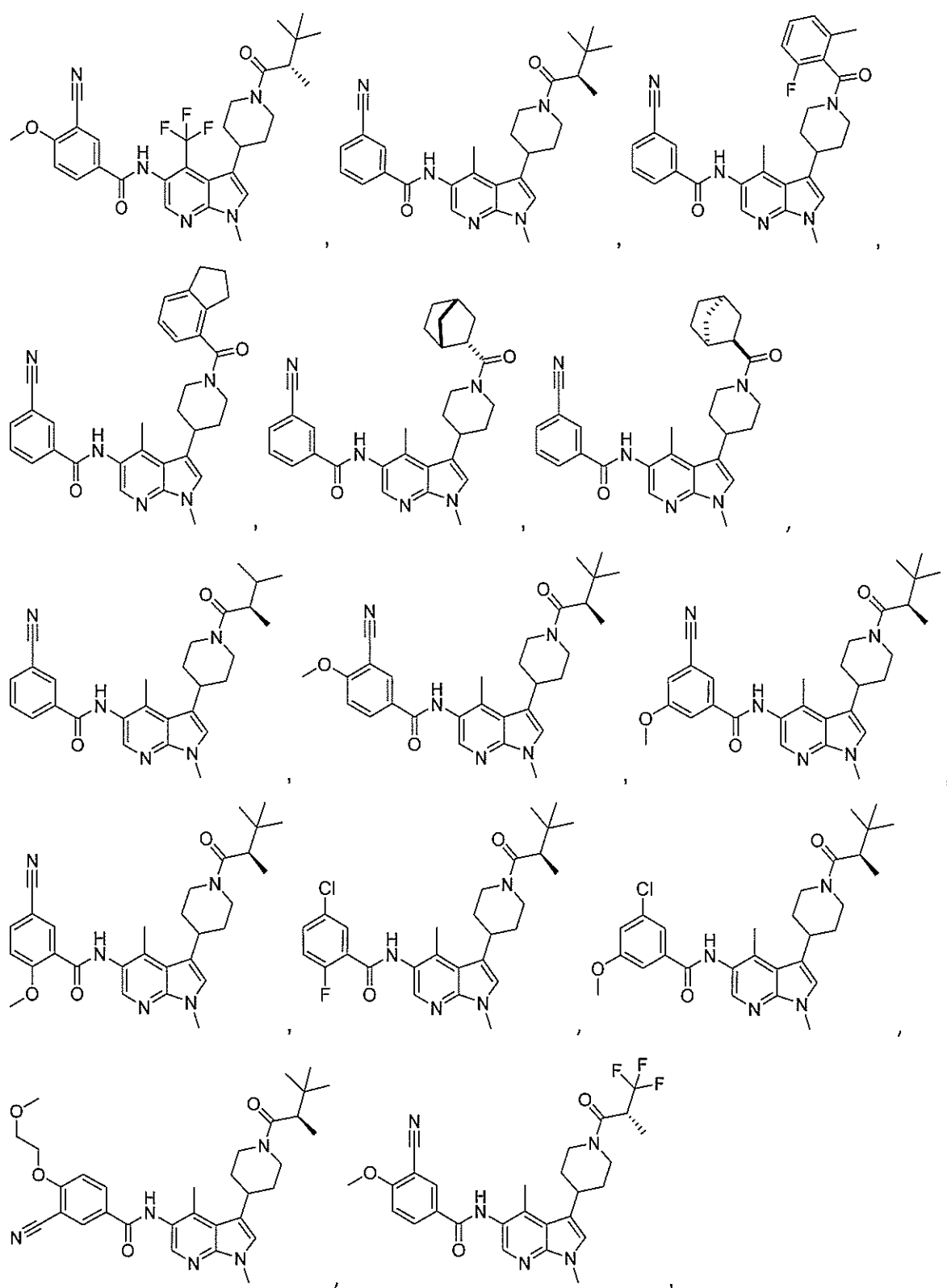


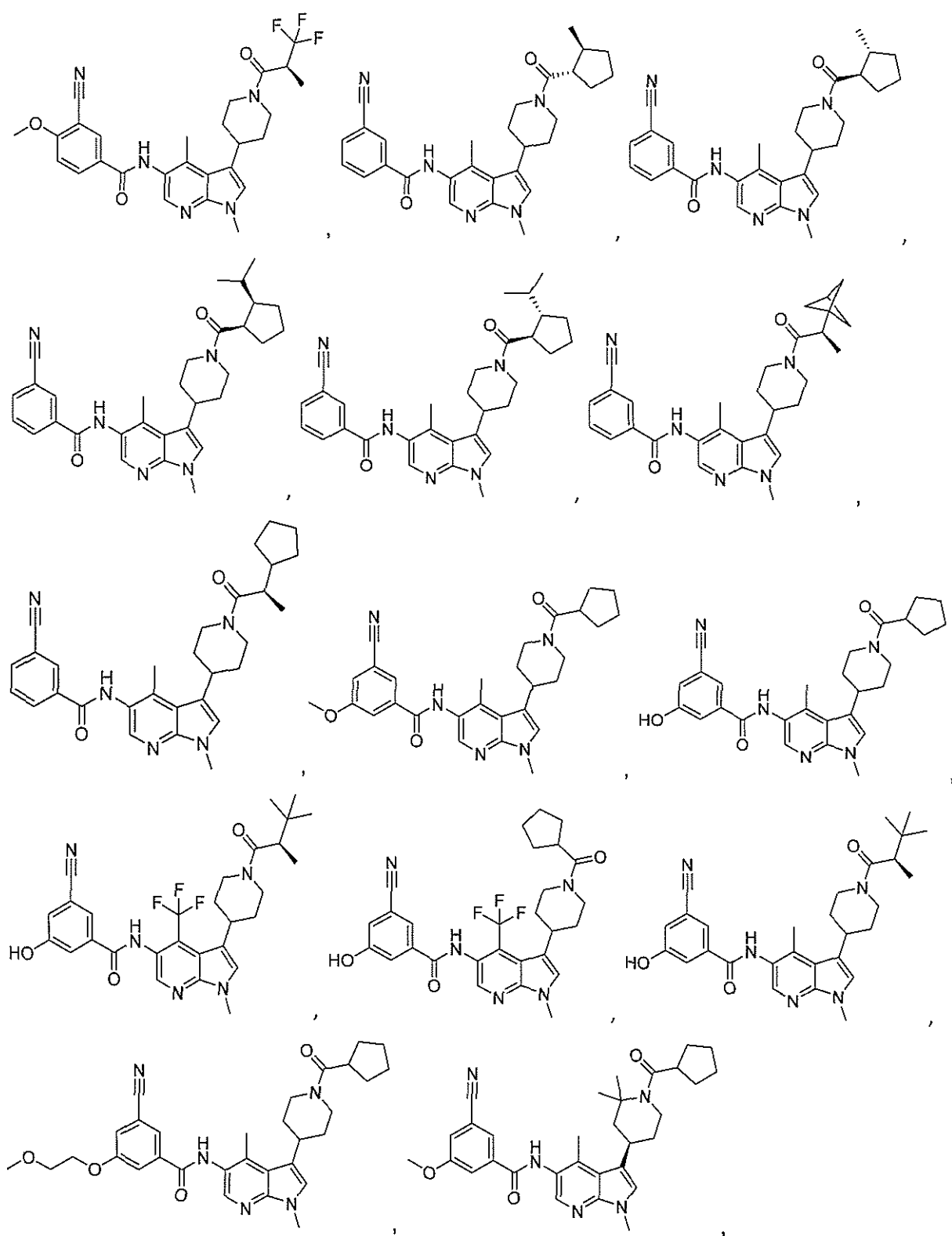
R² is (C₁-C₆)alkyl, (C₃-C₁₀)cycloalkyl, phenyl, tetrahydrothiophenyl, thietanyl or indanyl, optionally substituted with one, two, three, four or five substituents independently selected for each occurrence from the group consisting of -F, -Cl, -Br, -OH, (C₁-C₃)alkyl, (C₁-C₃)haloalkyl and (C₃-C₁₀)cycloalkyl.

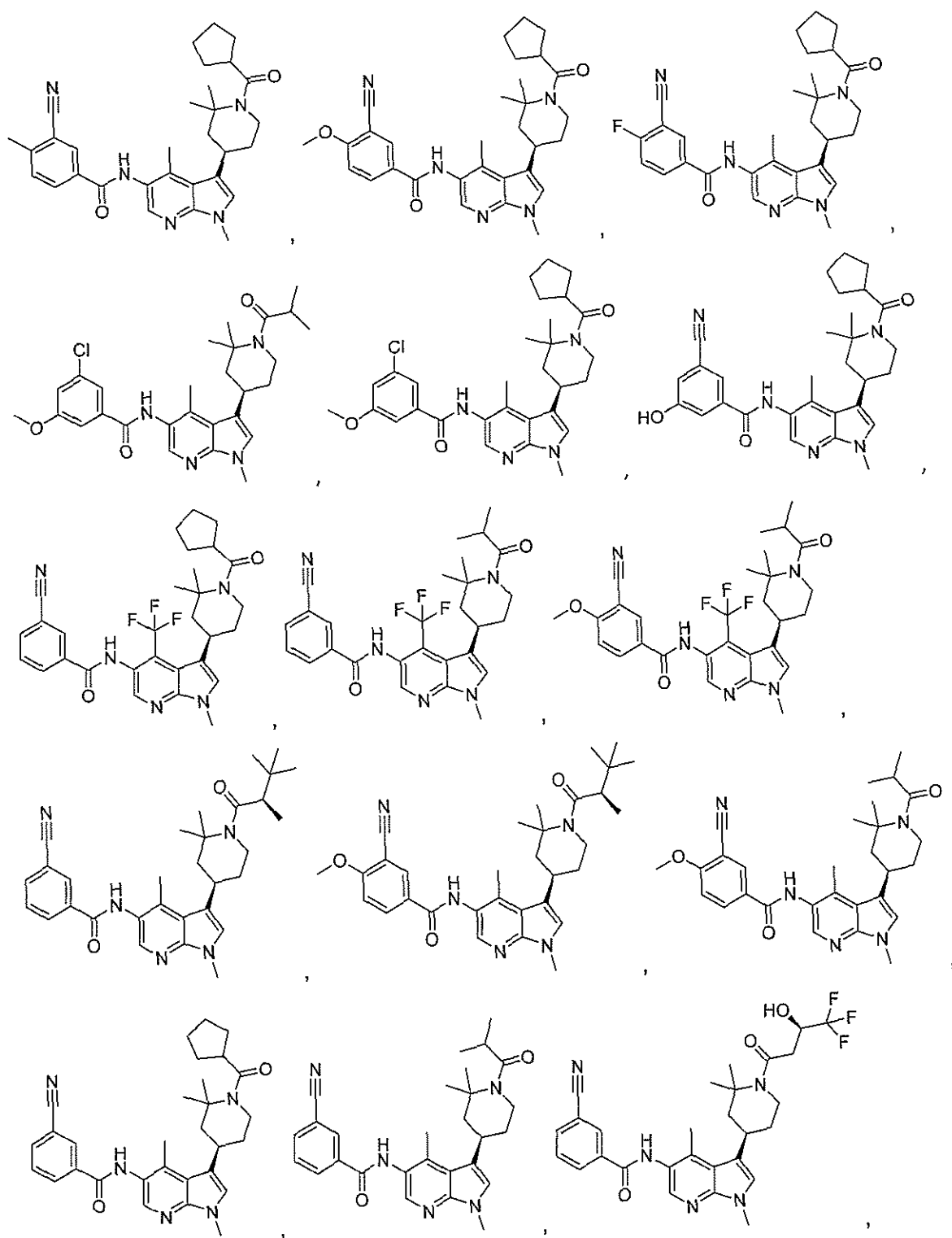
4. A compound selected from the group consisting of

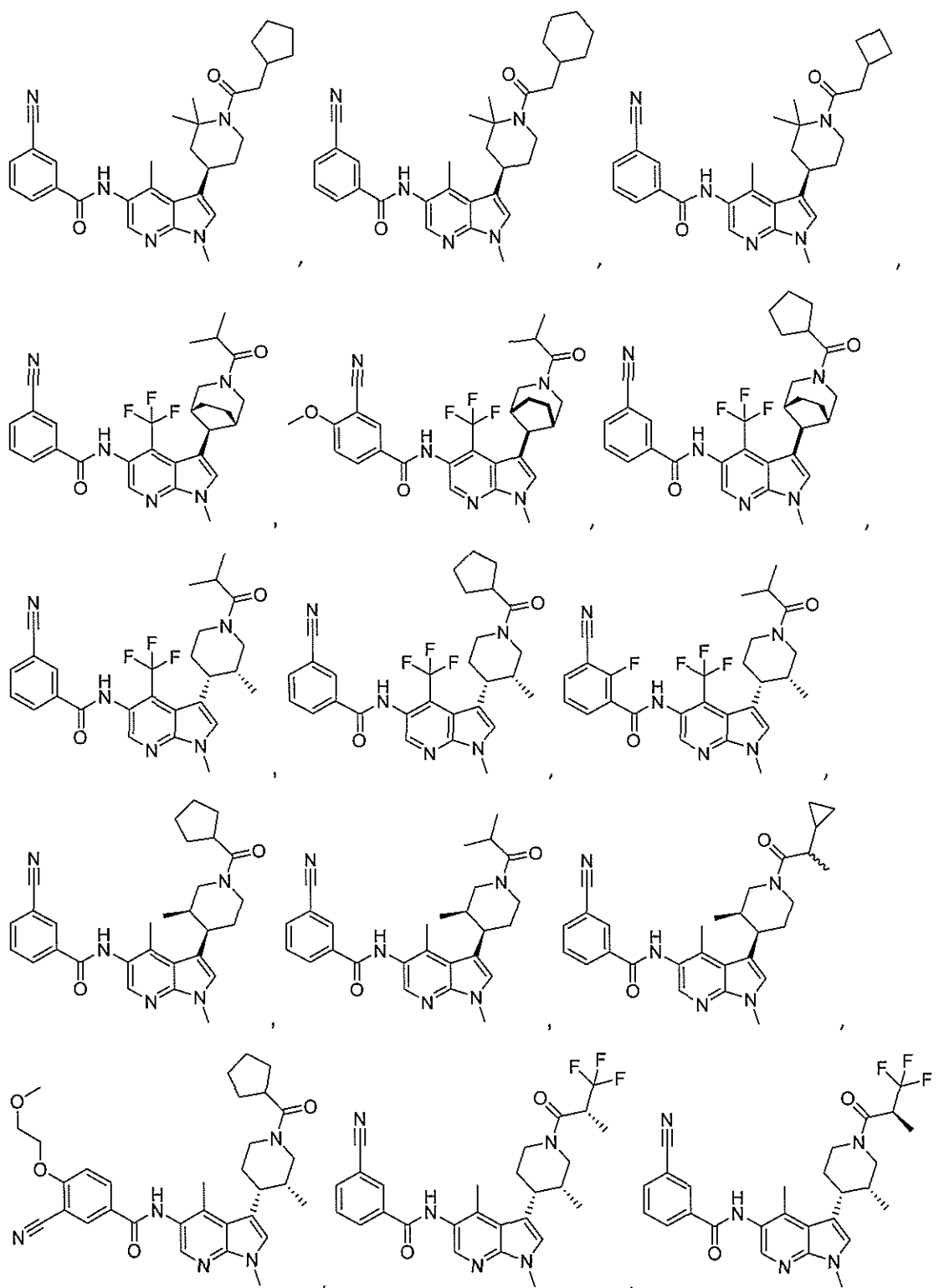


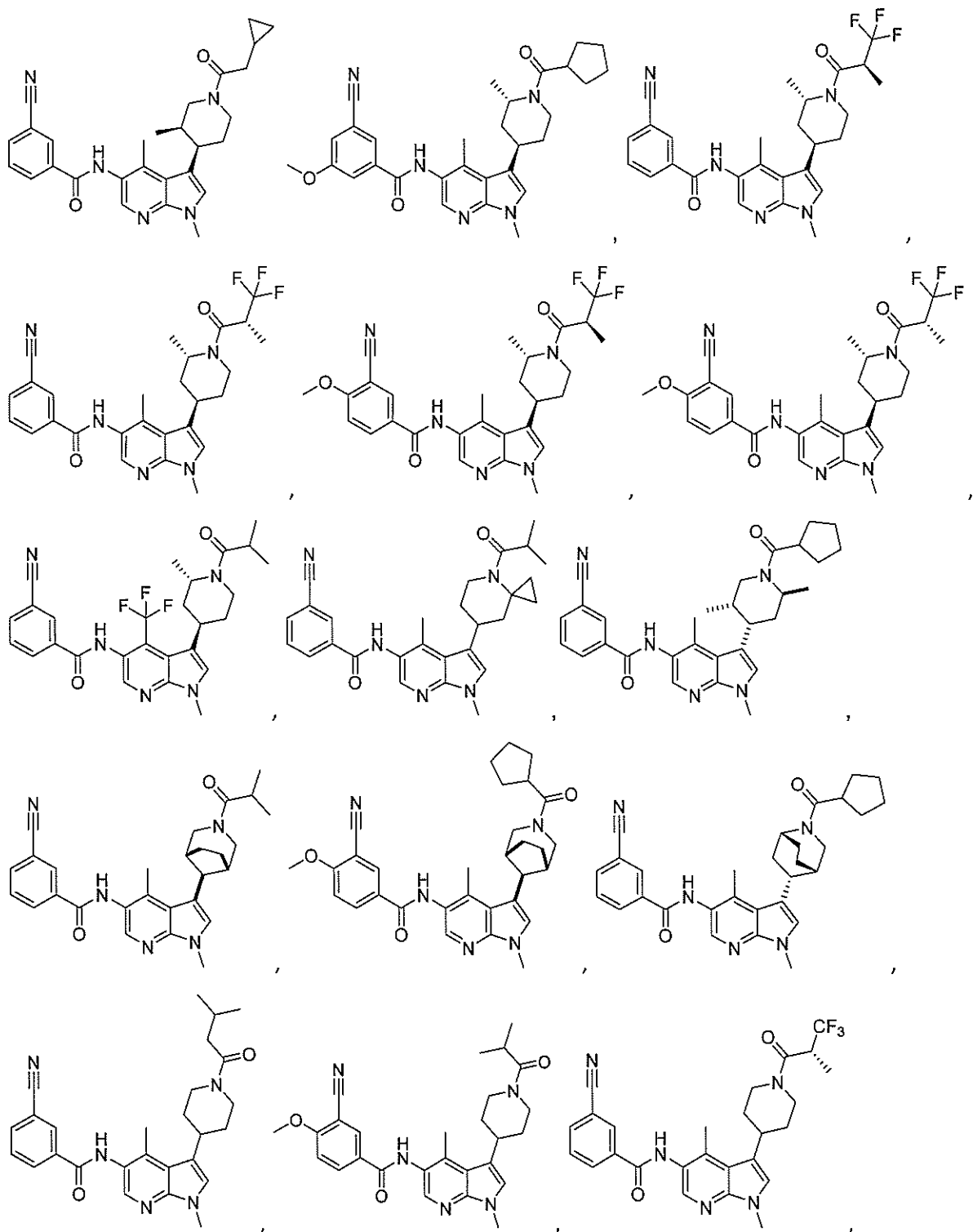


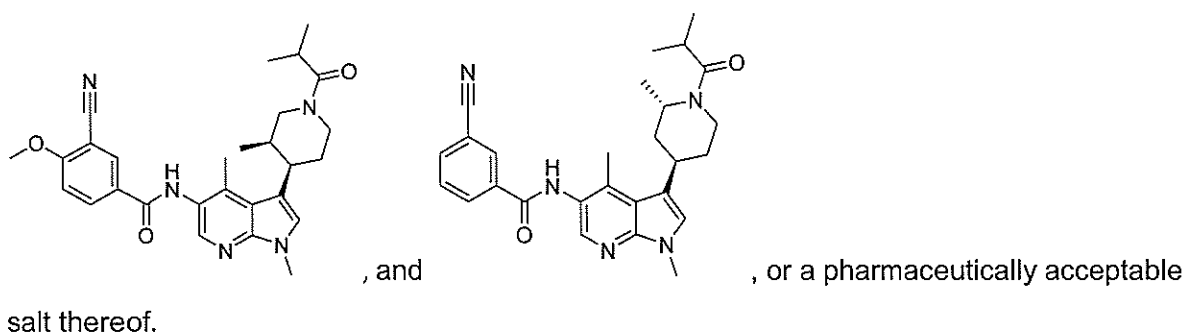




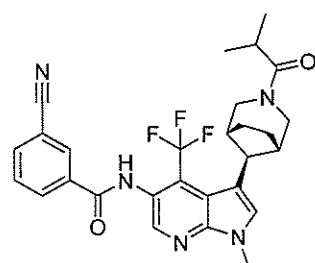






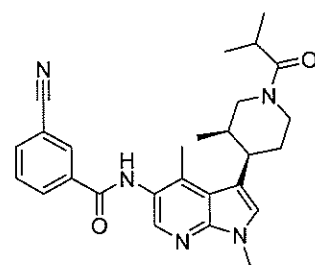


5. The compound of claim 1, wherein the compound is



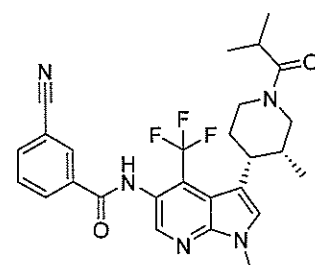
or a pharmaceutically acceptable salt thereof.

6. A compound that is



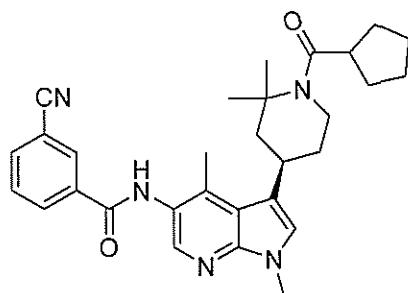
or a pharmaceutically acceptable salt thereof.

7. The compound of claim 1, wherein the compound is



or a pharmaceutically acceptable salt thereof.

8. A compound that is



or a pharmaceutically acceptable salt thereof.

9. A pharmaceutical composition comprising a compound according to any one of claims 1-8, or a pharmaceutically acceptable salt thereof, admixed with a pharmaceutically acceptable carrier, excipient or dilutant.