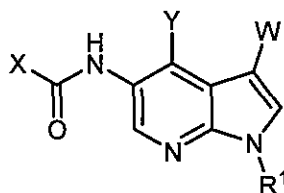


Claims

1. A compound of Formula I:



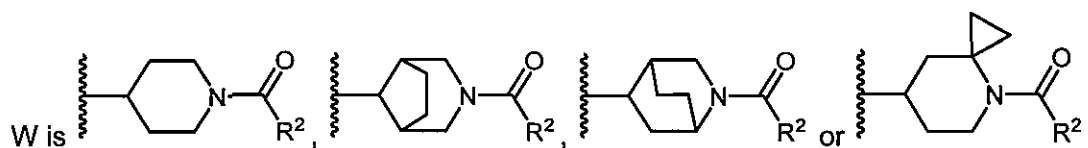
I

or a pharmaceutically acceptable salt, or pharmaceutically acceptable solvate thereof, wherein,

Y is $-\text{CF}_3$;

X is phenyl optionally substituted with one, two, three, four or five substituents independently selected from the group consisting of $-\text{CH}_3$, $-\text{CH}_2\text{CH}_3$, $-\text{CH}_2\text{OH}$, $-\text{OH}$, $-\text{OCH}_3$, $-\text{SCH}_3$, $-\text{OCH}_2\text{CH}_3$, $-\text{OCH}_2\text{CH}_2\text{OH}$, $-\text{OCH}_2\text{CH}_2\text{OCH}_3$, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$ and $-\text{CN}$;

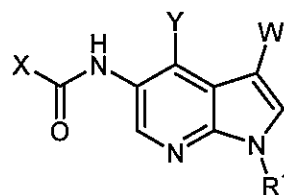
R^1 is $-\text{CH}_3$ or $-\text{CH}_2\text{CH}_3$;



each optionally substituted with one, two, three, four or five $-\text{CH}_3$; and

R^2 is $(\text{C}_1\text{-C}_6)$ alkyl, $(\text{C}_3\text{-C}_{10})$ 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 $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{OH}$, $(\text{C}_1\text{-C}_3)$ alkyl, $(\text{C}_1\text{-C}_3)$ haloalkyl and $(\text{C}_3\text{-C}_{10})$ cycloalkyl.

2. A compound of Formula I:



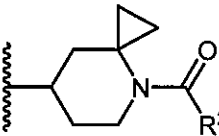
I

or a pharmaceutically acceptable salt, or pharmaceutically acceptable solvate thereof, wherein,

Y is $-\text{CH}_3$ or $-\text{CF}_3$;

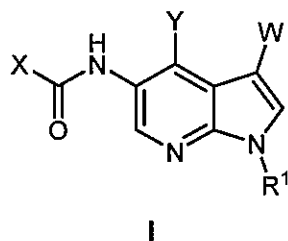
X is phenyl optionally substituted with one, two, three, four or five substituents independently selected from the group consisting of -CH₃, -CH₂CH₃, -CH₂OH, -OH, -OCH₃, -SCH₃, -OCH₂CH₃, -OCH₂CH₂OH, -OCH₂CH₂OCH₃, -F, -Cl, -Br and -CN;

R¹ is -CH₃ or -CH₂CH₃;

W is  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.

3. A compound of Formula I:

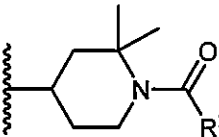


or a pharmaceutically acceptable salt, or pharmaceutically acceptable solvate thereof, wherein,

Y is -CH₃ or -CF₃;

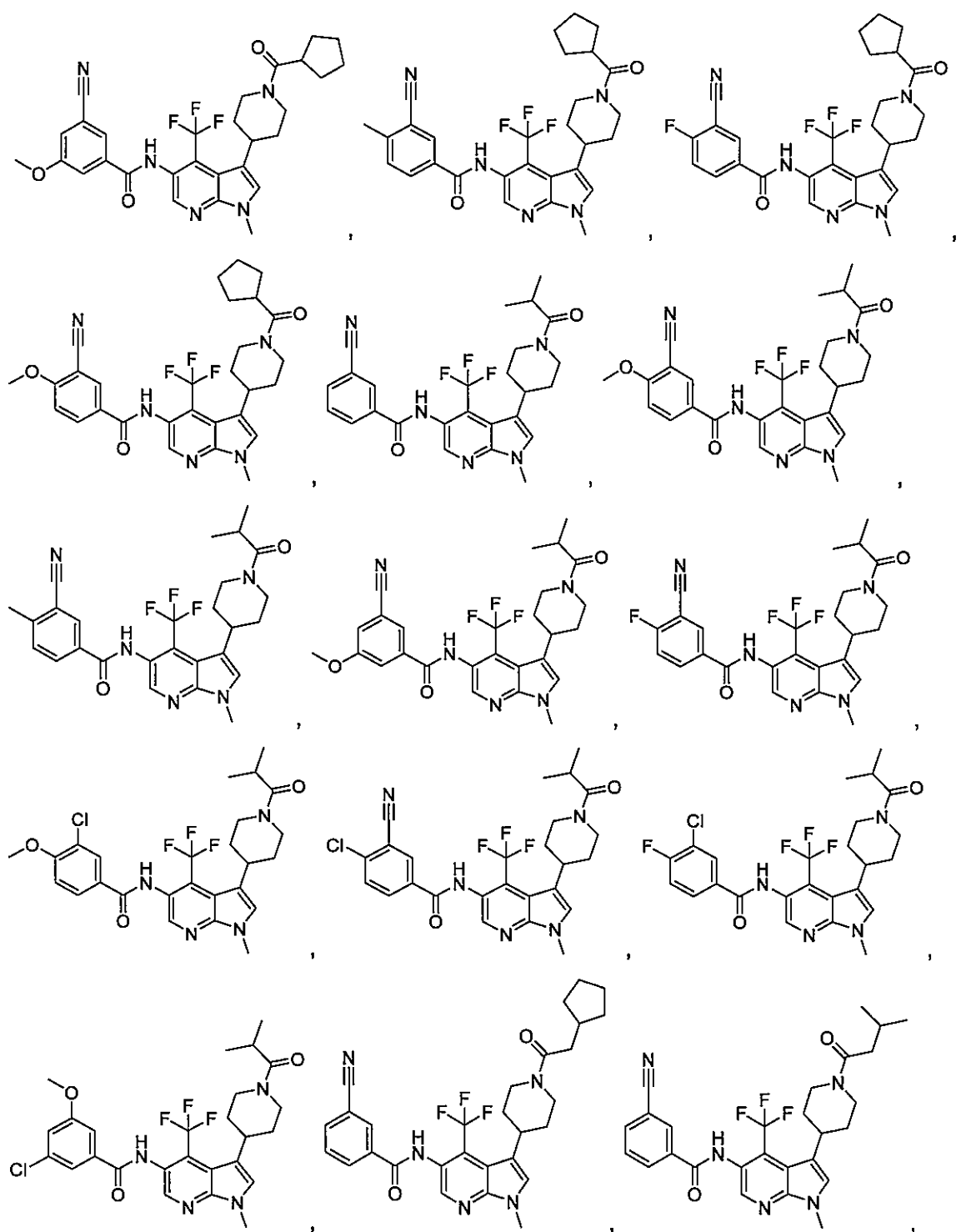
X is phenyl optionally substituted with one, two, three, four or five substituents independently selected from the group consisting of -CH₃, -CH₂CH₃, -CH₂OH, -OH, -OCH₃, -SCH₃, -OCH₂CH₃, -OCH₂CH₂OH, -OCH₂CH₂OCH₃, -F, -Cl, -Br and -CN;

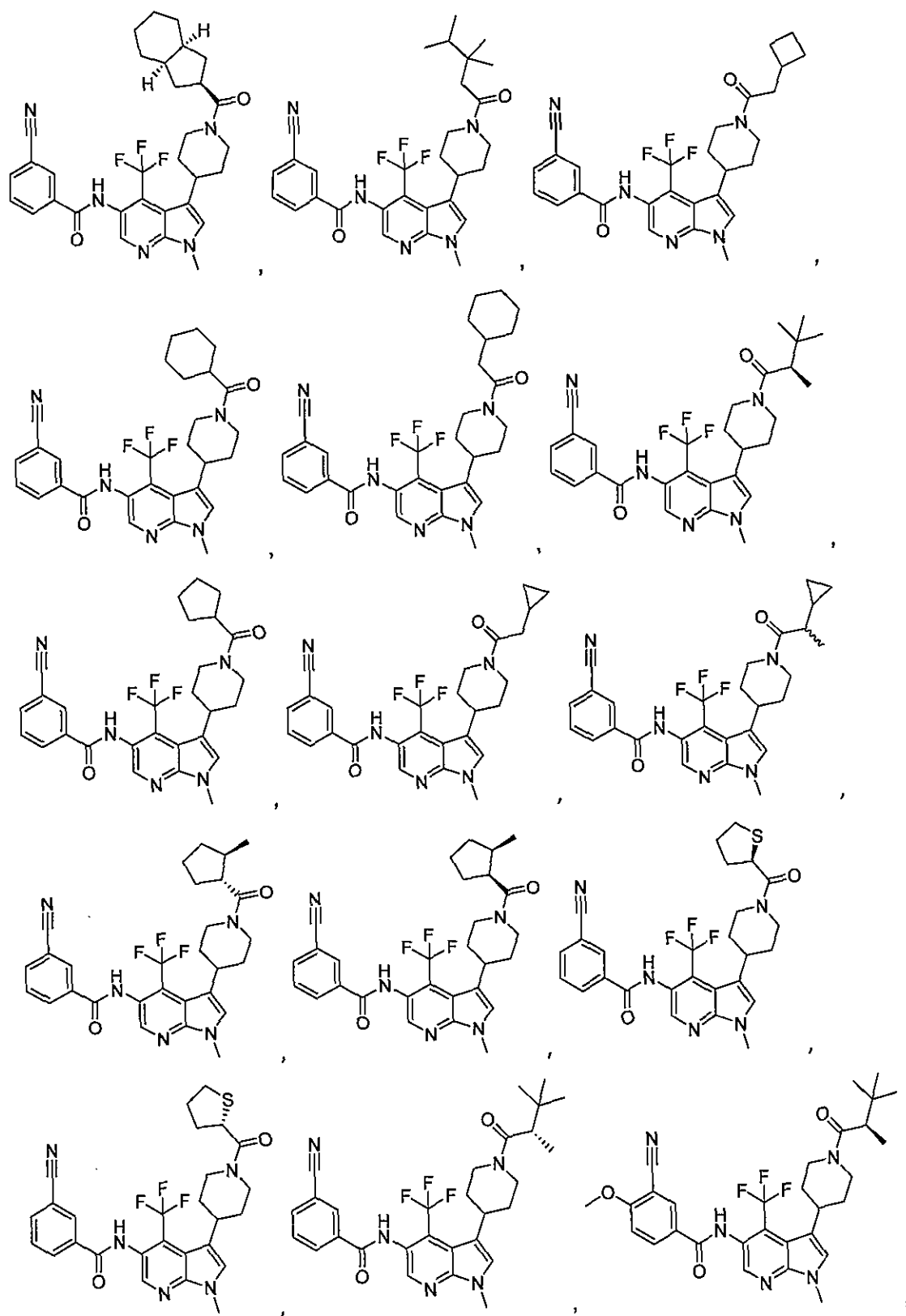
R¹ is -CH₃ or -CH₂CH₃;

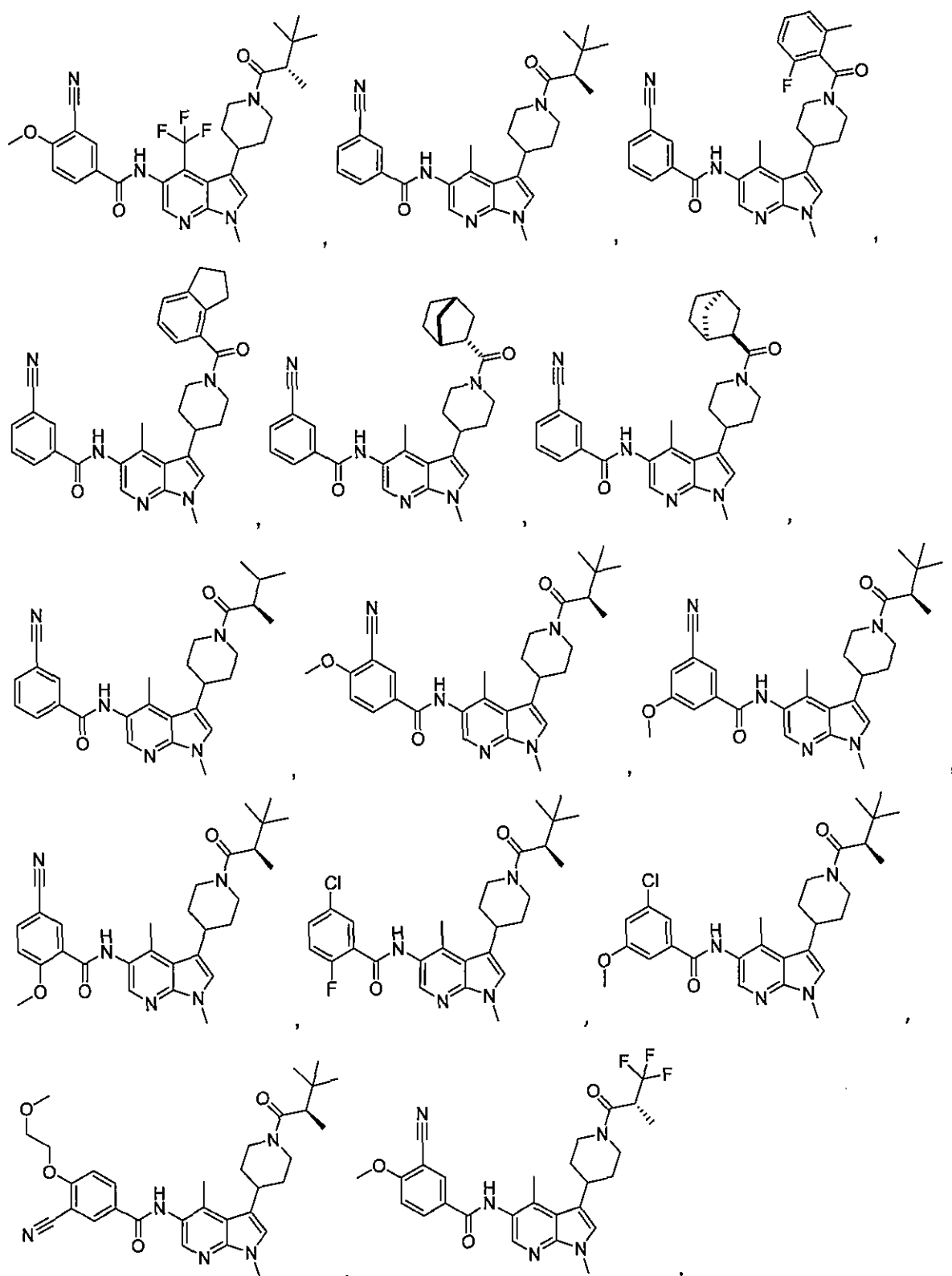
W is  ; and

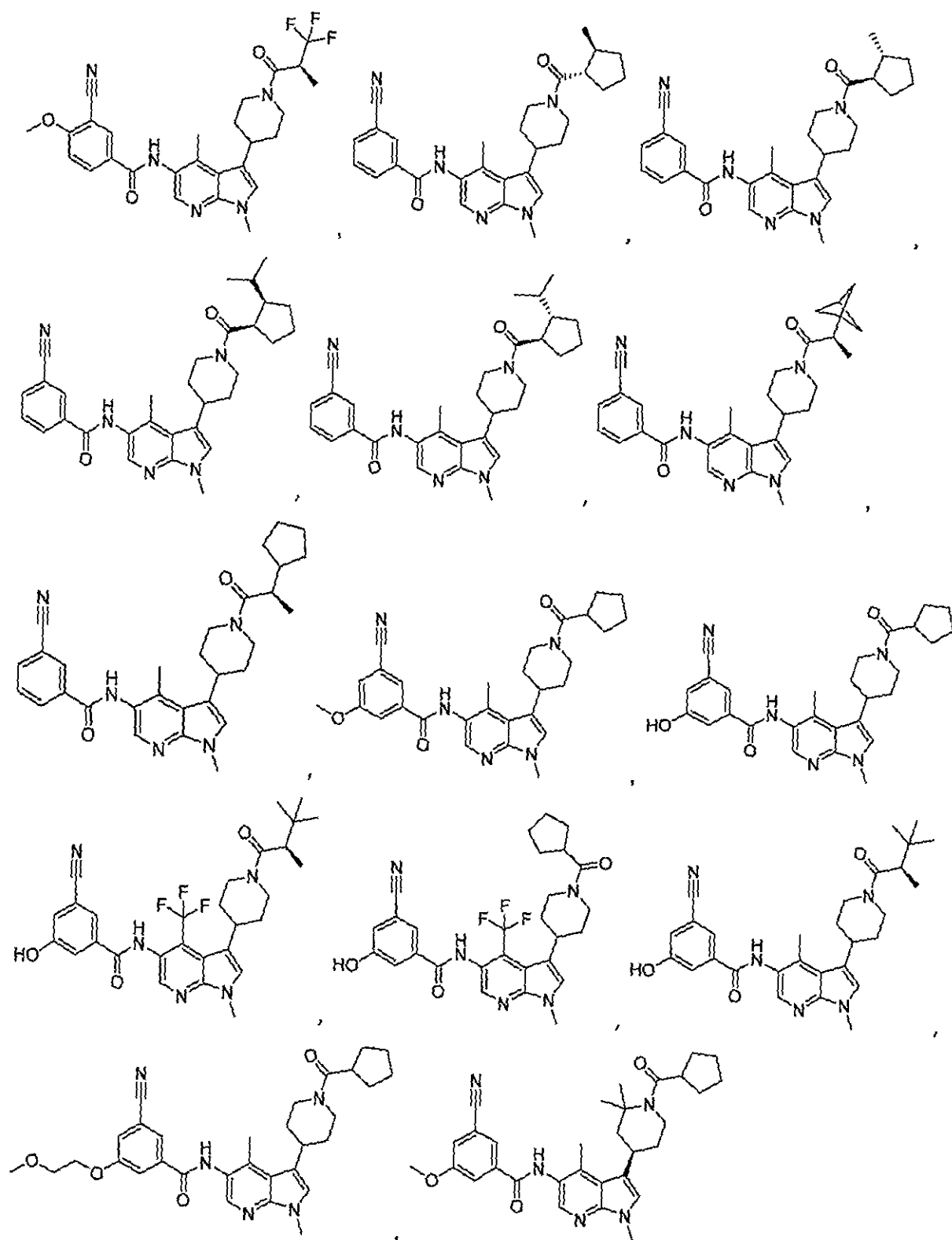
R² is (C₁-C₆)alkyl, (C₃-C₁₀)cycloalkyl, phenyl, tetrahydrothiophenyl, thietanyl or indanyl, optionally substituted with one, two, three, four or five substituents independently selected for each occurrence from the group consisting of -F, -Cl, -Br, -OH, (C₁-C₃)alkyl, (C₁-C₃)haloalkyl and (C₃-C₁₀)cycloalkyl.

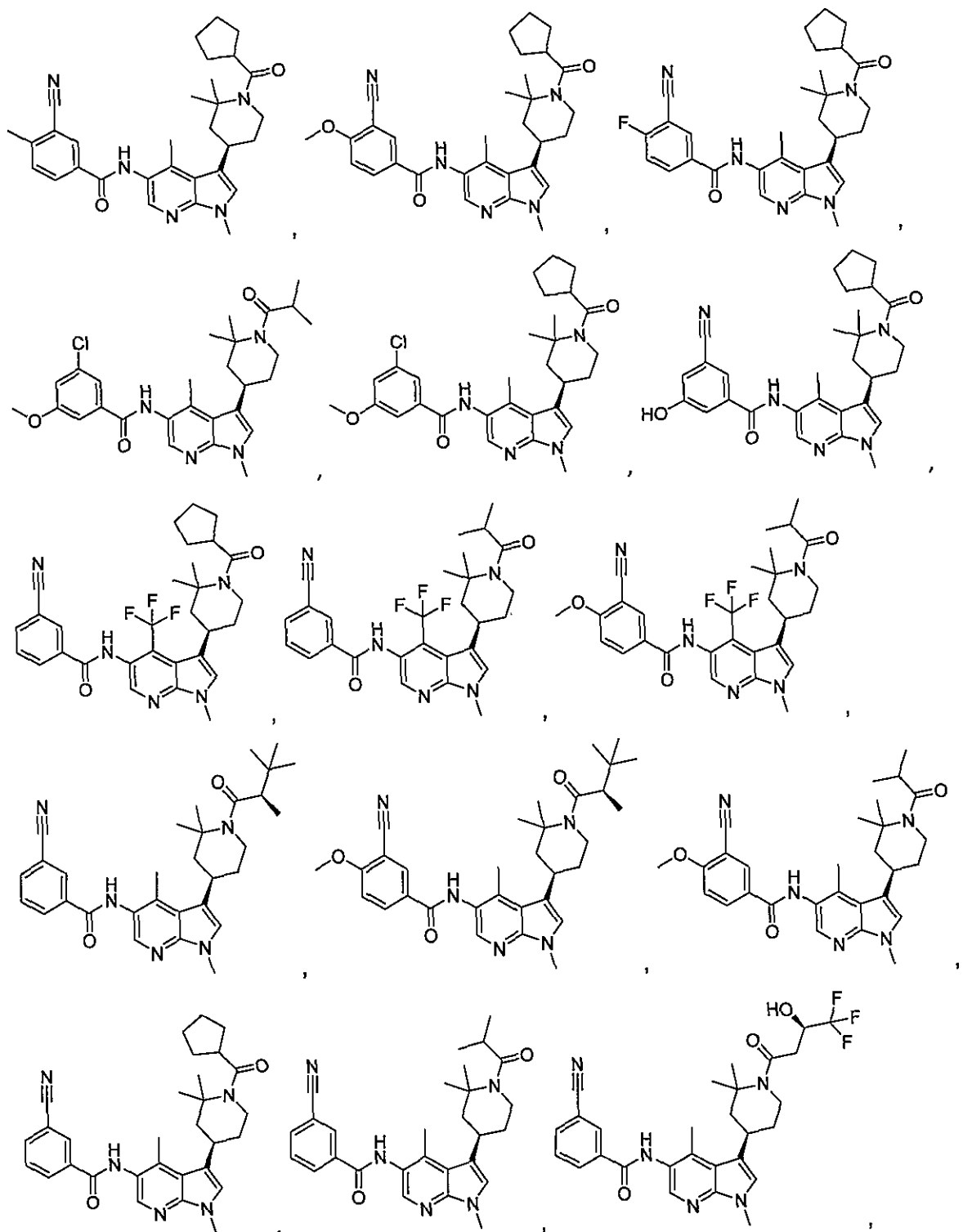
4. A compound selected from the group consisting of

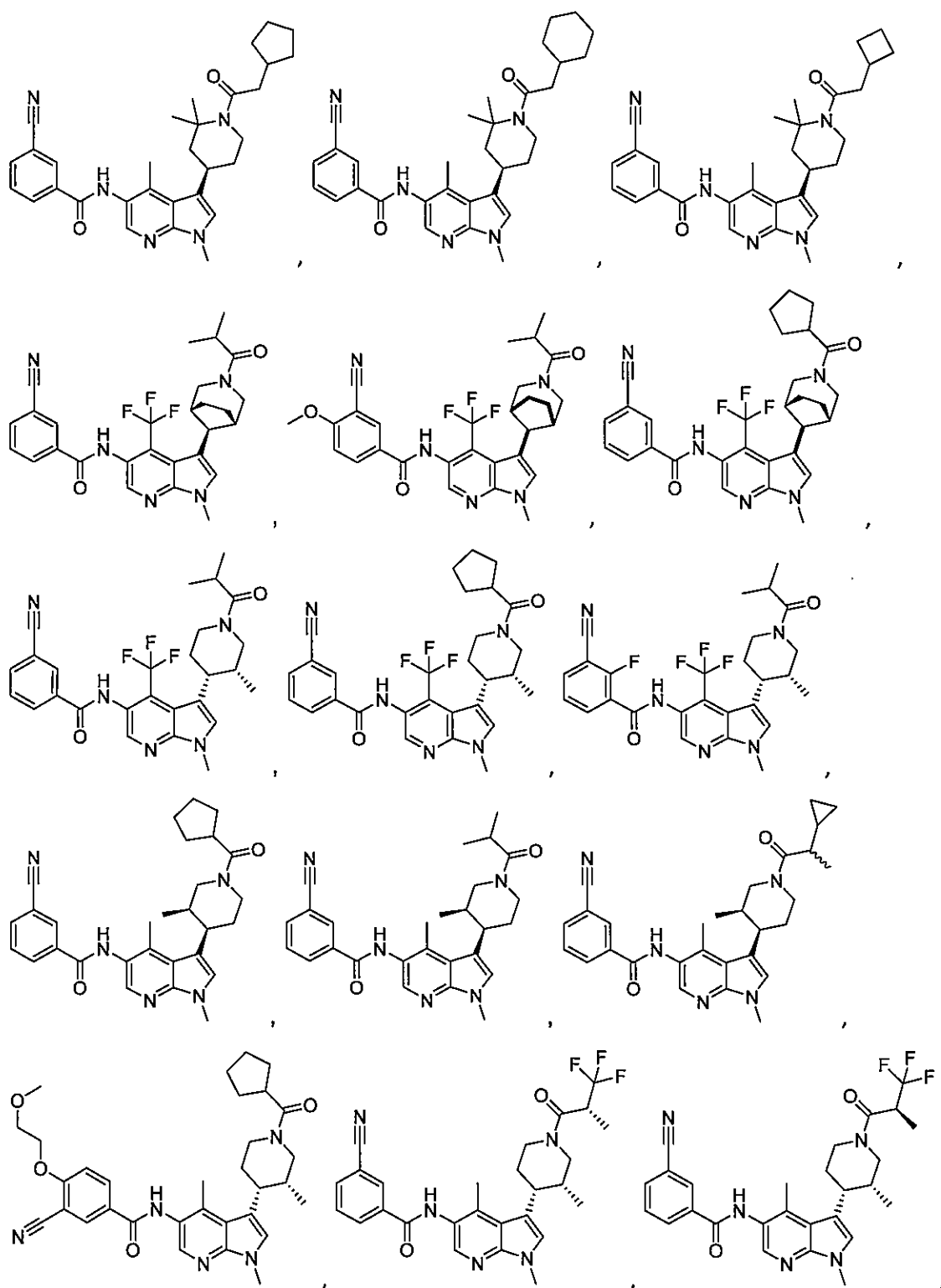


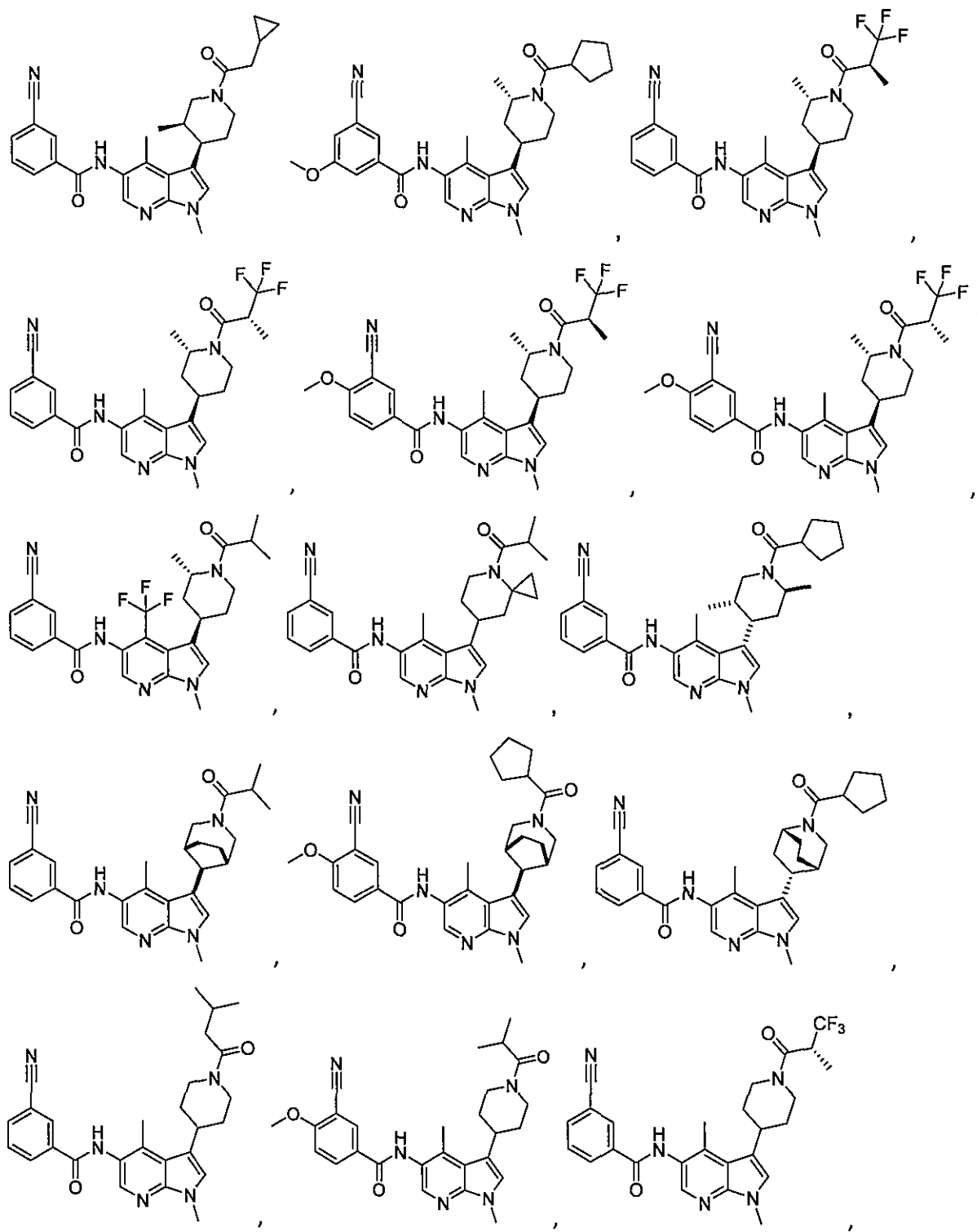


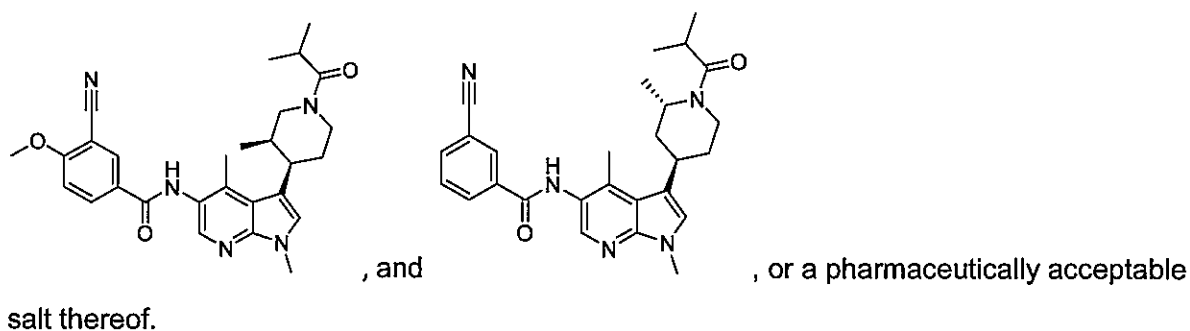




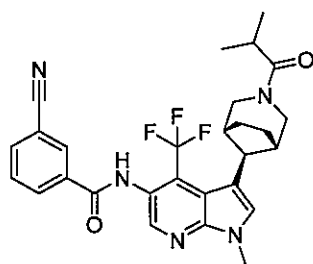






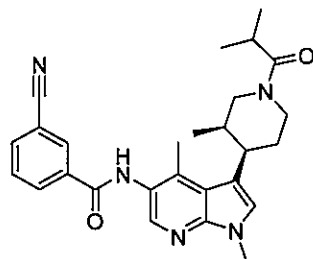


5. The compound of claim 1, wherein the compound is



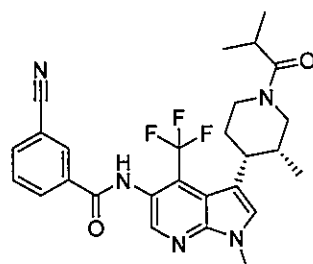
or a pharmaceutically acceptable salt thereof.

6. A compound that is



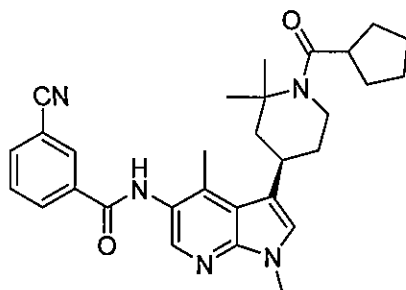
or a pharmaceutically acceptable salt thereof.

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.