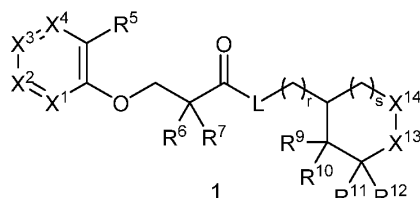


WHAT IS CLAIMED IS:

1. A compound of Formula 1,



or a pharmaceutically acceptable salt thereof in which:

X¹ is selected from N and CR¹;

X² is selected from N and CR²;

X³ is selected from N and CR³; and

X⁴ is selected from N and CR⁴, provided no more than two of X¹, X², X³, and X⁴ are N;

X¹³ is NR¹³ and X¹⁴ is CR¹⁵R¹⁶ or X¹³ is CH₂ and X¹⁴ is NR¹⁴;

L is selected from NR⁸ and O;

r is selected from 0 and 1;

s is selected from 0 and 1;

R¹, R², and R³ are each independently selected from

(a) hydrogen, halo, hydroxy, and cyano; and

(b) C₁₋₄ alkyl, C₃₋₆ cycloalkyl, and C₁₋₄ alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo;

R⁴ is selected from

(a) hydrogen, halo, hydroxy, and cyano; and

(b) C₁₋₄ alkyl, C₃₋₆ cycloalkyl, and C₁₋₄ alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo;

R⁵ is selected from

(a) hydrogen, halo, hydroxy, and cyano; and

(b) C₁₋₄ alkyl, C₃₋₆ cycloalkyl, and C₁₋₄ alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo, oxo, and phenyl which is substituted with 0 to 3 optional substituents independently selected from halo;
or

R⁴ and R⁵, together with the carbon atoms to which they are attached, form a cyclopent-1-en-1,2-diyl or a furan-2,3-diyl;

- R^6 and R^7 are each independently selected from halo and C_{1-3} alkyl, or R^6 and R^7 , together with the carbon atom to which they are attached, form a C_{3-4} cycloalkan-1,1-diyl;
- R^8 is selected from H and C_{1-4} alkyl;
- R^9 and R^{10} are each independently selected from
- (a) hydrogen, halo, hydroxy, and cyano;
 - (b) C_{1-4} alkyl, C_{3-6} cycloalkyl, and C_{1-4} alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo; and
 - (c) phenyl and C_{1-5} heteroaryl, each substituted with 0 to 3 optional substituents independently selected from halo, C_{1-4} alkyl, and C_{1-4} alkoxy, wherein the C_{1-5} heteroaryl substituent is a monocyclic ring with 5 to 6 ring members in which 1 to 4 ring members are heteroatoms, each of the heteroatoms independently selected from N, O, and S, provided no more than one of the ring members is O or S, and wherein the C_{1-4} alkyl and C_{1-4} alkoxy optional substituents on phenyl and C_{1-5} heteroaryl are each independently substituted with 0 to 3 optional substituents independently selected from halo; or
- R^9 and R^{10} , together with the carbon atom to which they are attached, form a C_{3-4} cycloalkan-1,1-diyl;
- R^{11} and R^{12} are each independently selected from
- (a) hydrogen, halo, hydroxy, and cyano; and
 - (b) C_{1-4} alkyl, C_{3-6} cycloalkyl, and C_{1-4} alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo; or
- R^{11} and R^{12} , together with the carbon atom to which they are attached, form a C_{3-4} cycloalkan-1,1-diyl;
- R^{13} and R^{14} are each independently selected from
- (a) hydrogen; and
 - (b) C_{1-4} alkyl, which is substituted with 0 to 3 optional substituents independently selected from cyano, oxo, and phenyl which is substituted with 0 to 3 optional substituents independently selected from halo; and
- R^{15} and R^{16} are each independently selected from
- (a) hydrogen, halo, hydroxy, and cyano; and
 - (b) C_{1-4} alkyl, C_{3-6} cycloalkyl, and C_{1-4} alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo; or

R^{15} and R^{16} , together with the carbon atom to which they are attached, form a C_{3-4} cycloalkan-1,1-diyl.

2. The compound or pharmaceutically acceptable salt according to claim 1, wherein X^1 is CR^1 , X^2 is CR^2 , X^3 is CR^3 , and X^4 is CR^4 .
3. The compound or pharmaceutically acceptable salt according to claim 2, wherein R^1 , R^2 , R^3 , and R^4 are each independently selected from
 - (a) hydrogen, halo, and cyano; and
 - (b) C_{1-4} alkyl, C_{3-6} cycloalkyl, and C_{1-4} alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo.
4. The compound or pharmaceutically acceptable salt according to claim 2, wherein R^1 , R^2 , R^3 , and R^4 are each independently selected from
 - (a) hydrogen, halo, and cyano; and
 - (b) C_{1-4} alkyl and C_{3-6} cycloalkyl, each substituted with 0 to 3 optional substituents independently selected from halo.
5. The compound or pharmaceutically acceptable salt according to claim 2, wherein R^1 , R^2 , R^3 , and R^4 are each independently selected from
 - (a) hydrogen, halo, and cyano; and
 - (b) methyl and cyclopropyl, each substituted with 0 to 3 optional substituents independently selected from halo.
6. The compound or pharmaceutically acceptable salt according to claim 2, wherein R^1 , R^2 , R^3 , and R^4 are each independently selected from hydrogen, halo, cyano, methyl, trifluoromethyl, and cyclopropyl.
7. The compound or pharmaceutically acceptable salt according to claim 1, wherein X^1 is N, X^2 is CR^2 , X^3 is CR^3 , and X^4 is CR^4 .
8. The compound or pharmaceutically acceptable salt according to claim 7, wherein R^2 , R^3 , and R^4 are each independently selected from
 - (a) hydrogen, halo, and cyano; and
 - (b) C_{1-4} alkyl, C_{3-6} cycloalkyl, and C_{1-4} alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo.

9. The compound or pharmaceutically acceptable salt according to claim 7, wherein R^2 , R^3 , and R^4 are each independently selected from
- (a) hydrogen, halo, and cyano; and
 - (b) C_{1-4} alkyl and C_{3-6} cycloalkyl, each substituted with 0 to 3 optional substituents independently selected from halo.
10. The compound or pharmaceutically acceptable salt according to claim 7, wherein R^2 , R^3 , and R^4 are each independently selected from
- (a) hydrogen, halo, and cyano; and
 - (b) methyl and cyclopropyl, each substituted with 0 to 3 optional substituents independently selected from halo.
11. The compound or pharmaceutically acceptable salt according to claim 7, wherein R^2 , R^3 , and R^4 are each independently selected from hydrogen, halo, cyano, methyl, trifluoromethyl, and cyclopropyl.
12. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 11, wherein R^5 is selected from
- (a) hydrogen, halo, hydroxy, and cyano; and
 - (b) C_{1-4} alkyl, C_{3-6} cycloalkyl, and C_{1-4} alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo, oxo, and phenyl which is substituted with 0 to 3 optional substituents independently selected from halo.
13. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 11, wherein R^5 is selected from
- (a) halo and cyano; and
 - (b) C_{1-4} alkyl, C_{3-6} cycloalkyl, and C_{1-4} alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo.
14. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 11, wherein R^5 is selected from
- (a) fluoro, chloro, bromo, and cyano; and
 - (b) methyl, ethyl, cyclopropyl, methoxy, and ethoxy, each substituted with 0 to 3 optional substituents independently selected from fluoro.

15. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 14, wherein R^6 and R^7 are each independently selected from fluoro and methyl or together with the carbon atom to which they are attached form a cyclopropan-1,1-diyl or a cyclobutan-1,1-diyl.
16. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 14, wherein each of R^6 and R^7 is methyl.
17. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 16, wherein L is NR^8 .
18. The compound or pharmaceutically acceptable salt according to claim 17, wherein R^8 is selected from hydrogen and methyl.
19. The compound or pharmaceutically acceptable salt according to claim 17, wherein R^8 is hydrogen.
20. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 19, wherein R^9 and R^{10} are each independently selected from
- (a) hydrogen, halo, hydroxy, and cyano;
 - (b) C_{1-4} alkyl, C_{3-6} cycloalkyl, and C_{1-4} alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo; and
 - (c) phenyl and C_{1-5} heteroaryl, each substituted with 0 to 3 optional substituents independently selected from halo, C_{1-4} alkyl, and C_{1-4} alkoxy, wherein the C_{1-5} heteroaryl substituent is a monocyclic ring with 5 to 6 ring members in which 1 to 4 ring members are heteroatoms, each of the heteroatoms independently selected from N, O, and S, provided no more than one of the ring members is O or S, and wherein the C_{1-4} alkyl and C_{1-4} alkoxy optional substituents on phenyl and C_{1-5} heteroaryl are each independently substituted with 0 to 3 optional substituents independently selected from halo.
21. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 19, wherein R^9 and R^{10} are each independently selected from
- (a) hydrogen and halo;
 - (b) C_{1-4} alkyl and C_{1-4} alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo; and

- (c) phenyl and C₁₋₅ heteroaryl, each substituted with 0 to 3 optional substituents independently selected from halo and C₁₋₄ alkyl, wherein the C₁₋₅ heteroaryl substituent is a monocyclic ring with 5 to 6 ring members in which 1 or 2 ring members are heteroatoms, each of the heteroatoms being N, and wherein the C₁₋₄ alkyl optional substituents on phenyl and C₁₋₅ heteroaryl are each independently substituted with 0 to 3 optional substituents independently selected from halo.
22. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 19, wherein R⁹ and R¹⁰ are each independently selected from
- (a) hydrogen and halo;
 - (b) C₁₋₄ alkyl and C₁₋₄ alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo; and
 - (c) phenyl, pyridinyl, and pyrazolyl, each substituted with 0 to 3 optional substituents independently selected from halo and C₁₋₄ alkyl which is substituted with 0 to 3 optional substituents independently selected from halo.
23. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 19, wherein R⁹ and R¹⁰ are selected from hydrogen, halo, and C₁₋₄ alkyl.
24. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 23, wherein R¹¹ and R¹² are each independently selected from
- (a) hydrogen, halo, hydroxy, and cyano; and
 - (b) C₁₋₄ alkyl, C₃₋₆ cycloalkyl, and C₁₋₄ alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo.
25. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 23, wherein R¹¹ and R¹² are each independently selected from hydrogen and C₁₋₄ alkyl.
26. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 25, wherein X¹³ is NR¹³ and X¹⁴ is CR¹⁵R¹⁶.
27. The compound or pharmaceutically acceptable salt according to claim 26, wherein R¹³ is selected from
- (a) hydrogen; and

- (b) C₁₋₄ alkyl, which is substituted with 0 to 3 optional substituents independently selected from cyano, oxo, and phenyl which is substituted with 0 to 3 optional substituents independently selected from halo.

28. The compound or pharmaceutically acceptable salt according to claim 26, wherein R¹³ is selected from hydrogen and methyl.

29. The compound or pharmaceutically acceptable salt according to claim 26, wherein R¹³ is methyl.

30. The compound or pharmaceutically acceptable salt according to any one of claims 26 to 29, wherein R¹⁵ and R¹⁶ are each independently selected from

(a) hydrogen, halo, hydroxy, and cyano; and

(b) C₁₋₄ alkyl, C₃₋₆ cycloalkyl, and C₁₋₄ alkoxy, each substituted with 0 to 3 optional substituents independently selected from halo.

31. The compound or pharmaceutically acceptable salt according to any one of claims 26 to 29, wherein R¹⁵ and R¹⁶ are each independently selected from hydrogen and C₁₋₄ alkyl which is substituted with 0 to 3 optional substituents independently selected from halo.

32. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 31, wherein r is 0.

33. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 32, wherein s is 0.

34. The compound or pharmaceutically acceptable salt according to any one of claims 1 to 32, wherein s is 1.

35. The compound according to claim 1, which is selected from the following compounds:

2,2-dimethyl-N-(1-methylpiperidin-4-yl)-3-((3-methylpyridin-2-yl)oxy)propanamide;

2,2-dimethyl-3-((3-methylpyridin-2-yl)oxy)-N-(1-methylpyrrolidin-3-yl)propanamide;

3-((3-cyanopyridin-2-yl)oxy)-2,2-dimethyl-N-(1-methylpiperidin-4-yl)propanamide;

(R)-3-((5-cyclopropylpyrimidin-4-yl)oxy)-2,2-dimethyl-N-(1-methylpyrrolidin-3-yl)propanamide;

3-((5-cyclopropylpyrimidin-4-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

3-((3-cyclopropyl-5-methylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

(*R*)-3-((3-cyclopropyl-5-methylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpyrrolidin-3-yl)propanamide;

3-((5-cyclopropyl-3-methylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

3-((3-cyclopropyl-6-methylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

(*R*)-3-((3-cyclopropyl-6-methylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpyrrolidin-3-yl)propanamide;

trans-N-(1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

cis-N-(1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-N-(1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

cis-N-(1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

trans-N-(4-isopropyl-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-N-(4-isopropyl-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

trans-N-(1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-methylpyridin-2-yl)oxy)propanamide;

trans-2,2-dimethyl-*N*-(1-methyl-4-(1-methyl-1*H*-pyrazol-4-yl)pyrrolidin-3-yl)-3-((3-methylpyridin-2-yl)oxy)propanamide;

trans-2,2-dimethyl-*N*-(1-methyl-4-(1-methyl-1*H*-pyrazol-4-yl)pyrrolidin-3-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-3-((3-cyclopropylpyridin-2-yl)oxy)-*N*-(1,3-dimethylpiperidin-4-yl)-2,2-dimethylpropanamide;

cis-3-((3-cyclopropylpyridin-2-yl)oxy)-*N*-(1,3-dimethylpiperidin-4-yl)-2,2-dimethylpropanamide;

trans-3-((3-cyclopropylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methyl-4-(1-methyl-1*H*-pyrazol-4-yl)pyrrolidin-3-yl)propanamide;

trans-*N*-(4-(1,3-dimethyl-1*H*-pyrazol-4-yl)-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-methylpyridin-2-yl)oxy)propanamide;

trans-3-((3-cyclopropylpyridin-2-yl)oxy)-*N*-(4-(1,3-dimethyl-1*H*-pyrazol-4-yl)-1-methylpyrrolidin-3-yl)-2,2-dimethylpropanamide;

trans-*N*-(4-(1,3-dimethyl-1*H*-pyrazol-4-yl)-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-*N*-(4-(1,5-dimethyl-1*H*-pyrazol-4-yl)-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-methylpyridin-2-yl)oxy)propanamide;

trans-3-((3-cyclopropylpyridin-2-yl)oxy)-*N*-(4-(1,5-dimethyl-1*H*-pyrazol-4-yl)-1-methylpyrrolidin-3-yl)-2,2-dimethylpropanamide;

trans-*N*-(4-(1,5-dimethyl-1*H*-pyrazol-4-yl)-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-*N*-(4-(1,3-dimethyl-1*H*-pyrazol-4-yl)-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

trans-*N*-(4-(1,5-dimethyl-1*H*-pyrazol-4-yl)-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

trans-*N*-(1,3-dimethylpiperidin-4-yl)-3-((3,5-dimethylpyridin-2-yl)oxy)-2,2-dimethylpropanamide;

cis-*N*-(1,3-dimethylpiperidin-4-yl)-3-((3,5-dimethylpyridin-2-yl)oxy)-2,2-dimethylpropanamide;

trans-*N*-(3-isopropyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

cis-*N*-(3-isopropyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)-*N*-(1,5,5-trimethylpyrrolidin-3-yl)propanamide;

(*R*)-2,2-dimethyl-*N*-((1-methylpyrrolidin-3-yl)methyl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

(*S*)-2,2-dimethyl-*N*-((1-methylpyrrolidin-3-yl)methyl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;
3-((3-chloropyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((3-fluoropyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((3,5-difluoropyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;
2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)-3-((4-(trifluoromethyl)pyridin-3-yl)oxy)propanamide;
2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-4-yl)oxy)propanamide;
3-((4-chloro-3-fluoropyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((3-chloro-5-fluoropyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((6-chloro-2-methylpyridin-3-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((3-cyclopropylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((5-chloro-3-methylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((3-methoxypyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)-3-((3-methylpyridin-4-yl)oxy)propanamide;
3-((5-cyano-3-methylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)-3-((2-(trifluoromethyl)pyridin-3-yl)oxy)propanamide;
(*R*)-3-((3-cyclopropylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpyrrolidin-3-yl)propanamide;
3-((3-ethoxypyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

trans-*N*-(4-(4-chlorophenyl)-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-methylpyridin-2-yl)oxy)propanamide;

trans-*N*-(4-(4-chlorophenyl)-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-2,2-dimethyl-*N*-(1-methyl-4-(6-(trifluoromethyl)pyridin-3-yl)pyrrolidin-3-yl)-3-((3-methylpyridin-2-yl)oxy)propanamide;

trans-3-((3-cyclopropylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methyl-4-(6-(trifluoromethyl)pyridin-3-yl)pyrrolidin-3-yl)propanamide;

trans-2,2-dimethyl-*N*-(1-methyl-4-(6-(trifluoromethyl)pyridin-3-yl)pyrrolidin-3-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-2,2-dimethyl-*N*-(1-methyl-4-(6-methylpyridin-3-yl)pyrrolidin-3-yl)-3-((3-methylpyridin-2-yl)oxy)propanamide;

trans-3-((3-cyclopropylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methyl-4-(6-methylpyridin-3-yl)pyrrolidin-3-yl)propanamide;

trans-2,2-dimethyl-*N*-(1-methyl-4-(6-methylpyridin-3-yl)pyrrolidin-3-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-(1-methylpiperidin-4-yl)-1-(((3-methylpyridin-2-yl)oxy)methyl)cyclopropane-1-carboxamide;

2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)-3-(2-(trifluoromethyl)phenoxy)propanamide;

(*R*)-2,2-dimethyl-*N*-(1-methylpyrrolidin-3-yl)-3-(2-(trifluoromethyl)phenoxy)propanamide;

3-(2-chlorophenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

3-(2-bromophenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

(*R*)-3-(2-chlorophenoxy)-2,2-dimethyl-*N*-(1-methylpyrrolidin-3-yl)propanamide;

(*R*)-3-(2-bromophenoxy)-2,2-dimethyl-*N*-(1-methylpyrrolidin-3-yl)propanamide;

3-(2-cyclopropylphenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

(*R*)-3-(2-cyclopropylphenoxy)-2,2-dimethyl-*N*-(1-methylpyrrolidin-3-yl)propanamide;

3-(2-ethylphenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

3-(4-cyano-2-(trifluoromethyl)phenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)-3-phenoxypropanamide;

3-(3-cyano-2-methylphenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)-3-(*o*-tolylloxy)propanamide;
3-(2-cyano-6-methylphenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-(2-ethoxyphenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((2,3-dihydro-1*H*-inden-4-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((3-cyclopropylpyrazin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
1-((2-chlorophenoxy)methyl)-*N*-(1-methylpiperidin-4-yl)cyclopropane-1-carboxamide;
3-((3-chloro-5-methylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((3,6-dimethylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((3,5-dimethylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((3-ethylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
(*R*)-3-((3-ethylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpyrrolidin-3-yl)propanamide;
3-((3-ethylpyrazin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
3-((3,4-dimethylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
1-(((3-cyclopropylpyridin-2-yl)oxy)methyl)-*N*-(1-methylpiperidin-4-yl)cyclobutane-1-carboxamide;
3-((3-(difluoromethyl)pyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
2,2-dimethyl-3-((3-methyl-5-(trifluoromethyl)pyridin-2-yl)oxy)-*N*-(1-methylpiperidin-4-yl)propanamide;
2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;
trans-N-(4-(4-chlorophenyl)-1-methylpyrrolidin-3-yl)-3-((3,5-dimethylpyridin-2-yl)oxy)-2,2-dimethylpropanamide;

trans-N-(4-(4-chlorophenyl)-1-methylpyrrolidin-3-yl)-3-((3-cyclopropylpyridin-2-yl)oxy)-2,2-dimethylpropanamide;

trans-N-(4-ethoxy-1-methylpyrrolidin-3-yl)-3-(2-ethylphenoxy)-2,2-dimethylpropanamide;

trans-3-(2-ethylphenoxy)-*N*-(4-methoxy-1-methylpyrrolidin-3-yl)-2,2-dimethylpropanamide;

3-(2-ethylphenoxy)-2,2-dimethyl-*N*-(1-methyl-4-phenylpyrrolidin-3-yl)propanamide;

3-((6-chloro-4-(trifluoromethyl)pyridazin-3-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

3-(furo[3,2-*c*]pyridin-4-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

3-((5-bromo-3-(trifluoromethyl)pyridin-2-yl)oxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

2,2-dimethyl-3-((5-methyl-3-(trifluoromethyl)pyridin-2-yl)oxy)-*N*-(1-methylpiperidin-4-yl)propanamide;

2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)-3-(2-(trifluoromethoxy)phenoxy)propanamide;

trans-N-(1-(cyanomethyl)-4-(6-methylpyridin-3-yl)pyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-(1-(2-(4-chlorophenyl)-2-oxoethyl)piperidin-4-yl)-3-(2-ethylphenoxy)-2,2-dimethylpropanamide;

N-((3*S*,4*S*)-1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-((3*R*,4*R*)-1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-((3*S*,4*S*)-1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

N-((3*R*,4*R*)-1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

trans-N-(1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-N-(3-ethyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

cis-*N*-(3-ethyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)-*N*-(1,3,3-trimethylpiperidin-4-yl)propanamide;

N-(3-fluoro-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(5-methyl-5-azaspiro[2.5]octan-8-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-(*trans*-3-ethyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

N-(*cis*-3-ethyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

N-((3*S*,4*S*)-3-fluoro-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-(3,3-difluoro-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(5-methyl-5-azaspiro[2.5]octan-8-yl)-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

N-(3-chloro-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-((3*S*,4*S*)-3-fluoro-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-methylpyridin-2-yl)oxy)propanamide;

3-((3,5-dimethylpyridin-2-yl)oxy)-*N*-((3*S*,4*S*)-3-fluoro-1-methylpiperidin-4-yl)-2,2-dimethylpropanamide;

3-((3-cyclopropylpyridin-2-yl)oxy)-*N*-((3*S*,4*S*)-3-fluoro-1-methylpiperidin-4-yl)-2,2-dimethylpropanamide;

N-((3*S*,4*S*)-3-fluoro-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

trans-*N*-(3-(1-(difluoromethyl)-1*H*-pyrazol-4-yl)-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-*N*-(3-(1-(difluoromethyl)-1*H*-pyrazol-4-yl)-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

trans-N-(3-ethyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-methylpyridin-2-yl)oxy)propanamide;

trans-3-((3,5-dimethylpyridin-2-yl)oxy)-*N*-(3-ethyl-1-methylpiperidin-4-yl)-2,2-dimethylpropanamide;

cis-N-(1,2-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-N-(4-fluoropyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-(4,4-difluoropyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide ;

2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)-*N*-(1,2,2-trimethylpiperidin-4-yl)propanamide;

trans-N-(1,5-dimethylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

cis-N-(1,5-dimethylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-N-(1,2-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

cis-N-(1,2-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

N-(2-ethyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(5-methyl-5-azaspiro[2.4]heptan-7-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-N-(1,2-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-((3*R*,4*R*)-1,3-dimethylpiperidin-4-yl)-3-((3,5-dimethylpyridin-2-yl)oxy)-2,2-dimethylpropanamide;

N-((3*S*,4*S*)-1,3-dimethylpiperidin-4-yl)-3-((3,5-dimethylpyridin-2-yl)oxy)-2,2-dimethylpropanamide;

N-((3*R*,4*R*)-1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-methylpyridin-2-yl)oxy)propanamide;

N-((3*S*,4*S*)-1,3-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-methylpyridin-2-yl)oxy)propanamide;

N-((3*R*,4*R*)-3-ethyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-((3*S*,4*S*)-3-ethyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-(2-ethyl-1-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

3-((3-cyclopropylpyridin-2-yl)oxy)-*N*-((2*R*,4*R*)-1,2-dimethylpiperidin-4-yl)-2,2-dimethylpropanamide;

(*R*)-2,2-dimethyl-*N*-(5-methyl-5-azaspiro[2.4]heptan-7-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-N-(2-ethyl-1-methylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

(*R*)-2,2-dimethyl-*N*-(5-methyl-5-azaspiro[2.4]heptan-7-yl)-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(5-methyl-5-azaspiro[3.4]octan-8-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-((2*S*,4*S*)-1,2-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

(*R*)-2,2-dimethyl-*N*-(5-methyl-5-azaspiro[2.4]heptan-7-yl)-3-((3-methylpyridin-2-yl)oxy)propanamide;

(*R*)-3-((3,5-dimethylpyridin-2-yl)oxy)-2,2-dimethyl-*N*-(5-methyl-5-azaspiro[2.4]heptan-7-yl)propanamide;

N,2,2-trimethyl-*N*-(*trans*-3-(*o*-tolyl)piperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N,2,2-trimethyl-*N*-(*cis*-3-(*o*-tolyl)piperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(*trans*-3-(*o*-tolyl)piperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-((3*S*,4*S*)-3-fluoropiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-(3,3-difluoropiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-((3*S*,4*S*)-3-fluoropiperidin-4-yl)-*N*,2,2-trimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(*trans*-1-methyl-3-(*o*-tolyl)piperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-*N*,2,2-trimethyl-*N*-(3-phenylpiperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(*cis*-1-methyl-3-phenylpiperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(*trans*-1-methyl-3-phenylpiperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

N-(*trans*-1,4-dimethylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(*cis*-1-methyl-3-phenylpiperidin-4-yl)-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(*trans*-1-methyl-3-phenylpiperidin-4-yl)-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

trans-3-(*o*-tolyl)piperidin-4-yl 2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanoate;

2,2-dimethyl-*N*-(*trans*-1-methyl-4-(1-methyl-1*H*-pyrazol-4-yl)piperidin-3-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

trans-3-(2-methylpyridin-3-yl)piperidin-4-yl 2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanoate;

trans-3-phenylpiperidin-4-yl 2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanoate;

N,2,2-trimethyl-*N*-(*trans*-3-(2-methylpyridin-3-yl)piperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

(*S*)-2,2-dimethyl-*N*-(5-methyl-5-azaspiro[2.4]heptan-7-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(4-methyl-4-azaspiro[2.5]octan-7-yl)-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

2,2-dimethyl-*N*-(4-methyl-4-azaspiro[2.5]octan-7-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;
N-((2*R*,4*R*)-1,2-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;
N-((2*R*,4*S*)-1,2-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;
N-((2*S*,4*S*)-1,2-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;
N-((2*R*,4*R*)-1,2-dimethylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;
N-(*trans*-1,4-dimethylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;
3-((3-cyclopropylpyridin-2-yl)oxy)-*N*-(*trans*-1,4-dimethylpyrrolidin-3-yl)-2,2-dimethylpropanamide;
N-(*trans*-1,4-dimethylpyrrolidin-3-yl)-2,2-dimethyl-3-((3-methylpyridin-2-yl)oxy)propanamide;
3-((3,5-dimethylpyridin-2-yl)oxy)-*N*-(*trans*-1,4-dimethylpyrrolidin-3-yl)-2,2-dimethylpropanamide;
trans-*N*-(1-(cyanomethyl)-3-methylpiperidin-4-yl)-2,2-dimethyl-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;
3-((3-(chlorodifluoromethoxy)pyridin-2-yl)oxy)-*N*-((3*S*,4*S*)-1,3-dimethylpiperidin-4-yl)-2,2-dimethylpropanamide;
trans-3-(2-chlorophenoxy)-*N*-(1,3-dimethylpiperidin-4-yl)-2,2-dimethylpropanamide;
3-(2-chlorophenoxy)-*N*-((3*S*,4*S*)-3-fluoropiperidin-4-yl)-2,2-dimethylpropanamide;
3-(2-chlorophenoxy)-*N*-((3*S*,4*S*)-3-fluoro-1-methylpiperidin-4-yl)-2,2-dimethylpropanamide;
trans-*N*-(1,3-dimethylpiperidin-4-yl)-2,2-difluoro-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide ;
3-(2-chlorophenoxy)-2,2-dimethyl-*N*-((3*S*,4*S*)-3-methylpiperidin-4-yl)propanamide;
cis-*N*-(1,2-dimethylpiperidin-4-yl)-2,2-difluoro-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;
trans-3-((3-cyclopropylpyridin-2-yl)oxy)-*N*-(1,3-dimethylpiperidin-4-yl)-2,2-difluoropropanamide;

3-(2-chlorophenoxy)-*N*-((3*S*,4*S*)-1,3-dimethylpiperidin-4-yl)-2,2-dimethylpropanamide;

3-((3-cyclopropylpyridin-2-yl)oxy)-*N*-((2*R*,4*R*)-1,2-dimethylpiperidin-4-yl)-2,2-difluoropropanamide;

3-(2-fluorophenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

3-(2-fluorophenoxy)-2,2-dimethyl-*N*-((3*S*,4*S*)-3-methylpiperidin-4-yl)propanamide;

2,2-difluoro-*N*-((3*S*,4*S*)-3-fluoropiperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

3-((3-cyclopropylpyridin-2-yl)oxy)-2,2-difluoro-*N*-((3*S*,4*S*)-3-fluoropiperidin-4-yl)propanamide;

N-(*trans*-1,3-dimethylpiperidin-4-yl)-2,2-difluoro-3-((3-methylpyridin-2-yl)oxy)propanamide;

N-(*trans*-1,3-dimethylpiperidin-4-yl)-2,2-difluoro-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

2,2-difluoro-*N*-((3*S*,4*S*)-3-fluoropiperidin-4-yl)-3-((3-methylpyridin-2-yl)oxy)propanamide;

N-((3*S*,4*S*)-1,3-dimethylpiperidin-4-yl)-3-(2-fluorophenoxy)-2,2-dimethylpropanamide;

3-(3-fluorophenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

3-(4-fluorophenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;

N-((2*R*,4*R*)-1,2-dimethylpiperidin-4-yl)-2,2-difluoro-3-((3-methylpyridin-2-yl)oxy)propanamide;

N-((2*R*,4*R*)-1,2-dimethylpiperidin-4-yl)-2,2-difluoro-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

2,2-difluoro-*N*-((3*S*,4*S*)-3-fluoropiperidin-4-yl)-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

2,2-difluoro-*N*-((3*S*,4*S*)-3-methylpiperidin-4-yl)-3-((3-(trifluoromethyl)pyridin-2-yl)oxy)propanamide;

3-((3-cyclopropylpyridin-2-yl)oxy)-2,2-difluoro-*N*-((3*S*,4*S*)-3-methylpiperidin-4-yl)propanamide;

2,2-difluoro-*N*-((3*S*,4*S*)-3-methylpiperidin-4-yl)-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;

3-((3-cyclopropylpyridin-2-yl)oxy)-2,2-difluoro-*N*-((2*R*,4*R*)-2-methylpiperidin-4-yl)propanamide;
2,2-difluoro-*N*-((2*R*,4*R*)-2-methylpiperidin-4-yl)-3-((3-(trifluoromethoxy)pyridin-2-yl)oxy)propanamide;
3-(2-chlorophenoxy)-2,2-dimethyl-*N*-(piperidin-4-yl)propanamide;
3-((3,5-dimethylpyridin-2-yl)oxy)-*N*-((3*R*,4*R*)-3-fluoropiperidin-4-yl)-2,2-dimethylpropanamide;
3-(2-chlorophenoxy)-2,2-difluoro-*N*-(1-methylpiperidin-4-yl)propanamide;
3-(2-chlorophenoxy)-2,2-difluoro-*N*-((3*S*,4*S*)-3-methylpiperidin-4-yl)propanamide;
3-(4-chlorophenoxy)-2,2-dimethyl-*N*-(1-methylpiperidin-4-yl)propanamide;
N-((2*R*,4*R*)-1,2-dimethylpiperidin-4-yl)-3-((3,5-dimethylpyridin-2-yl)oxy)-2,2-dimethylpropanamide; and

a pharmaceutically acceptable salt of any one of the aforementioned compounds.

36. A pharmaceutical composition comprising:
a compound or pharmaceutically acceptable salt as defined in any one of claims 1 to 35; and
a pharmaceutically acceptable excipient.
37. A compound or pharmaceutically acceptable salt as defined in any one of claims 1 to 35 for use as a medicament.
38. A compound or pharmaceutically acceptable salt as defined in any one of claims 1 to 35 for treatment of a disease, disorder or condition selected from Alzheimer's disease, depression, anxiety, schizophrenia, bipolar disorder, autism, epilepsy, pain, and hyperactivity disorder.
39. A method of treating a disease, disorder or condition in a subject, the method comprising administering to the subject a compound or pharmaceutically acceptable salt as defined in any one of claims 1 to 35, wherein the disease, disorder or condition is selected from Alzheimer's disease, depression, anxiety, schizophrenia, bipolar disorder, autism, epilepsy, pain, and hyperactivity disorder.
40. A combination comprising a compound or pharmaceutically acceptable salt as defined in any one of claims 1 to 35, and at least one additional pharmacologically active agent.

41. The combination according to claim 40, wherein the additional pharmacologically active agent is selected from beta-secretase inhibitors, gamma-secretase inhibitors, HMG-CoA reductase inhibitors, nonsteroidal anti-inflammatory drugs, vitamin E, anti-amyloid antibodies, antidepressants, antipsychotics, anxiolytics, and anticonvulsants.