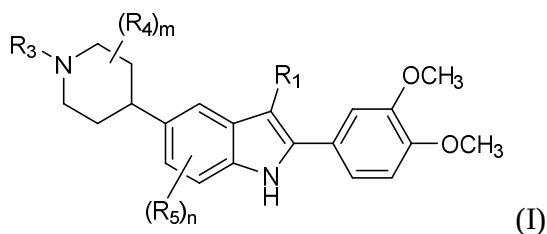


CLAIMS

1. A compound of Formula (I)



5 or a salt thereof, wherein:

R₁ is -CH₂CH₂CH₃, -CH(CH₃)₂, -CH₂CH(CH₃)₂, -CH₂CHF₂, -CH₂CF₃, -CH(CF₃)OH, -CH₂(cyclopropyl), or tetrahydropyranyl;

R₃ is:

(a) -L₁-A; or

- 10 (b) H, C₁₋₆ alkyl, C₁₋₃ fluoroalkyl, C₁₋₃ cyanoalkyl, C₁₋₆ hydroxyalkyl, C₁₋₃ hydroxy-fluoroalkyl, -CR_xR_xCR_x(OH)CR_x=CR_xR_x, -(CR_xR_x)₁₋₄O(C₁₋₃ alkyl), -(CR_xR_x)₁₋₄O(CR_xR_x)₁₋₃O(C₁₋₃ alkyl), -CH₂CH(OH)CH₂O(C₁₋₃ alkyl), -(CR_xR_x)₁₋₃S(C₁₋₃ alkyl), -(CH₂)₁₋₃C(O)OC(CH₃)₃, -(CR_xR_x)₀₋₃NR_xR_y, -(CR_xR_x)₀₋₃NR_x(C₁₋₄ hydroxyalkyl), -CH₂CH(OH)CH₂NR_xR_y,
- 15 -(CR_xR_x)₁₋₂NHC(O)(CR_xR_x)₁₋₄C≡CH, -C(O)H, -C(O)(C₁₋₆ alkyl), -C(O)(C₁₋₃ hydroxyalkyl), -C(O)(C₁₋₃ fluoroalkyl), -C(O)(C₁₋₃ chloroalkyl), -C(O)(C₁₋₃ cyanoalkyl), -(CR_xR_x)₀₋₃C(O)OH, -C(O)(CH₂)₀₋₂O(C₁₋₄ alkyl), -C(O)(CR_xR_x)₀₋₂O(CR_xR_x)₁₋₂O(C₁₋₃ alkyl), -C(O)CR_xR_xS(O)₂(C₁₋₃ alkyl), -C(O)CR_xR_xNR_xS(O)₂(C₁₋₃ alkyl), -C(O)CR_xR_xOC(O)(C₁₋₃ alkyl),
- 20 -C(O)(CR_xR_x)₀₋₃NR_yR_y, -C(O)(CR_xR_x)₀₋₁NR_x(C₁₋₃ cyanoalkyl), -C(O)(CR_xR_x)₀₋₂NR_y(C₁₋₆ hydroxyalkyl), -C(O)(CR_xR_x)₀₋₁NR_x(C₁₋₃ fluoroalkyl), -C(O)(CR_xR_x)₀₋₁NR_x(C₁₋₅ hydroxy-fluoroalkyl), -C(O)(CR_xR_x)₀₋₁NR_x(CH₂)₁₋₂O(C₁₋₃ hydroxyalkyl), -C(O)(CR_xR_x)₀₋₁NR_x(CH₂)₁₋₂NR_xC(O)(C₁₋₂ alkyl), -C(O)(CR_xR_x)₀₋₁NR_x((CR_xR_x)₁₋₂O(C₁₋₂ alkyl)), -C(O)CR_x(NH₂)(CR_xR_x)₁₋₄NR_xR_x,

- $-\text{C}(\text{O})\text{CR}_x(\text{NH}_2)(\text{CR}_x\text{R}_x)_{1-4}\text{NR}_x\text{C}(\text{O})\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-3}\text{NR}_x(\text{CH}_2)_{0-1}\text{C}(\text{O})(\text{C}_{1-3}$
 alkyl), $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-1}\text{NR}_x(\text{CH}_2)_{0-1}\text{C}(\text{O})(\text{C}_{1-3}$ cyanoalkyl),
 $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-1}\text{NR}_x(\text{CH}_2)_{1-2}\text{C}(\text{O})\text{NR}_y\text{R}_y$, $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{1-3}\text{C}(\text{O})\text{NR}_y\text{R}_y$,
 $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-1}\text{NR}_x(\text{CHR}_y(\text{CH}_2\text{OH}))$, $-(\text{CR}_x\text{R}_x)_{1-2}\text{C}(\text{O})\text{NR}_y\text{R}_y$,
 5 $-(\text{CR}_x\text{R}_x)_{1-2}\text{C}(\text{O})\text{NR}_y(\text{C}_{1-3}$ fluoroalkyl), $-(\text{CR}_x\text{R}_x)_{1-2}\text{C}(\text{O})\text{NR}_y(\text{C}_{1-4}$ hydroxyalkyl),
 $-(\text{CR}_x\text{R}_x)_{1-2}\text{C}(\text{O})\text{NR}_y(\text{C}_{1-3}$ cyanoalkyl), $-(\text{CR}_x\text{R}_x)_{1-2}\text{C}(\text{O})\text{NR}_x(\text{CH}_2)_{1-2}\text{O}(\text{C}_{1-3}$ alkyl),
 $-(\text{CR}_x\text{R}_x)_{1-2}\text{C}(\text{O})\text{NR}_x\text{CH}(\text{C}_{1-4}$ alkyl)(C_{1-3} hydroxyalkyl), $-(\text{CH}_2)_{1-2}\text{C}(\text{O})\text{NR}_x(\text{CH}_2)_{1-2}\text{C}(\text{O})\text{NR}_x\text{R}_x$,
 $-(\text{CH}_2)_{1-2}\text{C}(\text{O})\text{NR}_x(\text{CH}_2)_{1-2}\text{S}(\text{O})_2\text{OH}$,
 $-(\text{CH}_2)_{1-2}\text{C}(\text{O})\text{NR}_x(\text{CH}_2)_{1-2}\text{NR}_x\text{C}(\text{O})(\text{C}_{1-3}$ alkyl), $-(\text{CH}_2)_{1-2}\text{C}(\text{O})\text{NR}_x(\text{CH}_2)_{1-3}\text{NR}_x\text{R}_x$,
 10 $-(\text{CH}_2)_{1-2}\text{C}(\text{O})\text{N}(\text{CH}_2\text{CH}_3)(\text{CH}_2)_{1-3}\text{NR}_x\text{R}_x$, $-(\text{CH}_2)_{0-2}\text{S}(\text{O})_2(\text{C}_{1-4}$ alkyl),
 $-(\text{CH}_2)_{0-2}\text{S}(\text{O})_2(\text{C}_{1-3}$ fluoroalkyl), $-(\text{CH}_2)_{0-2}\text{S}(\text{O})_2\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})\text{C}(\text{O})\text{OH}$,
 $-\text{C}(\text{O})\text{C}(\text{O})\text{NR}_y\text{R}_y$, or $-\text{C}(\text{O})\text{C}(\text{O})\text{NR}_y(\text{CR}_x\text{R}_x)_{1-2}\text{NR}_y\text{R}_y$;
 L_1 is a bond, $-(\text{CR}_x\text{R}_x)_{1-2}-$, $-(\text{CR}_x\text{R}_x)_{1-2}\text{CR}_x(\text{OH})-$, $-(\text{CR}_x\text{R}_x)_{1-2}\text{O}-$, $-\text{CR}_x\text{R}_x\text{C}(\text{O})-$,
 $-(\text{CR}_x\text{R}_x)_2\text{NR}_x(\text{CR}_x\text{R}_x)_{0-1}-$, $-\text{CR}_x\text{R}_x\text{C}(\text{O})\text{NR}_x(\text{CR}_x\text{R}_x)_{0-4}-$, $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-3}-$,
 15 $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-2}\text{NR}_x(\text{CR}_x\text{R}_x)_{0-2}-$, $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-2}\text{N}(\text{C}_{1-2}$ hydroxyalkyl)($\text{CR}_x\text{R}_x)_{0-2}-$,
 $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-2}\text{NR}_x(\text{CR}_x\text{R}_x)_{1-2}\text{CR}_x(\text{OH})-$, $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{1-2}\text{C}(\text{O})\text{NR}_x-$,
 $-(\text{CR}_x\text{R}_x)_{0-2}\text{C}(\text{O})\text{NR}_x(\text{CR}_x\text{R}_x)_{1-2}\text{CR}_x(\text{OH})-$, $-(\text{CR}_x\text{R}_x)_{0-2}\text{C}(\text{O})\text{N}(\text{C}_{1-2}$
 hydroxyalkyl)($\text{CR}_x\text{R}_x)_{1-2}-$, $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-1}\text{O}-$, $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{1-2}\text{NHS}(\text{O})_2-$,
 $-\text{C}(\text{O})\text{CR}_x(\text{NH}_2)\text{CR}_x\text{R}_x-$, $-\text{C}(\text{O})\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-2}-$, $-\text{C}(\text{O})\text{NR}_x(\text{CR}_x\text{R}_x)_{1-2}-$, or $-\text{S}(\text{O})_2-$;
 20 A is 2-oxa-6-azaspiro[3,3]heptanyl, 4-oxaspiro[2.5]octanyl, 7-azaspiro[3.5]nonanyl,
 8-azabicyclo[3.2.1]octanyl, 8-oxa-3-azabicyclo[3.2.1]octanyl,
 9-azabicyclo[3.3.1]nonanyl, adamantanyl, azepanyl, azetidiny, C_{3-6} cycloalkyl,
 diazepanyl, dihydroinonyl, dihydropyrimidinonyl, dioxidoisothiazolidiny, dioxidothiazinanyl,
 dioxotetrahydrothiophenyl, dioxotetrahydrothiopyranyl,
 25 dioxothiomorpholinyl, furanyl, imidazolyl, imidazolidinonyl, indolyl, isoquinolinyl,
 isoxazolyl, morpholinyl, morpholinonyl, naphthalenyl, octahydrocyclopenta[b]pyranyl,
 oxazolidinonyl, oxadiazolyl, oxetanyl, oxazolyl, phenyl, piperidiny, piperidinonyl,
 piperaziny, piperazinonyl, pyraziny, pyrazolyl, pyridazinonyl, pyridinonyl, pyridiny,

pyrimidinyl, pyrrolidinonyl, pyrrolidinyl, pyrrolyl, quinolinyl, quinolizinonyl, tetrahydrofuranyl, tetrahydropyranyl, tetrazolyl, thiadiazolyl, thiazolyl, or triazolyl, each substituted with $-L_2-R_a$ and zero to 4 R_b ;

L_2 is a bond or $-CR_xR_x-$;

5 R_a is:

(a) H, F, Cl, $-CN$, $-OH$, C_{1-6} alkyl, C_{1-3} fluoroalkyl, C_{1-5} hydroxyalkyl, $-(CH_2)_{0-4}O(C_{1-3}$ alkyl), $-(CR_xR_x)_{1-3}S(C_{1-3}$ alkyl), $-(CR_xR_x)_{1-3}NHC(O)O(C_{1-4}$ alkyl), $-(CR_xR_x)_{1-3}NR_yR_y$, $-(CR_xR_x)_{1-3}C(O)NR_yR_y$, $-O(C_{1-3}$ fluoroalkyl), $-S(O)_2NR_xR_x$, $-O(CR_xR_x)_{1-3}NR_xR_x$, $-NHS(O)_2(C_{1-3}$ alkyl), $-NR_xR_x$, $-NR_x(C_{1-4}$ alkyl), $-NR_xC(O)(C_{1-4}$ alkyl), $-(CR_xR_x)_{0-3}C(O)OH$, $-C(O)(C_{1-5}$ alkyl), $-C(O)(C_{1-3}$ fluoroalkyl), $-C(O)CH_2N(C_{1-3}$ alkyl)₂, $-C(O)O(C_{1-4}$ alkyl), $-C(O)NH(C_{1-3}$ cyanoalkyl), $-C(O)NR_yR_y$, $-C(O)NR_xCH_2C(O)NR_xR_x$, or $-C(O)NR_xCH_2CH_2NHC(O)(C_{1-3}$ alkyl);

10

(b) C_{3-6} cycloalkyl or $-C(O)NH(C_{3-6}$ cycloalkyl), wherein each cycloalkyl is substituted with zero to 2 substituents independently selected from $-OH$, C_{1-3} alkyl, C_{1-3} hydroxyalkyl, C_{1-3} fluoroalkyl, and $-C(O)O(C_{1-3}$ alkyl); or

15

(c) A_1 , $-CH_2A_1$, $-C(O)A_1$, $-NR_xA_1$, or $-C(O)NR_xA_1$, wherein A_1 is furanyl, imidazolyl, indolyl, isoxazolyl, morpholinyl, octahydropyrrolo[3,4-c]pyrrolyl, oxazolyl, oxetanyl, phenyl, piperazinyl, piperidinyl, pyrazolyl, pyridinyl, pyrimidinyl, pyrrolidinyl, pyrrolyl, tetrahydrofuranyl, tetrahydropyranyl, thiadiazolyl, thiazolyl, thiophenyl, or triazolyl, each substituted with zero to three substituents independently selected from $-OH$, C_{1-3} alkyl, C_{1-3} hydroxyalkyl, $-C(O)(C_{1-2}$ alkyl), $-C(O)O(C_{1-3}$ alkyl), $-NR_xR_x$, phenyl, trifluoromethyl-phenyl, $-CH_2$ (bromophenyl), and $-CH_2CH_2$ (pyrrolidinyl);

20

each R_b is independently F, $-CH_3$, $-CF_3$, or $-OCH_3$;

each R_x is independently H or $-CH_3$;

25

each R_y is independently H or C_{1-6} alkyl;

R_z is H, C_{1-2} alkyl, or C_{1-2} fluoroalkyl;

each R_4 is independently F or $-OH$; or two R_4 attached to the same carbon atom form $=O$;

R_5 is F, Cl, $-CN$, C_{1-2} alkyl, C_{1-2} fluoroalkyl, or $-OCH_3$;

m is zero, 1, 2, 3, or 4; and

n is zero, 1, or 2.

2. The compound according to claim 1 or a salt thereof, wherein:

- 5 R_1 is $-\text{CH}_2\text{CH}_2\text{CH}_3$, $-\text{CH}(\text{CH}_3)_2$, $-\text{CH}_2\text{CH}(\text{CH}_3)_2$, $-\text{CH}_2\text{CHF}_2$, $-\text{CH}_2\text{CF}_3$, $-\text{CH}(\text{CF}_3)\text{OH}$,
 $-\text{CH}_2(\text{cyclopropyl})$, or tetrahydropyranyl;

R_3 is:

(a) $-\text{L}_1-\text{A}$; or

- (b) H, C_{1-6} alkyl, C_{1-3} fluoroalkyl, C_{1-5} hydroxyalkyl, $-(\text{CR}_x\text{R}_x)_{1-4}\text{O}(\text{C}_{1-3} \text{ alkyl})$,
10 $-(\text{CR}_x\text{R}_x)_{1-3}\text{S}(\text{C}_{1-3} \text{ alkyl})$, $-(\text{CH}_2)_{1-3}\text{C}(\text{O})\text{OC}(\text{CH}_3)_3$, $-(\text{CH}_2)_{1-3}\text{NR}_x\text{R}_x$,
 $-\text{CHR}_x\text{CH}_2\text{NR}_x(\text{C}_{1-4} \text{ alkyl})$, $-(\text{CH}_2)_{1-3}\text{NR}_x(\text{C}_{1-3} \text{ alkyl})$,
 $-(\text{CR}_x\text{R}_x)_{1-2}\text{NHC}(\text{O})(\text{CH}_2)_{1-4}\text{C}\equiv\text{CH}$, $-\text{S}(\text{O})_2\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})\text{O}(\text{C}_{1-4} \text{ alkyl})$, $-\text{C}(\text{O})(\text{C}_{1-3} \text{ alkyl})$,
 $-\text{C}(\text{O})(\text{C}_{1-3} \text{ hydroxyalkyl})$, $-\text{C}(\text{O})(\text{C}_{1-3} \text{ fluoroalkyl})$, $-\text{C}(\text{O})(\text{C}_{1-3} \text{ chloroalkyl})$,
 $-\text{C}(\text{O})\text{CH}_2\text{O}(\text{C}_{1-3} \text{ alkyl})$, $-\text{C}(\text{O})\text{CH}_2\text{S}(\text{O})_2(\text{C}_{1-3} \text{ alkyl})$, $-\text{C}(\text{O})(\text{C}_{1-3} \text{ hydroxyalkyl})$,
15 $-\text{C}(\text{O})\text{CR}_x\text{R}_x\text{OC}(\text{O})(\text{C}_{1-3} \text{ alkyl})$, $-\text{C}(\text{O})(\text{tetrahydrofuranyl})$, $-\text{C}(\text{O})(\text{tetrahydropyranyl})$,
 $-\text{C}(\text{O})(\text{piperidiny})$, $-\text{C}(\text{O})(\text{ethoxypiperidiny})$, $-\text{C}(\text{O})\text{NR}_x(\text{C}_{1-3} \text{ cyanoalkyl})$,
 $-\text{C}(\text{O})\text{NR}_x(\text{C}_{1-3} \text{ alkyl})$, $-\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CH}_2\text{NR}_x\text{C}(\text{O})(\text{C}_{1-2} \text{ alkyl})$,
 $-\text{C}(\text{O})\text{N}(\text{C}_{1-3} \text{ alkyl})_2$, $-\text{C}(\text{O})\text{NR}_x\text{C}(\text{R}_x)_3$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-3} \text{ cyanoalkyl})$,
 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{C}(\text{O})(\text{C}_{1-3} \text{ alkyl})$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{C}(\text{O})\text{N}(\text{C}_{1-3} \text{ alkyl})_2$,
20 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-2} \text{ alkyl})$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-3} \text{ hydroxyalkyl})$,
 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-3} \text{ fluoroalkyl})$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{CH}_2\text{O}(\text{C}_{1-3} \text{ hydroxyalkyl})$,
 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{N}(\text{C}_{1-2} \text{ alkyl})_2$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{CH}(\text{CH}_2\text{OH})(\text{C}_{1-4} \text{ alkyl}))$,
 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-5} \text{ alkyl})$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-5} \text{ hydroxy-fluoroalkyl})$,
 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-6} \text{ hydroxyalkyl})$, $-\text{C}(\text{O})(\text{CH}_2)_{1-3}\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})(\text{CH}_2)_{1-3}\text{NR}_x(\text{C}_{1-3} \text{ alkyl})$,
25 $-\text{C}(\text{O})\text{CH}_2\text{CH}(\text{CH}_3)\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{R}_x$, $-\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{R}_x$,
 $-\text{CH}_2\text{C}(\text{O})\text{NR}_x(\text{C}_{1-3} \text{ cyanoalkyl})$, $-\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{R}_x$, $-\text{CH}_2\text{C}(\text{O})\text{NR}_x(\text{C}_{1-3} \text{ hydroxyalkyl})$,
 $-\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CH}_2\text{S}(\text{O})_2\text{OH}$, $-\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CH}_2\text{NR}_x\text{C}(\text{O})(\text{C}_{1-3} \text{ alkyl})$,
 $-\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_2\text{CH}_3)\text{CH}_2\text{CH}_2\text{NR}_x\text{R}_x$, $-\text{CH}_2\text{C}(\text{O})\text{NR}_x(\text{C}_{1-3} \text{ alkyl})$,

$-\text{CH}_2\text{C}(\text{O})\text{NR}_x(\text{CH}_2)_{1-3}\text{NR}_x\text{R}_x$, $-\text{CH}_2\text{C}(\text{O})\text{NR}_x(\text{C}_{1-5} \text{ alkyl})$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{C}(\text{O})(\text{C}_{1-3} \text{ cyanoalkyl})$, or $-\text{C}(\text{O})(\text{CH}_2)_{1-3}\text{NR}_x\text{C}(\text{O})(\text{C}_{1-3} \text{ alkyl})$;

L_1 is a bond, $-\text{CR}_x\text{R}_x-$, $-\text{CR}_x\text{R}_x\text{C}(\text{O})-$, $-(\text{CR}_x\text{R}_x)_2\text{NR}_x(\text{CR}_x\text{R}_x)_{0-1}-$,

$-\text{CR}_x\text{R}_x\text{C}(\text{O})\text{NR}_x(\text{CR}_x\text{R}_x)_{0-4}-$, $-\text{C}(\text{O})(\text{CR}_x\text{R}_x)_{0-3}-$, $-\text{C}(\text{O})\text{CR}_x\text{R}_x\text{NR}_x-$, $-\text{C}(\text{O})\text{CR}_x\text{R}_x\text{O}-$,
5 or $-\text{C}(\text{O})\text{NR}_x(\text{CR}_x\text{R}_x)_{1-2}-$;

A is a ring selected from adamantanyl, 8-azabicyclo[3.2.1]octanyl, azepanyl, C_{3-6} cycloalkyl, diazepanyl, furanyl, imidazolyl, indolyl, isoquinoliny, morpholiny, naphthalenyl, oxetanyl, phenyl, piperaziny, piperidiny, pyraziny, pyrazolyl, pyridiny, pyrrolidinonyl, pyrrolidiny, pyrrolyl, quinoliny, tetrahydropyranyl, tetrazolyl,
10 thiadiazolyl, and thiazolyl, each substituted with $-\text{L}_2-\text{R}_a$ and zero to 4 R_b ;

L_2 is a bond or $-\text{CR}_x\text{R}_x-$;

R_a is:

(a) H, $-\text{CN}$, $-\text{OH}$, C_{1-6} alkyl, C_{1-3} fluoroalkyl, C_{1-5} hydroxyalkyl, $-(\text{CH}_2)_{1-4}\text{O}(\text{C}_{1-3} \text{ alkyl})$,
 $-(\text{CR}_x\text{R}_x)_{1-3}\text{S}(\text{C}_{1-3} \text{ alkyl})$, $-(\text{CR}_x\text{R}_x)_{1-3}\text{NHC}(\text{O})\text{O}(\text{C}_{1-4} \text{ alkyl})$, $-(\text{CR}_x\text{R}_x)_{1-3}\text{NH}_2$,
15 $-(\text{CR}_x\text{R}_x)_{1-3}\text{NR}_x(\text{C}_{1-4} \text{ alkyl})$, $-\text{O}(\text{C}_{1-3} \text{ fluoroalkyl})$, $-\text{S}(\text{O})_2\text{NR}_x\text{R}_x$, $-\text{O}(\text{CR}_x\text{R}_x)_{1-3}\text{NR}_x\text{R}_x$,
 $-\text{NHS}(\text{O})_2(\text{C}_{1-3} \text{ alkyl})$, $-\text{NR}_x\text{R}_x$, $-\text{NR}_x(\text{C}_{1-4} \text{ alkyl})$, $-(\text{CR}_x\text{R}_x)_{1-3}\text{C}(\text{O})\text{OH}$, -
 $(\text{CR}_x\text{R}_x)_{1-3}\text{C}(\text{O})\text{NH}(\text{C}_{1-4} \text{ alkyl})$, $-\text{C}(\text{O})\text{OH}$, $-\text{C}(\text{O})(\text{C}_{1-4} \text{ alkyl})$, $-\text{C}(\text{O})\text{CH}_2\text{N}(\text{C}_{1-3} \text{ alkyl})_2$,
 $-\text{C}(\text{O})\text{O}(\text{C}_{1-4} \text{ alkyl})$, $-\text{C}(\text{O})\text{NH}(\text{C}_{1-3} \text{ cyanoalkyl})$, $-\text{C}(\text{O})\text{NR}_x(\text{C}_{1-4} \text{ alkyl})$, $-\text{C}(\text{O})\text{N}(\text{C}_{1-3} \text{ alkyl})_2$,
 $-\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{R}_x$, or $-\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CH}_2\text{NHC}(\text{O})(\text{C}_{1-3} \text{ alkyl})$;

(b) C_{3-6} cycloalkyl or $-\text{C}(\text{O})\text{NH}(\text{C}_{3-6} \text{ cycloalkyl})$, wherein each cycloalkyl is substituted
20 with zero to 2 substituents independently selected from $-\text{OH}$, C_{1-3} alkyl, C_{1-3} hydroxyalkyl, C_{1-3} fluoroalkyl, and $-\text{C}(\text{O})\text{O}(\text{C}_{1-3} \text{ alkyl})$; or

(c) A_1 , $-\text{CH}_2\text{A}_1$, $-\text{C}(\text{O})\text{A}_1$, or $-\text{C}(\text{O})\text{NHA}_1$, wherein A_1 is furanyl, imidazolyl, indolyl, isoxazolyl, octahydropyrrolo[3,4-c]pyrrolyl, oxazolyl, oxetanyl, phenyl, piperaziny, piperidiny, pyrazolyl, pyridiny, pyrimidiny, pyrrolidiny, pyrrolyl, tetrahydrofuranyl, tetrahydropyranyl, thiadiazolyl, thiazolyl, thiophenyl, or triazolyl, each substituted with
25 zero to three substituents independently selected from $-\text{OH}$, C_{1-3} alkyl, C_{1-3} hydroxyalkyl, $-\text{C}(\text{O})(\text{C}_{1-2} \text{ alkyl})$, $-\text{C}(\text{O})\text{O}(\text{C}_{1-3} \text{ alkyl})$, $-\text{NR}_x\text{R}_x$, phenyl, trifluoromethyl-

phenyl, $-\text{CH}_2(\text{bromophenyl})$, and $-\text{CH}_2\text{CH}_2(\text{pyrrolidinyl})$;

each R_4 is independently F or $-\text{OH}$; or two R_4 attached to the same carbon atom form $=\text{O}$;

R_5 is F, $-\text{CH}_3$, or $-\text{OCH}_3$;

R_b is $-\text{CH}_3$;

5 each R_x is independently H or $-\text{CH}_3$;

m is zero, 1, or 2; and

n is zero or 1.

3. The compound according to claim 1 or a salt thereof, wherein:

10 R_1 is $-\text{CH}_2\text{CH}_2\text{CH}_3$, $-\text{CH}(\text{CH}_3)_2$, $-\text{CH}_2\text{CH}(\text{CH}_3)_2$, $-\text{CH}_2\text{CHF}_2$, $-\text{CH}_2\text{CF}_3$, $-\text{CH}(\text{CF}_3)\text{OH}$,
 $-\text{CH}_2(\text{cyclopropyl})$, or tetrahydropyranyl;

R_3 is:

(a) $-\text{L}_1-\text{A}$; or

(b) H, C_{1-6} alkyl, C_{1-3} fluoroalkyl, C_{1-5} hydroxyalkyl, $-(\text{CH}_2)_{1-4}\text{OCH}_3$, $-(\text{CR}_x\text{R}_x)_{1-3}\text{SCH}_3$,
15 $-\text{CH}_2\text{CH}_2\text{NHC}(\text{O})\text{OC}(\text{CH}_3)_3$, $-\text{CH}_2\text{CH}_2\text{NR}_x\text{R}_x$, $-\text{CHR}_x\text{CH}_2\text{NR}_x(\text{C}_{1-4} \text{ alkyl})$,
 $-\text{CH}_2\text{CH}_2\text{CH}_2\text{NH}(\text{CH}_3)$, $-(\text{CH}_2)_{1-2}\text{NHC}(\text{O})(\text{CH}_2)_{1-4}\text{C}\equiv\text{CH}$, $-\text{S}(\text{O})_2\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})\text{O}(\text{C}_{1-4}$
alkyl), $-\text{C}(\text{O})(\text{C}_{1-2} \text{ alkyl})$, $-\text{C}(\text{O})(\text{C}_{1-3} \text{ hydroxyalkyl})$, $-\text{C}(\text{O})(\text{C}_{1-2} \text{ chloroalkyl})$, $-\text{C}(\text{O})(\text{C}_{1-3}$
3 fluoroalkyl), $-\text{C}(\text{O})\text{CH}_2\text{O}(\text{C}_{1-3} \text{ alkyl})$, $-\text{C}(\text{O})\text{CH}_2\text{S}(\text{O})_2(\text{C}_{1-2} \text{ alkyl})$,
 $-\text{C}(\text{O})\text{CR}_x\text{R}_x\text{OC}(\text{O})(\text{C}_{1-2} \text{ alkyl})$, $-\text{C}(\text{O})(\text{tetrahydrofuranyl})$, $-\text{C}(\text{O})(\text{ethoxypiperidinyl})$,
20 $-\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CN}$, $-\text{C}(\text{O})\text{NH}(\text{C}_{1-3} \text{ alkyl})$, $-\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{C}(\text{O})\text{NH}_2$,
 $-\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CH}_2\text{NHC}(\text{O})(\text{C}_{1-2} \text{ alkyl})$, $-\text{C}(\text{O})\text{N}(\text{C}_{1-3} \text{ alkyl})_2$, $-\text{C}(\text{O})\text{NR}_x\text{C}(\text{R}_x)_3$,
 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{CN}$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{C}(\text{O})(\text{C}_{1-2} \text{ alkyl})$,
 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{C}(\text{O})\text{N}(\text{C}_{1-2} \text{ alkyl})_2$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-2} \text{ alkyl})$,
 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{CH}_2\text{OH}$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-2} \text{ fluoroalkyl})$,
25 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{CH}_2\text{OCH}_2\text{CH}_2\text{OH}$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{N}(\text{C}_{1-2} \text{ alkyl})_2$,
 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{CH}(\text{CH}_2\text{OH})(\text{C}_{1-4} \text{ alkyl}))$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-4} \text{ hydroxyalkyl})$,
 $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-5} \text{ alkyl})$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{CH}_2\text{CHFC}(\text{CH}_3)_2\text{OH}$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x(\text{C}_{1-6}$
hydroxyalkyl), $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{NR}_x\text{R}_x$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{NR}_x\text{R}_x$,

- $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{NR}_x(\text{C}_{1-3} \text{ alkyl}), -\text{C}(\text{O})\text{CH}_2\text{CH}(\text{CH}_3)\text{NR}_x\text{R}_x,$
 $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{R}_x, -\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{R}_x, -\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CN},$
 $-\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{R}_x, -\text{CH}_2\text{C}(\text{O})\text{NR}_x(\text{C}_{1-3} \text{ hydroxyalkyl}),$
 $-\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CH}_2\text{S}(\text{O})_2\text{OH}, -\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CH}_2\text{NR}_x\text{R}_x,$
5 $-\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CH}_2\text{NR}_x\text{C}(\text{O})(\text{C}_{1-2} \text{ alkyl}), -\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_2\text{CH}_3)\text{CH}_2\text{CH}_2\text{NR}_x\text{R}_x,$
 $-\text{CH}_2\text{C}(\text{O})\text{NR}_x(\text{C}_{1-3} \text{ alkyl}), -\text{CH}_2\text{C}(\text{O})\text{NR}_x\text{CH}_2\text{CH}_2\text{CH}_2\text{NR}_x\text{R}_x, -\text{CH}_2\text{C}(\text{O})\text{NR}_x(\text{C}_{1-5}$
 $\text{alkyl}), -\text{C}(\text{O})\text{CH}_2\text{NR}_x\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{CN}, \text{ or } -\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{NR}_x\text{C}(\text{O})(\text{C}_{1-2} \text{ alkyl});$
- L_1 is a bond, $-\text{CHR}_x-$, $-\text{CHR}_x\text{C}(\text{O})-$, $-(\text{CH}_2)_2\text{NR}_x(\text{CH}_2)_{0-1}-$, $-\text{CH}_2\text{C}(\text{O})\text{NR}_x(\text{CH}_2)_{0-4}-$,
 $-\text{C}(\text{O})(\text{CH}_2)_{0-3}-$, $-\text{C}(\text{O})\text{CH}_2\text{NR}_x-$, $-\text{C}(\text{O})\text{CH}_2\text{O}-$, or $-\text{C}(\text{O})\text{NR}_x(\text{CH}_2)_{1-2}-$;
- 10 A is a ring selected from adamantanyl, azabicyclo[3.2.1]octanyl, azepanyl, C_{3-6} cycloalkyl,
diazepanyl, furanyl, imidazolyl, indolyl, isoquinolinyl, morpholinyl, naphthalenyl,
oxetanyl, phenyl, piperazinyl, piperidinyl, pyrazinyl, pyrazolyl, pyridinyl, pyrrolidinonyl,
pyrrolidinyl, pyrrolyl, quinolinyl, tetrahydropyranlyl, tetrazolyl, thiadiazolyl, and
thiazolyl, each substituted with $-\text{L}_2-\text{R}_a$ and zero to 4 R_b ;
- 15 and
- R_a is:
- (a) H , $-\text{CN}$, $-\text{OH}$, C_{1-6} alkyl, C_{1-3} fluoroalkyl, C_{1-5} hydroxyalkyl, $-(\text{CH}_2)_{1-4}\text{OCH}_3$,
 $-(\text{CR}_x\text{R}_x)_{1-3}\text{SCH}_3$, $-\text{CH}_2\text{CH}_2\text{NHC}(\text{O})\text{OC}(\text{CH}_3)_3$, $-\text{CH}_2\text{CH}_2\text{NH}_2$, $-\text{CHR}_x\text{CH}_2\text{NR}_x(\text{C}_{1-4}$
 $\text{alkyl}), -\text{CH}_2\text{CH}_2\text{CH}_2\text{NH}(\text{CH}_3)$, $-\text{OCF}_3$, $-\text{S}(\text{O})_2\text{NR}_x\text{R}_x$, $-\text{OCH}_2\text{CH}_2\text{CH}_2\text{NR}_x\text{R}_x$,
20 $-\text{NHS}(\text{O})_2(\text{C}_{1-2} \text{ alkyl}), -\text{NR}_x\text{R}_x$, $-\text{NR}_x(\text{C}_{1-4} \text{ alkyl}), -\text{CH}_2\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{OH}$,
 $-\text{CH}_2\text{C}(\text{O})\text{NH}(\text{C}_{1-3} \text{ alkyl}), -\text{C}(\text{O})\text{OH}$, $-\text{C}(\text{O})(\text{C}_{1-3} \text{ alkyl}), -\text{C}(\text{O})\text{CH}_2\text{N}(\text{C}_{1-3} \text{ alkyl})_2$,
 $-\text{C}(\text{O})\text{O}(\text{C}_{1-3} \text{ alkyl}), -\text{C}(\text{O})\text{NHCH}_2\text{CN}$, $-\text{C}(\text{O})\text{NR}_x(\text{C}_{1-4} \text{ alkyl}), -\text{C}(\text{O})\text{N}(\text{C}_{1-3} \text{ alkyl})_2$,
 $-\text{C}(\text{O})\text{NHCH}_2\text{C}(\text{O})\text{NR}_x\text{R}_x$, or $-\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{NHC}(\text{O})(\text{C}_{1-3} \text{ alkyl});$
- (b) C_{3-6} cycloalkyl or $-\text{C}(\text{O})\text{NH}(\text{C}_{3-6} \text{ cycloalkyl})$, wherein each cycloalkyl is substituted
25 with zero to 1 substituent selected from $-\text{OH}$, C_{1-3} alkyl, C_{1-3} hydroxyalkyl, C_{1-2}
fluoroalkyl, and $-\text{C}(\text{O})\text{O}(\text{C}_{1-3} \text{ alkyl})$; or
- (c) A_1 , $-\text{CH}_2\text{A}_1$, $-\text{C}(\text{O})\text{A}_1$, or $-\text{C}(\text{O})\text{NHA}_1$, wherein A_1 is furanyl, imidazolyl, indolyl,
isoxazolyl, octahydropyrrolo[3,4-c]pyrrolyl, oxazolyl, oxetanyl, phenyl, piperazinyl,

piperidinyl, pyrazolyl, pyridinyl, pyrimidinyl, pyrrolidinyl, pyrrolyl, tetrahydrofuranyl, tetrahydropyranyl, thiadiazolyl, thiazolyl, thiophenyl, or triazolyl, each substituted with zero to three substituents independently selected from $-OH$, C_{1-3} alkyl, C_{1-3} hydroxyalkyl, $-C(O)(C_{1-2} \text{ alkyl})$, $-C(O)O(C_{1-3} \text{ alkyl})$, $-NR_xR_x$, phenyl, trifluoromethyl-phenyl, $-CH_2(\text{bromophenyl})$, and $-CH_2CH_2(\text{pyrrolidinyl})$.

4. The compound according to claim 1 or a salt thereof, wherein:

R_1 is $-CH_2CH_2CH_3$, $-CH(CH_3)_2$, $-CH_2CH(CH_3)_2$, $-CH_2CHF_2$, $-CH_2CF_3$, $-CH(CF_3)OH$, $-CH_2(\text{cyclopropyl})$, or tetrahydropyranyl;

10 R_3 is:

(a) $-L_1-A$; or

(b) H , $-CH_3$, $-CH_2CH(CH_3)_2$, $-CH_2CH(CH_3)CF_3$, $-CH_2CH_2CH_2OH$, $-CH(CH_3)CH_2OH$, $-CH_2CH_2NHC(O)OC(CH_3)_3$, $-CH_2CH_2NH_2$, $-CH_2CH_2NH(CH_3)$, $-CH_2CH_2N(CH_3)_2$, $-CH_2CH_2N(CH_3)(CH_2CH_3)$, $-CH_2CH_2N(CH_3)(CH_2CH(CH_3)_2)$, $-CH_2CH_2CH_2NH(CH_3)$, $-CH(CH_3)CH_2N(CH_3)_2$, $-CH_2CH_2NHC(O)(CH_2)_4C\equiv CH$, $-C(O)OC(CH_3)_3$, $-C(O)CH_3$, $-C(O)CH_2OH$, $-C(O)CH_2Cl$, $-C(O)CH_2CF_3$, $-C(O)CH_2CH_2CF_3$, $-C(O)CH_2OCH(CH_3)_2$, $-C(O)CH_2S(O)_2CH_3$, $-C(O)CH_2CH(CH_3)OH$, $-C(O)C(CH_3)_2OC(O)CH_3$, $-C(O)NHCH_2CH_3$, $-C(O)NHCH(CH_3)_2$, $-C(O)CH_2NHCH_3$, $-C(O)CH_2NHCH_2CN$, $-C(O)CH_2NHCH_2C(O)CH_3$, $-C(O)CH_2N(CH_3)CH_2C(O)N(CH_2CH_3)_2$, $-C(O)CH_2N(CH_3)_2$, $-C(O)CH_2NHCH_2CH_3$, $-C(O)CH_2NHCH_2CH_2OH$, $-C(O)CH_2NHCH_2CF_3$, $-C(O)CH_2NHCH_2CH_2OCH_2CH_2OH$, $-C(O)CH_2N(CH_3)CH_2CH_2C(O)N(CH_2CH_3)_2$, $-C(O)CH_2NHCH(CH_2OH)CH_2CH_3$, $-C(O)CH_2NHCH(CH_2OH)_2$, $-C(O)CH_2N(CH_3)CH_2CH_2OH$, $-C(O)CH_2NHCH_2CH(CH_3)OH$, $-C(O)CH_2NHCH_2CH(OH)CH_2OH$, $-C(O)CH_2N(CH_3)CH(CH_3)_2$, $-C(O)CH_2NHCH_2CH(CH_3)_2$, $-C(O)CH_2NHCH(CH_3)CH(CH_3)_2$, $-C(O)CH_2NHCH_2CH_2CH(CH_3)OH$, $-C(O)CH_2NH(CH_2OH)(CH_2CH_2CH_2CH_3)$, $-C(O)CH_2NHCH(CH_2OH)CH_2CH(CH_3)_2$, $-C(O)CH_2NH(CH(CH_2CH_3)_2)$,

- $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CHFC}(\text{CH}_3)_2\text{OH}$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{C}(\text{CH}_3)_2\text{CH}_2\text{OH}$,
 $-\text{C}(\text{O})\text{CH}_2\text{NHCH}(\text{CH}_2\text{OH})\text{CH}_2\text{CH}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$,
 $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{NHCH}(\text{CH}_3)_2$,
 $-\text{C}(\text{O})\text{CH}_2\text{CH}(\text{CH}_3)\text{NH}_2$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{NH}_2$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_3)_2$,
5 $-\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_3)_2$, $-\text{CH}_2\text{C}(\text{O})\text{NHC}(\text{CH}_3)_2$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CN}$,
 $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{C}(\text{O})\text{NH}_2$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{OH}$,
 $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{S}(\text{O})_2\text{OH}$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$,
 $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{NHC}(\text{O})\text{CH}_3$, $-\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_3)\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$,
 $-\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_2\text{CH}_3)\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}(\text{CH}_3)_2$,
10 $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{CH}_2\text{OH}$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$,
 $-\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_3)\text{CH}_2\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}(\text{CH}_3)(\text{CH}_2\text{CH}_3)$,
 $-\text{CH}_2\text{C}(\text{O})\text{NHCH}(\text{CH}_2\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{NHC}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{CN}$,
 $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{NHC}(\text{O})\text{CH}_3$, or $-\text{CH}_2\text{CH}_2\text{NHC}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2\text{C}\equiv\text{CH}$;
 L_1 is a bond, $-\text{CHR}_x-$, $-\text{CH}_2\text{C}(\text{O})-$, $-(\text{CH}_2)_2\text{NR}_x(\text{CH}_2)_{0-1}-$, $-\text{CH}_2\text{C}(\text{O})\text{NR}_x(\text{CH}_2)_{0-4}-$,
15 $-\text{C}(\text{O})(\text{CH}_2)_{0-3}-$, $-\text{C}(\text{O})\text{CH}_2\text{NH}-$, $-\text{C}(\text{O})\text{CH}_2\text{O}-$, or $-\text{C}(\text{O})\text{NH}(\text{CH}_2)_{1-2}-$;
A is a ring selected from adamantanyl, azabicyclo[3.2.1]octanyl, azepanyl, cyclohexyl,
cyclopentyl, cyclopropyl, diazepanyl, furanyl, imidazolyl, indolyl, isoquinolinyl,
morpholinyl, naphthalenyl, oxetanyl, phenyl, piperazinyl, piperidiny, pyrazinyl,
pyrazolyl, pyridinyl, pyrrolidinonyl, pyrrolidinyl, pyrrolyl, quinolinyl,
20 tetrahydropyranlyl, tetrazolyl, thiadiazolyl, and thiazolyl, each substituted with $-\text{L}_2-\text{R}_a$
and zero to 4 R_b ;
 L_2 is a bond or $-\text{CHR}_x-$;
 R_a is:
(a) H, $-\text{CN}$, $-\text{OH}$, C_{1-6} alkyl, $-\text{CF}_3$, $-\text{CH}_2\text{CH}_2\text{CF}_3$, $-\text{CH}_2\text{OH}$, $-\text{C}(\text{CH}_3)_2\text{OH}$,
25 $-\text{CH}_2\text{CH}_2\text{CH}(\text{CH}_3)\text{OH}$, $-\text{CH}_2\text{C}(\text{CH}_3)_2\text{CH}_2\text{OH}$, $-\text{CH}_2\text{CH}(\text{OH})\text{CH}_2\text{OH}$,
 $-\text{CH}(\text{OH})\text{CH}_2\text{CH}_2\text{CH}_2\text{OH}$, $-\text{CH}_2\text{CH}_2\text{CH}_2\text{SCH}_3$, $-\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2\text{OCH}_3$,
 $-\text{CH}_2\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{OH}$, $-\text{OCF}_3$, $-\text{OCH}_2\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{NH}_2$, $-\text{N}(\text{CH}_3)_2$,
 $-\text{N}(\text{CH}_3)(\text{CH}_2\text{CH}(\text{CH}_3)_2)$, $-\text{NH}(\text{CH}_2\text{CH}(\text{CH}_3)_2)$, $-\text{S}(\text{O})_2\text{NH}_2$, $-\text{NHS}(\text{O})_2\text{CH}_3$,

$-\text{CH}_2\text{C}(\text{O})\text{NHCH}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{OH}$, $-\text{C}(\text{O})\text{CH}_3$, $-\text{C}(\text{O})\text{OCH}_2\text{CH}_3$, $-\text{C}(\text{O})\text{CH}_2\text{N}(\text{CH}_3)_2$,
 $-\text{C}(\text{O})\text{NHCH}_2\text{CN}$, $-\text{C}(\text{O})\text{NHCH}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{N}(\text{C}_{1-2}\text{ alkyl})_2$, $-\text{C}(\text{O})\text{N}(\text{CH}_3)\text{CH}(\text{CH}_3)_2$,
 $-\text{C}(\text{O})\text{NHCH}_2\text{C}(\text{O})\text{NH}_2$, or $-\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{NHC}(\text{O})\text{CH}_3$;

(b) C_{3-6} cycloalkyl or $-\text{C}(\text{O})\text{NH}(\text{C}_{3-6}\text{ cycloalkyl})$, wherein each cycloalkyl is substituted
5 with zero to 1 substituent selected from $-\text{OH}$, $-\text{CH}_3$, $-\text{CH}_2\text{OH}$, $-\text{CF}_3$, and
 $-\text{C}(\text{O})\text{OCH}_2\text{CH}_3$; or

(c) A_1 , $-\text{CH}_2\text{A}_1$, $-\text{C}(\text{O})\text{A}_1$, or $-\text{C}(\text{O})\text{NHA}_1$, wherein A_1 is furanyl, imidazolyl, indolyl,
isoxazolyl, octahydropyrrolo[3,4-c]pyrrolyl, oxazolyl, oxetanyl, phenyl, piperazinyl,
piperidinyl, pyrazolyl, pyridinyl, pyrimidinyl, pyrrolidinyl, pyrrolyl, tetrahydrofuranyl,
10 tetrahydropyranlyl, thiadiazolyl, thiazolyl, thiophenyl, or triazolyl, each substituted with
zero to three substituents independently selected from $-\text{OH}$, $-\text{CH}_3$, $-\text{CH}(\text{CH}_3)_2$,
 $-\text{CH}_2\text{CH}_2\text{OH}$, $-\text{C}(\text{O})\text{CH}_3$, $-\text{C}(\text{O})\text{OCH}_2\text{CH}_3$, $-\text{N}(\text{CH}_3)_2$, phenyl, trifluoromethyl-phenyl,
 $-\text{CH}_2(\text{bromophenyl})$, and $-\text{CH}_2\text{CH}_2(\text{pyrrolidinyl})$; and

each R_4 is independently F or $-\text{OH}$; or two R_4 attached to the same carbon atom form $=\text{O}$.

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5. The compound according to claim 1 or a salt thereof, wherein:

R_3 is $-\text{L}_1-\text{A}$.

6. The compound according to claim 1 or a salt thereof, wherein:

20 R_3 is H, $-\text{CH}_3$, $-\text{CH}_2\text{CH}(\text{CH}_3)_2$, $-\text{CH}_2\text{CH}(\text{CH}_3)\text{CF}_3$, $-\text{CH}_2\text{CH}_2\text{CH}_2\text{OH}$, $-\text{CH}(\text{CH}_3)\text{CH}_2\text{OH}$,
 $-\text{CH}_2\text{CH}_2\text{NHC}(\text{O})\text{OC}(\text{CH}_3)_3$, $-\text{CH}_2\text{CH}_2\text{NH}_2$, $-\text{CH}_2\text{CH}_2\text{NH}(\text{CH}_3)$,
 $-\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)(\text{CH}_2\text{CH}_3)$, $-\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)(\text{CH}_2\text{CH}(\text{CH}_3)_2)$, $-\text{CH}_2\text{CH}_2\text{CH}_2\text{NH}(\text{CH}_3)$,
 $-\text{CH}(\text{CH}_3)\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{OC}(\text{CH}_3)_3$, $-\text{C}(\text{O})\text{CH}_3$, $-\text{C}(\text{O})\text{CH}_2\text{OH}$, $-\text{C}(\text{O})\text{CH}_2\text{Cl}$,
 $-\text{C}(\text{O})\text{CH}_2\text{CF}_3$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CF}_3$, $-\text{C}(\text{O})\text{CH}_2\text{OCH}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{S}(\text{O})_2\text{CH}_3$,
25 $-\text{C}(\text{O})\text{CH}_2\text{CH}(\text{CH}_3)\text{OH}$, $-\text{C}(\text{O})\text{C}(\text{CH}_3)_2\text{OC}(\text{O})\text{CH}_3$, $-\text{C}(\text{O})\text{NHCH}_2\text{CH}_3$, $-\text{C}(\text{O})\text{NHCH}(\text{CH}_3)_2$,
 $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_3$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CN}$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{C}(\text{O})\text{CH}_3$,
 $-\text{C}(\text{O})\text{CH}_2\text{N}(\text{CH}_3)\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_2\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CH}_3$,
 $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CH}_2\text{OH}$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CF}_3$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CH}_2\text{OCH}_2\text{CH}_2\text{OH}$,

- $-\text{C}(\text{O})\text{CH}_2\text{N}(\text{CH}_3)\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_2\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{NH}(\text{CH}_2\text{OH})\text{CH}_2\text{CH}_3$,
 $-\text{C}(\text{O})\text{CH}_2\text{NHCH}(\text{CH}_2\text{OH})_2$, $-\text{C}(\text{O})\text{CH}_2\text{N}(\text{CH}_3)\text{CH}_2\text{CH}_2\text{OH}$,
 $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CH}(\text{CH}_3)\text{OH}$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CH}(\text{OH})\text{CH}_2\text{OH}$,
 $-\text{C}(\text{O})\text{CH}_2\text{N}(\text{CH}_3)\text{CH}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CH}(\text{CH}_3)_2$,
5 $-\text{C}(\text{O})\text{CH}_2\text{NHCH}(\text{CH}_3)\text{CH}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CH}_2\text{CH}(\text{CH}_3)\text{OH}$,
 $-\text{C}(\text{O})\text{CH}_2\text{NH}(\text{CH}_2\text{OH})(\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_3)$, $-\text{C}(\text{O})\text{CH}_2\text{NH}(\text{CH}_2\text{OH})\text{CH}_2\text{CH}(\text{CH}_3)_2$,
 $-\text{C}(\text{O})\text{CH}_2\text{NH}(\text{CH}(\text{CH}_2\text{CH}_3)_2)$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{CHFC}(\text{CH}_3)_2\text{OH}$,
 $-\text{C}(\text{O})\text{CH}_2\text{NHCH}_2\text{C}(\text{CH}_3)_2\text{CH}_2\text{OH}$, $-\text{C}(\text{O})\text{CH}_2\text{NHCH}(\text{CH}_2\text{OH})\text{CH}_2\text{CH}(\text{CH}_3)_2$,
 $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{NHCH}(\text{CH}_3)_2$,
10 $-\text{C}(\text{O})\text{CH}_2\text{CH}(\text{CH}_3)\text{NH}_2$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{NH}_2$, $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_3)_2$,
 $-\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_3)_2$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CN}$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{C}(\text{O})\text{NH}_2$,
 $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{OH}$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{S}(\text{O})_2\text{OH}$,
 $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{NHC}(\text{O})\text{CH}_3$,
 $-\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_3)\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_2\text{CH}_3)\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$,
15 $-\text{CH}_2\text{C}(\text{O})\text{NHCH}(\text{CH}_3)_2$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{CH}_2\text{OH}$,
 $-\text{CH}_2\text{C}(\text{O})\text{NHCH}_2\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$, $-\text{CH}_2\text{C}(\text{O})\text{N}(\text{CH}_3)\text{CH}_2\text{CH}_2\text{CH}_2\text{N}(\text{CH}_3)_2$,
 $-\text{CH}_2\text{C}(\text{O})\text{NHCH}(\text{CH}_3)(\text{CH}_2\text{CH}_3)$, $-\text{CH}_2\text{C}(\text{O})\text{NHCH}(\text{CH}_2\text{CH}_3)_2$,
 $-\text{C}(\text{O})\text{CH}_2\text{NHC}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{CN}$, or $-\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{CH}_2\text{NHC}(\text{O})\text{CH}_3$.
- 20 7. The compound according to claim 1 or a salt thereof, wherein said compound is selected from 2-(3,4-dimethoxyphenyl)-5-(1'-isobutyl-[1,4'-bipiperidin]-4-yl)-3-propyl-1H-indole-di-trifluoroacetic acid (121); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(piperidin-4-yl)-1H-indole hydrochloride (194); 2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-N-methylethan-1-amine (195); 5-([1,4'-bipiperidin]-4-yl)-2-(3,4-dimethoxyphenyl)-3-
25 isopropyl-1H-indole dihydrochloride (197); 5-(1'-(cyclopropylmethyl)-[1,4'-bipiperidin]-4-yl)-2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indole di trifluoroacetic acid (198); 2-(3,4-dimethoxyphenyl)-5-(1'-isobutyl-[1,4'-bipiperidin]-4-yl)-3-isopropyl-1H-indole (199); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(1'-methyl-[1,4'-bipiperidin]-4-yl)-1H-indole (200); 2-

(3,4-dimethoxyphenyl)-3-isopropyl-5-(1'-isopropyl-[1,4'-bipiperidin]-4-yl)-1H-indole (201); 5-(1'-cyclopentyl-[1,4'-bipiperidin]-4-yl)-2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indole (202); 2-((2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-oxoethyl)(methyl)amino)-N,N-diethylacetamide, TFA (203); 3-((2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)(methyl) amino)-N,N-diethylpropanamide (204); 2-(3,4-dimethoxyphenyl)-3-isopropyl-6-methyl-5-(piperidin-4-yl)-1H-indole (205); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(1'-isopropyl-[1,4'-bipiperidin]-4-yl)-6-methyl-1H-indole di trifluoroacetic acid (206); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(1'-isopropyl-[1,4'-bipiperidin]-4-yl)-6-methyl-1H-indole-ditrifluoroacetic acid (207); 2-(4-(2-(3,4-dimethoxyphenyl)-6-fluoro-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-N-methylethan-1-amine (208); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(dimethylamino)ethan-1-one (209); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((2-hydroxypropyl)amino)ethan-1-one (210); 2-([1,4'-bipiperidin]-1'-yl)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (211); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((2-hydroxyethyl) amino)ethan-1-one (212); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-((1-(hydroxymethyl)cyclopentyl)amino)ethan-1-one (213); 2-((2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)amino) acetonitrile (214); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((4-(dimethylamino)cyclohexyl)amino)ethan-1-one (215); 2-((3-(tert-butyl)-1H-pyrazol-5-yl)amino)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (216); 2-(((3s,5s,7s)-adamantan-1-yl)amino)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (217); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(((1r,3s,5R,7S)-3-hydroxyadamantan-1-yl)amino)ethan-1-one (218); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((3-hydroxy-2,2-dimethylpropyl)amino)ethan-1-one (219); 2-((1,3-dihydroxypropan-2-yl)amino)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (220); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-

yl)piperidin-1-yl)-2-((2,2,2-trifluoroethyl) amino)ethan-1-one (221); 2-(4-acetyl-1,4-diazepan-1-yl)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (222); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-morpholinoethan-1-one (223); 1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)-N,N-diethylpiperidine-3-carboxamide (224); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(2-(((S)-1-methylpyrrolidin-2-yl)methyl) piperidin-1-yl)ethan-1-one (225); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(hydroxymethyl)piperidin-1-yl)ethan-1-one (226); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(isopropyl(methyl)amino) ethan-1-one (227); 2-((2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-oxoethyl)amino)acetamide (228); 2-(4-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)piperazin-1-yl)-N-isopropylacetamide (229); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-((2-(2-hydroxyethoxy)ethyl)amino)ethan-1-one (230); (S)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((1-hydroxy-4-methylpentan-2-yl)amino)ethan-1-one (231); 2-(((1r,4r)-4-aminocyclohexyl)amino)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (232); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((1-hydroxybutan-2-yl)amino)ethan-1-one (233); (R)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((1-hydroxy-4-methylpentan-2-yl)amino) ethan-1-one (234); (S)-2-((2,3-dihydroxypropyl)amino)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (235); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((2-hydroxycyclopentyl) amino)ethan-1-one (236); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-((3-hydroxybutyl)amino)ethan-1-one (237); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(4-hydroxypiperidin-1-yl) ethan-1-one (238); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(4-methyl-1,4-diazepan-1-yl)ethan-1-one (239); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(4-isopropylpiperazin-1-yl)ethan-1-one (240); 1-(4-(2-(3,4-dimethoxyphenyl)-3-

isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(4-(4-methylpiperazin-1-yl)piperidin-1-yl)ethan-1-one (241); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((2-hydroxyethyl)(methyl)amino)ethan-1-one (242); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(4-(hydroxymethyl)piperidin-1-yl)ethan-1-one (243); 1-(4-

5 (2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(4-(dimethylamino)piperidin-1-yl)ethan-1-one (244); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((4-(2-hydroxypropan-2-yl)cyclohexyl)amino)ethan-1-one (245); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((4-hydroxycyclohexyl) amino)ethan-1-one (246); 1-(2-(4-(2-

10 (3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)-N,N-diethylpiperidine-3-carboxamide (247 and 248); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(hydroxymethyl)piperidin-1-yl)ethan-1-one (249 and 250); (R)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((2-fluoro-3-hydroxy-3-methylbutyl)amino)ethan-1-one (251); N-(2-(4-(2-(3,4-

15 dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)hex-5-ynamide (252); (S)-3-amino-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)butan-1-one (253); (S)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-3-hydroxybutan-1-one (254 and 255); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(methylamino)ethan-1-one (256); 1-(4-(2-(3,4-

20 dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-4-(isopropylamino)butan-1-one (257); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(1-((1-methyl-1H-imidazol-5-yl)methyl)piperidin-4-yl)-1H-indole (258); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(1-((1-methyl-1H-pyrazol-4-yl)methyl)piperidin-4-yl)-1H-indole (259); N-(3-((4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)methyl)phenyl)

25 methanesulfonamide (260); 2-((4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)methyl)thiazole (261); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(1-((1-methyl-1H-imidazol-2-yl)methyl)piperidin-4-yl)-1H-indole (262); 5-(1-(4-(1H-imidazol-1-yl)benzyl)piperidin-4-yl)-2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indole (263); 3-((4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)methyl)benzonitrile (264);

2-(3,4-dimethoxyphenyl)-5-(1-(2-ethylbutyl)piperidin-4-yl)-3-isopropyl-1H-indole (265); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(1-((2-methyl-1H-imidazol-5-yl)methyl)piperidin-4-yl)-1H-indole (266); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(1-((4-methyl-1H-imidazol-5-yl)methyl)piperidin-4-yl)-1H-indole (267); 5-(1-((1H-imidazol-5-yl)methyl)piperidin-4-yl)-2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indole (268); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(1-(3-(trifluoromethyl)benzyl)piperidin-4-yl)-1H-indole (269); 4-((4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)methyl)-N,N-dimethylaniline (270); 2-(3,4-dimethoxyphenyl)-3-isopropyl-5-(1-(4-(trifluoromethoxy)benzyl)piperidin-4-yl)-1H-indole (271); 1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)acetyl)-N,N-diethylpiperidine-3-carboxamide (272); 2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-N-(1-isopropylpiperidin-4-yl)acetamide (273); 2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-1-(4-methylpiperazin-1-yl)ethan-1-one (274); 2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-1-(3-(hydroxymethyl)piperidin-1-yl)ethan-1-one (275); 2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-N-(2-hydroxyethyl)acetamide (276); 2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-N-(3-hydroxypropyl)acetamide (277); (S)-2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-1-(3-hydroxypyrrolidin-1-yl)ethan-1-one (278); 2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-N-((1r,4r)-4-(2-hydroxypropan-2-yl)cyclohexyl)acetamide (279); 1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)-N-(1-isopropylpiperidin-4-yl)piperidine-3-carboxamide (280 and 281); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(3-hydroxypyrrolidine-1-carbonyl)piperidin-1-yl)ethan-1-one (282 and 283); (S)-2-(3-(4-acetylpiperazine-1-carbonyl)piperidin-1-yl)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (284); (S)-1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)-N-isopropylpiperidine-3-carboxamide (285); (S)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(4-isopropylpiperazine-1-carbonyl)piperidin-1-yl)ethan-1-one (286); (S)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(pyrrolidine-1-

carbonyl)piperidin-1-yl)ethan-1-one (287); (S)-1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)-N-isopropyl-N-methylpiperidine-3-carboxamide (288); (S)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-(3-(4-(dimethylamino)piperidine-1-carbonyl)piperidin-1-yl)ethan-1-one (289); (S)-1-
 5 (4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(4-(3-(trifluoromethyl)phenyl)piperazine-1-carbonyl)piperidin-1-yl)ethan-1-one (290); (S)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(4-(2-hydroxyethyl)piperidine-1-carbonyl)piperidin-1-yl)ethan-1-one (291); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((3S)-3-(3-
 10 hydroxypiperidine-1-carbonyl)piperidin-1-yl)ethan-1-one (292); (S)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(4-(2-(pyrrolidin-1-yl)ethyl)piperazine-1-carbonyl)piperidin-1-yl)ethan-1-one (293); (S)-2-(3-(4-(4-bromobenzyl)piperazine-1-carbonyl)piperidin-1-yl)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (294); (S)-N-(2-acetamidoethyl)-1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl) piperidine-3-carboxamide (295); (S)-1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)-N-(1-(hydroxymethyl)cyclopentyl)piperidine-3-carboxamide (296); (S)-N-(cyanomethyl)-1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)piperidine-3-carboxamide (297); (S)-N-(2-amino-2-oxoethyl)-
 15 1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-oxoethyl)piperidine-3-carboxamide (298); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((3S)-3-(octahydropyrrolo[3,4-c]pyrrole-2-carbonyl)piperidin-1-yl)ethan-1-one (299); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((3S)-3-(2,5-dimethylpyrrolidine-1-carbonyl)piperidin-1-yl)ethan-1-one (300); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((S)-3-((S)-3-hydroxypyrrolidine-1-carbonyl)piperidin-1-yl)ethan-1-one (301); (R)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(4-isopropylpiperazine-1-carbonyl)piperidin-1-yl)ethan-1-one (302); (R)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(pyrrolidine-1-carbonyl)piperidin-1-yl)ethan-1-one (303);

(R)-1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)-N-isopropyl-N-methylpiperidine-3-carboxamide (304); (R)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-(3-(4-(dimethylamino)piperidine-1-carbonyl)piperidin-1-yl)ethan-1-one (305); (R)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(4-(2-hydroxyethyl)piperidine-1-carbonyl)piperidin-1-yl)ethan-1-one (306); (R)-1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-(3-(4-(2-(pyrrolidin-1-yl)ethyl)piperazine-1-carbonyl)piperidin-1-yl)ethan-1-one (307); (R)-N-(cyanomethyl)-1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)piperidine-3-carboxamide (308); (R)-N-(2-amino-2-oxoethyl)-1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)piperidine-3-carboxamide (309); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-((3R)-3-(octahydropyrrolo[3,4-c]pyrrole-2-carbonyl)piperidin-1-yl) ethan-1-one (310); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((3R)-3-(2,5-dimethylpyrrolidine-1-carbonyl)piperidin-1-yl)ethan-1-one (311); 2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-N,N-dimethylacetamide (312); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-hydroxyethanone (313); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-1-yl)-2-isopropoxyethan-1-one (314); 1-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)-2-((1-methylpiperidin-4-yl) oxy)ethan-1-one (315); tert-butyl 5-(1-(tert-butoxycarbonyl)-3-hydroxypiperidin-4-yl)-2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indole-1-carboxylate (316); 4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-3-ol 2,2,2-trifluoroacetate (317 and 318); 2-(3,4-dimethoxyphenyl)-5-(3-fluoropiperidin-4-yl)-3-isopropyl-1H-indole (319 and 320); 2-(3,4-dimethoxyphenyl)-5-(3-fluoropiperidin-4-yl)-3-isopropyl-1H-indole (321, 322, 323, and 324); 5-(3,3-difluoropiperidin-4-yl)-2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indole (325); 4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl) piperidin-2-one (326); 3-(2,2-difluoroethyl)-2-(3,4-dimethoxyphenyl)-5-(piperidin-4-yl)-1H-indole (328); 2-(3,4-dimethoxyphenyl)-5-(piperidin-4-yl)-3-(2,2,2-trifluoroethyl)-1H-indole (329); 1-(4-(2-(3,4-dimethoxyphenyl)-3-

(2,2,2-trifluoroethyl)-1H-indol-5-yl) piperidin-1-yl)-3-(piperidin-1-yl)propan-1-one (330); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-4-(1H-imidazol-1-yl)butan-1-one (331); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-(1H-indol-3-yl)ethan-1-one (332); 4-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-4-oxobutanamide (333); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-(1H-imidazol-4-yl)ethan-1-one (334); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-(pyridin-3-yl)ethan-1-one (335); (S)-1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidine-1-carbonyl)pyrrolidin-1-yl)ethan-1-one (336); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-3-(1H-indol-1-yl)propan-1-one (337); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-(4-(dimethylamino)phenyl)ethan-1-one (338); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-3,3,3-trifluoropropan-1-one (339); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-3-(dimethylamino)propan-1-one (340); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-(1H-tetrazol-5-yl)ethan-1-one (341); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-(pyrazin-2-yl)ethan-1-one (342); N-(4-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-4-oxobutyl)acetamide (343); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-3-(1H-pyrrol-1-yl)propan-1-one (344); 1-(4-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidine-1-carbonyl)piperidin-1-yl)ethan-1-one (345); 4-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N,N-dimethyl-4-oxobutanamide (346); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-(methylsulfonyl)ethan-1-one (347); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-3-(pyridin-3-yl)propan-1-one (348); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-(pyridin-3-yloxy)ethan-1-one (349); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-

yl)piperidin-1-yl)-4,4,4-trifluorobutan-1-one (350); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-4-(dimethylamino)butan-1-one (351); (S)-1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-3-hydroxybutan-1-one (352); (R)-1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-3-hydroxybutan-1-one (353); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-(dimethylamino)ethan-1-one (354); 2-(3,4-dimethoxyphenyl)-5-(1-methylpiperidin-4-yl)-3-(2,2,2-trifluoroethyl)-1H-indole (355); 4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)-N-ethylpiperidine-1-carboxamide (356); 4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)-N-phenethylpiperidine-1-carboxamide (357); 1-([1,4'-bipiperidin]-1'-yl)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (359); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-(2-(1-methylpyrrolidin-2-yl)ethyl) acetamide (360); 1-(4-acetypiperazin-1-yl)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)ethan-1-one (361); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl) piperidin-1-yl)-1-(4-isopropylpiperazin-1-yl)ethan-1-one (362); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-1-(4-(4-methylpiperazin-1-yl)piperidin-1-yl)ethan-1-one (363); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl) piperidin-1-yl)-1-(4-(dimethylamino)piperidin-1-yl) ethan-1-one (364); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl) piperidin-1-yl)-N-(3-(dimethylamino)propyl)-N-methylacetamide (365); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-1-(4-(pyridin-4-yl)piperazin-1-yl) ethan-1-one (366); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl) piperidin-1-yl)-N-(2-(dimethylamino)ethyl)-N-methylacetamide (367); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-1-(4-(pyrimidin-2-yl)piperazin-1-yl)ethan-1-one (368); 1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)acetyl) piperidine-4-carboxamide (369); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-methyl-N-(1-methylpyrrolidin-3-yl)acetamide (370); (R)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-

yl)piperidin-1-yl)-1-(2-(hydroxymethyl)pyrrolidin-1-yl)ethan-1-one (371); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-(2-(dimethylamino)ethyl)-N-ethylacetamide (372); (R)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-1-(3-(dimethylamino)pyrrolidin-1-yl) ethan-1-one (373); (S)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-1-(3-(dimethylamino)pyrrolidin-1-yl)ethan-1-one (374); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-(2-(pyrrolidin-1-yl)ethyl)acetamide (375); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-(3-(2-oxopyrrolidin-1-yl)propyl) acetamide (376); N-(sec-butyl)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)acetamide (377); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-(pentan-3-yl)acetamide (378); N-(2-acetamidoethyl)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)acetamide (379); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-(2-(dimethylamino)ethyl)acetamide (380); N-(cyanomethyl)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)acetamide (381); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-(3-(dimethylamino)propyl)acetamide (382); N-(3-(1H-imidazol-1-yl)propyl)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)acetamide (383); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-(4-sulfamoylphenethyl)acetamide (384); N-(2-amino-2-oxoethyl)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)acetamide (385); 2-(2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)acetamido)ethane-1-sulfonic acid (386); N-(5-(tert-butyl)-1H-pyrazol-3-yl)-2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)acetamide (387); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-(4-(pyrrolidin-1-yl)butyl)acetamide (388); 2-(3,4-dimethoxyphenyl)-5-(piperidin-4-yl)-3-(tetrahydro-2H-pyran-4-yl)-1H-indole (389); 1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-(tetrahydro-2H-pyran-4-yl)-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)-N,N-

diethylpiperidine-3-carboxamide (391 and 392); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(tetrahydro-2H-pyran-4-yl)-1H-indol-5-yl)piperidin-1-yl)-2-(dimethylamino)ethan-1-one (393); (R)-1-(4-(2-(3,4-dimethoxyphenyl)-3-(tetrahydro-2H-pyran-4-yl)-1H-indol-5-yl)piperidin-1-yl)-2-((3-methylbutan-2-yl)amino)ethan-1-one (394); 1-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoro-1-hydroxyethyl)-1H-indol-5-yl)piperidin-1-yl)-2-(dimethylamino)ethan-1-one (395); 1-(4-(3-(2,2-difluoroethyl)-2-(3,4-dimethoxyphenyl)-1H-indol-5-yl)piperidin-1-yl)-2-(dimethylamino)ethan-1-one (399); 2-(4-(3-(2,2-difluoroethyl)-2-(3,4-dimethoxyphenyl)-1H-indol-5-yl)piperidin-1-yl)-N,N-dimethylacetamide (400); 2-(4-(3-(2,2-difluoroethyl)-2-(3,4-dimethoxyphenyl)-1H-indol-5-yl)piperidin-1-yl)-N-methylacetamide (401); N-(2-(4-(2-(3,4-dimethoxyphenyl)-3-isopropyl-1H-indol-5-yl)piperidin-1-yl)ethyl)hept-6-ynamide (411); 2-(3,4-dimethoxyphenyl)-5-(1'-isopropyl-[1,4'-bipiperidin]-4-yl)-3-(2,2,2-trifluoroethyl)-1H-indole (412); 2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-N-methylethanamine (413); tert-butyl (2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)(methyl)carbamate (414); 1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)-N,N-diethylpiperidine-3-carboxamide (418, 420-421); and 1-(2-(4-(2-(3,4-dimethoxyphenyl)-3-(2,2,2-trifluoroethyl)-1H-indol-5-yl)piperidin-1-yl)-2-oxoethyl)-N,N-dimethylpiperidine-3-carboxamide (419).

8. A pharmaceutical composition comprising a compound according to any one of claims 1 to 7 or a pharmaceutically-acceptable salt thereof; and a pharmaceutically acceptable carrier.

9. A compound according to any one of claims 1 to 7 or composition according to claim 8 for use in therapy.

10. A compound according to any one of claim 1 to 7 or composition according to claim 8 for use in the treatment of inflammatory or autoimmune disease.

11. The compound or composition for use according to claim 10, wherein said autoimmune disease or chronic inflammatory disease is selected from systemic lupus erythematosus (SLE), rheumatoid arthritis, multiple sclerosis (MS), and Sjögren's syndrome.