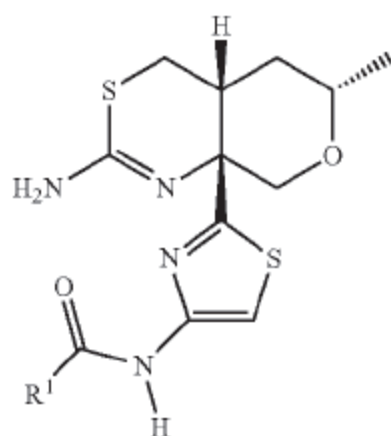


We claim:

1. A compound of Formula I



I

wherein

R^1 is selected from the group consisting of:

phenyl optionally substituted with one to three R^2 ;

C_{3-9} cycloalkyl optionally substituted with one to three R^2 ;
and

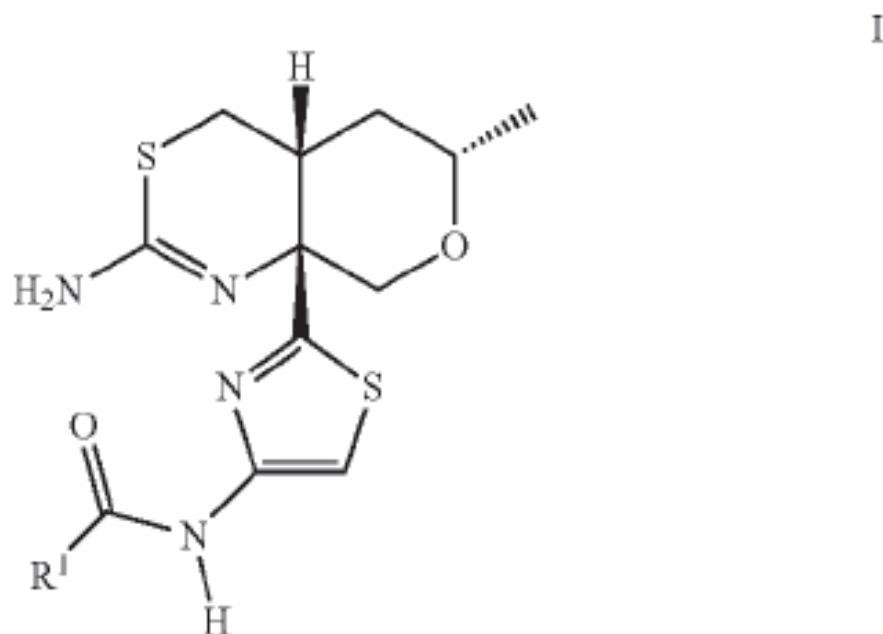
a 5- to 10-membered heteroaryl, having one to four heteroatoms independently selected from N, O or S, wherein at least one of the heteroatoms is N and wherein said N is optionally substituted with R^3 ; and wherein said 5- to 10-membered heteroaryl is optionally substituted on carbon with one to three R^2 ;

R^2 at each occurrence is independently selected from the group consisting of halogen, hydroxy, cyano, C_{1-6} alkyl, C_{1-6} alkoxy, C_{3-6} alkenyl, C_{3-6} alkenyloxy, C_{3-6} alkynyl, C_{3-6} alkynyloxy, C_{1-6} alkoxy- C_{1-6} alkyl, C_{3-6} cycloalkoxy, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl- C_{1-6} alkoxy, C_{3-6} cycloalkyl- C_{1-6} alkoxy, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl- C_{1-6} alkyl; wherein said C_{1-6} alkyl, C_{1-6} alkoxy, C_{3-6} alkenyl, C_{3-6} alkenyloxy, C_{3-6} alkynyl, C_{3-6} alkynyloxy, C_{1-6} alkoxy- C_{1-6} alkyl, C_{3-6} cycloalkoxy, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl- C_{1-6} alkyl, C_{3-6} cycloalkyl- C_{1-6} alkoxy, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl- C_{1-6} alkyl are each optionally substituted with one to three substituents independently selected from fluoro, chloro, hydroxy, cyano, methyl, fluoromethyl, difluoromethyl, trifluoromethyl, methoxy, fluoromethoxy, difluoromethoxy and trifluoromethoxy; or two R^2 groups taken together can be a C_{3-5} alkylene; and

R^3 is selected from the group consisting of hydrogen, C_{1-6} alkyl, C_{3-6} alkenyl, C_{3-6} alkynyl, C_{1-6} alkoxy- C_{1-6} alkyl, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl- C_{1-6} alkyl, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl- C_{1-6} alkyl; wherein said C_{1-6} alkyl, C_{3-6} alkenyl, C_{3-6} alkynyl, C_{1-6} alkoxy- C_{1-6} alkyl, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl- C_{1-6} alkyl, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl- C_{1-6} alkyl are each optionally substituted with one to three substituents independently selected from fluoro, chloro, hydroxy, cyano, methyl, fluoromethyl, difluoromethyl, trifluoromethyl, methoxy, fluoromethoxy, difluoromethoxy and trifluoromethoxy; or R^3 and R^2 taken together can be a C_{3-5} alkylene;

or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

2. The compound of claim 1 of Formula I



wherein

R¹ is selected from the group consisting of:
phenyl optionally substituted with one to three R²;
C₃₋₆cycloalkyl optionally substituted with one to three R²;
and

- a 5- to 10-membered heteroaryl, having one to four heteroatoms independently selected from N, O or S, wherein at least one of the heteroatoms is N and wherein said N is optionally substituted with R³; and wherein said 5- to 10-membered heteroaryl is optionally substituted on carbon with one to three R²;

R^2 at each occurrence is independently selected from the group consisting of halogen, hydroxy, cyano, C_{1-6} alkyl, C_{1-6} alkoxy, C_{3-6} alkenyl, C_{3-6} alkenyloxy, C_{3-6} alkynyl, C_{3-6} alkynyloxy, C_{1-6} alkoxy- C_{1-6} alkyl, C_{3-6} cycloalkoxy, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl- C_{1-6} alkyl, C_{3-6} cycloalkyl- C_{1-6} alkoxy, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl- C_{1-6} alkyl; wherein said C_{1-6} alkyl, C_{1-6} alkoxy, C_{3-6} alkenyl, C_{3-6} alkenyloxy, C_{3-6} alkynyl, C_{3-6} alkynyloxy, C_{1-6} alkoxy- C_{1-6} alkyl, C_{3-6} cycloalkoxy, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl- C_{1-6} alkyl, C_{3-6} cycloalkyl- C_{1-6} alkoxy, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl- C_{1-6} alkyl are each optionally substituted with one to three substituents independently selected from fluoro, chloro, hydroxy, methyl, fluoromethyl, difluoromethyl, trifluoromethyl, methoxy, fluoromethoxy, difluoromethoxy and trifluoromethoxy; or two R^2 groups taken together can be a C_{3-5} alkylene; and

R^3 is selected from the group consisting of hydrogen, C_{1-6} alkyl, C_{3-6} alkenyl, C_{3-6} alkynyl, C_{1-6} alkoxy- C_{1-6} alkyl, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl- C_{1-6} alkyl, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl- C_{1-6} alkyl; wherein said C_{1-6} alkyl, C_{3-6} alkenyl, C_{3-6} alkynyl, C_{1-6} alkoxy- C_{1-6} alkyl, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl- C_{1-6} alkyl, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl- C_{1-6} alkyl are each optionally substituted with one to three substituents independently selected from fluoro, chloro, hydroxy, methyl, fluoromethyl, difluoromethyl, trifluoromethyl, methoxy, fluoromethoxy, difluoromethoxy and trifluoromethoxy; or R^3 and R^2 taken together can be a C_{3-5} alkylene;

or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

3. The compound of claim 2 wherein

R¹ is a 5-membered heteroaryl selected from the group consisting of pyrazolyl, imidazolyl, oxazolyl, isoxazolyl, oxadiazolyl and triazolyl; each optionally sub-

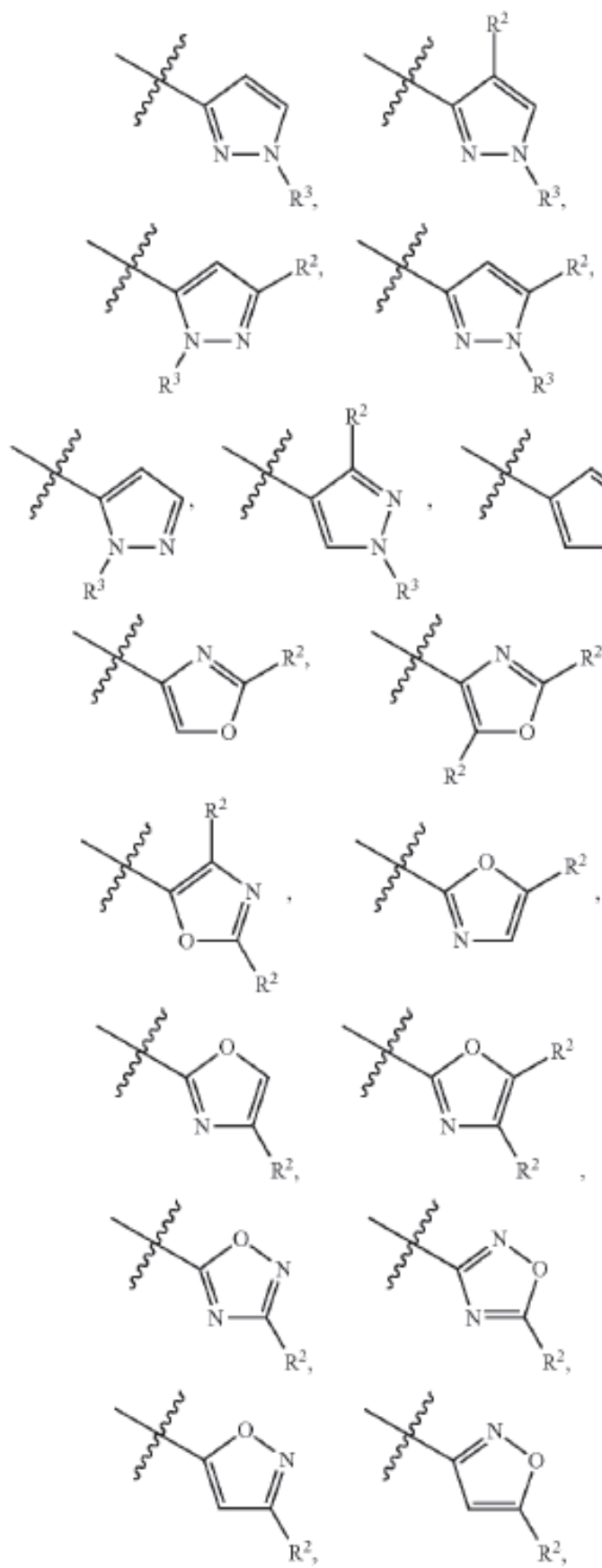
stituted on carbon with one to two R²; and wherein said pyrazolyl, imidazolyl and triazolyl are substituted on N with R³;

R² at each occurrence is independently selected from the group consisting of halogen, C₁₋₃alkyl, C₃₋₆cycloalkyl, and C₁₋₃alkoxy-C₁₋₃alkyl; wherein said C₁₋₃alkyl is optionally substituted with one to three fluoro; and

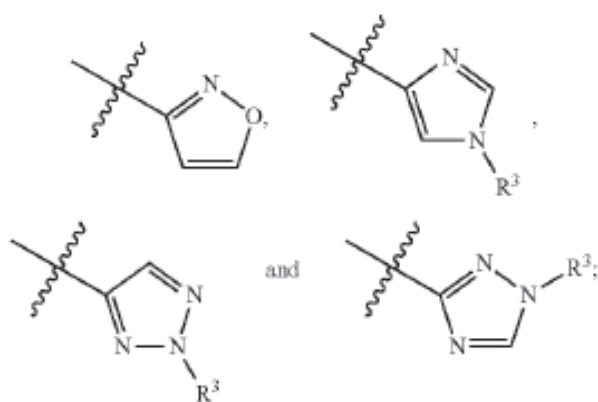
R³ is C₁₋₃alkyl or C₃₋₆cycloalkyl, wherein said C₁₋₃alkyl is optionally substituted with one to three fluoro;

or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

4. The compound of claim 3 wherein R¹ is selected from the group consisting of



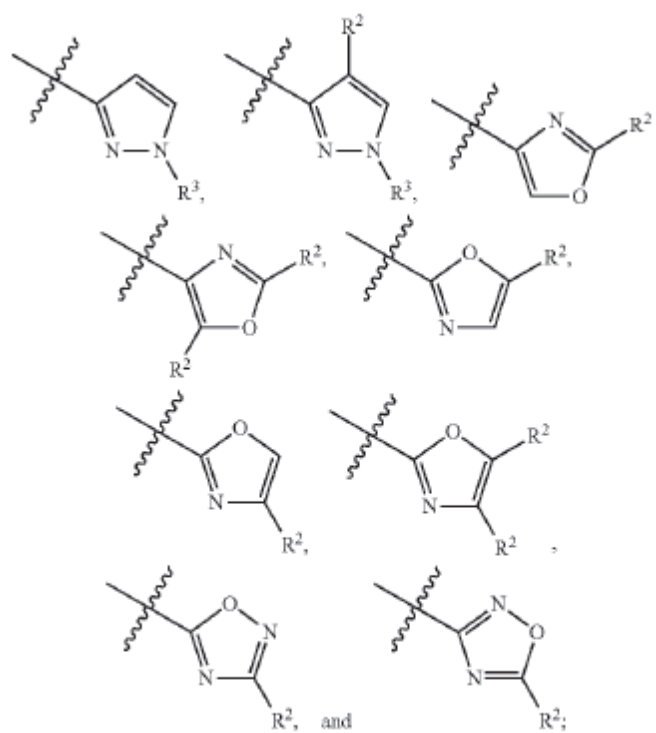
-continued



R^2 at each occurrence is independently selected from the group consisting of chloro, methyl, ethyl, isopropyl, isobutyl, fluoromethyl, difluoromethyl, trifluoromethyl, cyclopropyl, cyclobutyl and methoxymethyl; and

R^3 is methyl, ethyl, isopropyl, difluoromethyl, cyclopropyl or cyclobutyl; or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

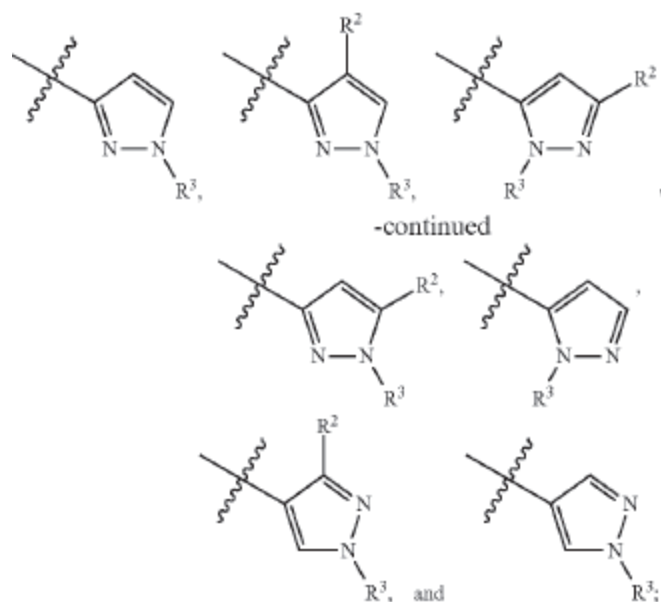
5. The compound according to claim 4 wherein R^1 is selected from the group consisting of



or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

6. The compound according to claim 4 wherein

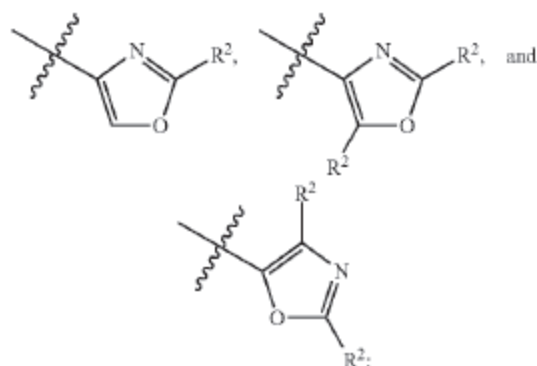
R^1 is selected from the group consisting of



or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

7. The compound according to claim 4 wherein

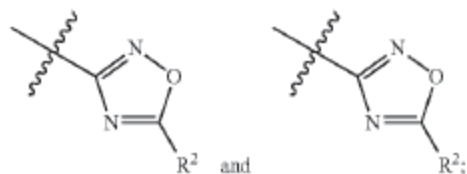
R^1 is selected from the group consisting of



or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

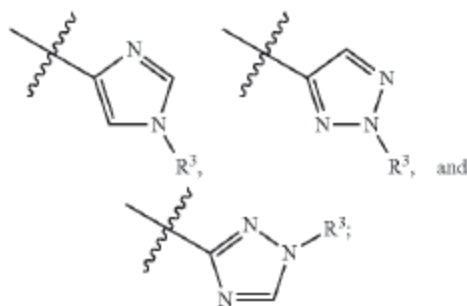
8. The compound according to claim 4 wherein

R^1 is selected from the group consisting of



or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

9. The compound according to claim 4 wherein R^1 is selected from the group consisting of



or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

10. The compound according to claim 2 wherein

R^1 is a 6-membered heteroaryl selected from the group consisting of pyridinyl, pyridonyl, pyrimidinyl, pyridazinyl and pyrazinyl; each optionally substituted on carbon with one to two R^2 ; and wherein said pyridonyl is substituted on N with R^3 ;

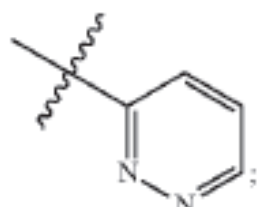
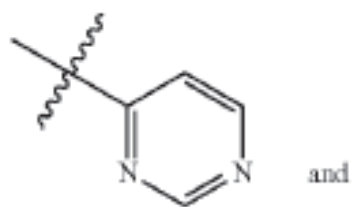
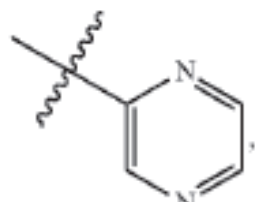
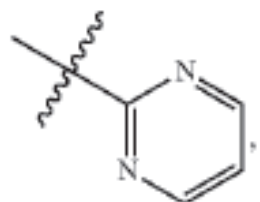
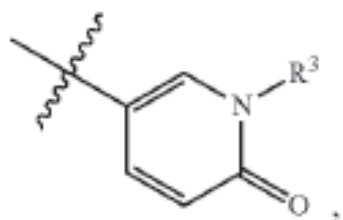
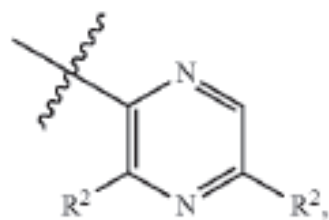
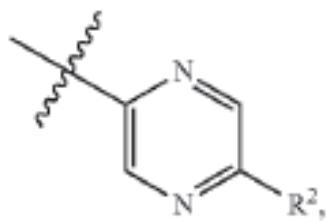
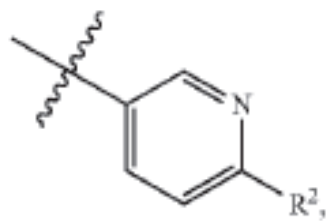
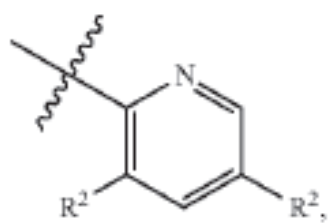
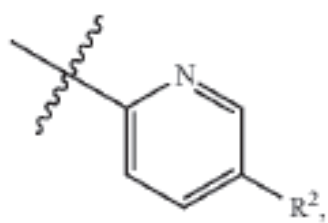
R^2 at each occurrence is independently selected from the group consisting of halogen, hydroxy, cyano, C_{1-6} alkyl, C_{1-6} alkoxy, C_{1-6} alkoxy- C_{1-6} alkyl, C_{3-6} cycloalkyl, C_{3-6} alkynyloxy, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl- C_{1-6} alkyl; wherein said C_{1-6} alkyl, C_{1-6} alkoxy, C_{1-6} alkoxy- C_{1-6} alkyl, C_{3-6} cycloalkyl, C_{3-6} alkynyloxy, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl- C_{1-6} alkyl are optionally substituted with one to three fluoro or hydroxy; and

R^3 is C_{1-3} alkyl;

or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

11. The compound according to claim 10 wherein

R^1 is selected from the group consisting of

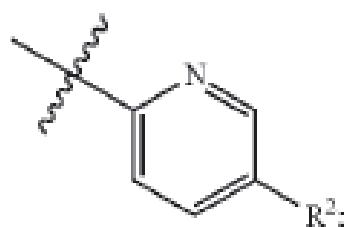


and

R^2 at each occurrence is independently selected from the group consisting of fluoro, chloro, cyano, methyl, fluoromethyl, difluoromethyl, trifluoromethyl, hydroxy, hydroxymethyl, methoxy, fluoromethoxy, difluoromethoxy, trifluoromethoxy, difluoroethoxy, methoxydifluoroethyl, difluoropropoxy, butynyloxy and cyclopropyl; and R^3 is methyl;

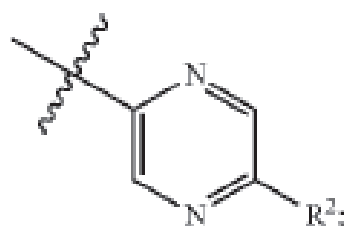
or a tautomer thereof or pharmaceutically acceptable salt of said compound or tautomer.

12. The compound according to claim 11 wherein R^1 is



or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

13. The compound according to claim 11 wherein R^1 is



or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

14. The compound according to claim 2 wherein

R¹ is phenyl optionally substituted with one to two R², or C₃₋₆-cycloalkyl optionally substituted with one to two R²; or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

15. The compound according to claim 2 selected from the group consisting of:

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(difluoromethoxy)pyridine-2-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(difluoromethoxy)pyrazine-2-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-cyanopyridine-2-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-chloropyridine-2-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-fluoropyridine-2-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-methoxypyrazine-2-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(fluoromethyl)pyrazine-2-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(2,2-difluoroethoxy)pyrazine-2-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-fluoro-3-methylpyridine-2-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(difluoromethyl)pyrazine-2-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(trifluoromethyl)pyrazine-2-carboxamide;

- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(trifluoromethyl)pyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-3,5-difluoropyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-methoxypyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(trifluoromethoxy)pyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(but-2-yn-1-yloxy)pyrazine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(difluoromethyl)pyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(fluoromethoxy)pyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}pyrazine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}pyridazine-3-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-1-methyl-6-oxo-1,6-dihydropyridine-3-carboxamide;

- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(2,2-difluoroethoxy)pyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(1,1-difluoroethoxy)pyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(difluoromethoxy)-3-methylpyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-methoxy-3-methylpyrazine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-methoxy-3-methylpyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}pyrimidine-4-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-methylpyrazine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}pyrimidine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a, 5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-cyclopropylpyrazine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}pyrazolo[1,5-a]pyridine-2-carboxamide;
- N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-6-(difluoromethoxy)pyridine-3-carboxamide;

N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(cyclopropylmethoxy)pyrazine-2-carboxamide;
 N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(cyclopropyloxy)pyrazine-2-carboxamide;
 N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-([(2R)-2-fluoropropyl]oxy)pyrazine-2-carboxamide;
 N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-([(2S)-2-fluoropropyl]oxy)pyrazine-2-carboxamide; and
 N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(2,2-difluoropropoxy)pyrazine-2-carboxamide;
 or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

16. The compound N-{2-[(4aR,6S,8aR)-2-Amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(difluoromethoxy)pyridine-2-carboxamide; or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

17. The compound N-{2-[(4aR,6S,8aR)-2-Amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(difluoromethoxy)pyrazine-2-carboxamide; or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

18. The compound N-{2-[(4aR,6S,8aR)-2-Amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-chloropyridine-2-carboxamide; or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

19. The compound N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-

yl]-1,3-thiazol-4-yl}-5-fluoropyridine-2-carboxamide; or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

20. The compound N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(difluoromethoxy)-3-methylpyridine-2-carboxamide; or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

21. The compound N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-5-(difluoromethoxy)-3-methylpyrazine-2-carboxamide; or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

22. The compound N-{2-[(4aR,6S,8aR)-2-amino-6-methyl-4,4a,5,6-tetrahydropyrano[3,4-d][1,3]thiazin-8a(8H)-yl]-1,3-thiazol-4-yl}-3-chloro-5-(difluoromethoxy)pyridine-2-carboxamide; or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer.

23. A pharmaceutical composition comprising a therapeutically effective amount of a compound of claim 1, or a tautomer thereof or a pharmaceutically acceptable salt of said compound or tautomer, and a pharmaceutically acceptable vehicle, diluent or carrier.