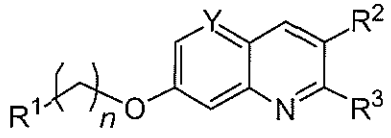
CLAIMS

1. A chemical entity, wherein the chemical entity is selected from the group consisting of compounds of Formula (I) and pharmaceutically acceptable salts of compounds of Formula (I)



Formula (I) , wherein:

n is 1 or 2;

Y is CH or N;

R^1 is a pyridine substituted with $-CF_3$, or phenyl substituted only in the *meta* and *para* positions with a total of one, two, or three R^a members;

each R^a is independently selected from the group consisting of halo, $-C_{1-4}$alkyl, $-CF_3$, $-NO_2$, and $-OC_{1-4}$alkyl;

R^2 is selected from the group consisting of $-C(R^b)_2R^c$ or $-CO-R^d$;

each R^b is independently selected from the group consisting of -H, -F, and $-C_{1-3}$alkyl, or optionally two R^b members are taken together with the carbon to which they are attached to form a C_{3-6}cycloalkyl ring;

R^c is selected from the group consisting of -F, $-NH_2$, -OH, $-OC_{1-3}$alkyl, $-CH_2OH$, -CN, $-CO_2-C_{1-4}$alkyl, $-CO-NHR^e$, and $-C(CH_3)_2OH$; provided that when at least one R^b is -F then R^c is not -F;

R^d is selected from the group consisting of $-CH_3$, $-OC_{1-4}$alkyl, $-NHR^e$, and $-NHCH_2CH_2N(R^e)_2$;

each R^e is independently -H or $-CH_3$; and

R^3 is selected from the group consisting of -H, $-CH_3$, -OH, and $-CF_3$.

2. A chemical entity as in claim 1, wherein *n* is 1.

3. A chemical entity as in claim 1, wherein *n* is 2.

4. A chemical entity as in claim 1, wherein Y is CH.

5. A chemical entity as in claim 1, wherein Y is N.

6. A chemical entity as in claim 1, wherein R^1 is 2-(trifluoromethyl)pyridin-4-yl or 6-(trifluoromethyl)pyridin-2-yl.

7. A chemical entity as in claim 1, wherein R^1 is phenyl substituted only in the *meta* and *para* positions with a total of, two, or three R^a members independently selected from the group consisting of halo, $-CF_3$, $-CH_3$, $-OCH_3$, and $-NO_2$.

8. A chemical entity as in claim 1, wherein R^1 is selected from the group consisting of 3-chlorophenyl, 3-fluorophenyl, 3-nitrophenyl, 3-methylphenyl, 3-methoxyphenyl, 3-(trifluoromethyl)phenyl, 3-chloro-4-fluorophenyl, 3,4-difluorophenyl, 3-chloro-5-fluorophenyl, 3,5-difluorophenyl, 3-fluoro-5-(trifluoromethyl)phenyl, 3,4,5-trifluorophenyl, 4-chlorophenyl, 4-fluorophenyl, 4-trifluoromethyl)phenyl, 4-fluoro-3-(trifluoromethyl)phenyl, 4-nitrophenyl, 4-methoxyphenyl, 2-(trifluoromethyl)pyridin-4-yl, and 6-(trifluoromethyl)pyridin-2-yl.

9. A chemical entity as in claim 1, wherein R^2 is $-(CR^b)_2R^c$.

10. A chemical entity as in claim 1, wherein R^b is independently selected from the group consisting of -H, -F and $-CH_3$.

11. A chemical entity as in claim 1, wherein two R^b members are taken together with the carbon to which they are attached to form a cyclopropyl, cyclobutyl, cyclopentyl, or cyclohexyl ring.

12. A chemical entity as in claim 1, wherein R^c is selected from the group consisting of -F, $-NH_2$, -OH, $-OCH_3$, $-CH_2OH$, -CN, $-CO_2-C_{1-4}$alkyl, $-CO-NHR^e$, and $-C(CH_3)_2OH$.

13. A chemical entity as in claim 1,wherein R^2 is selected from the group consisting of $-CH_2NH_2$, $-CH_2OH$, $-CH_2CH_2OH$, $-CH_2OCH_3$, $-CH_2CN$, $-CH_2(C=O)OCH_3$,

$-CH_2(C{=}O)OCH_2CH_3$, $-CH_2(C{=}O)NH_2$, $-CH_2(CH_3)_2OH$, $-CH(OH)CH_3$, $-C(CH_3)_2OH$, $-C(CH_3)_2CH_2OH$, $-C(CH_3)_2(C{=}O)NH_2$, $-OCH_2CH_3$, and $-CF(CH_3)_2$.

14. A chemical entity as in claim 1, wherein R^2 is $-CO-R^d$.

15. A chemical entity as in claim 1, wherein R^d is selected from the group consisting of $-CH_3$, $-OC_{1\text{-}4}$alkyl, $-NH_2$, $-NH(CH_3)$, $-NHCH_2CH_2NH(CH_3)$ and $-NHCH_2CH_2N(CH_3)_2$.

16. A chemical entity as in claim 1, wherein R^2 is selected from the group consisting of $-(C{=}O)CH_3$, $-C(=O)OCH_3$, $-C(=O)OCH_2CH_3$, $-(C{=}O)NH_2$, $-(C{=}O)NHCH_3$, $-(C{=}O)N(CH_3)_2$, $-C{=}O)NHCH_2CH_2NH_2$, $-(C{=}O)NHCH_2CH_2NHCH_3$, and $-(C{=}O)NHCH_2CH_2N(CH_3)_2$.

17. A chemical entity as in claim 1, wherein R^3 is H or $-CH_3$.

18. A chemical entity as in claim 1, wherein R^3 is $-CF_3$ or -OH.

19. A chemical entity selected from the group consisting of:

Ethyl 7-[(4-chlorophenyl)methoxy]quinoline-3-carboxylate;

Ethyl 7-[(4-fluorophenyl)methoxy]quinoline-3-carboxylate;

Ethyl 7-[(3-chlorophenyl)methoxy]quinoline-3-carboxylate;

Ethyl 7-[(3-fluorophenyl)methoxy]quinoline-3-carboxylate;

Ethyl 7-((3-(trifluoromethyl)benzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((4-(trifluoromethyl)benzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-methylbenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((4-methylbenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-methoxybenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((4-methoxybenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((4-nitrobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((4-fluoro-3-(trifluoromethyl)benzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-fluoro-5-(trifluoromethyl)benzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3,4-difluorobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3,5-difluorobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3,4,5-trifluorobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-chloro-4-fluorobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-nitrobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-chloro-5-fluorobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-chlorobenzyl)oxy)-2-methylquinoline-3-carboxylate;

Ethyl 7-((3-fluorobenzyl)oxy)-2-methylquinoline-3-carboxylate;

Ethyl 7-((4-fluorobenzyl)oxy)-2-methylquinoline-3-carboxylate;

Ethyl 7-((3-chloro-4-fluorobenzyl)oxy)-2-methylquinoline-3-carboxylate;

Ethyl 7-(3-fluorophenethoxy)quinoline-3-carboxylate;

Ethyl 7-(3-chlorophenethoxy)quinoline-3-carboxylate;

Methyl 7-((3-chlorobenzyl)oxy)quinoline-3-carboxylate;

2-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)propan-2-ol;

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)propan-2-ol;

2-(7-((3,5-Difluorobenzyl)oxy)quinolin-3-yl)propan-2-ol;

2-(7-((3,4,5-Trifluorobenzyl)oxy)quinolin-3-yl)propan-2-ol;

2-(7-((4-Fluorobenzyl)oxy)-2-methylquinolin-3-yl)propan-2-ol;

2-(7-((3-Chlorobenzyl)oxy)-2-methylquinolin-3-yl)propan-2-ol;

2-(7-((3-Chloro-4-fluorobenzyl)oxy)-2-methylquinolin-3-yl)propan-2-ol;

2-(7-((3-Fluorobenzyl)oxy)-2-methylquinolin-3-yl)propan-2-ol;

(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)methanol;

(7-((3,5-Difluorobenzyl)oxy)quinolin-3-yl)methanol;

(7-((3,4,5-Trifluorobenzyl)oxy)quinolin-3-yl)methanol;

2-(7-(3-Fluorophenethoxy)quinolin-3-yl)propan-2-ol;

2-(7-(3-Chlorophenethoxy)quinolin-3-yl)propan-2-ol;

7-((3-Chlorobenzyl)oxy)-3-(methoxymethyl)quinoline;

7-((4-Fluorobenzyl)oxy)-3-(2-fluoropropan-2-yl)quinoline;

1-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)ethanone;

1-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)ethanol;

(R)-1-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)ethanol;

(S)-1-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)ethanol;

7-((3-Chlorobenzyl)oxy)-N-methylquinoline-3-carboxamide;

N-(2-Aminoethyl)-7-((3-chlorobenzyl)oxy)quinoline-3-carboxamide;

7-((3-Chlorobenzyl)oxy)-N-(2-(methylamino)ethyl)quinoline-3-carboxamide;

7-((3-Chlorobenzyl)oxy)-N-(2-(dimethylamino)ethyl)quinoline-3-carboxamide;

(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)methanamine;

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)acetamide;

Ethyl 2-(7-((3-fluorobenzyl)oxy)quinolin-3-yl)acetate;

Ethyl 2-(7-((4-fluorobenzyl)oxy)quinolin-3-yl)acetate;

Methyl 2-(7-((3-chlorobenzyl)oxy)quinolin-3-yl)acetate;

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)acetamide;

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)ethanol;

2-(7-((3-Fluorobenzyl)oxy)quinolin-3-yl)acetamide;

2-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)acetamide;

Ethyl 2-(7-((3-chlorobenzyl)oxy)quinolin-3-yl)-2-methylpropanoate;

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)-2-methylpropan-1-ol;

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)-2-methylpropanamide;

2-(7-((2-(Trifluoromethyl)pyridin-4-yl)methoxy)quinolin-3-yl)acetamide;

1-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)-2-methylpropan-2-ol;

7-((3-Chlorobenzyl)oxy)quinoline-3-carboxamide;

2-{7-[(4-Fluorophenyl)methoxy]-1,5-naphthyridin-3-yl}propan-2-ol;

7-[(4-Fluorophenyl)methoxy]-3-(2-hydroxypropan-2-yl)quinolin-1-ium-1-olate;

7-[(3-Fluorophenyl)methoxy]-1,5-naphthyridine-3-carboxamide;

7-((4-Fluorobenzyl)oxy)-1,5-naphthyridine-3-carboxamide;

7-((3-Chlorobenzyl)oxy)-1,5-naphthyridine-3-carboxamide;

7-((3,4-Difluorobenzyl)oxy)-1,5-naphthyridine-3-carboxamide;

7-[(3-Chloro-4-fluorophenyl)methoxy]-1,5-naphthyridine-3-carboxamide;

7-[(3-Chlorophenyl)methoxy]-N-methyl-1,5-naphthyridine-3-carboxamide;

(7-((3,4-Difluorobenzyl)oxy)-1,5-naphthyridin-3-yl)methanol;

2-{7-[(3,4-Difluorophenyl)methoxy]-1,5-naphthyridin-3-yl}acetonitrile;

3-[(4-Fluorophenyl)methoxy]-7-(2-fluoropropan-2-yl)-1,5-naphthyridine;

2-{7-[(3-Fluorophenyl)methoxy]-1,5-naphthyridin-3-yl}propan-2-ol;

2-{7-[(3,4-Difluorophenyl)methoxy]-1,5-naphthyridin-3-yl}propan-2-ol;

2-{7-[(3-Chloro-4-fluorophenyl)methoxy]-1,5-naphthyridin-3-yl}propan-2-ol;

2-{7-[(3-Chlorophenyl)methoxy]-1,5-naphthyridin-3-yl}propan-2-ol;

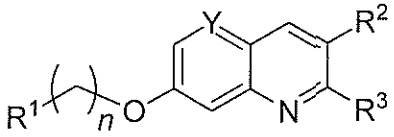
2-(7-((3,4-Difluorobenzyl)oxy)-1,5-naphthyridin-3-yl)acetamide;

2-(7-((6-(Trifluoromethyl)pyridin-2-yl)methoxy)-1,5-naphthyridin-3-yl)acetamide;

2-(7-((3-Chlorobenzyl)oxy)-2-oxo-1,2,3,4-tetrahydroquinolin-3-yl)acetamide;

2-(7-((3-Chlorobenzyl)oxy)-2-oxo-1,2-dihydroquinolin-3-yl)acetamide; and pharmaceutically acceptable salts thereof.

20. A pharmaceutical composition comprising an effective amount of at least one chemical entity selected from the compounds of Formula (I),



Formula (I) , wherein:

n is 1 or 2;

Y is CH or N;

R^1 is a pyridine substituted with $-CF_3$, or phenyl substituted only in the *meta* and *para* positions with a total of one, two, or three R^a members;

each R^a is independently selected from the group consisting of halo, $-C_{1-4}$alkyl, CF_3, $-NO_2$, and $-OC_{1-4}$alkyl;

R^2 is selected from the group consisting of $-C(R^b)_2R^c$ or $-CO-R^d$;

each R^b is independently selected from the group consisting of -H, -F, and $-C_{1-3}$alkyl, or optionally two R^b members are taken together with the carbon to which they are attached to form a C_{3-6}cycloalkyl ring;

R^c is selected from the group consisting of -F, $-NH_2$, -OH, $-OC_{1-3}$alkyl, $-CH_2OH$, -CN, $-CO_2-C_{1-4}$alkyl, $-CO-NHR^e$, and $-C(CH_3)_2OH$; provided that when at least one R^b is -F then R^c is not -F;

R^d is selected from the group consisting of $-CH_3$, $-OC_{1-4}$alkyl, $-NHR^e$, and $-NHCH_2CH_2N(R^e)_2$;

each R^e is independently -H or $-CH_3$;

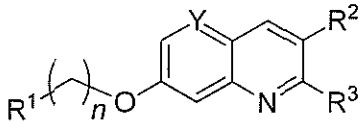
R^3 is selected from the group consisting of -H, $-CH_3$, -OH, and $-CF_3$,

and pharmaceutically acceptable salts of compounds of Formula (I).

21. A pharmaceutical composition comprising an effective amount of at least one chemical entity of claim 1.

22. A pharmaceutical composition comprising an effective amount of at least one chemical entity of claim 19.

23. A method for modulating one or more monoamine oxidase (MAO) enzymes, comprising exposing *in-vitro* MAO-A or MAO-B to an effective amount of at least one chemical entity selected from compounds of Formula (I)



Formula (I) , wherein:

n is 1 or 2;

Y is CH or N;

R^1 is a pyridine substituted with $-CF_3$, or phenyl substituted only in the *meta* and *para* positions with a total of one, two, or three R^a members;

each R^a is independently selected from the group consisting of halo, $-C_{1-4}$alkyl, CF_3, $-NO_2$, and $-OC_{1-4}$alkyl;

R^2 is selected from the group consisting of $-C(R^b)_2R^c$ or $-CO-R^d$;

each R^b is independently selected from the group consisting of -H, -F, and $-C_{1-3}$alkyl, or optionally two R^b members are taken together with the carbon to which they are attached to form a C_{3-6}cycloalkyl ring;

R^c is selected from the group consisting of -F, $-NH_2$, -OH, $-OC_{1-3}$alkyl, $-CH_2OH$, -CN, $-CO_2-C_{1-4}$alkyl, $-CO-NHR^e$, and $-C(CH_3)_2OH$; provided that when at least one R^b is -F then R^c is not -F;

R^d is selected from the group consisting of $-CH_3$, $-OC_{1-4}$alkyl, $-NHR^e$, and $-NHCH_2CH_2N(R^e)_2$;

each R^e is independently -H or $-CH_3$;

R^3 is selected from the group consisting of -H, $-CH_3$, -OH, and $-CF_3$,

and pharmaceutically acceptable salt of compounds of Formula (I).

24. A method for modulating one or more monoamine oxidase (MAO) enzymes, comprising exposing *in-vitro* MAO-A or MAO-B to an effective amount of at least one chemical entity of Claim 19.

25. A chemical entity as in claim 19, wherein the chemical entity is selected from the group consisting of:

Ethyl 7-[(4-chlorophenyl)methoxy]quinoline-3-carboxylate;

Ethyl 7-[(4-fluorophenyl)methoxy]quinoline-3-carboxylate;

Ethyl 7-[(3-chlorophenyl)methoxy]quinoline-3-carboxylate;

Ethyl 7-[(3-fluorophenyl)methoxy]quinoline-3-carboxylate;

Ethyl 7-((3-(trifluoromethyl)benzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((4-(trifluoromethyl)benzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-methylbenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((4-methylbenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-methoxybenzyl)oxy)quinoline-3-carboxylate; and

Ethyl 7-((4-methoxybenzyl)oxy)quinoline-3-carboxylate.

26. A chemical entity as in claim 19, wherein the chemical entity is selected from the group consisting of:

Ethyl 7-((4-nitrobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((4-fluoro-3-(trifluoromethyl)benzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-fluoro-5-(trifluoromethyl)benzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3,4-difluorobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3,5-difluorobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3,4,5-trifluorobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-chloro-4-fluorobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-nitrobenzyl)oxy)quinoline-3-carboxylate;

Ethyl 7-((3-chloro-5-fluorobenzyl)oxy)quinoline-3-carboxylate; and

Ethyl 7-((3-chlorobenzyl)oxy)-2-methylquinoline-3-carboxylate.

27. A chemical entity as in claim 19, wherein the chemical entity is selected from the group consisting of:

Ethyl 7-((3-fluorobenzyl)oxy)-2-methylquinoline-3-carboxylate;

Ethyl 7-((4-fluorobenzyl)oxy)-2-methylquinoline-3-carboxylate;

Ethyl 7-((3-chloro-4-fluorobenzyl)oxy)-2-methylquinoline-3-carboxylate;

Ethyl 7-(3-fluorophenethoxy)quinoline-3-carboxylate;

Ethyl 7-(3-chlorophenethoxy)quinoline-3-carboxylate;

Methyl 7-((3-chlorobenzyl)oxy)quinoline-3-carboxylate;

2-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)propan-2-ol;

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)propan-2-ol;

2-(7-((3,5-Difluorobenzyl)oxy)quinolin-3-yl)propan-2-ol; and

2-(7-((3,4,5-Trifluorobenzyl)oxy)quinolin-3-yl)propan-2-ol.

28. A chemical entity as in claim 19, wherein the chemical entity is selected from the group consisting of:

2-(7-((4-Fluorobenzyl)oxy)-2-methylquinolin-3-yl)propan-2-ol;

2-(7-((3-Chlorobenzyl)oxy)-2-methylquinolin-3-yl)propan-2-ol;

2-(7-((3-Chloro-4-fluorobenzyl)oxy)-2-methylquinolin-3-yl)propan-2-ol;

2-(7-((3-Fluorobenzyl)oxy)-2-methylquinolin-3-yl)propan-2-ol;

(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)methanol;

(7-((3,5-Difluorobenzyl)oxy)quinolin-3-yl)methanol;

(7-((3,4,5-Trifluorobenzyl)oxy)quinolin-3-yl)methanol;

2-(7-(3-Fluorophenethoxy)quinolin-3-yl)propan-2-ol;

2-(7-(3-Chlorophenethoxy)quinolin-3-yl)propan-2-ol; and

7-((3-Chlorobenzyl)oxy)-3-(methoxymethyl)quinoline.

29. A chemical entity as in claim 19, wherein the chemical entity is selected from the group consisting of:

7-((4-Fluorobenzyl)oxy)-3-(2-fluoropropan-2-yl)quinoline;

1-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)ethanone;

1-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)ethanol;

(R)-1-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)ethanol;

(S)-1-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)ethanol;

7-((3-Chlorobenzyl)oxy)-N-methylquinoline-3-carboxamide;

N-(2-Aminoethyl)-7-((3-chlorobenzyl)oxy)quinoline-3-carboxamide;

7-((3-Chlorobenzyl)oxy)-N-(2-(methylamino)ethyl)quinoline-3-carboxamide;

7-((3-Chlorobenzyl)oxy)-N-(2-(dimethylamino)ethyl)quinoline-3-carboxamide; and

(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)methanamine.

30. A chemical entity as in claim 19, wherein the chemical entity is selected from the group consisting of:

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)acetamide;

Ethyl 2-(7-((3-fluorobenzyl)oxy)quinolin-3-yl)acetate;

Ethyl 2-(7-((4-fluorobenzyl)oxy)quinolin-3-yl)acetate;

Methyl 2-(7-((3-chlorobenzyl)oxy)quinolin-3-yl)acetate;

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)acetamide;

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)ethanol;

2-(7-((3-Fluorobenzyl)oxy)quinolin-3-yl)acetamide;

2-(7-((4-Fluorobenzyl)oxy)quinolin-3-yl)acetamide;

Ethyl 2-(7-((3-chlorobenzyl)oxy)quinolin-3-yl)-2-methylpropanoate; and

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)-2-methylpropan-1-ol.

31. A chemical entity as in claim 19, wherein the chemical entity is selected from the group consisting of:

2-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)-2-methylpropanamide;

2-(7-((2-(Trifluoromethyl)pyridin-4-yl)methoxy)quinolin-3-yl)acetamide;

1-(7-((3-Chlorobenzyl)oxy)quinolin-3-yl)-2-methylpropan-2-ol;

7-((3-Chlorobenzyl)oxy)quinoline-3-carboxamide;

2-{7-[(4-Fluorophenyl)methoxy]-1,5-naphthyridin-3-yl}propan-2-ol;

7-[(4-Fluorophenyl)methoxy]-3-(2-hydroxypropan-2-yl)quinolin-1-ium-1-olate;

7-[(3-Fluorophenyl)methoxy]-1,5-naphthyridine-3-carboxamide;

7-((4-Fluorobenzyl)oxy)-1,5-naphthyridine-3-carboxamide;

7-((3-Chlorobenzyl)oxy)-1,5-naphthyridine-3-carboxamide;

7-((3,4-Difluorobenzyl)oxy)-1,5-naphthyridine-3-carboxamide;

7-[(3-Chloro-4-fluorophenyl)methoxy]-1,5-naphthyridine-3-carboxamide; and

7-[(3-Chlorophenyl)methoxy]-N-methyl-1,5-naphthyridine-3-carboxamide.

32. A chemical entity as in claim 19, wherein the chemical entity is selected from the group consisting of:

(7-((3,4-Difluorobenzyl)oxy)-1,5-naphthyridin-3-yl)methanol;

2-{7-[(3,4-Difluorophenyl)methoxy]-1,5-naphthyridin-3-yl}acetonitrile;

3-[(4-Fluorophenyl)methoxy]-7-(2-fluoropropan-2-yl)-1,5-naphthyridine;

2-{7-[(3-Fluorophenyl)methoxy]-1,5-naphthyridin-3-yl}propan-2-ol;

2-{7-[(3,4-Difluorophenyl)methoxy]-1,5-naphthyridin-3-yl}propan-2-ol;

2-{7-[(3-Chloro-4-fluorophenyl)methoxy]-1,5-naphthyridin-3-yl}propan-2-ol;

2-{7-[(3-Chlorophenyl)methoxy]-1,5-naphthyridin-3-yl}propan-2-ol;

2-(7-((3,4-Difluorobenzyl)oxy)-1,5-naphthyridin-3-yl)acetamide;

2-(7-((6-(Trifluoromethyl)pyridin-2-yl)methoxy)-1,5-naphthyridin-3-yl)acetamide;

2-(7-((3-Chlorobenzyl)oxy)-2-oxo-1,2,3,4-tetrahydroquinolin-3-yl)acetamide; and

2-(7-((3-Chlorobenzyl)oxy)-2-oxo-1,2-dihydroquinolin-3-yl)acetamide.

33. A chemical entity as in claim 1, wherein the chemical entity is a compound of Formula (I).

34. A chemical entity as in claim 1, wherein the chemical entity is a pharmaceutically acceptable salt of a compound of Formula (I).