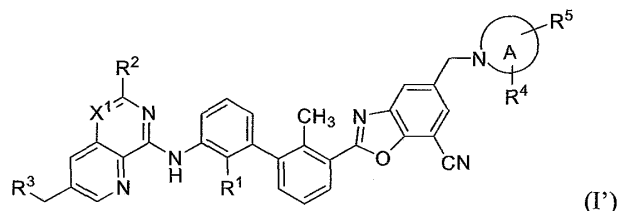


What is claimed is:

1. A compound of Formula (I'):



or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein:

ring A is azetidiny, pyrrolidiny or piperidiny;

X¹ is CH or N;

R¹ is methyl or halo;

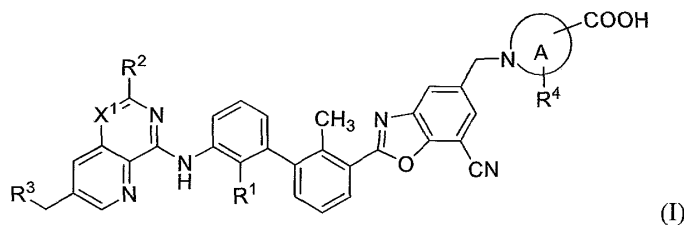
R² is C₁₋₄ alkyl, C₁₋₄ alkoxy, C₁₋₄ haloalkyl, C₁₋₄ haloalkoxy, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl-C₁₋₂ alkyl-, OH, NH₂, -NH-C₁₋₄ alkyl, -N(C₁₋₄ alkyl)₂, 4- to 6-membered heterocycloalkyl or 4- to 6-membered heterocycloalkyl-C₁₋₂ alkyl-, wherein the 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl-C₁₋₂ alkyl- each has one or two heteroatoms as ring members selected from O and N, and wherein the C₁₋₄ alkyl, C₁₋₄ alkoxy, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl-C₁₋₂ alkyl-, -NH-C₁₋₄ alkyl, -N(C₁₋₄ alkyl)₂, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl-C₁₋₂ alkyl- of R² are each optionally substituted with 1 or 2 substituents independently selected from halo, CN and OH;

R³ is selected from (R)-3-hydroxy-3-methylpyrrolidin-1-yl, (S)-3-hydroxy-3-methylpyrrolidin-1-yl, (R)-3-hydroxypyrrolidin-1-yl, (S)-3-hydroxypyrrolidin-1-yl, (R)-2-hydroxy-2-methyl-ethylamino, (S)-2-hydroxy-2-methyl-ethylamino, (R)-2-hydroxy-1-methyl-ethylamino and (S)-2-hydroxy-1-methyl-ethylamino; and

R⁴ is H or C₁₋₃ alkyl; and

R⁵ is C(O)OH, C(O)N(CH₃)₂, C(O)NH(CH₃), or C(O)NH(CH₂)₂C(O)OH.

2. A compound of Formula (I):



or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein:

ring A is azetidiny, pyrrolidinyl or piperidinyl;

X¹ is CH or N;

R¹ is methyl or halo;

R² is C₁₋₄ alkyl, C₁₋₄ alkoxy, C₁₋₄ haloalkyl, C₁₋₄ haloalkoxy, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl-C₁₋₂ alkyl-, OH, NH₂, -NH-C₁₋₄ alkyl, -N(C₁₋₄ alkyl)₂, 4- to 6-membered heterocycloalkyl or 4- to 6-membered heterocycloalkyl-C₁₋₂ alkyl-, wherein the 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl-C₁₋₂ alkyl- each has one or two heteroatoms as ring members selected from O and N, and wherein the C₁₋₄ alkyl, C₁₋₄ alkoxy, C₃₋₆ cycloalkyl, C₃₋₆ cycloalkyl-C₁₋₂ alkyl-, -NH-C₁₋₄ alkyl, -N(C₁₋₄ alkyl)₂, 4- to 6-membered heterocycloalkyl and 4- to 6-membered heterocycloalkyl-C₁₋₂ alkyl- of R² are each optionally substituted with 1 or 2 substituents independently selected from halo, CN and OH;

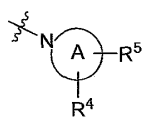
R³ is selected from (R)-3-hydroxy-3-methylpyrrolidin-1-yl, (S)-3-hydroxy-3-methylpyrrolidin-1-yl, (R)-3-hydroxypyrrolidin-1-yl, (S)-3-hydroxypyrrolidin-1-yl, (R)-2-hydroxy-2-methyl-ethylamino, (S)-2-hydroxy-2-methyl-ethylamino, (R)-2-hydroxy-1-methyl-ethylamino and (S)-2-hydroxy-1-methyl-ethylamino; and

R⁴ is H or C₁₋₃ alkyl.

3. The compound of claim 1 or 2, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein ring A is pyrrolidinyl.

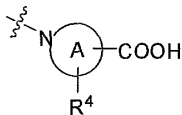
4. The compound of claim 1 or 2, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein ring A is piperidinyl.

5. The compound of claim 1 or 2, or a pharmaceutically acceptable salt or a



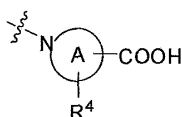
stereoisomer thereof, wherein is selected from 4-carboxypiperidin-1-yl, 3-carboxypyrrolidin-1-yl, 3-methyl-3-carboxypyrrolidin-1-yl, 4-(N,N-dimethylaminocarbonyl)piperidin-1-yl, 4-(N-methylaminocarbonyl)piperidin-1-yl, and 4-(2-carboxyethylaminocarbonyl)piperidin-1-yl, wherein the wavy line indicates the point of attachment to the rest of the molecule.

6. The compound of claim 1 or 2, or a pharmaceutically acceptable salt or a



stereoisomer thereof, wherein R^4 is selected from 4-carboxypiperidin-1-yl, 3-carboxypyrrolidin-1-yl, and 3-methyl-3-carboxypyrrolidin-1-yl, wherein the wavy line indicates the point of attachment to the rest of the molecule.

7. The compound of claim 1 or 2, or a pharmaceutically acceptable salt or a



stereoisomer thereof, wherein R^4 is selected from 4-carboxypiperidin-1-yl, (R)-3-carboxypyrrolidin-1-yl, (S)-3-carboxypyrrolidin-1-yl, (R)-3-methyl-3-carboxypyrrolidin-1-yl and (S)-3-methyl-3-carboxypyrrolidin-1-yl, wherein the wavy line indicates the point of attachment to the rest of the molecule.

8. The compound of any one of claims 1-7, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein X^1 is N.

9. The compound of any one of claims 1-7, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein X^1 is CH.

10. The compound of any one of claims 1-9, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^1 is CH_3 or Cl.

11. The compound of any one of claims 1-9, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^1 is CH_3 .

12. The compound of any one of claims 1-11, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is C_{1-4} alkyl, C_{1-4} alkoxy, C_{1-4} haloalkyl, C_{1-4} haloalkoxy, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl- C_{1-2} alkyl-, OH, NH_2 , -NH- C_{1-4} alkyl, -N(C_{1-4} alkyl) $_2$, 1-azetidiny, azetidin-1-ylmethyl, 1-pyrrolidinyl, pyrrolidin-1-ylmethyl, 1-piperidinyl, or piperidin-1-ylmethyl, wherein the C_{1-4} alkyl, C_{1-4} alkoxy, C_{3-6} cycloalkyl, C_{3-6} cycloalkyl- C_{1-2} alkyl-, -NH- C_{1-4} alkyl, -N(C_{1-4} alkyl) $_2$, 1-azetidiny, azetidin-1-ylmethyl, 1-pyrrolidinyl, pyrrolidin-1-

ylmethyl, 1-piperidinyl and piperidin-1-ylmethyl of R^2 are each optionally substituted with 1 or 2 substituents independently selected from halo, CN and OH.

13. The compound of claim **12**, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is methyl, ethyl, isopropyl, methoxy, ethoxy, CF_3 , CHF_2 , CFH_2 , OCF_3 , $OCHF_2$, OCH_2F , cyclopropyl, cyclobutyl, cyclohexyl, cyclopropylmethyl, cyclobutylmethyl, cyclohexylmethyl, OH, NH_2 , $NHCH_3$, $N(CH_3)_2$, 1-azetidiny, azetidin-1-ylmethyl, 1-pyrrolidinyl, pyrrolidin-1-ylmethyl, 1-piperidinyl or piperidin-1-ylmethyl, wherein the methyl, ethyl, isopropyl, methoxy, ethoxy, cyclopropyl, cyclobutyl, cyclohexyl, cyclopropylmethyl, cyclobutylmethyl, cyclohexylmethyl, $NHCH_3$, $N(CH_3)_2$, 1-azetidiny, azetidin-1-ylmethyl, 1-pyrrolidinyl, pyrrolidin-1-ylmethyl, 1-piperidinyl and piperidin-1-ylmethyl of R^2 are each optionally substituted with 1 or 2 substituents independently selected from F, Cl, Br, CN and OH.

14. The compound of any one of claims **1-11**, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is C_{1-4} alkyl or C_{1-4} haloalkyl, each of which is optionally substituted with 1 or 2 substituents independently selected from F, Cl, Br, CN and OH.

15. The compound of any one of claims **1-11**, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is C_{1-4} alkyl or C_{1-4} haloalkyl.

16. The compound of any one of claims **1-11**, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is CH_3 , CF_3 , CHF_2 , $CH(CH_3)_2$, NH_2 , cyclopropyl, or CH_2OH .

17. The compound of any one of claims **1-11**, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is CH_3 , CF_3 , CHF_2 or $CH(CH_3)_2$.

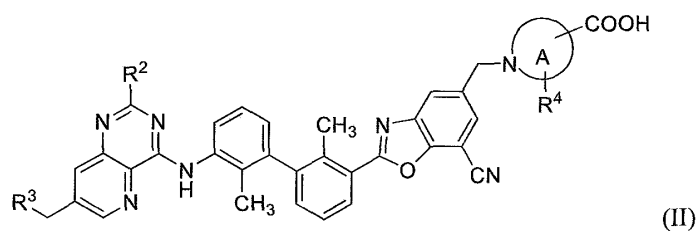
18. The compound of any one of claims **1-11**, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is CH_3 .

19. The compound of any one of claims **1-11**, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is CF_3 or CHF_2 .

20. The compound of any one of claims 1-11, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is $\text{CH}(\text{CH}_3)_2$.
21. The compound of any one of claims 1-11, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is NH_2 .
22. The compound of any one of claims 1-11, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is cyclopropyl.
23. The compound of any one of claims 1-11, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^2 is CH_2OH .
24. The compound of any one of claims 1-23, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^3 is (R)-3-hydroxy-3-methylpyrrolidin-1-yl or (S)-3-hydroxy-3-methylpyrrolidin-1-yl.
25. The compound of any one of claims 1-23, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^3 is (R)-3-hydroxypyrrolidin-1-yl or (S)-3-hydroxypyrrolidin-1-yl.
26. The compound of any one of claims 1-23, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^3 is (R)-2-hydroxy-2-methyl-ethylamino or (S)-2-hydroxy-2-methyl-ethylamino.
27. The compound of any one of claims 1-23, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^3 is (R)-2-hydroxy-1-methyl-ethylamino or (S)-2-hydroxy-1-methyl-ethylamino.
28. The compound of any one of claims 1-27, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^4 is H or CH_3 .
29. The compound of any one of claims 1-27, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R^4 is H.

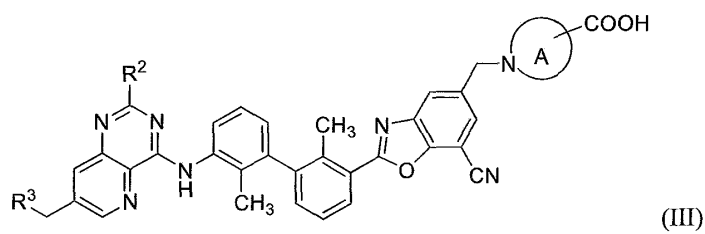
30. The compound of any one of claims 1-27, or a pharmaceutically acceptable salt or a stereoisomer thereof, wherein R⁴ is CH₃.

31. The compound of claim 1 having Formula II:



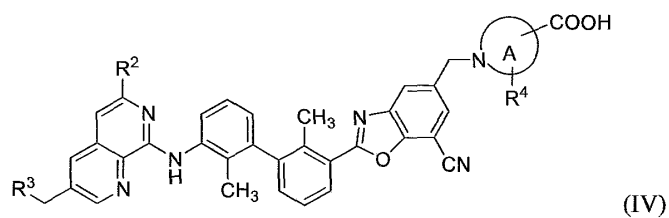
or a pharmaceutically acceptable salt or a stereoisomer thereof.

32. The compound of claim 1 having Formula III:



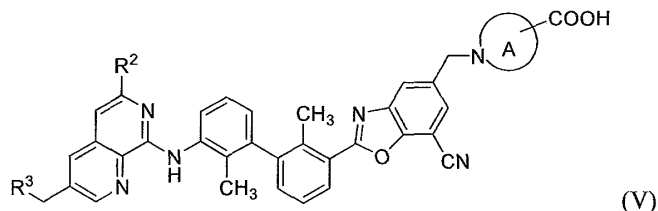
or a pharmaceutically acceptable salt or a stereoisomer thereof.

33. The compound of claim 1 having Formula IV:



or a pharmaceutically acceptable salt or a stereoisomer thereof.

34. The compound of claim 1 having Formula V:



or a pharmaceutically acceptable salt or a stereoisomer thereof.

35. The compound of claim 1, wherein the compound is selected from:

(R)-1-((7-cyano-2-(3'-(7-((3-hydroxypyrrolidin-1-yl)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-((3-hydroxy-3-methylpyrrolidin-1-yl)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(S)-1-((7-cyano-2-(3'-(7-((3-hydroxy-3-methylpyrrolidin-1-yl)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(S)-1-((7-cyano-2-(3'-(7-((1-hydroxypropan-2-ylamino)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(S)-1-((7-cyano-2-(3'-(7-((2-hydroxypropylamino)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((R)-3-hydroxypyrrolidin-1-yl)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)pyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((R)-3-hydroxy-3-methylpyrrolidin-1-yl)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)pyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((S)-3-hydroxy-3-methylpyrrolidin-1-yl)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)pyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((S)-1-hydroxypropan-2-ylamino)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)pyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((S)-2-hydroxypropylamino)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)pyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((R)-3-hydroxypyrrolidin-1-yl)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-3-methylpyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((R)-3-hydroxy-3-methylpyrrolidin-1-yl)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-3-methylpyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((S)-3-hydroxy-3-methylpyrrolidin-1-yl)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-3-methylpyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((S)-1-hydroxypropan-2-ylamino)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-3-methylpyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((S)-2-hydroxypropylamino)methyl)-2-methylpyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-3-methylpyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-((3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-(((R)-3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)pyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-(((R)-3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-3-methylpyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-((3-hydroxypyrrolidin-1-yl)methyl)-2-(trifluoromethyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(7-(((R)-3-hydroxypyrrolidin-1-yl)methyl)-2-(trifluoromethyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)pyrrolidine-3-carboxylic acid; and

(R)-1-((7-cyano-2-(3'-(7-(((R)-3-hydroxypyrrolidin-1-yl)methyl)-2-(trifluoromethyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-3-methylpyrrolidine-3-carboxylic acid,

or a pharmaceutically acceptable salt or a stereoisomer thereof.

36. The compound of claim 1, wherein the compound is selected from:

(S)-1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-(((R)-3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)pyrrolidine-3-carboxylic acid;

(S)-1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-(((R)-3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-3-methylpyrrolidine-3-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-((3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-4-methylpiperidine-4-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-((3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-N,N-dimethylpiperidine-4-carboxamide;

(R)-1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-((3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-N-methylpiperidine-4-carboxamide;

(R)-3-(1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-((3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxamido)propanoic acid;

(R)-1-((7-cyano-2-(3'-(2-cyclopropyl-7-(((R)-3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)pyrrolidine-3-carboxylic acid;

(R)-1-((2-(3'-(2-amino-7-((3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)-7-cyanobenzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-((3-hydroxy-3-methylpyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(S)-1-((7-cyano-2-(3'-(2-(difluoromethyl)-7-((3-hydroxy-3-methylpyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(2-(hydroxymethyl)-7-((3-hydroxypyrrolidin-1-yl)methyl)pyrido[3,2-d]pyrimidin-4-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)piperidine-4-carboxylic acid;

(R)-1-((7-cyano-2-(3'-(3-(((R)-3-hydroxypyrrolidin-1-yl)methyl)-6-methyl-1,7-naphthyridin-8-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)pyrrolidine-3-carboxylic acid; and

(R)-1-((7-cyano-2-(3'-(6-(difluoromethyl)-3-(((R)-3-hydroxypyrrolidin-1-yl)methyl)-1,7-naphthyridin-8-ylamino)-2,2'-dimethylbiphenyl-3-yl)benzo[d]oxazol-5-yl)methyl)-3-methylpyrrolidine-3-carboxylic acid,

or a pharmaceutically acceptable salt or a stereoisomer thereof.

37. A pharmaceutical composition comprising a compound of any one of claims **1-36**, or a pharmaceutically acceptable salt or a stereoisomer thereof, and a pharmaceutically acceptable excipient or carrier.

38. A method of inhibiting PD-1/PD-L1 interaction, said method comprising administering to a patient a compound of any one of claims **1-36**, or a pharmaceutically acceptable salt or a stereoisomer thereof.

39. A method of treating a disease or disorder associated with inhibition of PD-1/PD-L1 interaction, said method comprising administering to a patient in need thereof a therapeutically effective amount of a compound of any one of claims **1-36**, or a pharmaceutically acceptable salt or a stereoisomer thereof.

40. The method of claim **39**, wherein the disease or disorder is an infection disease, inflammation, autoimmune disease, cancer, or neurodegenerative disorder

41. A method of enhancing, stimulating and/or increasing the immune response in a patient, said method comprising administering to the patient in need thereof a therapeutically effective amount of a compound of any one of claims **1-36**, or a pharmaceutically acceptable salt or a stereoisomer thereof.