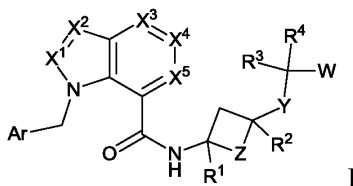


CLAIMS

WHAT IS CLAIMED IS:

1. A compound of Formula I:



or a pharmaceutically acceptable salt, solvate, solvate of the salt, hydrate, a single stereoisomer, a mixture of stereoisomers, a racemic mixture of stereoisomers, or prodrug thereof, wherein:

Ar is an aryl or a heteroaryl, wherein said aryl and said heteroaryl are each optionally substituted with 1 to 3 substituents independently selected from:

- (a) C₁-C₆ alkyl,
- (b) C₃-C₇ cycloalkyl,
- (c) heterocycle,
- (d) aryl,
- (e) heteroaryl,
- (f) halogen,
- (g) CN,
- (h) OR^b,
- (i) N(R^b)C(=O)R^c,
- (j) C(=O)N(R^b)(R^c),
- (k) S(=O)_mR^b,
- (l) S(=O)₂N(R^b)(R^c),
- (m) N(R^b)S(=O)₂R^c,
- (n) SF₅; and
- (o) C₁-C₆ haloalkyl

W is selected from:

- (a) C(=O)OR⁵,
- (b) C(=O)NHOH,
- (c) S(=O)₂NHR^b,
- (d) S(=O)₂NHC(=O)R^b,
- (e) NHC(=O)NHSO₂R^b,
- (f) 1*H*-tetrazole,
- (g) 1,2,4-oxadiazol-5(4*H*)one,

- (h) 1,2,4-thiadiazol-5(4*H*)one,
- (i) 1,2,4-oxadiazole-5(4*H*)-thione,
- (j) 1,2,4-triazole-5(4*H*)-one,
- (k) tetrazol-5(4*H*)-one, and
- (l) $C(=O)NHS(=O)_2R^b$;

X^1 , X^2 , X^3 , X^4 , and X^5 are each independently N or CR^a , wherein not more than 2 of X^1 , X^2 , X^3 , X^4 , and X^5 are N;

Y is selected from:

- (a) a bond,
- (b) $(CH_2)_n$ wherein 1 to 4 hydrogen atoms may be replaced by R^a ,
- (c) O, and
- (d) NR^b ;

Z is $(CH_2)_n$, wherein 1 to 4 hydrogen atoms may be replaced by R^a ;

R^1 and R^2 are independently selected from:

- (a) H,
- (b) C_1 - C_6 alkyl,
- (c) C_3 - C_6 cycloalkyl, and
- (d) C_1 - C_6 haloalkyl;

wherein R^2 is not H; or

R^1 and R^2 taken together represent $-(CH_2)_n$ -, $-(CH_2)_nO(CH_2)_p$ -, $-(CH_2)_nNR^b(CH_2)_p$ -, or $-(CH_2)_nS(=O)_m(CH_2)_p$;

R^3 and R^4 are independently selected from:

- (a) H,
- (b) C_1 - C_6 alkyl,
- (c) C_3 - C_6 cycloalkyl,
- (d) aryl,
- (e) heteroaryl,
- (f) halogen,
- (g) C_1 - C_6 haloalkyl; or

R^3 and R^4 taken together represent $-(CH_2)_n$ -, $-(CH_2)_nO(CH_2)_p$ -, $-(CH_2)_nNR^b(CH_2)_p$ -, or $-(CH_2)_nS(=O)_m(CH_2)_p$;

or R^1 , R^2 , R^3 and R^4 above are selected as follows:

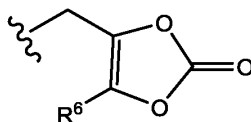
R^1 is selected from H, C_1 - C_6 alkyl, C_3 - C_6 cycloalkyl, and C_1 - C_6 haloalkyl;

R^3 and R^2 taken together represent $(CH_2)_n$, $(CH_2)_nO(CH_2)_p$, $(CH_2)_nNR^b(CH_2)_p$, or $(CH_2)_nS(=O)_m(CH_2)_p$; and

R⁴ is selected from H, C₁-C₆ alkyl, C₃-C₆ cycloalkyl, aryl, heteroaryl, halogen and C₁-C₆ haloalkyl;

R⁵ is selected from

- (a) H,
- (b) C₁-C₆ alkyl,
- (c) aryl,
- (d) aralkyl,
- (e) CH(R⁷)OC(=O)R⁸,
- (f) CH(R⁷)OC(=O)OR⁸, and
- (g) a (5-alkyl-2-oxo-1,3-dioxolen-4-yl)methyl group having the following formula:



wherein R⁶ is C₁-C₆ alkyl;

R⁷ is hydrogen or C₁-C₆ alkyl;

R⁸ is C₁-C₆ alkyl or C₃-C₆ -cycloalkyl;

R^a is selected from:

- (a) H,
- (b) C₁-C₆ alkyl,
- (c) halogen,
- (d) aryl,
- (e) OR^b,
- (f) cyano,
- (g) heteroaryl,
- (h) C₃-C₆ cycloalkyl, and
- (i) C₁-C₆ haloalkyl;

R^{a'} is selected from:

- (a) cyano,
- (b) C₁-C₆ alkyl,
- (c) halogen,
- (d) aryl,
- (e) OR^b,
- (f) heteroaryl,

(g) C₃-C₆ cycloalkyl, and

(h) C₁-C₆ haloalkyl;

R^b and R^c are independently selected from:

(a) H,

(b) C₁-C₆ alkyl,

(c) aryl,

(d) heteroaryl,

(e) C₃-C₆ cycloalkyl, and

(f) C₁-C₆ haloalkyl; or

R^b and R^c taken together with the N to which they are both attached form a 3- to 6-membered heterocycle optionally having an additional heteroatom selected from N, O and S;

m is 0, 1, or 2;

n is 1, 2 or 3; and

p is 1, 2 or 3.

2. A compound of Claim 1 wherein Ar is aryl or heteroaryl optionally substituted with 1 to 3 substituents independently selected from:

(a) halogen,

(b) cyano,

(c) C₁-C₆ alkyl,

(d) SF₅,

(e) C₁-C₆ haloalkyl,

(f) OR^b wherein R^b is C₁-C₆ alkyl, aryl, heteroaryl, C₃-C₆ cycloalkyl or C₁-C₆ haloalkyl;

(g) heterocycle,

(h) aryl, and

(i) heteroaryl.

3. A compound of Claim 1 or Claim 2 wherein Ar is phenyl.

4. A compound any one of Claims 1 to 3 wherein each of X¹, X², X³, X⁴ and X⁵ is independently C-R^a, or one of X¹, X², X³, X⁴ and X⁵ is N, and the others are each independently C-R^a.

5. A compound of Claim 4 wherein each of X¹, X², X³, X⁴ and X⁵ is independently C-R^a.

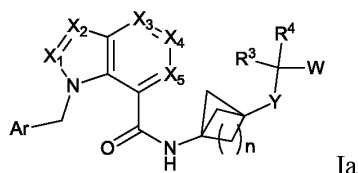
6. A compound of any of Claims 1 to 5 wherein R^a is H or a halogen atom.

7. A compound of any of Claims 1 to 6 wherein W is selected from the group consisting of:

(a) CO₂H; and

(b) 1H-tetrazole.

8. A compound of any of Claims 1 to 7 wherein Z is -CH₂-.
9. A compound of any of Claims 1 to 8 wherein Y is a bond or -CH₂-.
10. A compound of any of Claims 1 to 9 wherein R¹ and R² taken together represent -CH₂-, -CH₂CH₂- or -CH₂CH₂CH₂-.
11. A compound of any of Claims 1 to 8 wherein Y is -CH₂-, and R³ and R² taken together represent -CH₂-, -CH₂CH₂- or -CH₂CH₂CH₂-.
12. A compound of Claim 1 having the Formula Ia:



or a pharmaceutically acceptable salt, solvate, solvate of the salt, hydrate, a single stereoisomer, a mixture of stereoisomers, a racemic mixture of stereoisomers, or prodrug thereof, wherein Ar is phenyl optionally substituted with 1 to 3 substituents independently selected from:

- (a) halogen,
- (b) cyano,
- (c) C₁-C₆ alkyl,
- (d) SF₅,
- (e) C₁-C₆ haloalkyl,
- (f) OR^b wherein R^b is C₁-C₆ alkyl, aryl, heteroaryl, C₃-C₆ cycloalkyl or C₁-C₆ haloalkyl,
- (g) heterocycle,
- (h) aryl, and
- (i) heteroaryl;

each X¹, X², X³, X⁴ and X⁵ is independently C-R^a, or one of X¹, X², X³, X⁴ and X⁵ is N, and the others are each independently C-R^a;

W is selected from:

- (a) CO₂H and
- (b) 1H-tetrazole;

Y is a bond or -CH₂-;

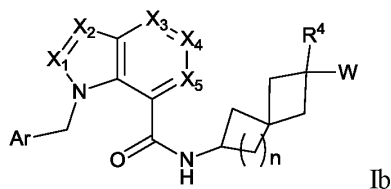
n is 1 or 2;

R³ and R⁴ are independently selected:

- (a) H,
- (b) C₁-C₃ alkyl, and
- (c) C₁-C₃ haloalkyl;

R^a is selected from:

- (a) H and
 - (b) halogen.
13. A compound of any of Claims 1 to 12 wherein Y is a bond, and n is 1.
14. A compound of Claim 1 having the Formula Ib:



or a pharmaceutically acceptable salt, solvate, solvate of the salt, hydrate, a single stereoisomer, a mixture of stereoisomers, a racemic mixture of stereoisomers, or prodrug thereof, wherein

Ar is phenyl optionally substituted with 1 to 3 substituents independently selected from:

- (a) halogen,
- (b) cyano,
- (c) C₁-C₆ alkyl,
- (d) SF₅,
- (e) C₁-C₆ haloalkyl,
- (f) OR^b wherein R^b is C₁-C₆ alkyl, aryl, heteroaryl, C₃-C₆ cycloalkyl or C₁-C₆ haloalkyl,
- (g) heterocycle,
- (h) aryl, and
- (i) heteroaryl;

each X¹, X², X³, X⁴ and X⁵ is independently C-R^a, or one of X¹, X², X³, X⁴ and X⁵ is N, and the others are each independently C-R^a;

W is selected from :

- (a) CO₂H and
- (b) 1H-tetrazole;

n is 1 or 2;

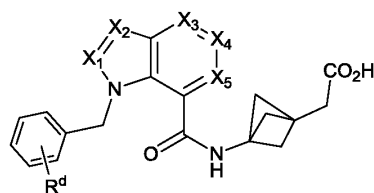
R⁴ is selected from :

- (a) H,
- (b) C₁-C₃ alkyl, and
- (c) C₁-C₃ haloalkyl;

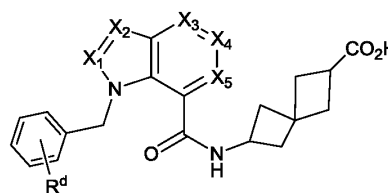
R^a is selected from :

- (a) H and

- (b) halogen.
15. A compound of Claim 14 wherein n is 1.
16. A compound of Claim 1 having the Formula Ic or Id:



Ic



Id

or a pharmaceutically acceptable salt, solvate, solvate of the salt, hydrate, a single stereoisomer, a mixture of stereoisomers, a racemic mixture of stereoisomers, or prodrug thereof, wherein

each X^1 , X^2 , X^3 , X^4 and X^5 is independently $C-R^a$, or one of X^1 , X^2 , X^3 , X^4 and X^5 is N, and the others are each independently $C-R^a$;

R^a is selected from :

- (a) H and
- (b) halogen;

R^d is selected from :

- (a) CN,
- (b) C_1 - C_3 alkyl,
- (c) SF_5 ,
- (d) C_1 - C_3 haloalkyl,
- (e) OR^b wherein R^b is C_1 - C_6 alkyl, aryl, heteroaryl, C_3 - C_6 cycloalkyl or C_1 - C_6 haloalkyl,
- (f) heterocycle,
- (g) aryl, and
- (h) heteroaryl.

17. A compound of any of Claims 1 to 16 wherein each X^1 , X^2 , X^3 , X^4 and X^5 is CH; or one of X^1 , X^2 , X^3 , X^4 and X^5 is $C-R^a$ and the others are CH, and R^a is halogen.

18. A compound of Claim 1 selected from :

2-(3-(1-(4-cyanobenzyl)-1*H*-indole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)acetic acid;

2-(3-(1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)acetic acid;

2-(3-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[3,2-*b*]pyridine-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)acetic acid;

2-(3-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[3,2-*c*]pyridine-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)acetic acid;

2-(3-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[2,3-*c*]pyridine-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)acetic acid;

2-(3-(1-(4-(trifluoromethyl)benzyl)-1*H*-indazole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)acetic acid;

2-(3-(1-((4-(trifluoromethyl)phenyl)methyl-*d*₂)-1*H*-indole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)acetic acid;

6-(1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*R*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*S*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

6-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[3,2-*b*]pyridine-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*R*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[3,2-*b*]pyridine-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*S*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[3,2-*b*]pyridine-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

6-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[3,2-*c*]pyridine-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*R*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[3,2-*c*]pyridine-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*S*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[3,2-*c*]pyridine-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

6-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[2,3-*c*]pyridine-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*R*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[2,3-*c*]pyridine-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*S*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-pyrrolo[2,3-*c*]pyridine-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

6-(1-(4-(trifluoromethyl)benzyl)-1*H*-indazole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*R*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-indazole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*S*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-indazole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

6-(1-(4-(trifluoromethyl)benzyl)-1*H*-benzo[*d*]imidazole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*R*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-benzo[*d*]imidazole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*S*)-6-(1-(4-(trifluoromethyl)benzyl)-1*H*-benzo[*d*]imidazole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

3-(3-(1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)propanoic acid;

3-(3-methyl-3-(1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)cyclobutyl)propanoic acid;

cis-3-(3-methyl-3-(1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)cyclobutyl)propanoic acid;

N-(3-(2-oxo-2-(phenylsulfonamido)ethyl)bicyclo[1.1.1]pentan-1-yl)-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamide;

N-(3-((3-(phenylsulfonyl)ureido)methyl)bicyclo[1.1.1]pentan-1-yl)-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamide;

N-(3-((1*H*-tetrazol-5-yl)methyl)bicyclo[1.1.1]pentan-1-yl)-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamide;

2-(4-(1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)bicyclo[2.1.1]hexan-1-yl)acetic acid;

6-(4-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*R*)-6-(4-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*S*)-6-(4-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

6-(5-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*R*)-6-(5-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

(*S*)-6-(5-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

6-(5-chloro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*R*)-6-(5-chloro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*S*)-6-(5-chloro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
6-(6-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*R*)-6-(6-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*S*)-6-(6-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
6-(1-(4-cyanobenzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*R*)-6-(1-(4-cyanobenzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*S*)-6-(1-(4-cyanobenzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
6-(1-(4-(difluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*R*)-6-(1-(4-(difluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*S*)-6-(1-(4-(difluoromethyl)benzyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
2-(4-(1-(4-(difluoromethyl)benzyl)-1*H*-indole-7-carboxamido)bicyclo[2.1.1]hexan-1-yl)acetic acid;
2-(3-(4-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)acetic acid;
2-(4-(4-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)bicyclo[2.1.1]hexan-1-yl)acetic acid;
2-(4-(1-((4-(pentafluorothiol)phenyl)methyl)-1*H*-indole-7-carboxamido)bicyclo[2.1.1]hexan-1-yl)acetic acid;
6-(1-((4-(pentafluorothiol)phenyl)methyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*R*)-6-(1-((4-(pentafluorothiol)phenyl)methyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*S*)-6-(1-((4-(pentafluorothiol)phenyl)methyl)-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;

6-(1-(4-(difluoromethyl)benzyl)-4-fluoro-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*R*)-6-(1-(4-(difluoromethyl)benzyl)-4-fluoro-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
(*S*)-6-(1-(4-(difluoromethyl)benzyl)-4-fluoro-1*H*-indole-7-carboxamido)spiro[3.3]heptane-2-carboxylic acid;
2-(3-(1-(4-(difluoromethyl)benzyl)-4-fluoro-1*H*-indole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)acetic acid;
2-(3-(5-fluoro-1-(4-(trifluoromethyl)benzyl)-1*H*-indole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl)acetic acid;
2-(3-(1-([1,1'-biphenyl]-4-ylmethyl)-1*H*-indole-7-carboxamido)bicyclo[1.1.1] pentan-1-yl)acetic acid;
2-(3-(1-([1,1'-biphenyl]-4-ylmethyl)-4-fluoro-1*H*-indole-7-carboxamido)bicyclo[1.1.1] pentan-1-yl)acetic acid;
2-(3-(1-([1,1'-biphenyl]-4-ylmethyl)-1*H*-indazole-7-carboxamido)bicyclo[1.1.1] pentan-1-yl)acetic acid;
2-(3-(1-(4-(trifluoromethoxy)benzyl)-1*H*-indazole-7-carboxamido)bicyclo[1.1.1] pentan-1-yl)acetic acid;
2-(3-(4-fluoro-1-(4-iodobenzyl)-1*H*-indole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl) acetic acid;
2-(3-(4-fluoro-1-(4-(pyridine-4-yl)benzyl)-1*H*-indole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl) acetic acid; and
2-(3-(4-fluoro-1-(4-morpholinobenzyl)-1*H*-indole-7-carboxamido)bicyclo[1.1.1]pentan-1-yl) acetic acid;
or a pharmaceutically acceptable salt, solvate, solvate of the salt or prodrug thereof.

19. A pharmaceutical composition comprising a compound of any of Claims 1 to 18, or a pharmaceutically acceptable salt, solvate, solvate of the salt, hydrate, a single stereoisomer, a mixture of stereoisomers, a racemic mixture of stereoisomers, or prodrug thereof, and a pharmaceutically acceptable carrier.

20. A method for the treatment of cancer comprising administering to a patient in need thereof a compound of any of Claims 1 to 18 or the pharmaceutical composition of Claim 19.

21. The method of Claim 20 wherein said cancer is selected from glioblastoma, bone cancer, head and neck cancer, melanoma, basal cell carcinoma, squamous cell carcinoma, adenocarcinoma, oral cancer, esophageal cancer, gastric cancer, intestinal cancer, colon cancer,

bladder cancer, hepatocellular carcinoma, renal cell carcinoma, pancreatic cancer, ovarian cancer, cervical cancer, lung cancer, breast cancer, and prostate cancer.

22. The method of Claim 20 or Claim 21 wherein said cancer is selected from colon cancer, bladder cancer, hepatocellular carcinoma, pancreatic cancer, ovarian cancer, cervical cancer, lung cancer, breast cancer, and prostate cancer.

23. The method of any of Claims 20 to 22 wherein said treatment further comprises an additional agent selected from an anti-PD-1 antibody and an anti-PD-L1 antibody.