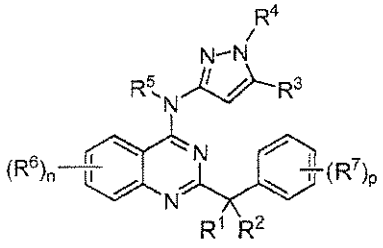
What is claimed is:

1. A compound having formula (I):



(I)

or pharmaceutically acceptable salts, solvates or hydrates thereof, wherein

R^1 and R^2 are selected from (i), (ii), (iii), (iv) and (v) as follows:

(i) R^1 and R^2 together form =O, =S, =NR^9 or =$CR^{10}R^{11}$;

(ii) R^1 and R^2 are both -OR^8, or R^1 and R^2, together with the carbon atom to which they are attached, form dioxacycloalkyl;

(iii) R^1 is hydrogen or halo; and R^2 is halo;

(iv) R^1 is alkyl, alkenyl, alkynyl, cycloalkyl or aryl, wherein the alkyl, alkenyl, alkynyl, cycloalkyl or aryl is optionally substituted with one or more substitutents selected from halo, cyano, alkyl, -R^xOR^w, -$R^xS(O)_qR^v$, -$R^xNR^yR^z$ and -$C(O)OR^w$; and R^2 is hydrogen, halo or -OR^8; and

(v) R^1 is halo, deutero, -OR^{12}, -$NR^{13}R^{14}$, or -$S(O)_qR^{15}$; and R^2 is hydrogen, deutero, alkyl, alkenyl, alkynyl, cycloalkyl or aryl, wherein the alkyl, alkenyl, alkynyl, cycloalkyl or aryl, is optionally substituted with one or more substitutents selected from halo, cyano, alkyl, -R^xOR^w, -$R^xS(O)_qR^v$ and -$R^xNR^yR^z$;

R^3 is hydrogen, halo, alkyl, cyano, haloalkyl, cycloalkyl, cycloalkylalkyl, hydroxy or alkoxy;

R^4 and R^5 are each independently hydrogen or alkyl;

each R^6 is independently selected from halo, alkyl, alkenyl, alkynyl, haloalkyl, cycloalkyl, -R^xOR^{18}, -$R^xNR^{19}R^{20}$, and -$R^xS(O)_qR^v$;

each R^7 is independently halo, alkyl, haloalkyl or -R^xOR^w;

R^8 is alkyl, alkenyl or alkynyl;

R^9 is hydrogen, alkyl, haloalkyl, hydroxy, alkoxy or amino;

R^{10} is hydrogen or alkyl;

R^{11} is hydrogen, alkyl, haloalkyl or -$C(O)OR^8$;

R^{12} is selected from hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl, heteroaralkyl, $-C(O)R^v$, $-C(O)OR^w$ and $-C(O)NR^yR^z$, wherein the alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl or heteroaralkyl are each optionally substituted with one or more substituents independently selected from halo, oxo, alkyl, hydroxy, alkoxy, amino and alkylthio;

R^{13} and R^{14} are selected as follows:

(i) R^{13} is hydrogen or alkyl; and R^{14} is selected from hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl, heteroaralkyl, alkoxy, $-C(O)R^v$, $-C(O)OR^w$, $-C(O)NR^yR^z$ and $-S(O)_qR^v$, wherein the alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl or heteroaralkyl are each optionally substituted with one or more substituents independently selected from halo, oxo, alkyl, hydroxy, alkoxy, amino and alkylthio; or

(ii) R^{13} and R^{14}, together with the nitrogen atom to which they are attached, form heterocyclyl or heteroaryl wherein the heterocyclyl or heteroaryl is optionally substituted with one or more substituents independently selected from halo, alkyl, hydroxy, alkoxy, amino and alkylthio, and wherein the heterocyclyl is also optionally substituted with oxo;

R^{15} is alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl, heteroaralkyl, $-C(O)NR^yR^z$ or $-NR^yR^z$, wherein the alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl or heteroaralkyl are each optionally substituted with one or more substituents independently selected from halo, oxo, alkyl, hydroxy, alkoxy, amino and alkylthio;

R^{18} is hydrogen, alkyl, haloalkyl, hydroxyC_{2-6}alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl or heteroarylalkyl; wherein R^{18} is optionally substituted with 1 to 3 groups Q^1, each Q^1 independently selected from alkyl, hydroxyl, halo, haloalkyl, alkoxy, aryloxy, alkoxyalkyl, alkoxycarbonyl, alkoxysulfonyl, carboxyl, cycloalkyl, heterocyclyl, aryl, heteroaryl, haloaryl and amino;

R^{19} and R^{20} are selected as follows:

(i) R^{19} and R^{20} are each independently hydrogen or alkyl; or

(ii) R^{19} and R^{20}, together with the nitrogen atom to which they are attached, form a heterocyclyl or heteroaryl which is optionally substituted with 1 to 2 groups each independently selected from halo, alkyl, haloalkyl, hydroxyl and alkoxy;

each R^x is independently alkylene or a direct bond;

R^y is hydrogen, alkyl, alkenyl or alkynyl;

R^w is independently hydrogen, alkyl, alkenyl, alkynyl or haloalkyl;

R^y and R^z are selected as follows:

(i) R^y and R^z are each independently hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl or haloalkyl;

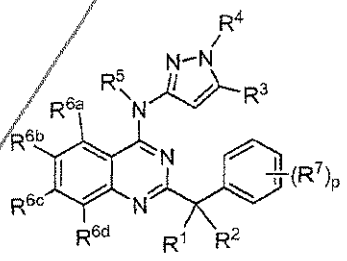
(ii) R^y and R^z, together with the nitrogen atom to which they are attached, form a heterocyclyl or heteroaryl which is optionally substituted with 1 to 2 groups each independently selected from halo, alkyl, haloalkyl, hydroxyl and alkoxy;

n is 0-4;

p is 0-5; and

each q is independently 0, 1 or 2.

2. The compound of claim 1 having formula (II)



(II)

or a pharmaceutically acceptable salt, solvate or hydrate thereof, wherein

R^1 and R^2 are selected from (i), (ii), (iii), (iv) and (v) as follows:

(i) R^1 and R^2 together form =O, =S, $=NR^9$ or $=CR^{10}R^{11}$;

(ii) R^1 and R^2, are both $-OR^8$, or R^1 and R^2, together with the carbon atom to which they are attached, form dioxacycloalkyl;

(iii) R^1 is hydrogen or halo, and R^2 is halo;

(iv) R^1 is alkyl, alkenyl, alkynyl, cycloalkyl or aryl, wherein the alkyl, alkenyl, alkynyl, cycloalkyl or aryl is optionally substituted with one or more substitutents selected from halo, alkyl, $-R^xOR^w$, $-R^xS(O)_qR^v$ and $-R^xNR^yR^z$, and R^2 is hydrogen, halo or $-OR^8$; and

(v) R^1 is halo, $-OR^{12}$, $-NR^{13}R^{14}$, $-S(O)_qR^{15}$ or $-R^{17}C(O)OR^{18}$, and R^2 is hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl or aryl, wherein the alkyl, alkenyl, alkynyl, cycloalkyl or aryl is optionally substituted with one or more substitutents selected from halo, alkyl, $-R^xOR^w$, $-R^xS(O)_qR^v$ and $-R^xNR^yR^z$;

R^3 is hydrogen, alkyl or, cycloalkyl;

R^4 and R^5 are each independently hydrogen or alkyl;

R^{6a}, R^{6b}, R^{6c}, and R^{6d} are each independently selected from halo, alkyl, haloalkyl and $-R^xOR^{18}$;

each R^7 is independently selected from halo, alkyl, haloalkyl or $-R^xOR^w$;

R^8 is alkyl, alkenyl or alkynyl;

R^9 is hydrogen, alkyl, haloalkyl, hydroxy, alkoxy or amino;

R^{10} is hydrogen or alkyl;

R^{11} is hydrogen, alkyl, haloalkyl or $-C(O)OR^8$;

each R^{12} is independently hydrogen, alkyl, haloalkyl, hydroxyalkyl, alkoxyalkyl, aminoalkyl, thioalkyl, heterocyclylalkyl or $-C(O)NR^yR^z$;

R^{13} and R^{14} are selected as follows:

(i) R^{13} is hydrogen or alkyl, and R^{14} is selected from hydrogen, alkyl, haloalkyl, hydroxyalkyl, alkoxyalkyl, aminoalkyl, thioalkyl, heterocycylalkyl, $-C(O)OR^v$, $-C(O)OR^w$, $-C(O)NR^yR^z$ and $-S(O)_qR^v$; or

(ii) R^{13} and R^{14}, together with the nitrogen atom to which they are attached, form heterocyclyl optionally substituted with one more more substituents independently selected from halo, oxo, alkyl, hydroxy, alkoxy, amino and alkylthio;

R^{15} is selected from hydrogen, alkyl, haloalkyl, hydroxyalkyl, alkoxyalkyl, aminoalkyl, thioalkyl, heterocycylalkyl, $-C(O)NR^yR^z$ and $-NR^yR^z$;

R^{18} is hydrogen, alkyl, haloalkyl, hydroxyalkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl or heteroarylalkyl; wherein R^{18} is optionally substituted with 1 to 3 groups Q^1, where

each Q^1 independently selected from alkyl, hydroxyl, halo, haloalkyl, alkoxy, aryloxy, alkoxyalkyl, alkoxycarbonyl, carboxyl, cycloalkyl, heterocyclyl, aryl, heteroaryl, haloaryl and amino;

R^v is alkyl, alkenyl or alkynyl;

each R^x is independently alkylene or a direct bond;

R^w is independently hydrogen or alkyl;

R^y and R^z are selected as follows:

(i) R^y and R^z are each independently hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl or haloalkyl; or

(ii) R^y and R^z, together with the nitrogen atom to which they are attached, form a heterocyclyl or heteroaryl which is optionally substituted with 1 to 2 groups each independently selected from halo, alkyl, haloalkyl, hydroxyl and alkoxy; and

each q is independently 0, 1 or 2.

3. The compound of claim 1 or 2, wherein R^1 and R^2 together form =O

4. The compound of claim 1 or 2, wherein R^1 is hydrogen or fluoro, and R^2 is fluoro.

5. The compound of claims 1 or 2, wherein R^1 is hydroxyl, methoxy, amino or methoxycarbonylamino, and R^2 is hydrogen, methyl, phenyl or fluorophenyl.

6. The compound of any of claims 1-5, wherein R^3 is hydrogen, alkyl or alkoxy.

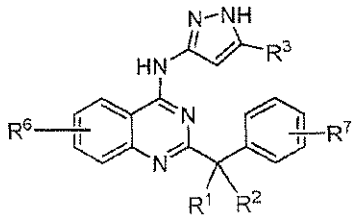
7. The compound of any of claims 1 or 3-6, wherein each R^6 is independently selected from hydrogen, halo, alkyl, hydroxyl and alkoxy.

8. The compound of any of claims 1-7, wherein R^7 is halo.

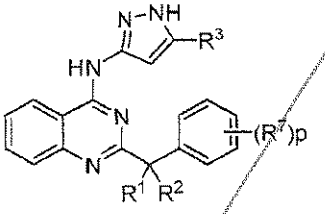
9. The compound of any of claims 1 or 3-7, wherein n is 0 or 1.

10. The compound of any of claims 1-9, wherein p is 1 or 2.

11. The compound of claim 1 having formula (III) or (IIIa)



(III) or



(IIIa)

or a pharmaceutically acceptable salt thereof, wherein:

R^1 and R^2 are selected from (i) and (ii) as follows:

(i) R^1 and R^2 are both alkoxy or R^1 and R^2, together form =O; and

(ii) R^1 is $-OR^{12}$ or $-NR^{13}R^{14}$, and R^2 is hydrogen, alkyl, aryl or haloaryl;

R^{12} is selected from hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl, heteroaralkyl, $-C(O)R^v$, $-C(O)OR^w$ and $-C(O)NR^yR^z$, wherein the alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl or heteroaralkyl are each optionally substituted with one or more substituents independently selected from halo, oxo, alkyl, hydroxy, alkoxy, amino and alkylthio;

R^{13} and R^{14} are selected as follows:

(i) R^{13} is hydrogen or alkyl and R^{14} is selected from hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl, heteroaralkyl, alkoxy, $-C(O)R^v$, $-C(O)OR^w$, $-C(O)NR^yR^z$ and $-S(O)_qR^v$ wherein the alkyl, alkenyl, alkynyl, cycloalkyl, cycloalkylalkyl, heterocyclyl, heterocyclylalkyl, aryl, aralkyl, heteroaryl or heteroaralkyl are each optionally substituted with one or more substituents independently selected from halo, oxo, alkyl, hydroxy, alkoxy, amino and alkylthio; or

(ii) R^{13} and R^{14}, together with the nitrogen atom to which they are attached, form heterocyclyl or heteroaryl wherein the heterocyclyl or heteroaryl is optionally substituted with one or more substituents independently selected from halo, alkyl,

hydroxy, alkoxy, amino and alkylthio and wherein the heterocyclyl is also optionally substituted with oxo;

R^3 is hydrogen or alkyl;

each R^6 is independently selected from halo, alkyl, haloalkyl and $-R^xOR^{18}$; where R^{18} is hydrogen, alkyl, haloalkyl, hydroxyalkyl or heterocyclyl; wherein R^{18} is optionally substituted with group Q^1, where Q^1 is selected from hydroxyl, alkoxy, alkoxycarbonyl, carboxyl, heterocyclyl and amino;

each R^7 is independently halo, alkyl, haloalkyl, hydroxyl or alkoxy;

R^v is hydrogen, alkyl, alkenyl or alkynyl;

R^w is independently hydrogen, alkyl, alkenyl, alkynyl or haloalkyl;

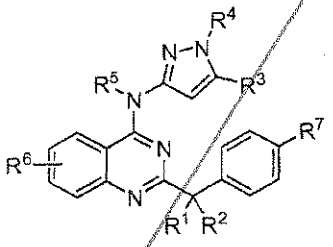
R^y and R^z are selected as follows:

(i) R^y and R^z are each independently hydrogen, alkyl, alkenyl, alkynyl, cycloalkyl or haloalkyl; or

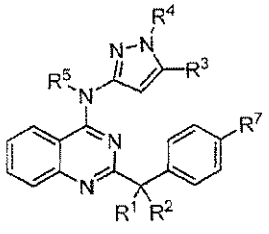
(ii) R^y and R^z, together with the nitrogen atom to which they are attached, form a heterocyclyl or heteroaryl which is optionally substituted with 1 to 2 groups each independently selected from halo, alkyl, haloalkyl, hydroxyl and alkoxy;

and q is 0, 1 or 2.

12. The compound of claim 1 having formula (IV) or (IVa)



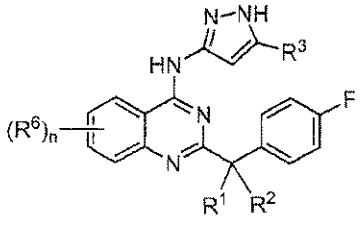
(IV) or



(IVa)

or a pharmaceutically acceptable salt thereof.

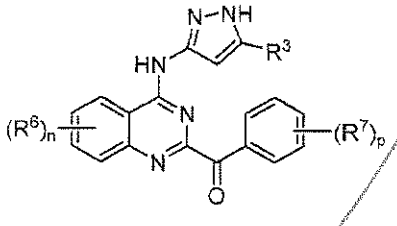
13. The compound of claim 1 having formula (VI)



(VI)

or a pharmaceutically acceptable salt thereof.

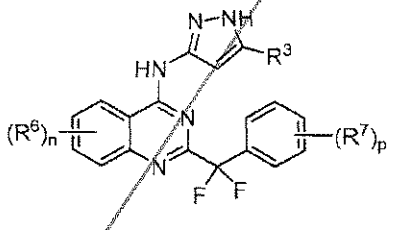
14. The compound of claim 1 having formula (VII)



(VII)

or a pharmaceutically acceptable salt thereof.

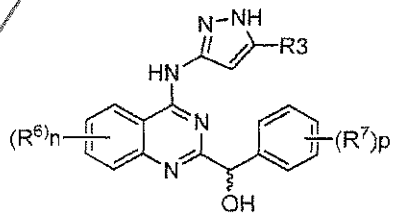
15. The compound of claim 1 having formula (VIII)



(VIII)

or a pharmaceutically acceptable salt thereof.

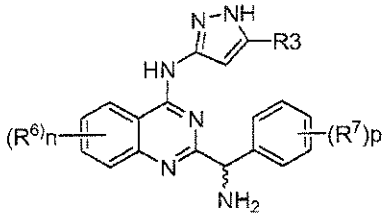
16. The compound of claim 1 having formula (IX)



(IX)

or a pharmaceutically acceptable salt thereof.

17. The compound of claim 1 having formula (X)



(X)

or a pharmaceutically acceptable salt thereof.

18. The compound of claim 1 selected from:

(4-chloroquinazolin-2-yl)(3-fluorophenyl)methanone;

(4-(1H-pyrazol-3-ylamino)quinazolin-2-yl)(3-fluorophenyl)methanone;

(4-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanone;

(4-(1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanone;

(4-(1H-pyrazol-3-ylamino)quinazolin-2-yl)(2-methoxyphenyl) methanone;

(4-(1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanol;

2-(fluoro(4-fluorophenyl)methyl)-N-(1H-pyrazol-3-yl)quinazolin-4-amine;

2-(difluoro(4-fluorophenyl)methyl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;

2-(difluoro(4-fluorophenyl)methyl)-N-(1H-pyrazol-3-yl)quinazolin-4-amine;

N-(5-cyclopropyl-1H-pyrazol-3-yl)-2-(difluoro(4-fluorophenyl)methyl) quinazolin-4-amine;

3-(2-(4-fluorobenzoyl)quinazolin-4-ylamino)-1H-pyrazole-5-carbonitrile;

(4-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;

2-((4-fluorophenyl)(methoxy)methyl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;

2-(amino(4-fluorophenyl)methyl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;

3-(2-((4-fluorophenyl)(hydroxy)methyl)quinazolin-4-ylamino)-1H-pyrazole-5-carbonitrile;

(5-fluoro-4-(5-methyl-1H-pyrazol-3-ylamino) quinazolin-2-yl) (4-fluorophenyl) methanol;

(4-fluorophenyl) (4-(5-methyl-1H-pyrazol-3-ylamino)-7-(trifluoromethyl) quinazolin-2-yl) methanone;

(4-fluorophenyl) (4-(5-methyl-1H-pyrazol-3-ylamino)-7-(trifluoromethyl) quinazolin-2-yl);

(7-fluoro-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanone;

2-(difluoro(4-fluorophenyl)methyl)-7-fluoro-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;

2-(difluoro(4-fluorophenyl)methyl)-7-fluoro-N-(1H-pyrazol-3-yl) quinazolin-4-amine;

(4-(1H-pyrazol-3-ylamino)-7-iodoquinazolin-2-yl)(4-fluorophenyl)methanone;

(4-(1H-pyrazol-3-ylamino)-7-iodoquinazolin-2-yl)(4-fluorophenyl)methanol;

(4-fluorophenyl)(7-methyl-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanone;

(4-fluorophenyl)(7-methyl-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;

2-(difluoro(4-fluorophenyl)methyl)-7-methyl-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;

2-(difluoro(4-fluorophenyl)methyl)-7-methyl-N-(1H-pyrazol-3-yl)quinazolin-4-amine;

(4-(1H-pyrazol-3-ylamino)-7-methoxyquinazolin-2-yl)(4-fluorophenyl)methanone;

(4-(1H-pyrazol-3-ylamino)-7-methoxyquinazolin-2-yl)(4-fluorophenyl)methanol;

(4-fluorophenyl)(7-methoxy-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanone;

(4-fluorophenyl)(7-methoxy-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;

2-(difluoro(4-fluorophenyl)methyl)-7-methoxy-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;

2-(difluoro(4-fluorophenyl)methyl)-7-methoxy-N-(1H-pyrazol-3-yl)quinazolin-4-amine;

2-(difluoro(4-fluorophenyl)methyl)-8-fluoro-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;

(4-(1H-pyrazol-3-ylamino)-8-methoxyquinazolin-2-yl)(4-fluorophenyl)methanone;

2-((4-fluorophenyl)(hydroxy)methyl)-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-7-ol;

(4-fluorophenyl)(7-hydroxy-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanone;

(4-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)-7-(2-morpholinoethoxy)quinazolin-2-yl)methanol;

2-(2-((4-fluorophenyl)(hydroxy)methyl)-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-7-yloxy)ethanol;

3-(2-((4-fluorophenyl)(hydroxy)methyl)-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-7-yloxy)propan-1-ol;

(4-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)-7-(piperidin-4-yloxy)quinazolin-2-yl)methanol;

(4-fluorophenyl)(7-(2-methoxyethoxy)-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;

tert-butyl 2-(2-((4-fluorophenyl)(hydroxy)methyl)-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-7-yloxy)acetate;

2-(2-((4-fluorophenyl)(hydroxy)methyl)-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-7-yloxy)acetic acid;

{(4-fluoro-phenyl)-[4-(5-methyl-1H-pyrazol-3-ylamino)-quinazolin-2-yl]-methyl}-carbamic acid methyl ester;

bis-(4-fluoro-phenyl)-[4-(5-methyl-1H-pyrazol-3-ylamino)-quinazolin-2-yl]-methanol;

methyl (4-fluorophenyl)(4-(5-methyl-4H-pyrazol-3-ylamino) quinazolin-2-yl) methylcarbamate;

(4-fluorophenyl)(8-methyl-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;

(7-fluoro-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanol;

(4-(1H-pyrazol-3-ylamino)quinazolin-2-yl)bis(4-fluorophenyl)methanol;

(2-(difluoro(4-fluorophenyl)methyl)-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-7-yl)methanol;

2-(difluoro(4-fluorophenyl)methyl)-N-(5-methyl-1H-pyrazol-3-yl)-7-(methylsulfonylmethyl)quinazolin-4-amine;
2-(Difluoro(4-fluorophenyl)methyl)-7-(ethoxymethyl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;
(7-chloro-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanol;
(6-fluoro-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanone-(6-fluoro-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanol;
(4-(1H-pyrazol-3-ylamino)-6-fluoroquinazolin-2-yl)(4-fluorophenyl)methanol;
(7-bromo-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanone;
(7-bromo-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanol;
(4-(1H-pyrazol-3-ylamino)-7-bromoquinazolin-2-yl)(4-fluorophenyl)methanol;
2-(2-(4-fluorophenyl)-1,3-dioxolan-2-yl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;
(8-fluoro-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanone;
(8-fluoro-4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanol;
(2-methoxyphenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanone;
(2-methoxyphenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;
(3-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;
N-((4-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methyl)formamide;
(3,4-difluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;
(3-chloro-4-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;
3-(4-fluorophenyl)-3-(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)propanenitrile;

2-((cyclopropylamino)(4-fluorophenyl)methyl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;
2-(1-(4-fluorophenyl)-2-(methylsulfonyl)ethyl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;
2-(3-amino-1-(4-fluorophenyl)propyl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;
(4-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol-1-*d*;
(4-fluorophenyl)(4-(5-methoxy-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanone;
(4-(5-ethyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanol;
(4-Fluorophenyl)(4-(5-methoxy-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;
(4-fluoro-3-methoxyphenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanone;
(4-fluoro-3-hydroxyphenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanone; and
(2-fluoro-5-(hydroxy(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methyl)phenol acetate.

19. A pharmaceutical composition comprising a compound of any of claims 1-18 and a pharmaceutically acceptable carrier, diluent or excipient.

20. A method for treatment of a JAK modulated disease comprising administering a therapeutically effective amount of a compound of any of claims 1-18.

21. A method for treatment of a JAK2 modulated disease comprising administering a therapeutically effective amount of a compound of any of claims 1-18.

22. The method of claim 21, wherein JAK2 is wild type or mutant JAK2.

23. The method of claim 20, wherein the disease is cancer, myeloproliferative disorder, inflammation or autoimmune disease.

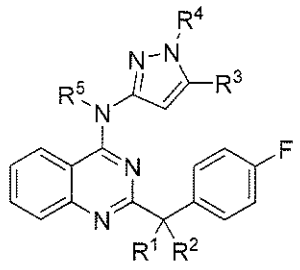
24. The method of claim 20 or 21, further comprising administering a second pharmaceutical agent selected from anti-proliferative agent, anti-inflammatory agent, immunomodulatory agent and immunosuppressive agent.

25. The compound of any of claims 1-18 for use in treatment of a JAK modulated disease.

26. Use of the compound of any of claims 1-18 for preparation of a medicament for treating a JAK modulated disease.

What is claimed is:

1. A compound having formula (Va)



(Va)

or a pharmaceutically acceptable salt thereof, where

R^1 and R^2 are selected from (i), (ii), and (iii) as follows:

(i) R^1 and R^2 together form =O;

(ii) R^1 is hydrogen or halo, and R^2 is halo; and

(iii) R^1 is hydroxyl, $C_{1\text{-}6}$alkoxy, amino, -NHCH(O) or $C_{1\text{-}6}$alkoxycarbonylamino, and R^2 is hydrogen or $C_{1\text{-}6}$alkyl;

R^3 is hydrogen or $C_{1\text{-}6}$alkyl;

R^4 is hydrogen; and

R^5 is hydrogen.

2. The compound of claim 1, wherein R^1 and R^2 together form =O.

3. The compound of claim 1 or 2, wherein R^1 is hydrogen or fluoro, and R^2 is fluoro.

4. The compound of claim 1 or 2, wherein R^1 is hydroxyl, methoxy, amino or methoxycarbonylamino, and R^2 is hydrogen.

5. The compound of claim 1 selected from:

(4-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanone;

(4-(1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanone;

(4-(1H-pyrazol-3-ylamino)quinazolin-2-yl)(2-methoxyphenyl) methanone;

(4-(1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanol;

2-(fluoro(4-fluorophenyl)methyl)-N-(1H-pyrazol-3-yl)quinazolin-4-amine;

2-(difluoro(4-fluorophenyl)methyl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;

2-(difluoro(4-fluorophenyl)methyl)-N-(1H-pyrazol-3-yl)quinazolin-4-amine;

(4-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methanol;

2-((4-fluorophenyl)(methoxy)methyl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;

2-(amino(4-fluorophenyl)methyl)-N-(5-methyl-1H-pyrazol-3-yl)quinazolin-4-amine;

3-(2-((4-fluorophenyl)(hydroxy)methyl)quinazolin-4-ylamino)-1H-pyrazole-5-carbonitrile;
{(4-fluoro-phenyl)-[4-(5-methyl-1H-pyrazol-3-ylamino)-quinazolin-2-yl]-methyl}-carbamic acid methyl ester;
N-((4-fluorophenyl)(4-(5-methyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)methyl)formamide; and
(4-(5-ethyl-1H-pyrazol-3-ylamino)quinazolin-2-yl)(4-fluorophenyl)methanol.

6. A pharmaceutical composition comprising a compound of any of claims 1-5 and a pharmaceutically acceptable carrier or excipient.

7. Use of a therapeutically effective amount of a compound of any of claims 1-5 in the manufacture of a medicament for treatment of a JAK modulated disease.

8. Use of a therapeutically effective amount of a compound of any of claims 1-5 in the manufacture of a medicament for treatment of a JAK2 modulated disease.

9. The use of claim 8, wherein JAK2 is wild type or mutant JAK2.

10. The use of claim 7, wherein the disease is cancer, myeloproliferative disorder, inflammation or autoimmune disease.

11. The use of claim 7 or 8, wherein the medicament further comprises a second pharmaceutical agent selected from anti-proliferative agent, anti-inflammatory agent, and immunomodulatory agent.

12. The compound of any of claims 1-5 for use in treatment of a JAK modulated disease.

13. Use of the compound of any of claims 1-5 for preparation of a medicament for treating a JAK modulated disease.