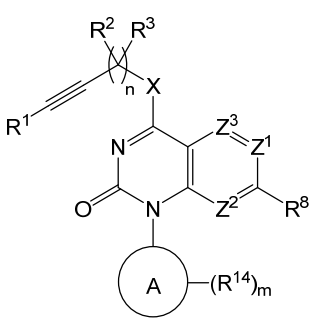
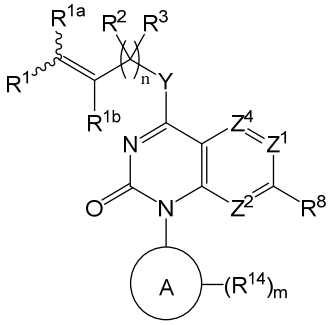
CLAIMS

WHAT IS CLAIMED IS:

1. A compound of Formula (I) or Formula (II), or a pharmaceutically acceptable salt thereof:



Formula (I);



Formula (II);

wherein:

A is selected from C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-10}heteroaryl;

Z^1 is CR^7 or N;

Z^2 is CR^9 or N;

Z^3 is CR^6 or N;

Z^4 is CR^{6a} or N;

X is selected from $-N(R^4)-$, -O-, and $-C(R^5)(R^{5a})-$;

Y is selected from $-N(R^{4a})-$, -O-, and $-C(R^5)(R^{5a})-$;

R^1 is selected from hydrogen, halogen, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, C_{1-9}heteroaryl, $-OR^{10}$, $-SR^{10}$, $-SF_5$, $-N(R^{10})(R^{11})$, $-C(O)OR^{10}$, $-OC(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)OR^{13}$, $-N(R^{12})S(O)_2R^{13}$, $-C(O)R^{13}$, $-S(O)R^{13}$, $-OC(O)R^{13}$, $-C(O)N(R^{10})(R^{11})$, $-C(O)C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)R^{13}$, $-S(O)_2R^{13}$, $-S(O)_2N(R^{10})(R^{11})-$, $-N=S(=O)(R^{13})_2$, $-S(=O)(=NH)N(R^{10})(R^{11})$, $-S(=O)(=NH)C(R^{10})(R^{11})$, $-S(=O)(=NR^{13})R^{13}$, $-CH_2C(O)N(R^{10})(R^{11})$, $-CH_2N(R^{12})C(O)R^{13}$, $-CH_2S(O)_2R^{13}$, $-CH_2S(O)_2N(R^{10})(R^{11})$, $-Si(C_{1-6}alkyl)_3$, and $-P(O)(R^{10})_2$, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from R^{15a};

R^{1a} and R^{1b} are independently selected from hydrogen, halogen, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl; wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from R^{15a};

each R^2 and each R^3 are each independently selected from hydrogen, halogen, oxo, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, C_{1-9}heteroaryl, $-OR^{10}$, $-SR^{10}$, $-SF_5$, $-N(R^{10})(R^{11})$, $-C(O)OR^{10}$, $-OC(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)OR^{13}$, $-N(R^{12})S(O)_2R^{13}$, $-C(O)R^{13}$, $-S(O)R^{13}$, $-OC(O)R^{13}$,

-C(O)N(R^{10})(R^{11}), -C(O)C(O)N(R^{10})(R^{11}), -N(R^{12})C(O)R^{13}, -S(O)$_2$$R^{13}$, -S(O)$_2$N($R^{10}$)($R^{11}$)-, -N=S(=O)($R^{13}$)$_2$, -S(=O)(=NH)N($R^{10}$)($R^{11}$), -S(=O)(=NH)C($R^{10}$)($R^{11}$), -S(=O)(=N$R^{13}$)$R^{13}$, -CH$_2$C(O)N($R^{10}$)($R^{11}$), -CH$_2$N($R^{12}$)C(O)$R^{13}$, -CH$_2$S(O)$_2$$R^{13}$, -CH$_2$S(O)$_2$N($R^{10}$)($R^{11}$), -Si($C_{1-6}$alkyl)$_3$, and -P(O)($R^{10}$)$_2$, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from R^{15b}; or R^2 and R^3, together with the carbon to which they are attached, form a C_{3-6}cycloalkyl or C_{2-9}heterocycloalkyl;

R^4 is selected from hydrogen, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from halogen, -CN, hydroxy, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{1-6}alkoxy, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl; or R^4 and an R^3 are combined to form a C_{2-9}heterocycloalkyl optionally substituted with one, two, or three groups selected from halogen, -CN, hydroxy, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{1-6}alkoxy, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl;

R^{4a} is selected from hydrogen, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from halogen, -CN, hydroxy, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{1-6}alkoxy, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl; or R^{4a} and an R^3 are combined to form a C_{2-9}heterocycloalkyl optionally substituted with one, two, or three groups selected from halogen, -CN, hydroxy, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{1-6}alkoxy, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl;

R^5 and R^{5a} are independently selected from hydrogen, halogen, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, C_{1-9}heteroaryl, -OR^{10}, -SR^{10}, -SF_5, -N(R^{10})(R^{11}), -C(O)OR^{10}, -OC(O)N(R^{10})(R^{11}), -N(R^{12})C(O)N(R^{10})(R^{11}), -N(R^{12})C(O)OR^{13}, -N(R^{12})S(O)$_2$$R^{13}$, -C(O)$R^{13}$, -S(O)$R^{13}$, -OC(O)$R^{13}$, -C(O)N($R^{10}$)($R^{11}$), -C(O)C(O)N($R^{10}$)($R^{11}$), -N($R^{12}$)C(O)$R^{13}$, -S(O)$_2$$R^{13}$, -S(O)$_2$N($R^{10}$)($R^{11}$)-, -N=S(=O)($R^{13}$)$_2$, -S(=O)(=NH)N($R^{10}$)($R^{11}$), -S(=O)(=NH)C($R^{10}$)($R^{11}$), -S(=O)(=N$R^{13}$)$R^{13}$, -CH$_2$C(O)N($R^{10}$)($R^{11}$), -CH$_2$N($R^{12}$)C(O)$R^{13}$, -CH$_2$S(O)$_2$$R^{13}$, -CH$_2$S(O)$_2$N($R^{10}$)($R^{11}$), -Si($C_{1-6}$alkyl)$_3$, and -P(O)($R^{10}$)$_2$, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from halogen, -CN, hydroxy, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{1-6}alkoxy, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl; or R^5 and an R^3 are combined to form a C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, or C_{2-9}heteroaryl, wherein C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with

one, two, or three groups selected from halogen, -CN, hydroxy, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{1-6}alkoxy, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl;

R^6, R^7, R^8, and R^9 are independently selected from hydrogen, halogen, -CN, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, C_{1-9}heteroaryl, $-OR^{10}$, $-SR^{10}$, $-SF_5$, $-N(R^{10})(R^{11})$, $-C(O)OR^{10}$, $-OC(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)OR^{13}$, $-N(R^{12})S(O)_2R^{13}$, $-C(O)R^{13}$, $-S(O)R^{13}$, $-OC(O)R^{13}$, $-C(O)N(R^{10})(R^{11})$, $-C(O)C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)R^{13}$, $-S(O)_2R^{13}$, $-S(O)_2N(R^{10})(R^{11})$-, $-N=S(=O)(R^{13})_2$, $-S(=O)(=NH)N(R^{10})(R^{11})$, $-S(=O)(=NH)C(R^{10})(R^{11})$, $-S(=O)(=NR^{13})R^{13}$, $-CH_2C(O)N(R^{10})(R^{11})$, $-CH_2N(R^{12})C(O)R^{13}$, $-CH_2S(O)_2R^{13}$, $-CH_2S(O)_2N(R^{10})(R^{11})$, $-Si(C_{1-6}alkyl)_3$, and $-P(O)(R^{10})_2$, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from R^{15c};

R^{6a} is selected from hydrogen, halogen, -CN, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{1-9}heteroaryl, $-OR^{10}$, $-SR^{10}$, $-SF_5$, $-N(R^{10})(R^{11})$, $-C(O)OR^{10}$, $-OC(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)OR^{13}$, $-N(R^{12})S(O)_2R^{13}$, $-C(O)R^{13}$, $-S(O)R^{13}$, $-OC(O)R^{13}$, $-C(O)N(R^{10})(R^{11})$, $-C(O)C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)R^{13}$, $-S(O)_2R^{13}$, $-S(O)_2N(R^{10})(R^{11})$-, $-N=S(=O)(R^{13})_2$, $-S(=O)(=NH)N(R^{10})(R^{11})$, $-S(=O)(=NH)C(R^{10})(R^{11})$, $-S(=O)(=NR^{13})R^{13}$, $-CH_2C(O)N(R^{10})(R^{11})$, $-CH_2N(R^{12})C(O)R^{13}$, $-CH_2S(O)_2R^{13}$, $-CH_2S(O)_2N(R^{10})(R^{11})$, $-Si(C_{1-6}alkyl)_3$, and $-P(O)(R^{10})_2$, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from R^{15c};

each R^{10} is independently selected from hydrogen, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from halogen, -CN, hydroxy, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{1-6}alkoxy, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl;

each R^{11} is independently selected from hydrogen, C_{1-6}alkyl, and C_{1-6}haloalkyl; or R^{10} and R^{11}, together with the nitrogen to which they are attached, form a C_{2-9}heterocycloalkyl;

each R^{12} is independently selected from hydrogen, C_{1-6}alkyl, and C_{1-6}haloalkyl;

each R^{13} is independently selected C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from halogen, -CN, hydroxy, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{1-6}alkoxy, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl;

each R^{14} is independently selected from halogen, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, C_{1-9}heteroaryl, $-OR^{10}$, $-SR^{10}$, $-SF_5$, -

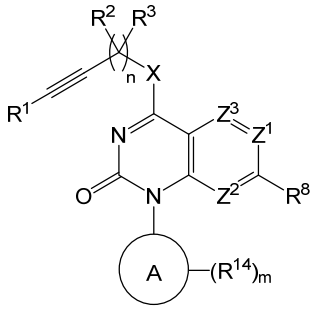
$N(R^{10})(R^{11})$, $-C(O)OR^{10}$, $-OC(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)OR^{13}$, $-N(R^{12})S(O)_2R^{13}$, $-C(O)R^{13}$, $-S(O)R^{13}$, $-OC(O)R^{13}$, $-C(O)N(R^{10})(R^{11})$, $-C(O)C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)R^{13}$, $-S(O)_2R^{13}$, $-S(O)_2N(R^{10})(R^{11})$-, $-N=S(=O)(R^{13})_2$, $-S(=O)(=NH)N(R^{10})(R^{11})$, $-S(=O)(=NH)C(R^{10})(R^{11})$, $-S(=O)(=NR^{13})R^{13}$, $-CH_2C(O)N(R^{10})(R^{11})$, $-CH_2N(R^{12})C(O)R^{13}$, $-CH_2S(O)_2R^{13}$, $-CH_2S(O)_2N(R^{10})(R^{11})$, $-Si(C_{1-6}alkyl)_3$, and $-P(O)(R^{10})_2$, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, C_{2-9}heterocycloalkyl, C_{6-10}aryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups selected from R^{15d};

each R^{15a}, R^{15b}, R^{15c}, and R^{15d} are each independently selected from halogen, oxo, -CN, C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-10}cycloalkyl, $-CH_2-C_{3-6}$cycloalkyl, C_{2-9}heterocycloalkyl, $-CH_2-C_{2-9}$heterocycloalkyl, C_{6-10}aryl, $-CH_2-C_{6-10}$aryl, C_{1-9}heteroaryl, $-CH_2-C_{1-9}$heteroaryl, $-OR^{10}$, $-SR^{10}$, $-SF_5$, $-N(R^{10})(R^{11})$, $-C(O)OR^{10}$, $-OC(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)OR^{13}$, $-N(R^{12})S(O)_2R^{13}$, $-C(O)R^{13}$, $-S(O)R^{13}$, $-OC(O)R^{13}$, $-C(O)N(R^{10})(R^{11})$, $-C(O)C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)R^{13}$, $-S(O)_2R^{13}$, $-S(O)_2N(R^{10})(R^{11})$-, $-N=S(=O)(R^{13})_2$, $-S(=O)(=NH)N(R^{10})(R^{11})$, $-S(=O)(=NH)C(R^{10})(R^{11})$, $-S(=O)(=NR^{13})R^{13}$, $-CH_2C(O)N(R^{10})(R^{11})$, $-CH_2N(R^{12})C(O)R^{13}$, $-CH_2S(O)_2R^{13}$, $-CH_2S(O)_2N(R^{10})(R^{11})$, $-Si(C_{1-6}alkyl)_3$, and $-P(O)(R^{10})_2$, wherein C_{1-6}alkyl, C_{2-6}alkenyl, C_{2-6}alkynyl, C_{3-6}cycloalkyl, $-CH_2-C_{3-10}$cycloalkyl, C_{2-9}heterocycloalkyl, $-CH_2-C_{2-9}$heterocycloalkyl, C_{6-10}aryl, $-CH_2-C_{6-10}$aryl, $-CH_2-C_{1-9}$heteroaryl, and C_{1-9}heteroaryl are optionally substituted with one, two, or three groups independently selected from halogen, oxo, -CN, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{1-6}alkoxy, C_{1-6}haloalkoxy, $-OR^{10}$, $-SR^{10}$, $-SF_5$, $-N(R^{10})(R^{11})$, $-C(O)OR^{10}$, $-OC(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)OR^{13}$, $-N(R^{12})S(O)_2R^{13}$, $-C(O)R^{13}$, $-S(O)R^{13}$, $-OC(O)R^{13}$, $-C(O)N(R^{10})(R^{11})$, $-C(O)C(O)N(R^{10})(R^{11})$, $-N(R^{12})C(O)R^{13}$, $-S(O)_2R^{13}$, $-S(O)_2N(R^{10})(R^{11})$-, $-N=S(=O)(R^{13})_2$, $-S(=O)(=NH)N(R^{10})(R^{11})$, $-S(=O)(=NH)C(R^{10})(R^{11})$, $-S(=O)(=NR^{13})R^{13}$, $-CH_2C(O)N(R^{10})(R^{11})$, $-CH_2N(R^{12})C(O)R^{13}$, $-CH_2S(O)_2R^{13}$, $-CH_2S(O)_2N(R^{10})(R^{11})$ and $-P(O)(R^{10})_2$;

m is 0, 1, 2, 3, 4, or 5; and

n is 0, 1, 2, 3, 4, 5, or 6.

2. The compound of claim 1, or a pharmaceutically acceptable salt thereof, having the structure of Formula (I):

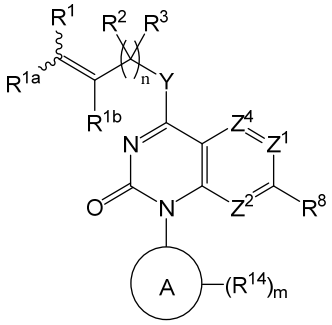


Formula (I).

3. The compound of claim 1 or claim 2, or a pharmaceutically acceptable salt thereof, wherein Z^3 is CR^6; R^6 is selected from hydrogen and $-OR^{10}$; and R^{10} is C_{1-6}alkyl.

4. The compound of any one of claims 1-3, or a pharmaceutically acceptable salt thereof, wherein X is $-N(R^4)-$ and R^4 is hydrogen or C_{1-6}alkyl.

5. The compound of claim 1, or a pharmaceutically acceptable salt thereof, having the structure of Formula (II):



Formula (II).

6. The compound of claim 1 or claim 5, or a pharmaceutically acceptable salt thereof, wherein Z^4 is CR^{6a}; R^{6a} is selected from hydrogen and $-OR^{10}$; and R^{10} is C_{1-6}alkyl.

7. The compound of claim 5 or 6, or a pharmaceutically acceptable salt thereof, wherein R^{1a} and R^{1b} are independently selected from hydrogen and unsubstituted C_{1-6}alkyl.

8. The compound of any one of claims 5-7, or a pharmaceutically acceptable salt thereof, wherein Y is $-N(R^{4a})-$ and R^{4a} is hydrogen or C_{1-6}alkyl.

9. The compound of any one of claims 1-8, or a pharmaceutically acceptable salt thereof, wherein each R^2 is independently selected from hydrogen and C_{1-6}alkyl and each R^3 is independently selected from hydrogen and C_{1-6}alkyl.

10. The compound of any one of claims 1-9, or a pharmaceutically acceptable salt thereof, wherein R^2 and R^3, together with the carbon to which they are attached, form a C_{3-6}cycloalkyl.

11. The compound of any one of claims 1-10, or a pharmaceutically acceptable salt thereof, wherein n is 1.

12. The compound of any one of claims 1-11, or a pharmaceutically acceptable salt thereof, wherein Z^1 is CR^7; R^7 is hydrogen, halogen, C_{1-6}alkyl, or C_{1-6}haloalkyl; Z^2 is CR^9; and R^9 is hydrogen, halogen, C_{1-6}alkyl, or C_{1-6}haloalkyl.

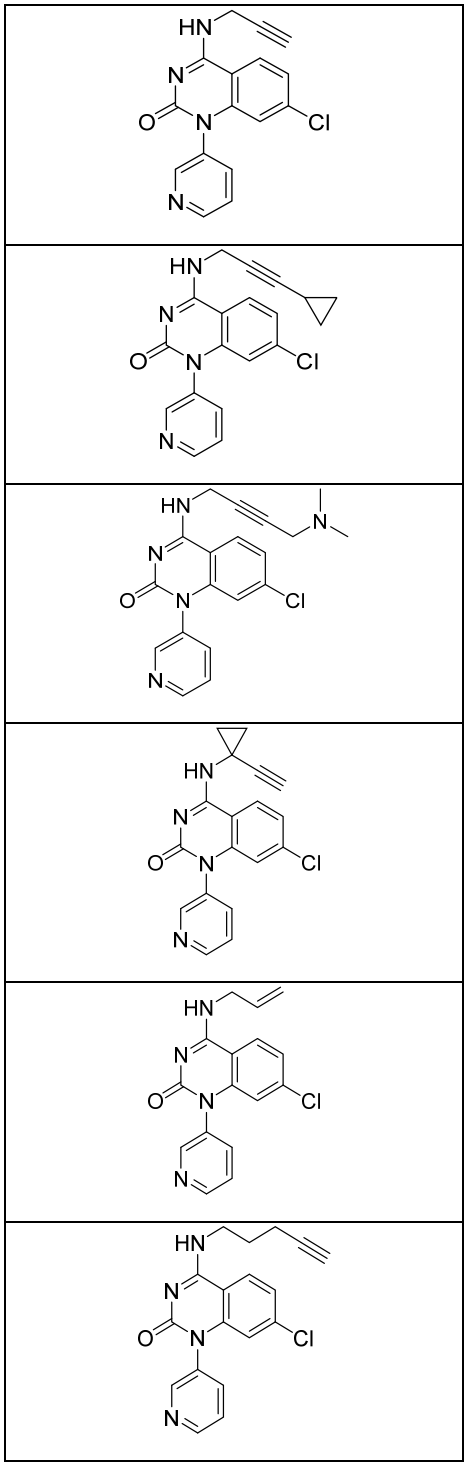
13. The compound of any one of claims 1-12, or a pharmaceutically acceptable salt thereof, wherein  is pyridyl.

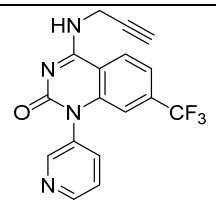
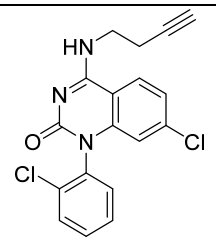
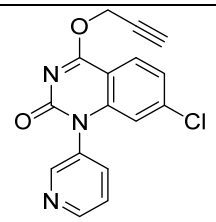
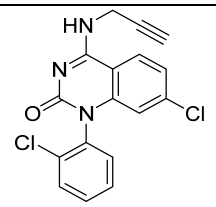
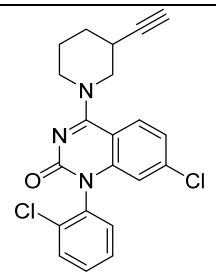
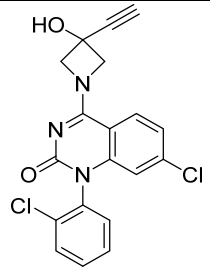
14. The compound of any one of claims 1-13, or a pharmaceutically acceptable salt thereof, wherein m is 0 or 1 and each R^{14} is independently selected from halogen and C_{1-6}alkyl.

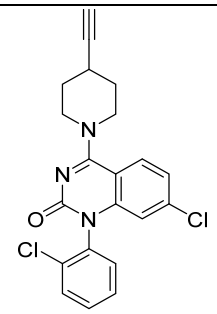
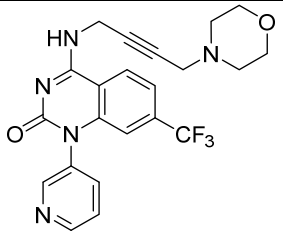
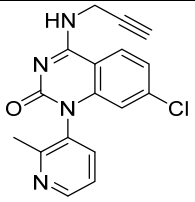
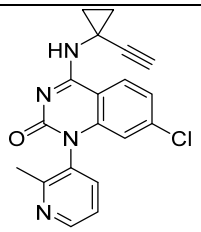
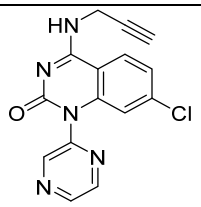
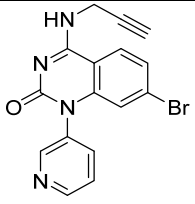
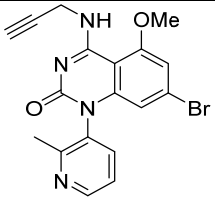
15. The compound of any one of claims 1-14, or a pharmaceutically acceptable salt thereof, wherein R^8 is hydrogen, halogen or C_{1-6}haloalkyl.

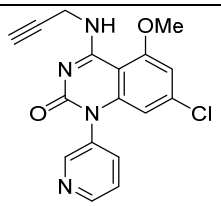
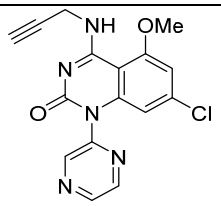
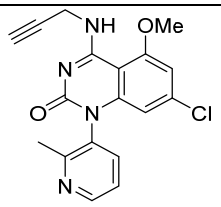
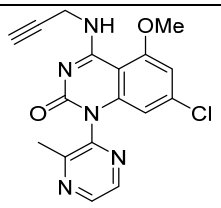
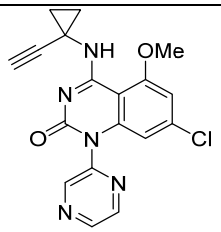
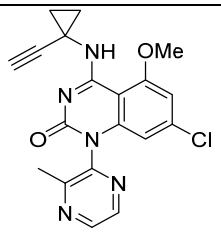
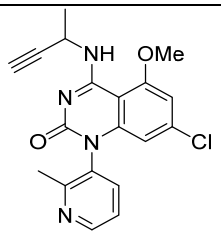
16. The compound of any one of claims 1-15, or a pharmaceutically acceptable salt thereof, wherein R^1 is selected from hydrogen and unsubstituted C_{3-6}cycloalkyl.

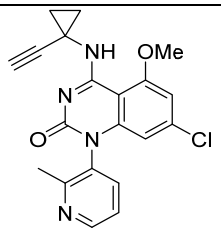
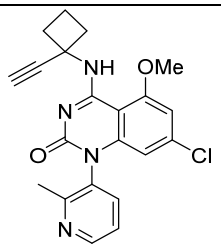
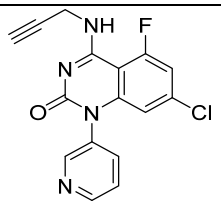
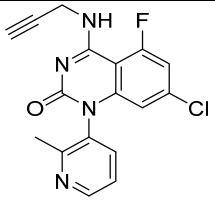
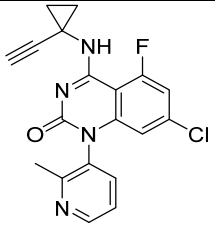
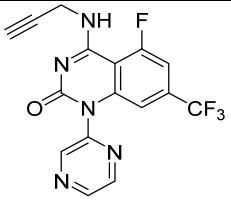
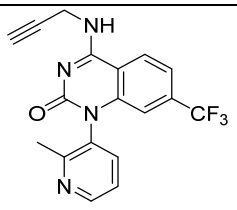
17. The compound of claim 1, or a pharmaceutically acceptable salt thereof, wherein the compound is a compound selected from the following compounds:

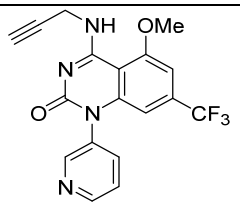
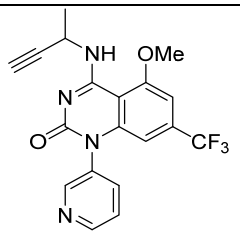
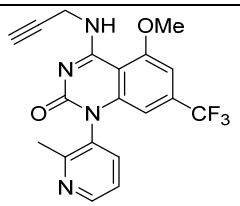
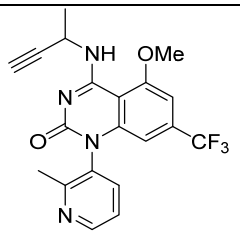
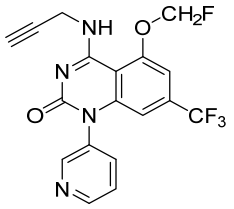
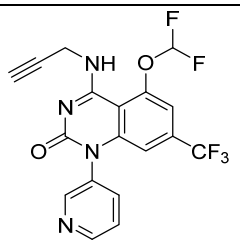
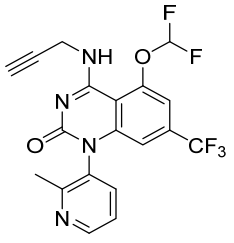


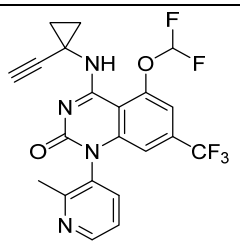
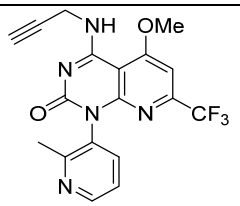
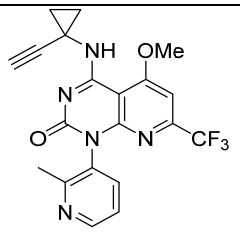
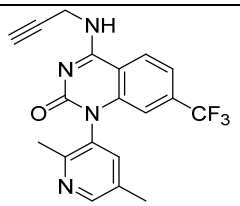
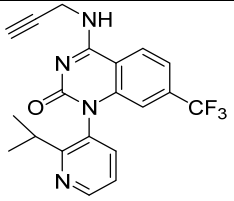
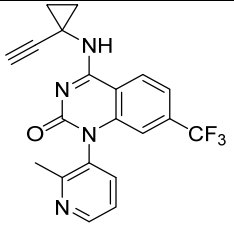
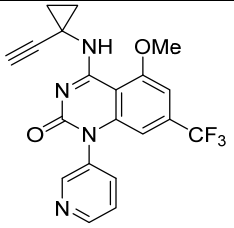







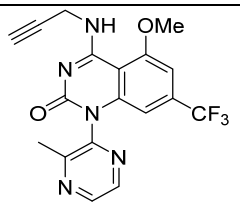
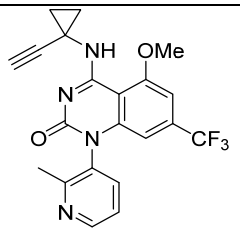
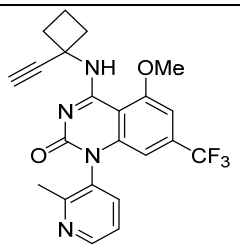
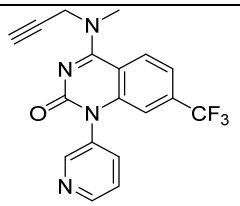
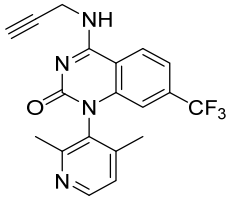
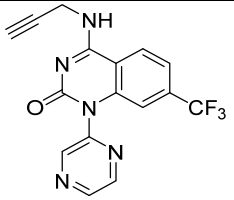
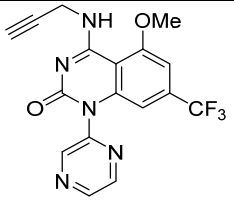








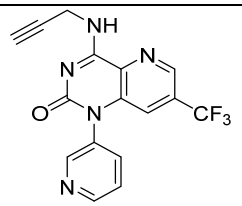
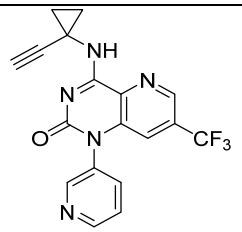
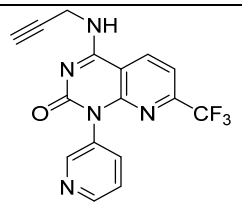
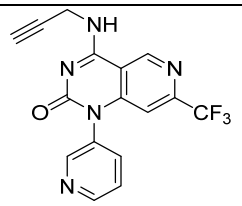
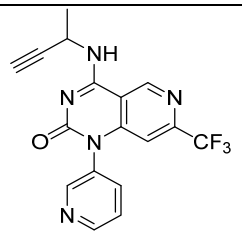
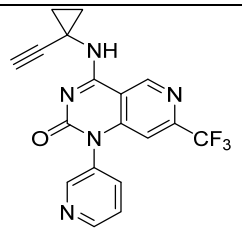
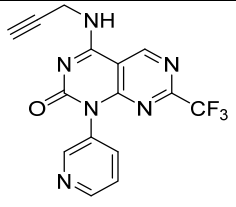








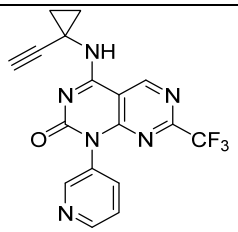
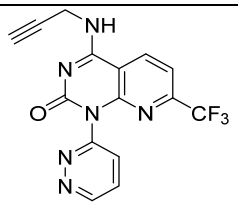
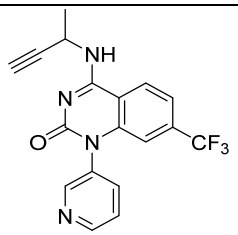
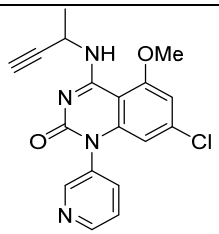
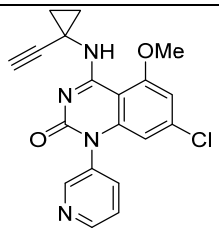
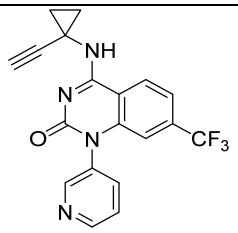
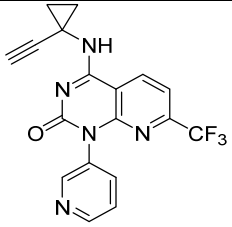








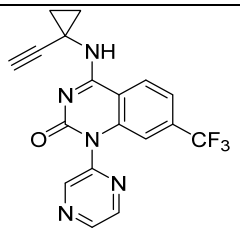
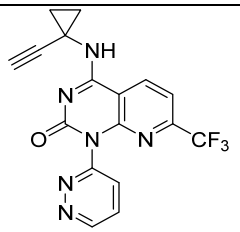
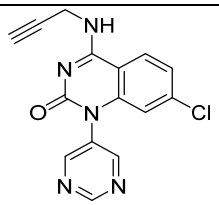
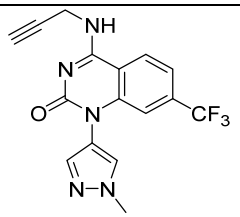
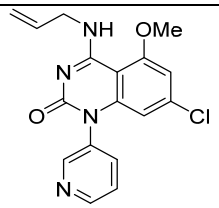
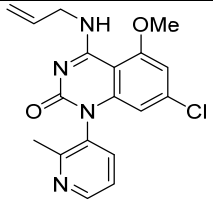
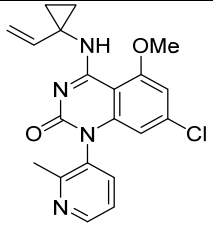








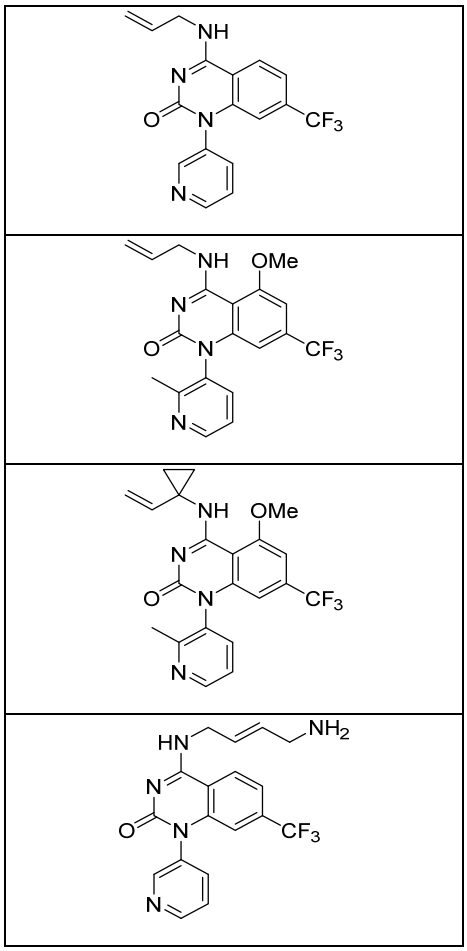








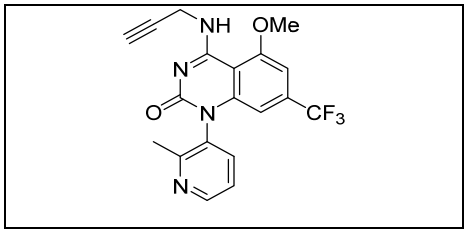










or a pharmaceutically acceptable salt thereof.

18. The compound of claim 1, or a pharmaceutically acceptable salt thereof, wherein the compound is:



or a pharmaceutically acceptable salt thereof.

19. A pharmaceutical composition comprising a compound of any one of claims 1-18, or a pharmaceutically acceptable salt thereof, and at least one pharmaceutically acceptable excipient.

20. Use of a compound of any one of claims 1-18, or a pharmaceutically acceptable salt thereof, in the manufacture of a medicament for treating cancer in a mammal.

21. A compound of any one of claims 1-18, or a pharmaceutically acceptable salt thereof, for use as a medicament.