



- (51) **International Patent Classification:**
A61K 31/4162 (2006.01)
- (21) **International Application Number:**
PCT/US2015/048680
- (22) **International Filing Date:**
4 September 2015 (04.09.2015)
- (25) **Filing Language:** English
- (26) **Publication Language:** English
- (30) **Priority Data:**
62/047,371 8 September 2014 (08.09.2014) US
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- (81) **Designated States** (unless otherwise indicated, for every kind of national protection available): AE, AG, AL, AM, AO, AT, AU, AZ, BA, BB, BG, BH, BN, BR, BW, BY,

BZ, CA, CH, CL, CN, CO, CR, CU, CZ, DE, DK, DM, DO, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, GT, HN, HR, HU, ID, IL, IN, IR, IS, JP, KE, KG, KN, KP, KR, KZ, LA, LC, LK, LR, LS, LU, LY, MA, MD, ME, MG, MK, MN, MW, MX, MY, MZ, NA, NG, NI, NO, NZ, OM, PA, PE, PG, PH, PL, PT, QA, RO, RS, RU, RW, SA, SC, SD, SE, SG, SK, SL, SM, ST, SV, SY, TH, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, ZA, ZM, ZW.

- (84) **Designated States** (unless otherwise indicated, for every kind of regional protection available): ARIPO (BW, GH, GM, KE, LR, LS, MW, MZ, NA, RW, SD, SL, ST, SZ, TZ, UG, ZM, ZW), Eurasian (AM, AZ, BY, KG, KZ, RU, TJ, TM), European (AL, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HR, HU, IE, IS, IT, LT, LU, LV, MC, MK, MT, NL, NO, PL, PT, RO, RS, SE, SI, SK, SM, TR), OAPI (BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, KM, ML, MR, NE, SN, TD, TG).

Declarations under Rule 4.17:

- as to applicant's entitlement to apply for and be granted a patent (Rule 4.17(ii))
- as to the applicant's entitlement to claim the priority of the earlier application (Rule 4.17(iii))

Published:

- with international search report (Art. 21(3))
- before the expiration of the time limit for amending the claims and to be republished in the event of receipt of amendments (Rule 48.2(h))



WO 2016/040184 A1

(54) **Title:** 3-(3H-IMIDAZO[4,5-B]PYRIDIN-2-YL)-1H-PYRAZOLO[3,4-C]PYRIDINE AND THERAPEUTIC USES THEREOF

(57) **Abstract:** Azaindazole compounds for treating various diseases and pathologies are disclosed. More particularly, the present invention concerns the use of an azaindazole compound or analogs thereof, in the treatment of disorders characterized by the activation of Wnt pathway signaling (e.g., cancer, abnormal cellular proliferation, angiogenesis, fibrotic disorders, bone or cartilage diseases, and osteoarthritis), the modulation of cellular events mediated by Wnt pathway signaling, as well as genetic diseases and neurological conditions/disorders/diseases due to mutations or dysregulation of the Wnt pathway and/or of one or more of Wnt signaling components. Also provided are methods for treating Wnt-related disease states.

3-(3H-IMIDAZO[4,5-B]PYRIDIN-2-YL)-1H-PYRAZOLO[3,4-C]PYRIDINE AND THERAPEUTIC USES THEREOF

RELATED APPLICATIONS

[001] This application claims the benefit of U.S. Provisional Application No. 62/047,371, filed September 8, 2014, which is incorporated herein by reference in its entirety.

BACKGROUND

Technical Field

[002] This disclosure relates to inhibitors of one or more proteins in the Wnt pathway, including inhibitors of one or more Wnt proteins, and compositions comprising the same. More particularly, it concerns the use of an azaindazole compound or salts or analogs thereof, in the treatment of disorders characterized by the activation of Wnt pathway signaling (e.g., cancer, abnormal cellular proliferation, angiogenesis, fibrotic disorders, bone or cartilage diseases, and osteoarthritis), the modulation of cellular events mediated by Wnt pathway signaling, as well as genetic diseases and neurological conditions/disorders/diseases due to mutations or dysregulation of the Wnt pathway and/or of one or more of Wnt signaling components. Also provided are methods for treating Wnt-related disease states.

Background

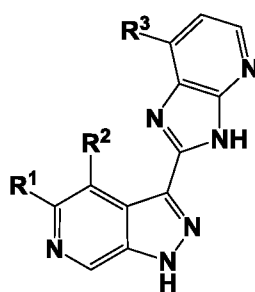
[003] The Wnt growth factor family includes more than 10 genes identified in the mouse and at least 19 genes identified in the human. Members of the Wnt family of signaling molecules mediate many short-and long-range patterning processes during invertebrate and vertebrate development. The Wnt signaling pathway is known for its role in the inductive interactions that regulate growth and differentiation, and it also plays roles in the homeostatic maintenance of post-embryonic tissue integrity. Wnt stabilizes cytoplasmic β -catenin, which stimulates the expression of genes including c-myc, c jun, fra-1, and cyclin D1. In addition, misregulation of Wnt signaling can cause developmental defects and is implicated in the genesis of several human cancers. The Wnt pathway has also been implicated in the maintenance of stem or progenitor cells in a growing list of adult tissues including skin, blood, gut, prostate, muscle, and the nervous system.

SUMMARY

[004] The present disclosure provides methods and reagents, involving contacting a cell with an agent, such as an azaindazole compound, in a sufficient amount to antagonize a Wnt activity, e.g., to reverse or control an aberrant growth state or correct a genetic disorder due to mutations in Wnt signaling components.

[005] Some embodiments disclosed herein include Wnt inhibitors containing an azaindazole core. Other embodiments disclosed herein include pharmaceutical compositions and methods of treatment using these compounds.

[006] One embodiment disclosed herein includes a compound having the structure of Formula I:



I

as well as prodrugs and pharmaceutically acceptable salts thereof.

[007] In some embodiments of Formula (I):

R^1 is selected from the group consisting of $-\text{heteroaryl}(R^4)_q$ and $-\text{heterocyclyl}(R^5)_h$;

R^2 is selected from the group consisting of H and halide;

R^3 is selected from the group consisting of H, $-\text{heteroaryl}(R^6)_q$, $-\text{heterocyclyl}(R^7)_h$, and $-\text{aryl}(R^8)_k$;

each R^4 is one substituent attached to the heteroaryl and is independently selected from the group consisting of halide, $-(C_{1-6} \text{ alkyl})$, $-(C_{1-4} \text{ alkylene})_p$ heterocyclyl(R^9) $_h$, $-(C_{1-4} \text{ alkylene})_p$ carbocyclyl(R^{10}) $_j$, $-(C_{1-4} \text{ alkylene})_p$ aryl(R^{11}) $_k$, $-\text{NHC}(=\text{O})R^{12}$, $-\text{NR}^{13}R^{14}$, $-(C_{1-6} \text{ alkylene})NR^{15}R^{16}$, and $-\text{OR}^{22}$;

each R^5 is one substituent attached to the heterocyclyl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-\text{CF}_3$, and $-\text{CN}$;

each R^6 is one substituent attached to the heteroaryl and is independently selected from the group consisting of $-(C_{1-6} \text{ alkyl})$, halide, $-\text{CF}_3$, $-\text{OCH}_3$, $-\text{CN}$, and $-\text{C}(=\text{O})R^{17}$;

each R^7 is one substituent attached to the heterocyclyl and is independently selected from the group consisting of $-(C_{1-6} \text{ alkyl})$, halide, $-\text{CF}_3$, $-\text{CN}$, and $-\text{OCH}_3$;

each R⁸ is one substituent attached to the aryl and is independently selected from the group consisting of $-(C_{1-6} \text{ alkyl})$, halide, $-CF_3$, $-CN$, $-OCH_3$, $-(C_{1-6} \text{ alkylene})_pNHSO_2R^{17}$, $-NR^{13}(C_{1-6} \text{ alkylene})NR^{13}R^{14}$, $-(C_{1-6} \text{ alkylene})_pNR^{13}R^{14}$, and $-OR^{25}$;

each R⁹ is one substituent attached to the heterocyclyl and is independently selected from the group consisting of amino, $-(C_{1-4} \text{ alkyl})$, halide, $-CF_3$, and $-CN$;

each R¹⁰ is one substituent attached to the carbocyclyl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-CF_3$, and $-CN$;

each R¹¹ is one substituent attached to the aryl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-CF_3$, and $-CN$;

each R¹² is independently selected from the group consisting of $-(C_{1-9} \text{ alkyl})$, $-$ heteroaryl(R¹⁸)_q, $-$ aryl(R¹⁹)_k, $-CH_2$ aryl(R¹⁹)_k, $-$ carbocyclyl(R²⁰)_j, $-CH_2$ carbocyclyl(R²⁰)_j, $-(C_{1-4} \text{ alkylene})_pNR^{23}R^{24}$, $-$ heterocyclyl(R²¹)_h, and $-CH_2$ heterocyclyl(R²¹)_h;

each R¹³ is independently selected from the group consisting of H and $-(C_{1-6} \text{ alkyl})$;

each R¹⁴ is independently selected from the group consisting of H, $-(C_{1-6} \text{ alkyl})$, $-CH_2$ aryl(R¹⁹)_k, and $-CH_2$ carbocyclyl(R²⁰)_j;

each R¹⁵ is independently selected from the group consisting of H and $-(C_{1-6} \text{ alkyl})$;

each R¹⁶ is independently selected from the group consisting of H, $-(C_{1-6} \text{ alkyl})$, $-CH_2$ aryl(R¹⁹)_k, and $-CH_2$ carbocyclyl(R²⁰)_j;

each R¹⁷ is a $-(C_{1-6} \text{ alkyl})$;

each R¹⁸ is one substituent attached to the heteroaryl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-CF_3$, and $-CN$;

each R¹⁹ is one substituent attached to the aryl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-CF_3$, and $-CN$;

each R²⁰ is one substituent attached to the carbocyclyl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-CF_3$, and $-CN$;

each R²¹ is one substituent attached to the heterocyclyl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-CF_3$, and $-CN$;

each R²² is independently selected from the group consisting of H, $-(C_{1-6} \text{ alkyl})$, $-(C_{1-4} \text{ alkylene})_p$ heterocyclyl(R²¹)_h, $-(C_{1-4} \text{ alkylene})_p$ carbocyclyl(R²⁰)_j, $-(C_{1-4} \text{ alkylene})_p$ aryl(R¹⁹)_k, and $-(C_{1-6} \text{ alkylene})_pNR^{23}R^{24}$;

each R²³ is independently selected from the group consisting of H and $-(C_{1-6} \text{ alkyl})$;

each R²⁴ is independently selected from the group consisting of H and $-(C_{1-6} \text{ alkyl})$;

each R²⁵ is independently selected from the group consisting of H, $-(C_{1-6} \text{ alkyl})$, $-(C_{1-4} \text{ alkylene})_p$ heterocyclyl(R²¹)_h, and $-(C_{1-6} \text{ alkylene})_pNR^{23}R^{24}$;

each p is independently 0 or 1;
each q is independently 0 to 4;
each h is independently 0 to 10;
each k is independently 0 to 5; and
each j is independently 0 to 12.

[008] Some embodiments include stereoisomers and pharmaceutically acceptable salts of a compound of Formula (I).

[009] Some embodiments include pro-drugs of a compound of Formula (I).

[010] Some embodiments of the present disclosure include pharmaceutical compositions comprising a compound of Formula (I) and a pharmaceutically acceptable carrier, diluent, or excipient.

[011] Other embodiments disclosed herein include methods of inhibiting one or more members of the Wnt pathway, including one or more Wnt proteins by administering to a patient affected by a disorder or disease in which aberrant Wnt signaling is implicated, such as cancer and other diseases associated with abnormal angiogenesis, cellular proliferation, cell cycling and mutations in Wnt signaling components, a compound according to Formula (I). Accordingly, the compounds and compositions provided herein can be used to treat cancer, to reduce or inhibit angiogenesis, to reduce or inhibit cellular proliferation and correct a genetic disorder due to mutations in Wnt signaling components.

[012] Non-limiting examples of diseases which can be treated with the compounds and compositions provided herein include a variety of cancers, diabetic retinopathy, pulmonary fibrosis, rheumatoid arthritis, sepsis, ankylosing spondylitis, psoriasis, scleroderma, mycotic and viral infections, osteochondrodysplasia, Alzheimer's disease, lung disease, bone/osteoporotic (wrist, spine, shoulder and hip) fractures, articular cartilage (chondral) defects, degenerative disc disease (or intervertebral disc degeneration), polyposis coli, osteoporosis-pseudoglioma syndrome, familial exudative vitreoretinopathy, retinal angiogenesis, early coronary disease, tetra-Amelia syndrome, Müllerian-duct regression and virilization, SERKAL syndrome, diabetes mellitus type 2, Fuhrmann syndrome, Al-Awadi/Raas-Rothschild/Schinzel phocomelia syndrome, odontonycho-dermal dysplasia, obesity, split-hand/foot malformation, caudal duplication syndrome, tooth agenesis, Wilms tumor, skeletal dysplasia, focal dermal hypoplasia, autosomal recessive anonychia, neural tube defects, alpha-thalassemia (ATRX) syndrome, fragile X syndrome, ICF syndrome, Angelman syndrome, Prader-Willi syndrome, Beckwith-Wiedemann Syndrome, Norrie disease, and Rett syndrome.

[013] Some embodiments of the present disclosure include methods to prepare compounds of Formula (I).

[014] It is to be understood that both the foregoing general description and the following detailed description are exemplary and explanatory only and are not restrictive of the invention, as claimed.

DETAILED DESCRIPTION

[015] Provided herein are compositions and methods for inhibiting one or more members of the Wnt pathway, including one or more Wnt proteins. Other Wnt inhibitors and methods for using the same are disclosed in U.S. Application Ser. Nos. 12/852,706; 12/968,505; 13/552,188; 13/800,963; 13/855,874; 13/887,177 13/938,691; 13/938,692; 14/019,103; 14/019,147; 14/019,940; 14/149,948; 14/178,749; 14/331,427; and 14/334,005; and U.S. Provisional Application Ser. Nos. 61/232,603; 61/288,544; 61/305,459; 61/620,107; 61/642,915; and 61/750,221, all of which are incorporated by reference in their entirety herein.

[016] Some embodiments provided herein relate to a method for treating a disease or disorder including, but not limited to, cancers, diabetic retinopathy, pulmonary fibrosis, rheumatoid arthritis, sepsis, ankylosing spondylitis, psoriasis, scleroderma, mycotic and viral infections, bone and cartilage diseases, Alzheimer's disease, lung disease, osteoarthritis, bone/osteoporotic (wrist, spine, shoulder and hip) fractures, articular cartilage (chondral) defects, degenerative disc disease (or intervertebral disc degeneration), polyposis coli, bone density and vascular defects in the eye (Osteoporosis-pseudoglioma Syndrome, OPPG) and other eye diseases or syndromes associated with defects or damaged photoreceptors, familial exudative vitreoretinopathy, retinal angiogenesis, early coronary disease, tetra-amelia, Müllerian-duct regression and virilization, SERKAL syndrome, type II diabetes, Fuhrmann syndrome, Al-Awadi/Raas-Rothschild/Schinzel phocomelia syndrome, odonto-onycho-dermal dysplasia, obesity, split-hand/foot malformation, caudal duplication, tooth agenesis, Wilms tumor, skeletal dysplasia, focal dermal hypoplasia, autosomal recessive anonychia, neural tube defects, alpha-thalassemia (ATRX) syndrome, fragile X syndrome, ICF syndrome, Angelman's syndrome, Prader-Willi syndrome, Beckwith-Wiedemann Syndrome, Norrie disease, and Rett syndrome.

[017] In some embodiments, non-limiting examples of bone and cartilage diseases which can be treated with the compounds and compositions provided herein include bone spur (osteophytes), craniosynostosis, fibrodysplasia ossificans progressive, fibrous dysplasia, giant cell tumor of bone, hip labral tear, meniscal tears, bone/osteoporotic (wrist, spine, shoulder and hip) fractures, articular cartilage (chondral) defects, degenerative disc disease (or intervertebral disc

degeneration), osteochondritis dissecans, osteochondroma (bone tumor), osteopetrosis, relapsing polychondritis, and Salter-Harris fractures.

[018] In some embodiments, pharmaceutical compositions are provided that are effective for treatment of a disease of an animal, e.g., a mammal, caused by the pathological activation or mutations of the Wnt pathway. The composition includes a pharmaceutically acceptable carrier and a compound as described herein.

Definitions

[019] Unless defined otherwise, all technical and scientific terms used herein have the same meaning as is commonly understood by one of ordinary skill in the art to which this disclosure belongs. All patents, applications, published applications, and other publications are incorporated by reference in their entirety. In the event that there is a plurality of definitions for a term herein, those in this section prevail unless stated otherwise.

[020] As used herein, “alkyl” means a branched, or straight chain chemical group containing only carbon and hydrogen, such as methyl, ethyl, n-propyl, iso-propyl, n-butyl, iso-butyl, sec-butyl, tert-butyl, n-pentyl, iso-pentyl, sec-pentyl and neo-pentyl. Alkyl groups can either be unsubstituted or substituted with one or more substituents. Alkyl groups can be saturated or unsaturated (e.g., containing $-C=C-$ or $-C\equiv C-$ subunits), at one or several positions. In some embodiments, alkyl groups include 1 to 9 carbon atoms (for example, 1 to 6 carbon atoms, 1 to 4 carbon atoms, or 1 to 2 carbon atoms).

[021] As used herein, “alkylene” means a bivalent branched, or straight chain chemical group containing only carbon and hydrogen, such as methylene, ethylene, n-propylene, iso-propylene, n-butylene, iso-butylene, sec-butylene, tert-butylene, n-pentylene, iso-pentylene, sec-pentylene and neo-pentylene. Alkylene groups can either be unsubstituted or substituted with one or more substituents. Alkylene groups can be saturated or unsaturated (e.g., containing $-C=C-$ or $-C\equiv C-$ subunits), at one or several positions. In some embodiments, alkylene groups include 1 to 9 carbon atoms (for example, 1 to 6 carbon atoms, 1 to 4 carbon atoms, or 1 to 2 carbon atoms).

[022] As used herein, “carbocyclyl” means a cyclic ring system containing only carbon atoms in the ring system backbone, such as cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl, and cyclohexenyl. Carbocyclyls may include multiple fused rings. Carbocyclyls may have any degree of saturation provided that at least one ring in the ring system is not aromatic. Carbocyclyl groups can either be unsubstituted or substituted with one or more substituents. In some embodiments, carbocyclyl groups include 3 to 10 carbon atoms, for example, 3 to 6 carbon atoms.

[023] As used herein, “lower alkyl” means a subset of alkyl having 1 to 3 carbon atoms, which is linear or branched. Examples of lower alkyls include methyl, ethyl, n-propyl and isopropyl. Likewise, radicals using the terminology “lower” refer to radicals having 1 to about 3 carbons in the alkyl portion of the radical.

[024] As used herein, “aryl” means a mono-, bi-, tri- or polycyclic group with only carbon atoms present in the ring backbone having 5 to 14 ring atoms, alternatively 5, 6, 9, or 10 ring atoms; and having 6, 10, or 14 pi electrons shared in a cyclic array; wherein at least one ring in the system is aromatic. Aryl groups can either be unsubstituted or substituted with one or more substituents. Examples of aryl include phenyl, naphthyl, tetrahydronaphthyl, 2,3-dihydro-1H-indenyl, and others. In some embodiments, the aryl is phenyl.

[025] As used herein, “arylalkyl” means an aryl-alkyl- group in which the aryl and alkyl moieties are as previously described. In some embodiments, arylalkyl groups contain a C₁-alkyl moiety. Exemplary arylalkyl groups include benzyl and 2-phenethyl.

[026] As used herein, the term “heteroaryl” means a mono-, bi-, tri- or polycyclic group having 5 to 14 ring atoms, alternatively 5, 6, 9, or 10 ring atoms; and having 6, 10, or 14 pi electrons shared in a cyclic array; wherein at least one ring in the system is aromatic, and at least one ring in the system contains one or more heteroatoms independently selected from the group consisting of N, O, and S. Heteroaryl groups can either be unsubstituted or substituted with one or more substituents. Examples of heteroaryl include thienyl, pyridinyl, furyl, oxazolyl, oxadiazolyl, pyrrolyl, imidazolyl, triazolyl, thiodiazolyl, pyrazolyl, isoxazolyl, thiadiazolyl, pyranyl, pyrazinyl, pyrimidinyl, pyridazinyl, triazinyl, thiazolyl benzothienyl, benzoxadiazolyl, benzofuranyl, benzimidazolyl, benzotriazolyl, cinnolinyl, indazolyl, indolyl, isoquinolinyl, isothiazolyl, naphthyridinyl, purinyl, thienopyridinyl, pyrido[2,3-*d*]pyrimidinyl, pyrrolo[2,3-*b*]pyridinyl, quinazolinyl, quinolinyl, thieno[2,3-*c*]pyridinyl, pyrazolo[3,4-*b*]pyridinyl, pyrazolo[3,4-*c*]pyridinyl, pyrazolo[4,3-*c*]pyridine, pyrazolo[4,3-*b*]pyridinyl, tetrazolyl, chromane, 2,3-dihydrobenzo[*b*][1,4]dioxine, benzo[*d*][1,3]dioxole, 2,3-dihydrobenzofuran, 2,3-dihydrobenzo[*b*][1,4]oxathiine, and others. In some embodiments, the heteroaryl is selected from thienyl, pyridinyl, furyl, pyrazolyl, imidazolyl, pyranyl, pyrazinyl, and pyrimidinyl.

[027] As used herein, “halo”, “halide” or “halogen” is a chloro, bromo, fluoro, or iodo atom radical. In some embodiments, a halo is a chloro, bromo or fluoro. For example, a halide can be fluoro.

[028] As used herein, “haloalkyl” means a hydrocarbon substituent, which is a linear or branched, alkyl, alkenyl or alkynyl substituted with one or more chloro, bromo, fluoro, and/or iodo atom(s). In some embodiments, a haloalkyl is a fluoroalkyls, wherein one or more of the

hydrogen atoms have been substituted by fluoro. In some embodiments, haloalkyls are of 1 to about 3 carbons in length (e.g., 1 to about 2 carbons in length or 1 carbon in length). The term “haloalkylene” means a diradical variant of haloalkyl, and such diradicals may act as spacers between radicals, other atoms, or between a ring and another functional group.

[029] As used herein, “heterocyclyl” means a nonaromatic cyclic ring system comprising at least one heteroatom in the ring system backbone. Heterocyclyls may include multiple fused rings. Heterocyclyls may be substituted or unsubstituted with one or more substituents. In some embodiments, heterocycles have 5-7 members. In six membered monocyclic heterocycles, the heteroatom(s) are selected from one to three of O, N or S, and wherein when the heterocycle is five membered, it can have one or two heteroatoms selected from O, N, or S. Examples of heterocyclyl include aziriny, aziridinyl, azetidiny, oxetanyl, thietanyl, 1,4,2-dithiazolyl, dihydropyridinyl, 1,3-dioxanyl, 1,4-dioxanyl, 1,3-dioxolanyl, morpholinyl, thiomorpholinyl, piperazinyl, pyranyl, pyrrolidinyl, tetrahydrofuryl, tetrahydropyridinyl, oxazinyl, thiazinyl, thiinyl, thiazolidinyl, isothiazolidinyl, oxazolidinyl, isoxazolidinyl, piperidinyl, pyrazolidinyl, imidazolidinyl, thiomorpholinyl, and others. In some embodiments, the heterocyclyl is selected from azetidiny, morpholinyl, piperazinyl, pyrrolidinyl, and tetrahydropyridinyl.

[030] As used herein, “monocyclic heterocyclyl” means a single nonaromatic cyclic ring comprising at least one heteroatom in the ring system backbone. Heterocyclyls may be substituted or unsubstituted with one or more substituents. In some embodiments, heterocycles have 5-7 members. In six membered monocyclic heterocycles, the heteroatom(s) are selected from one to three of O, N or S, and wherein when the heterocycle is five membered, it can have one or two heteroatoms selected from O, N, or S. Examples of heterocyclyl include aziriny, aziridinyl, azetidiny, oxetanyl, thietanyl, 1,4,2-dithiazolyl, dihydropyridinyl, 1,3-dioxanyl, 1,4-dioxanyl, 1,3-dioxolanyl, morpholinyl, thiomorpholinyl, piperazinyl, pyranyl, pyrrolidinyl, tetrahydrofuryl, tetrahydropyridinyl, oxazinyl, thiazinyl, thiinyl, thiazolidinyl, isothiazolidinyl, oxazolidinyl, isoxazolidinyl, piperidinyl, pyrazolidinyl, imidazolidinyl, thiomorpholinyl, and others.

[031] The term “substituted” refers to moieties having substituents replacing a hydrogen on one or more non-hydrogen atoms of the molecule. It will be understood that “substitution” or “substituted with” includes the implicit proviso that such substitution is in accordance with permitted valence of the substituted atom and the substituent, and that the substitution results in a stable compound, e.g., which does not spontaneously undergo transformation such as by rearrangement, cyclization, elimination, etc. Substituents can include, for example, $-(C_{1-9} \text{ alkyl})$ optionally substituted with one or more of hydroxyl, $-NH_2$, $-NH(C_{1-3} \text{ alkyl})$, and $-N(C_{1-3} \text{ alkyl})_2$; $-(C_{1-9} \text{ haloalkyl})$; a halide; a hydroxyl; a carbonyl [such as $-C(O)OR$,

and -C(O)R]; a thiocarbonyl [such as -C(S)OR, -C(O)SR, and -C(S)R]; -(C₁₋₉ alkoxy) optionally substituted with one or more of halide, hydroxyl, -NH₂, -NH(C₁₋₃ alkyl), and -N(C₁₋₃ alkyl)₂; -OPO(OH)₂; a phosphonate [such as -PO(OH)₂ and -PO(OR')₂]; -OPO(OR')R''; -NRR'; -C(O)NRR'; -C(NR)NR'R''; -C(NR')R''; a cyano; a nitro; an azido; -SH; -S-R; -OSO₂(OR); a sulfonate [such as -SO₂(OH) and -SO₂(OR)]; -SO₂NR'R''; and -SO₂R; in which each occurrence of R, R' and R'' are independently selected from H; -(C₁₋₉ alkyl); C₆₋₁₀ aryl optionally substituted with from 1-3R'''; 5-10 membered heteroaryl having from 1-4 heteroatoms independently selected from N, O, and S and optionally substituted with from 1-3 R'''; C₃₋₇ carbocyclyl optionally substituted with from 1-3 R'''; and 3-8 membered heterocyclyl having from 1-4 heteroatoms independently selected from N, O, and S and optionally substituted with from 1-3 R'''; wherein each R''' is independently selected from -(C₁₋₆ alkyl), -(C₁₋₆ haloalkyl), a halide (e.g., F), a hydroxyl, -C(O)OR, -C(O)R, -(C₁₋₆ alkoxy), -NRR', -C(O)NRR', and a cyano, in which each occurrence of R and R' is independently selected from H and -(C₁₋₆ alkyl). In some embodiments, the substituent is selected from -(C₁₋₆ alkyl), -(C₁₋₆ haloalkyl), a halide (e.g., F), a hydroxyl, -C(O)OR, -C(O)R, -(C₁₋₆ alkoxy), -NRR', -C(O)NRR', and a cyano, in which each occurrence of R and R' is independently selected from H and -(C₁₋₆ alkyl).

[032] As used herein, when two groups are indicated to be “linked” or “bonded” to form a “ring”, it is to be understood that a bond is formed between the two groups and may involve replacement of a hydrogen atom on one or both groups with the bond, thereby forming a carbocyclyl, heterocyclyl, aryl, or heteroaryl ring. The skilled artisan will recognize that such rings can and are readily formed by routine chemical reactions. In some embodiments, such rings have from 3-7 members, for example, 5 or 6 members.

[033] The skilled artisan will recognize that some structures described herein may be resonance forms or tautomers of compounds that may be fairly represented by other chemical structures, even when kinetically, the artisan recognizes that such structures are only a very small portion of a sample of such compound(s). Such compounds are clearly contemplated within the scope of this disclosure, though such resonance forms or tautomers are not represented herein.

[034] The compounds provided herein may encompass various stereochemical forms. The compounds also encompass diastereomers as well as optical isomers, e.g., mixtures of enantiomers including racemic mixtures, as well as individual enantiomers and diastereomers, which arise as a consequence of structural asymmetry in certain compounds. Separation of the individual isomers or selective synthesis of the individual isomers is accomplished by application of various methods which are well known to practitioners in the art. Unless otherwise indicated, when a disclosed compound is named or depicted by a structure without specifying the

stereochemistry and has one or more chiral centers, it is understood to represent all possible stereoisomers of the compound.

[035] The term “administration” or “administering” refers to a method of providing a dosage of a compound or pharmaceutical composition to a vertebrate or invertebrate, including a mammal, a bird, a fish, or an amphibian, where the method is, e.g., orally, subcutaneously, intravenously, intralymphatic, intranasally, topically, transdermally, intraperitoneally, intramuscularly, intrapulmonarily, vaginally, rectally, ontologically, neuro-otologically, intraocularly, subconjunctivally, via anterior eye chamber injection, intravitreally, intraperitoneally, intrathecally, intracisternally, intrapleurally, via wound irrigation, intrabuccally, intra-abdominally, intra-articularly, intra-aurally, intrabronchially, intracapsularly, intrameningeally, via inhalation, via endotracheal or endobronchial instillation, via direct instillation into pulmonary cavities, intraspinally, intrasynovially, intrathoracically, via thoracostomy irrigation, epidurally, intratympanically, intracisternally, intravascularly, intraventricularly, intraosseously, via irrigation of infected bone, or via application as part of any admixture with a prosthetic device. The method of administration can vary depending on various factors, e.g., the components of the pharmaceutical composition, the site of the disease, the disease involved, and the severity of the disease.

[036] A “diagnostic” as used herein is a compound, method, system, or device that assists in the identification or characterization of a health or disease state. The diagnostic can be used in standard assays as is known in the art.

[037] The term “mammal” is used in its usual biological sense. Thus, it specifically includes humans, cattle, horses, monkeys, dogs, cats, mice, rats, cows, sheep, pigs, goats, and non-human primates, but also includes many other species.

[038] The term “pharmaceutically acceptable carrier”, “pharmaceutically acceptable diluent” or “pharmaceutically acceptable excipient” includes any and all solvents, co-solvents, complexing agents, dispersion media, coatings, isotonic and absorption delaying agents and the like which are not biologically or otherwise undesirable. The use of such media and agents for pharmaceutically active substances is well known in the art. Except insofar as any conventional media or agent is incompatible with the active ingredient, its use in the therapeutic compositions is contemplated. Supplementary active ingredients can also be incorporated into the compositions. In addition, various adjuvants such as are commonly used in the art may be included. These and other such compounds are described in the literature, e.g., in the Merck Index, Merck & Company, Rahway, NJ. Considerations for the inclusion of various components in pharmaceutical compositions are described, e.g., in Gilman *et al.* (Eds.) (2010); Goodman and Gilman’s: The Pharmacological Basis of Therapeutics, 12th Ed., The McGraw-Hill Companies.

[039] The term “pharmaceutically acceptable salt” refers to salts that retain the biological effectiveness and properties of the compounds provided herein and, which are not biologically or otherwise undesirable. In many cases, the compounds provided herein are capable of forming acid and/or base salts by virtue of the presence of amino and/or carboxyl groups or groups similar thereto. Many such salts are known in the art, for example, as described in WO 87/05297. Pharmaceutically acceptable acid addition salts can be formed with inorganic acids and organic acids. Inorganic acids from which salts can be derived include, for example, hydrochloric acid, hydrobromic acid, sulfuric acid, nitric acid, phosphoric acid, and the like. Organic acids from which salts can be derived include, for example, acetic acid, propionic acid, glycolic acid, pyruvic acid, oxalic acid, maleic acid, malonic acid, succinic acid, fumaric acid, tartaric acid, citric acid, benzoic acid, cinnamic acid, mandelic acid, methanesulfonic acid, ethanesulfonic acid, p-toluenesulfonic acid, salicylic acid, and the like. Pharmaceutically acceptable base addition salts can be formed with inorganic and organic bases. Inorganic bases from which salts can be derived include, for example, sodium, potassium, lithium, ammonium, calcium, magnesium, iron, zinc, copper, manganese, aluminum, and the like; particularly preferred are the ammonium, potassium, sodium, calcium, and magnesium salts. Organic bases from which salts can be derived include, for example, primary, secondary, and tertiary amines, substituted amines including naturally occurring substituted amines, cyclic amines, basic ion exchange resins, and the like, specifically such as isopropylamine, trimethylamine, diethylamine, triethylamine, tripropylamine, and ethanolamine.

[040] “Solvate” refers to the compound formed by the interaction of a solvent and a compound as provided herein or a salt thereof. Suitable solvates are pharmaceutically acceptable solvates including hydrates.

[041] “Patient” as used herein, means a human or a non-human mammal, e.g., a dog, a cat, a mouse, a rat, a cow, a sheep, a pig, a goat, a non-human primate, or a bird, e.g., a chicken, as well as any other vertebrate or invertebrate. In some embodiments, the patient is a human.

[042] A “therapeutically effective amount” or “pharmaceutically effective amount” of a compound as provided herein is one which is sufficient to achieve the desired physiological effect and may vary according to the nature and severity of the disease condition, and the potency of the compound. “Therapeutically effective amount” is also intended to include one or more of the compounds of Formula I in combination with one or more other agents that are effective to treat the diseases and/or conditions described herein. The combination of compounds can be a synergistic combination. Synergy, as described, for example, by Chou and Talalay, *Advances in Enzyme Regulation* (1984), 22, 27-55, occurs when the effect of the compounds when administered in combination is greater than the additive effect of the compounds when administered alone as a

single agent. In general, a synergistic effect is most clearly demonstrated at sub-optimal concentrations of the compounds. It will be appreciated that different concentrations may be employed for prophylaxis than for treatment of an active disease. This amount can further depend upon the patient's height, weight, sex, age and medical history.

[043] A therapeutic effect relieves, to some extent, one or more of the symptoms of the disease.

[044] "Treat," "treatment," or "treating," as used herein refers to administering a compound or pharmaceutical composition as provided herein for therapeutic purposes. The term "therapeutic treatment" refers to administering treatment to a patient already suffering from a disease thus causing a therapeutically beneficial effect, such as ameliorating existing symptoms, ameliorating the underlying metabolic causes of symptoms, postponing or preventing the further development of a disorder, and/or reducing the severity of symptoms that will or are expected to develop.

[045] "Drug-eluting" and/or controlled release as used herein refers to any and all mechanisms, e.g., diffusion, migration, permeation, and/or desorption by which the drug(s) incorporated in the drug-eluting material pass therefrom over time into the surrounding body tissue.

[046] "Drug-eluting material" and/or controlled release material as used herein refers to any natural, synthetic or semi-synthetic material capable of acquiring and retaining a desired shape or configuration and into which one or more drugs can be incorporated and from which incorporated drug(s) are capable of eluting over time.

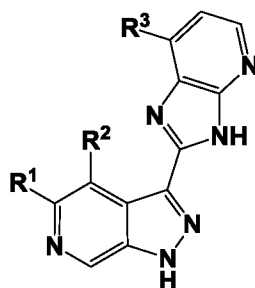
[047] "Elutable drug" as used herein refers to any drug or combination of drugs having the ability to pass over time from the drug-eluting material in which it is incorporated into the surrounding areas of the body.

[048] The term "comprising" as used herein is synonymous with "including," "containing," or "characterized by," and is inclusive or open-ended and does not exclude additional, unrecited elements or method steps.

Compounds

[049] The compounds and compositions described herein can be used as anti-proliferative agents, e.g., anti-cancer and anti-angiogenesis agents, and/or as inhibitors of the Wnt signaling pathway, e.g., for treating diseases or disorders associated with aberrant Wnt signaling. In addition, the compounds can be used as inhibitors of one or more kinases, kinase receptors, or kinase complexes. Such compounds and compositions are also useful for controlling cellular proliferation, differentiation, and/or apoptosis.

[050] Some embodiments of the present disclosure include compounds of Formula I:



I

or salts, pharmaceutically acceptable salts, or prodrugs thereof.

[051] In some embodiments, R^1 is selected from the group consisting of –pyridinyl(R^4) and –pyrimidinyl(R^5).

[052] In some embodiments, R^1 is selected from the group consisting of –heteroaryl(R^4)_q and –heterocyclyl(R^5)_h.

[053] In some embodiments, R^1 is selected from the group consisting of –piperidinyl(R^5)_h and –tetrahydropyridinyl(R^5)_h.

[054] In some embodiments, R^1 is selected from the group consisting of –pyridinyl(R^4)_q, –pyrimidinyl(R^4)_q, –pyrazinyl(R^4)_q, –pyrazolyl(R^4)_q, and –imidazolyl(R^4)_q.

[055] In some embodiments, R^2 is selected from the group consisting of H and halide.

[056] In some embodiments, R^3 is selected from the group consisting of –heteroaryl(R^6)_q, –heterocyclyl(R^7)_h, and –aryl(R^8)_k.

[057] In some embodiments, R^3 is selected from the group consisting of H, –heteroaryl(R^6)_q, –heterocyclyl(R^7)_h, and –aryl(R^8)_k.

[058] In some embodiments, R^3 is selected from the group consisting of –pyridinyl(R^6)_q, –imidazolyl(R^6)_q, –furanlyl(R^6)_q, –thiophenyl(R^6)_q, –piperidinyl(R^7)_h, –piperazinyl(R^7)_h, and –phenyl(R^8)_k.

[059] In some embodiments, R^4 is one substituent attached to the pyridinyl and is independently selected from the group consisting of H, halide, –(C₁₋₆ alkyl), –(C₁₋₄ alkylene)_p, heterocyclyl(R^9)_h, –(C₁₋₄ alkylene)_pcarbocyclyl(R^{10})_j, –(C₁₋₄ alkylene)_paryl(R^{11})_k, –NHC(=O) R^{12} , –NR¹³ R^{14} , and –(C₁₋₆ alkylene)NR¹⁵ R^{16} .

[060] In some embodiments, each R^4 is one substituent attached to the heteroaryl and is independently selected from the group consisting of halide, –(C₁₋₆ alkyl), –(C₁₋₄

alkylene)_pheterocyclyl(R⁹)_h, -(C₁₋₄ alkylene)_pcarbocyclyl(R¹⁰)_j, -(C₁₋₄ alkylene)_paryl(R¹¹)_k, -NHC(=O)R¹², -NR¹³R¹⁴, -(C₁₋₆ alkylene)NR¹⁵R¹⁶, and -OR²².

[061] In some embodiments, each R⁴ is one substituent attached to the heteroaryl and is independently selected from the group consisting of F, -Me, -Et, -(CH₂)heterocyclyl(R⁹)_h, -heterocyclyl(R⁹)_h, -(CH₂)carbocyclyl(R¹⁰)_j, -(CH₂)aryl(R¹¹)_k, -NHC(=O)(C₁₋₅ alkyl), -NHC(=O)phenyl(R¹⁹)_k, -NHC(=O)(CH₂)phenyl(R¹⁹)_k, -NHC(=O)carbocyclyl(R²⁰)_j, -NHC(=O)(CH₂)heterocyclyl(R²¹)_h, -NH₂, -N(C₁₋₃ alkyl)₂, -NH(C₁₋₄ alkyl), -(CH₂)N(C₁₋₃ alkyl)₂, -(CH₂)NH(C₁₋₄ alkyl), -OH, -O(C₁₋₃ alkyl), -Ocarbocyclyl(R²⁰)_j, -Oheterocyclyl(R²¹)_h, -O(CH₂CH₂)heterocyclyl(R²¹)_h, -O(CH₂CH₂)N(C₁₋₃ alkyl)₂, and -O(CH₂)phenyl(R¹⁹)_k.

[062] In some embodiments, R⁵ is one substituent attached to the pyrimidinyl and is independently selected from the group consisting of H, halide, -(C₁₋₆ alkyl), -(C₁₋₄ alkylene)_pheterocyclyl(R⁹)_h, -(C₁₋₄ alkylene)_pcarbocyclyl(R¹⁰)_j, -(C₁₋₄ alkylene)_paryl(R¹¹)_k, -NHC(=O)R¹², -NR¹³R¹⁴, and -(C₁₋₆ alkylene)NR¹⁵R¹⁶.

[063] In some embodiments, each R⁵ is one substituent attached to the heterocyclyl and is independently selected from the group consisting of -(C₁₋₄ alkyl), halide, -CF₃, and -CN.

[064] In some embodiments, each R⁶ is one substituent attached to the heteroaryl and is independently selected from the group consisting of H, -(C₁₋₆ alkyl), halide, -CF₃, -OCH₃, -CN, and -C(=O)R¹⁷.

[065] In some embodiments, each R⁶ is one substituent attached to the heteroaryl and is independently selected from the group consisting of -(C₁₋₆ alkyl), halide, -CF₃, -OCH₃, -CN, and -C(=O)R¹⁷.

[066] In some embodiments, each R⁶ is one substituent attached to the heteroaryl and is independently selected from the group consisting of -Me, -Et, F, -CF₃, -OCH₃, -CN, and -C(=O)(C₁₋₃ alkyl).

[067] In some embodiments, each R⁷ is one substituent attached to the heterocyclyl and is independently selected from the group consisting of H, -(C₁₋₆ alkyl), halide, -CF₃, -CN, and -OCH₃.

[068] In some embodiments, each R⁷ is one substituent attached to the heterocyclyl and is independently selected from the group consisting of -(C₁₋₆ alkyl), halide, -CF₃, -CN, and -OCH₃.

[069] In some embodiments, each R⁸ is one substituent attached to the aryl and is independently selected from the group consisting of H, -(C₁₋₆ alkyl), halide, -CF₃, -CN, -OCH₃, -(C₁₋₆ alkylene)_pNHSO₂R¹⁷, -NR¹³(C₁₋₆ alkylene)NR¹³R¹⁴, and -(C₁₋₆ alkylene)_pNR¹³R¹⁴.

[070] In some embodiments, each R⁸ is one substituent attached to the aryl and is independently selected from the group consisting of -(C₁₋₆ alkyl), halide, -CF₃, -CN, -OCH₃, -(C₁₋₆ alkylene)_pNHSO₂R¹⁷, -NR¹³(C₁₋₆ alkylene)NR¹³R¹⁴, -(C₁₋₆ alkylene)_pNR¹³R¹⁴, and -OR²⁵.

[071] In some embodiments, each R⁸ is one substituent attached to the aryl and is independently selected from the group consisting of -Me, -Et, F, -CF₃, -CN, -OCH₃, -(CH₂CH₂)NHSO₂(C₁₋₃ alkyl), -NH(CH₂CH₂)N(C₁₋₃ alkyl)₂, -OH, -O(C₁₋₃ alkyl), -O(CH₂CH₂)heterocyclyl(R²¹)_h, and -O(CH₂CH₂)N(C₁₋₃ alkyl)₂.

[072] In some embodiments, each R⁹ is one substituent attached to the heterocyclyl and is independently selected from the group consisting of H, -(C₁₋₄ alkyl), halide, -CF₃, and -CN.

[073] In some embodiments, each R⁹ is one substituent attached to the heterocyclyl and is independently selected from the group consisting of amino, -(C₁₋₄ alkyl), halide, -CF₃, and -CN.

[074] In some embodiments, each R⁹ is one substituent attached to the heterocyclyl and is independently selected from the group consisting of amino, Me, Et, F, Cl, and -CF₃.

[075] In some embodiments, each R¹⁰ is one substituent attached to the carbocyclyl and is independently selected from the group consisting of H, -(C₁₋₄ alkyl), halide, -CF₃, and -CN.

[076] In some embodiments, each R¹⁰ is one substituent attached to the carbocyclyl and is independently selected from the group consisting of -(C₁₋₄ alkyl), halide, -CF₃, and -CN.

[077] In some embodiments, each R¹⁰ is one substituent attached to the carbocyclyl and is independently selected from the group consisting of Me, Et, F, Cl, and -CF₃.

[078] In some embodiments, each R¹¹ is one substituent attached to the aryl and is independently selected from the group consisting of H, -(C₁₋₄ alkyl), halide, -CF₃, and -CN.

[079] In some embodiments, each R¹¹ is one substituent attached to the aryl and is independently selected from the group consisting of -(C₁₋₄ alkyl), halide, -CF₃, and -CN.

[080] In some embodiments, each R¹¹ is one substituent attached to the aryl and is independently selected from the group consisting of Me, Et, F, Cl, and -CF₃.

[081] In some embodiments, each R¹² is independently selected from the group consisting of -(C₁₋₉ alkyl), -heteroaryl(R¹⁸)_q, -aryl(R¹⁹)_k, -CH₂aryl(R¹⁹)_k, -carbocyclyl(R²⁰)_j, and -CH₂carbocyclyl(R²⁰)_j.

[082] In some embodiments, each R¹² is independently selected from the group consisting of -(C₁₋₉ alkyl), -heteroaryl(R¹⁸)_q, -aryl(R¹⁹)_k, -CH₂aryl(R¹⁹)_k, -carbocyclyl(R²⁰)_j, -CH₂carbocyclyl(R²⁰)_j, -(C₁₋₄ alkylene)_pNR²³R²⁴, -heterocyclyl(R²¹)_h, and -CH₂heterocyclyl(R²¹)_h.

[083] In some embodiments, each R^{12} is independently selected from the group consisting of $-(C_{1-5} \text{ alkyl})$, $-\text{phenyl}(R^{19})_k$, $-(CH_2)\text{phenyl}(R^{19})_k$, $-\text{carbocyclyl}(R^{20})_j$, $-(CH_2)\text{carbocyclyl}(R^{20})_j$, $-(CH_2)N(C_{1-3} \text{ alkyl})_2$, and $-(CH_2)\text{heterocyclyl}(R^{21})_h$.

[084] In some embodiments, each R^{13} is independently selected from the group consisting of H and $-(C_{1-6} \text{ alkyl})$.

[085] In some embodiments, each R^{13} is independently selected from the group consisting of H and $-(C_{1-3} \text{ alkyl})$.

[086] In some embodiments, each R^{14} is independently selected from the group consisting of H, $-(C_{1-6} \text{ alkyl})$, $-\text{CH}_2\text{aryl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$.

[087] In some embodiments, each R^{14} is independently selected from the group consisting of H, $-(C_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$.

[088] In some embodiments, each R^{15} is independently selected from the group consisting of H and $-(C_{1-6} \text{ alkyl})$.

[089] In some embodiments, each R^{15} is independently selected from the group consisting of H and $-(C_{1-3} \text{ alkyl})$.

[090] In some embodiments, each R^{16} is independently selected from the group consisting of H, $-(C_{1-6} \text{ alkyl})$, $-\text{CH}_2\text{aryl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$.

[091] In some embodiments, each R^{16} is independently selected from the group consisting of H, $-(C_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$.

[092] In some embodiments, each R^{17} is independently a $-(C_{1-6} \text{ alkyl})$.

[093] In some embodiments, each R^{17} is independently a $-(C_{1-3} \text{ alkyl})$.

[094] In some embodiments, each R^{18} is one substituent attached to the heteroaryl and is independently selected from the group consisting of H, $-(C_{1-4} \text{ alkyl})$, halide, $-\text{CF}_3$, and $-\text{CN}$.

[095] In some embodiments, each R^{18} is one substituent attached to the heteroaryl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-\text{CF}_3$, and $-\text{CN}$.

[096] In some embodiments, each R^{18} is one substituent attached to the heteroaryl and is independently selected from the group consisting of Me, Et, F, Cl, and $-\text{CF}_3$.

[097] In some embodiments, each R^{19} is one substituent attached to the aryl and is independently selected from the group consisting of H, $-(C_{1-4} \text{ alkyl})$, halide, $-\text{CF}_3$, and $-\text{CN}$.

[098] In some embodiments, each R^{19} is one substituent attached to the aryl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-\text{CF}_3$, and $-\text{CN}$.

[099] In some embodiments, each R^{19} is one substituent attached to the aryl and is independently selected from the group consisting of Me, Et, F, Cl, and $-\text{CF}_3$.

[0100] In some embodiments, each R^{20} is one substituent attached to the carbocyclyl and is independently selected from the group consisting of H, $-(C_{1-4} \text{ alkyl})$, halide, $-CF_3$, and $-CN$.

[0101] In some embodiments, each R^{20} is one substituent attached to the carbocyclyl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-CF_3$, and $-CN$.

[0102] In some embodiments, each R^{20} is one substituent attached to the carbocyclyl and is independently selected from the group consisting of Me, Et, F, Cl, and $-CF_3$.

[0103] In some embodiments, each R^{21} is one substituent attached to the heterocyclyl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-CF_3$, and $-CN$.

[0104] In some embodiments, each R^{21} is one substituent attached to the heterocyclyl and is independently selected from the group consisting of Me, Et, F, Cl, and $-CF_3$.

[0105] In some embodiments, R^{22} is selected from the group consisting of H, $-(C_{1-6} \text{ alkyl})$, $-(C_{1-4} \text{ alkylene})_p \text{heterocyclyl}(R^{21})_h$, $-(C_{1-4} \text{ alkylene})_p \text{carbocyclyl}(R^{20})_j$, $-(C_{1-4} \text{ alkylene})_p \text{aryl}(R^{19})_k$, and $-(C_{1-6} \text{ alkylene})_p \text{NR}^{23}R^{24}$.

[0106] In some embodiments, R^{22} is selected from the group consisting of H, $-Me$, $-Et$, $-iPr$, $-\text{heterocyclyl}(R^{21})_h$, $-(CH_2CH_2)\text{heterocyclyl}(R^{21})_h$, $-\text{carbocyclyl}(R^{20})_j$, $-(CH_2)\text{phenyl}(R^{19})_k$, and $-(CH_2CH_2)N(C_{1-3} \text{ alkyl})_2$.

[0107] In some embodiments, each R^{23} is independently selected from the group consisting of H and $-(C_{1-6} \text{ alkyl})$.

[0108] In some embodiments, each R^{23} is independently selected from the group consisting of Me and Et.

[0109] In some embodiments, each R^{24} is independently selected from the group consisting of H and $-(C_{1-6} \text{ alkyl})$.

[0110] In some embodiments, each R^{24} is independently selected from the group consisting of Me and Et.

[0111] In some embodiments, R^{25} is selected from the group consisting of H, $-(C_{1-6} \text{ alkyl})$, $-(C_{1-4} \text{ alkylene})_p \text{heterocyclyl}(R^{21})_h$, and $-(C_{1-6} \text{ alkylene})_p \text{NR}^{23}R^{24}$.

[0112] In some embodiments, R^{25} is selected from the group consisting of H, $-Me$, $-Et$, $-iPr$, $-(CH_2CH_2)\text{heterocyclyl}(R^{21})_h$, and $-(CH_2CH_2)N(C_{1-3} \text{ alkyl})_2$.

[0113] In some embodiments, each p is independently 0 or 1.

[0114] In some embodiments, each q is independently 1 to 4.

[0115] In some embodiments, each h is independently 1 to 10.

[0116] In some embodiments, each k is independently 1 to 5.

[0117] In some embodiments, each j is independently 1 to 12.

[0118] In some embodiments, each p is independently 0 or 1; in some embodiments, each p is 0; in some embodiments, each p is 1.

[0119] In some embodiments, each q is independently 0 to 4; in some embodiments, each q is 0; in some embodiments, each q is 1; in some embodiments, each q is 2; in some embodiments, each q is 3; in some embodiments, each q is 4.

[0120] In some embodiments, each h is independently 0 to 10; in some embodiments, each h is 0; in some embodiments, each h is 1; in some embodiments, each h is 2; in some embodiments, each h is 3; in some embodiments, each h is 4.

[0121] In some embodiments, each k is independently 0 to 5; in some embodiments, each k is 0; in some embodiments, each k is 1; in some embodiments, each k is 2; in some embodiments, each k is 3.

[0122] In some embodiments, each j is independently 0 to 12; in some embodiments, each j is 0; in some embodiments, each j is 1; in some embodiments, each j is 2; in some embodiments, each j is 3; in some embodiments, each j is 4.

[0123] In some embodiments, each R⁴ is one substituent attached to the heteroaryl and is selected from the group consisting of $-(C_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{heterocyclyl}(R^9)_h$, $-\text{NHC}(=\text{O})R^{12}$, $-\text{NR}^{13}R^{14}$, and $-\text{CH}_2\text{NR}^{15}R^{16}$.

[0124] In some embodiments, at least one R⁹ is halide.

[0125] In some embodiments, R¹² is selected from the group consisting of $-(C_{1-5} \text{ alkyl})$, $-\text{phenyl}(R^{19})_k$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{carbocyclyl}(R^{20})_j$.

[0126] In some embodiments, R¹³ and R¹⁴ are independently selected from H and $-(C_{1-5} \text{ alkyl})$.

[0127] In some embodiments, R¹⁵ and R¹⁶ are independently selected from H and $-(C_{1-5} \text{ alkyl})$.

[0128] In some embodiments, k is 1 or 2 and each R⁸ is independently a halide.

[0129] In some embodiments, k is 2, one R⁸ is halide and the other R⁸ is $-\text{CH}_2\text{NHSO}_2R^{17}$.

[0130] In some embodiments, R¹⁷ is $-(C_{1-3} \text{ alkyl})$.

[0131] In some embodiments, k is 2, one R⁸ is halide and the other R⁸ is $-\text{NHCH}_2\text{CH}_2\text{NR}^{13}R^{14}$.

[0132] In some embodiments, R¹³ and R¹⁴ are independently selected from H and $-(C_{1-3} \text{ alkyl})$.

[0133] In some embodiments, R³ is selected from the group consisting of $-\text{pyridinyl}(R^6)_q$, $-\text{imidazolyl}(R^6)_q$, $-\text{furanlyl}(R^6)_q$, and $-\text{thiophenyl}(R^6)_q$.

[0134] In some embodiments, q is 0 or 1, R⁶ is selected from the group consisting of halide, -(C₁₋₃ alkyl), and -C(=O)R¹⁷, wherein R¹⁷ is -(C₁₋₂ alkyl).

[0135] In some embodiments, R³ is selected from the group consisting of -piperidinylnyl(R⁷)_h and -piperazinyl(R⁷)_h.

[0136] In some embodiments, q is 1, and R⁷ is selected from the group consisting of H and -(C₁₋₃ alkyl).

[0137] In some embodiments, R² is H; in other embodiments, R² is halide, e.g. F.

[0138] In some embodiments, R¹ is -heteroaryl(R⁴)_q.

[0139] In some embodiments, R¹ is -pyridinyl(R⁴)_q.

[0140] In some embodiments, R¹ is -pyridin-3-yl(R⁴)_q.

[0141] In some embodiments, R¹ is -pyrimidinyl(R⁴)_q.

[0142] In some embodiments, R¹ is -pyrimidin-5-yl(R⁴)_q.

[0143] In some embodiments, R¹ is -pyrimidin-5-yl(R⁴)_q and q is 0.

[0144] In some embodiments, R¹ is -pyrazinyl(R⁴)_q.

[0145] In some embodiments, R¹ is -pyrazolyl(R⁴)_q.

[0146] In some embodiments, R¹ is -pyrazol-4-yl(R⁴)_q, q is 1, and R⁴ is Me.

[0147] In some embodiments, R¹ is -pyrazol-4-yl(R⁴)_q and q is 0.

[0148] In some embodiments, R¹ is -imidazolyl(R⁴)_q.

[0149] In some embodiments, R¹ is -imidazol-5-yl(R⁴)_q, q is 1, and R⁴ is Me.

[0150] In some embodiments, R¹ is -imidazol-5-yl(R⁴)_q, q is 2, and both R⁴ are Me.

[0151] In some embodiments, R¹ is -heterocyclyl(R⁵)_h.

[0152] In some embodiments, R¹ is -piperidinylnyl(R⁵)_h.

[0153] In some embodiments, R¹ is -piperidin-4-yl(R⁵)_h.

[0154] In some embodiments, R¹ is -piperidin-4-yl(R⁵)_h, and h is 0.

[0155] In some embodiments, R¹ is -tetrahydropyridinylnyl(R⁵)_h.

[0156] In some embodiments, R¹ is -1,2,3,6-tetrahydropyridinylnyl(R⁵)_h.

[0157] In some embodiments, R¹ is -1,2,3,6-tetrahydropyridinylnyl(R⁵)_h, and h is 0.

[0158] In some embodiments, R³ is H.

[0159] In some embodiments, R³ is -heteroaryl(R⁶)_q.

[0160] In some embodiments, R³ is -heterocyclyl(R⁷)_h.

[0161] In some embodiments, R³ is -piperidinylnyl(R⁷)_h.

[0162] In some embodiments, R³ is -piperazinyl(R⁷)_h.

[0163] In some embodiments, R³ is -aryl(R⁸)_k.

[0164] In some embodiments, R³ is -pyridinyl(R⁶)_q.

- [0165] In some embodiments, R^3 is $-\text{pyridin-3-yl}(R^6)_q$.
- [0166] In some embodiments, R^3 is $-\text{pyridin-4-yl}(R^6)_q$.
- [0167] In some embodiments, R^3 is $-\text{pyridin-5-yl}(R^6)_q$.
- [0168] In some embodiments, R^3 is $-\text{pyridin-3-yl}(R^6)_q$, q is 0.
- [0169] In some embodiments, R^3 is $-\text{pyridin-4-yl}(R^6)_q$, q is 0.
- [0170] In some embodiments, R^3 is $-\text{pyridin-5-yl}(R^6)_q$, q is 0.
- [0171] In some embodiments, R^3 is $-\text{imidazolyl}(R^6)_q$.
- [0172] In some embodiments, R^3 is $-\text{imidazol-1-yl}(R^6)_q$, q is 1, and R^6 is $-(C_{1-3} \text{ alkyl})$.
- [0173] In some embodiments, R^3 is $-\text{imidazol-1-yl}(R^6)_q$, q is 1, and R^6 is methyl.
- [0174] In some embodiments, R^3 is $-\text{furan-2-yl}(R^6)_q$.
- [0175] In some embodiments, R^3 is $-\text{furan-2-yl}(R^6)_q$.
- [0176] In some embodiments, R^3 is $-\text{furan-2-yl}(R^6)_q$ and q is 0.
- [0177] In some embodiments, R^3 is $-\text{furan-3-yl}(R^6)_q$.
- [0178] In some embodiments, R^3 is $-\text{furan-3-yl}(R^6)_q$ and q is 0.
- [0179] In some embodiments, R^3 is $-\text{thiophenyl}(R^6)_q$.
- [0180] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$.
- [0181] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$ and q is 0.
- [0182] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1 or 2, and each R^6 is independently a halide.
- [0183] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1 or 2, and R^6 is F.
- [0184] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1 or 2, and each R^6 is independently $-(C_{1-6} \text{ alkyl})$.
- [0185] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1 or 2, and each R^6 is independently $-(C_{1-2} \text{ alkyl})$.
- [0186] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1 or 2, and R^6 is methyl.
- [0187] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1 or 2, and R^6 is $-\text{CF}_3$.
- [0188] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1 or 2, and R^6 is CN.
- [0189] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1, and R^6 is $-\text{C}(=\text{O})R^{17}$.
- [0190] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1, R^6 is $-\text{C}(=\text{O})R^{17}$, and R^{17} is $-(C_{1-6} \text{ alkyl})$.
- [0191] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1, R^6 is $-\text{C}(=\text{O})R^{17}$, and R^{17} is $-(C_{1-4} \text{ alkyl})$.
- [0192] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1, R^6 is $-\text{C}(=\text{O})R^{17}$, and R^{17} is $-(C_{1-2} \text{ alkyl})$.

[0193] In some embodiments, R^3 is $-\text{thiophen-2-yl}(R^6)_q$, q is 1, R^6 is $-\text{C}(=\text{O})R^{17}$, and R^{17} is methyl.

[0194] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$.

[0195] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$ and q is 0.

[0196] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1 or 2, and each R^6 is independently halide.

[0197] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1 or 2, and R^6 is F.

[0198] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1 or 2, and each R^6 is independently $-(\text{C}_{1-6}$ alkyl).

[0199] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1 or 2, and each R^6 is independently $-(\text{C}_{1-2}$ alkyl).

[0200] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1 or 2, and R^6 is methyl.

[0201] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1 or 2, and R^6 is $-\text{CF}_3$.

[0202] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1 or 2, and R^6 is CN.

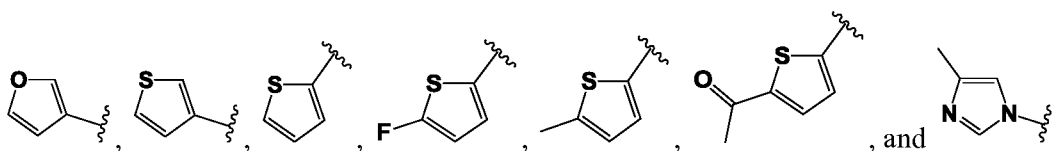
[0203] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1, and R^6 is $-\text{C}(=\text{O})R^{17}$.

[0204] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1, R^6 is $-\text{C}(=\text{O})R^{17}$, and R^{17} is $-(\text{C}_{1-4}$ alkyl).

[0205] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1, R^6 is $-\text{C}(=\text{O})R^{17}$, and R^{17} is $-(\text{C}_{1-2}$ alkyl).

[0206] In some embodiments, R^3 is $-\text{thiophen-3-yl}(R^6)_q$, q is 1, R^6 is $-\text{C}(=\text{O})R^{17}$, and R^{17} is methyl.

[0207] In some embodiments, R^3 is selected from the group consisting of:



[0208] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$.

[0209] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$ and k is 0.

[0210] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 1 or 2, and each R^8 is independently a halide.

[0211] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 1 or 2, and R^8 is F.

[0212] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 1, and R^8 is F.

[0213] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is a halide and the other R^8 is $-(\text{C}_{1-6}$ alkylene)_pNHSO₂R¹⁷.

[0214] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is a halide and the other R^8 is $-(C_{1-4} \text{ alkylene})_p\text{NHSO}_2R^{17}$, and p is 1.

[0215] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is a halide and the other R^8 is $-(C_{1-2} \text{ alkylene})_p\text{NHSO}_2R^{17}$, and p is 1.

[0216] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is a halide and the other R^8 is $-\text{CH}_2\text{NHSO}_2R^{17}$.

[0217] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is a halide and the other R^8 is $-\text{CH}_2\text{NHSO}_2R^{17}$, and R^{17} is $-(C_{1-4} \text{ alkyl})$.

[0218] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is a halide and the other R^8 is $-\text{CH}_2\text{NHSO}_2R^{17}$, and R^{17} is $-(C_{1-2} \text{ alkyl})$.

[0219] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is a halide and the other R^8 is $-\text{CH}_2\text{NHSO}_2R^{17}$, and R^{17} is methyl.

[0220] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is F and the other R^8 is $-\text{CH}_2\text{NHSO}_2R^{17}$, and R^{17} is $-(C_{1-2} \text{ alkyl})$.

[0221] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is F and the other R^8 is $-\text{CH}_2\text{NHSO}_2R^{17}$, and R^{17} is methyl.

[0222] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{NR}^{13}(C_{1-6} \text{ alkylene})\text{NR}^{13}R^{14}$.

[0223] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{NR}^{13}(C_{1-5} \text{ alkylene})\text{NR}^{13}R^{14}$.

[0224] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{NR}^{13}(C_{1-4} \text{ alkylene})\text{NR}^{13}R^{14}$.

[0225] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{NR}^{13}(C_{1-3} \text{ alkylene})\text{NR}^{13}R^{14}$.

[0226] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{NR}^{13}\text{CH}_2\text{CH}_2\text{NR}^{13}R^{14}$.

[0227] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, R^8 is halide and the other R^8 is $-\text{NHCH}_2\text{CH}_2\text{NR}^{13}R^{14}$, and R^{13} and R^{14} are independently selected from $-(C_{1-6} \text{ alkyl})$.

[0228] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{NHCH}_2\text{CH}_2\text{NR}^{13}R^{14}$, and R^{13} and R^{14} are independently selected from $-(C_{1-4} \text{ alkyl})$.

[0229] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{NHCH}_2\text{CH}_2\text{NR}^{13}R^{14}$, and R^{13} and R^{14} are independently selected from $-(C_{1-2} \text{ alkyl})$.

[0230] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{NHCH}_2\text{CH}_2\text{NR}^{13}R^{14}$, and R^{13} and R^{14} are both methyl.

[0231] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is F and the other R^8 is $-\text{NHCH}_2\text{CH}_2\text{NR}^{13}\text{R}^{14}$, and R^{13} and R^{14} are independently selected from $-(C_{1-2}\text{ alkyl})$.

[0232] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is F and the other R^8 is $-\text{NHCH}_2\text{CH}_2\text{NR}^{13}\text{R}^{14}$, and R^{13} and R^{14} are both methyl.

[0233] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{OCH}_2\text{CH}_2\text{NR}^{23}\text{R}^{24}$.

[0234] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{OCH}_2\text{CH}_2\text{NR}^{23}\text{R}^{24}$, and R^{23} and R^{24} are independently a $-(C_{1-2}\text{ alkyl})$.

[0235] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{OCH}_2\text{CH}_2\text{NR}^{23}\text{R}^{24}$, and R^{23} and R^{24} are both methyl.

[0236] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is F and the other R^8 is $-\text{OCH}_2\text{CH}_2\text{NR}^{23}\text{R}^{24}$, and R^{23} and R^{24} are both methyl.

[0237] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{CH}_2\text{NHSO}_2\text{R}^{17}$, and R^{17} is $-(C_{1-4}\text{ alkyl})$.

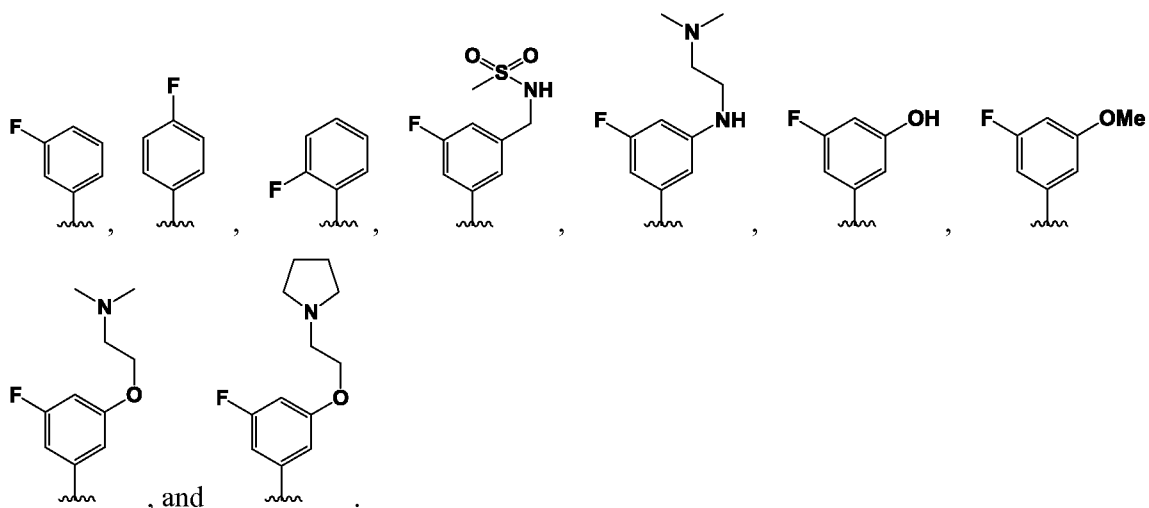
[0238] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{CH}_2\text{NHSO}_2\text{R}^{17}$, and R^{17} is $-(C_{1-2}\text{ alkyl})$.

[0239] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is halide and the other R^8 is $-\text{CH}_2\text{NHSO}_2\text{R}^{17}$, and R^{17} is methyl.

[0240] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is F and the other R^8 is $-\text{CH}_2\text{NHSO}_2\text{R}^{17}$, and R^{17} is $-(C_{1-2}\text{ alkyl})$.

[0241] In some embodiments, R^3 is $-\text{phenyl}(R^8)_k$, k is 2, one R^8 is F and the other R^8 is $-\text{CH}_2\text{NHSO}_2\text{R}^{17}$, and R^{17} is methyl.

[0242] In some embodiments, R^3 is selected from the group consisting of:



[0243] In some embodiments, R^3 is $-\text{piperidinyl}(R^7)_h$.

[0244] In some embodiments, R^3 is $-\text{piperidin-1-yl}(R^7)_h$.

[0245] In some embodiments, R^3 is $-\text{piperidin-1-yl}(R^7)_h$ and h is 0.

[0246] In some embodiments, R^3 is $-\text{piperidin-1-yl}(R^7)_h$, h is 1 or 2, and each R^7 is independently selected from a halide.

[0247] In some embodiments, R^3 is $-\text{piperazinyl}(R^7)_h$.

[0248] In some embodiments, R^3 is $-\text{piperazin-1-yl}(R^7)_h$.

[0249] In some embodiments, R^3 is $-\text{piperazin-1-yl}(R^7)_h$, h is 1, and R^7 is C_{1-3} alkyl.

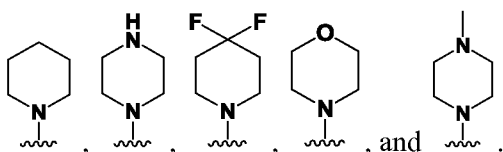
[0250] In some embodiments, R^3 is $-\text{piperazin-1-yl}(R^7)_h$, h is 1, and R^7 is methyl.

[0251] In some embodiments, R^3 is $-\text{morpholinyl}(R^7)_h$.

[0252] In some embodiments, R^3 is $-\text{morpholin-1-yl}(R^7)_h$.

[0253] In some embodiments, R^3 is $-\text{morpholin-1-yl}(R^7)_h$ and h is 0.

[0254] In some embodiments, R^3 is selected from the group consisting of:



[0255] In some embodiments, q is 0.

[0256] In some embodiments, at least one R^4 is a halide.

[0257] In some embodiments, at least one R^4 is a F.

[0258] In some embodiments, R^4 is F.

[0259] In some embodiments, at least one R^4 is $-(C_{1-6}$ alkyl).

[0260] In some embodiments, at least one R^4 is $-(C_{1-5}$ alkyl).

[0261] In some embodiments, at least one R^4 is $-(C_{1-4}$ alkyl).

[0262] In some embodiments, at least one R^4 is $-(C_{1-3}$ alkyl).

[0263] In some embodiments, at least one R^4 is $-(C_{1-2}$ alkyl).

[0264] In some embodiments, R^4 is a methyl.

[0265] In some embodiments, at least one R^4 is $-(C_{1-4}$ alkylene) $_p$ heterocyclyl(R^9) $_h$ and p is 0 or 1.

[0266] In some embodiments, at least one R^4 is $-(C_{1-3}$ alkylene) $_p$ heterocyclyl(R^9) $_h$ and p is 0 or 1.

[0267] In some embodiments, at least one R^4 is $-(C_{1-2}$ alkylene) $_p$ heterocyclyl(R^9) $_h$ and p is 0 or 1.

[0268] In some embodiments, at least one R^4 is $-\text{CH}_2\text{pyrrolidinyl}(R^9)_h$.

[0269] In some embodiments, at least one R^4 is $-\text{CH}_2\text{pyrrolidinyl}(R^9)_h$ and h is 0.

[0270] In some embodiments, R^4 is a $-\text{CH}_2\text{pyrrolidinyl}(R^9)_h$ and h is 0.

[0271] In some embodiments, at least one R⁴ is –CH₂pyrrolidinyl(R⁹)_h, h is 1 or 2, and at least one R⁹ is halide.

[0272] In some embodiments, at least one R⁴ is –CH₂pyrrolidinyl(R⁹)_h, h is 1 or 2, and at least one R⁹ is F.

[0273] In some embodiments, R⁴ is a –CH₂pyrrolidinyl(R⁹)_h, h is 1 or 2, and at least one R⁹ is halide.

[0274] In some embodiments, R⁴ is –CH₂pyrrolidinyl(R⁹)_h, h is 1 or 2, and at least one R⁹ is F.

[0275] In some embodiments, R⁴ is a –CH₂pyrrolidinyl(R⁹)_h, h is 1 or 2, and each R⁹ is F.

[0276] In some embodiments, at least one R⁴ is –CH₂piperidinyl(R⁹)_h.

[0277] In some embodiments, at least one R⁴ is –CH₂piperidinyl(R⁹)_h and h is 0.

[0278] In some embodiments, R⁴ is a –CH₂piperidinyl(R⁹)_h and h is 0.

[0279] In some embodiments, at least one R⁴ is –CH₂piperidinyl(R⁹)_h and at least one R⁹ is halide.

[0280] In some embodiments, at least one R⁴ is –CH₂piperidinyl(R⁹)_h and at least one R⁹ is F.

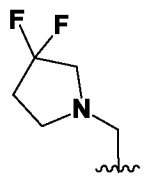
[0281] In some embodiments, at least one R⁴ is –CH₂piperidinyl(R⁹)_h, h is 1 or 2, and at least one R⁹ is halide.

[0282] In some embodiments, at least one R⁴ is –CH₂piperidinyl(R⁹)_h, h is 1 or 2, and at least one R⁹ is F.

[0283] In some embodiments, R⁴ is –CH₂piperidinyl(R⁹)_h, h is 1 or 2, and each R⁹ is a halide.

[0284] In some embodiments, R⁴ is –CH₂piperidinyl(R⁹)_h, h is 1 or 2, and each R⁹ is F.

[0285] In some embodiments, R⁴ is a –CH₂piperidinyl(R⁹)_h, h is 1 or 2, and each R⁹ is F.



[0286] In some embodiments, R⁴ is a

[0287] In some embodiments, at least one R⁴ is –(C₁₋₄ alkylene)_pcarbocyclyl(R¹⁰)_j.

[0288] In some embodiments, at least one R⁴ is –(C₁₋₄ alkylene)_pcarbocyclyl(R¹⁰)_j and j is 0 or 1.

- [0289] In some embodiments, at least one R^4 is $-(C_{1-3} \text{ alkylene})_p \text{ carbocyclyl}(R^{10})_j$ and j is 0 or 1.
- [0290] In some embodiments, at least one R^4 is $-(C_{1-2} \text{ alkylene})_p \text{ carbocyclyl}(R^{10})_j$ and j is 0 or 1.
- [0291] In some embodiments, at least one R^4 is $-\text{CH}_2 \text{ carbocyclyl}(R^{10})_j$.
- [0292] In some embodiments, R^4 is a $-\text{CH}_2 \text{ carbocyclyl}(R^{10})_j$.
- [0293] In some embodiments, at least one R^4 is $-(C_{1-4} \text{ alkylene})_p \text{ aryl}(R^{11})_k$ and k is 0 or 1.
- [0294] In some embodiments, at least one R^4 is $-(C_{1-3} \text{ alkylene})_p \text{ aryl}(R^{11})_k$ and k is 0 or 1.
- [0295] In some embodiments, at least one R^4 is $-(C_{1-2} \text{ alkylene})_p \text{ aryl}(R^{11})_k$ and k is 0 or 1.
- [0296] In some embodiments, at least one R^4 is $-\text{CH}_2 \text{ aryl}(R^{11})_k$.
- [0297] In some embodiments, at least one R^4 is $-\text{CH}_2 \text{ phenyl}(R^{11})_k$.
- [0298] In some embodiments, R^4 is a $-\text{CH}_2 \text{ phenyl}(R^{11})_k$.
- [0299] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$.
- [0300] In some embodiments, R^4 is a $-\text{NHC}(=\text{O})R^{12}$.
- [0301] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-9} \text{ alkyl})$.
- [0302] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-8} \text{ alkyl})$.
- [0303] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-7} \text{ alkyl})$.
- [0304] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-6} \text{ alkyl})$.
- [0305] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-5} \text{ alkyl})$.
- [0306] In some embodiments, R^4 is a $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-5} \text{ alkyl})$.
- [0307] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-4} \text{ alkyl})$.
- [0308] In some embodiments, R^4 is a $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-4} \text{ alkyl})$.
- [0309] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-3} \text{ alkyl})$.
- [0310] In some embodiments, R^4 is a $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-3} \text{ alkyl})$.
- [0311] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{1-2} \text{ alkyl})$.
- [0312] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{2-5} \text{ alkyl})$.
- [0313] In some embodiments, R^4 is a $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{2-5} \text{ alkyl})$.
- [0314] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-(C_{3-4} \text{ alkyl})$.
- [0315] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$ and R^{12} is $-\text{aryl}(R^{19})_k$.
- [0316] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})R^{12}$, R^{12} is $-\text{phenyl}(R^{19})_k$, and k is 0.

- [0317] In some embodiments, R^4 is a $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{phenyl}(\text{R}^{19})_k$, and k is 0.
- [0318] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$ and R^{12} is $-\text{CH}_2\text{aryl}(\text{R}^{19})_k$.
- [0319] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{CH}_2\text{phenyl}(\text{R}^{19})_k$, and k is 0.
- [0320] In some embodiments, R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{CH}_2\text{phenyl}(\text{R}^{19})_k$, and k is 0.
- [0321] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$ and R^{12} is $-\text{heteroaryl}(\text{R}^{18})_q$.
- [0322] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$ and R^{12} is $-\text{carbocyclyl}(\text{R}^{20})_j$.
- [0323] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{carbocyclyl}(\text{R}^{20})_j$, and j is 0.
- [0324] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{cyclopropyl}(\text{R}^{20})_j$, and j is 0.
- [0325] In some embodiments, R^4 is a $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{cyclopropyl}(\text{R}^{20})_j$, and j is 0.
- [0326] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{cyclobutyl}(\text{R}^{20})_j$, and j is 0.
- [0327] In some embodiments, R^4 is a $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{cyclobutyl}(\text{R}^{20})_j$, and j is 0.
- [0328] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{cyclopentyl}(\text{R}^{20})_j$, and j is 0.
- [0329] In some embodiments, R^4 is a $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{cyclopentyl}(\text{R}^{20})_j$, and j is 0.
- [0330] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{cyclohexyl}(\text{R}^{20})_j$, and j is 0.
- [0331] In some embodiments, R^4 is a $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{cyclohexyl}(\text{R}^{20})_j$, and j is 0.
- [0332] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{CH}_2\text{carbocyclyl}(\text{R}^{20})_j$, and j is 0.
- [0333] In some embodiments, at least one R^4 is $-\text{NHC}(=\text{O})\text{R}^{12}$, R^{12} is $-\text{CH}_2\text{cyclopropyl}(\text{R}^{20})_j$, and j is 0.
- [0334] In some embodiments, at least one R^4 is $-\text{NR}^{13}\text{R}^{14}$.

[0335] In some embodiments, at least one R^4 is $-NR^{13}R^{14}$, and R^{13} and R^{14} are independently selected from the group consisting of H and $-(C_{1-6}$ alkyl).

[0336] In some embodiments, at least one R^4 is $-NR^{13}R^{14}$, and R^{13} and R^{14} are independently selected from the group consisting of H and $-(C_{1-5}$ alkyl).

[0337] In some embodiments, at least one R^4 is $-NR^{13}R^{14}$, and R^{13} and R^{14} are independently selected from the group consisting of H and $-(C_{1-4}$ alkyl).

[0338] In some embodiments, at least one R^4 is $-NR^{13}R^{14}$, and R^{13} and R^{14} are independently selected from the group consisting of H and $-(C_{1-3}$ alkyl).

[0339] In some embodiments, at least one R^4 is $-NR^{13}R^{14}$, and R^{13} and R^{14} are independently selected from the group consisting of H and $-(C_{1-2}$ alkyl).

[0340] In some embodiments, at least one R^4 is $-NR^{13}R^{14}$, and R^{13} and R^{14} are independently selected from the group consisting of H and methyl.

[0341] In some embodiments, at least one R^4 is $-NH_2$.

[0342] In some embodiments, R^4 is a $-NH_2$.

[0343] In some embodiments, at least one R^4 is $-NHR^{14}$ and R^{14} is $-(C_{1-4}$ alkyl).

[0344] In some embodiments, at least one R^4 is $-NHR^{14}$ and R^{14} is $-(C_{1-3}$ alkyl).

[0345] In some embodiments, at least one R^4 is $-NHR^{14}$ and R^{14} is $-(C_{1-2}$ alkyl).

[0346] In some embodiments, R^4 is a $-NHR^{14}$ and R^{14} is $-(C_{1-2}$ alkyl).

[0347] In some embodiments, at least one R^4 is $-NHR^{14}$ and R^{14} is $-CH_2aryl(R^{19})_k$.

[0348] In some embodiments, at least one R^4 is $-NHR^{14}$, R^{14} is $-CH_2phenyl(R^{19})_k$, and k is 0.

[0349] In some embodiments, R^4 is $-NHR^{14}$, R^{14} is $-CH_2phenyl(R^{19})_k$, and k is 0.

[0350] In some embodiments, at least one R^4 is $-NHR^{14}$ and R^{14} is $-CH_2carbocyclyl(R^{20})_j$.

[0351] In some embodiments, at least one R^4 is $-NHR^{14}$, R^{14} is $-CH_2cyclopropyl(R^{20})_j$, and j is 0.

[0352] In some embodiments, R^4 is a $-NHR^{14}$, R^{14} is $-CH_2cyclopropyl(R^{20})_j$, and j is 0.

[0353] In some embodiments, at least one R^4 is $-NHR^{14}$, R^{14} is $-CH_2cyclobutyl(R^{20})_j$, and j is 0.

[0354] In some embodiments, R^4 is a $-NHR^{14}$, R^{14} is $-CH_2cyclobutyl(R^{20})_j$, and j is 0.

[0355] In some embodiments, at least one R^4 is $-NHR^{14}$, R^{14} is $-CH_2cyclopentyl(R^{20})_j$, and j is 0.

- [0356] In some embodiments, R^4 is a $-NHR^{14}$, R^{14} is $-CH_2cyclopentyl(R^{20})_j$, and j is 0.
- [0357] In some embodiments, at least one R^4 is $-NHR^{14}$, R^{14} is $-CH_2cyclohexyl(R^{20})_j$, and j is 0.
- [0358] In some embodiments, R^4 is a $-NHR^{14}$, R^{14} is $-CH_2cyclohexyl(R^{20})_j$, and j is 0.
- [0359] In some embodiments, at least one R^4 is $-(C_{1-6} alkylene)NR^{15}R^{16}$.
- [0360] In some embodiments, at least one R^4 is $-(C_{1-5} alkylene)NR^{15}R^{16}$.
- [0361] In some embodiments, at least one R^4 is $-(C_{1-4} alkylene)NR^{15}R^{16}$.
- [0362] In some embodiments, at least one R^4 is $-(C_{1-3} alkylene)NR^{15}R^{16}$.
- [0363] In some embodiments, at least one R^4 is $-(C_{1-2} alkylene)NR^{15}R^{16}$.
- [0364] In some embodiments, at least one R^4 is $-CH_2NR^{15}R^{16}$.
- [0365] In some embodiments, R^4 is a $-CH_2NR^{15}R^{16}$.
- [0366] In some embodiments, at least one R^4 is $-CH_2NR^{15}R^{16}$, and R^{15} and R^{16} are independently selected from the group consisting of H and $-(C_{1-6} alkyl)$.
- [0367] In some embodiments, at least one R^4 is $-CH_2NR^{15}R^{16}$, and R^{15} and R^{16} are independently selected from the group consisting of H and $-(C_{1-5} alkyl)$.
- [0368] In some embodiments, at least one R^4 is $-CH_2NR^{15}R^{16}$, and R^{15} and R^{16} are independently selected from the group consisting of H and $-(C_{1-4} alkyl)$.
- [0369] In some embodiments, at least one R^4 is $-CH_2NR^{15}R^{16}$, and R^{15} and R^{16} are independently selected from the group consisting of H and $-(C_{1-3} alkyl)$.
- [0370] In some embodiments, at least one R^4 is $-CH_2NR^{15}R^{16}$, and R^{15} and R^{16} are independently selected from the group consisting of H and $-(C_{1-2} alkyl)$.
- [0371] In some embodiments, at least one R^4 is $-CH_2NR^{15}R^{16}$, and R^{15} and R^{16} are independently selected from the group consisting of H and methyl.
- [0372] In some embodiments, R^4 is a $-CH_2NR^{15}R^{16}$, and R^{15} and R^{16} are independently selected from the group consisting of H and methyl.
- [0373] In some embodiments, at least one R^4 is $-CH_2NH_2$.
- [0374] In some embodiments, R^4 is a $-CH_2NH_2$.
- [0375] In some embodiments, at least one R^4 is $-CH_2NMe_2$.
- [0376] In some embodiments, R^4 is $-CH_2NMe_2$.
- [0377] In some embodiments, at least one R^4 is $-CH_2NHR^{16}$ and R^{16} is $-(C_{1-4} alkyl)$.
- [0378] In some embodiments, at least one R^4 is $-CH_2NHR^{16}$ and R^{16} is $-(C_{1-3} alkyl)$.
- [0379] In some embodiments, at least one R^4 is $-CH_2NHR^{16}$ and R^{16} is $-(C_{1-2} alkyl)$.

- [0380] In some embodiments, R^4 is a $-\text{CH}_2\text{NHR}^{16}$ and R^{16} is $-(\text{C}_{1-2} \text{ alkyl})$.
- [0381] In some embodiments, at least one R^4 is $-\text{CH}_2\text{NHR}^{16}$ and R^{16} is $-\text{CH}_2\text{aryl}(\text{R}^{19})_k$.
- [0382] In some embodiments, at least one R^4 is $-\text{CH}_2\text{NHR}^{16}$, R^{16} is $-\text{CH}_2\text{phenyl}(\text{R}^{19})_k$, and k is 0.
- [0383] In some embodiments, R^4 is a $-\text{CH}_2\text{NHR}^{16}$, R^{16} is $-\text{CH}_2\text{phenyl}(\text{R}^{19})_k$, and k is 0.
- [0384] In some embodiments, at least one R^4 is $-\text{CH}_2\text{NHR}^{16}$ and R^{16} is $-\text{CH}_2\text{carbocyclyl}(\text{R}^{20})_j$.
- [0385] In some embodiments, at least one R^4 is $-\text{CH}_2\text{NHR}^{16}$, R^{16} is $-\text{CH}_2\text{cyclopropyl}(\text{R}^{20})_j$, and j is 0.
- [0386] In some embodiments, R^4 is a $-\text{CH}_2\text{NHR}^{16}$, R^{16} is $-\text{CH}_2\text{cyclopropyl}(\text{R}^{20})_j$, and j is 0.
- [0387] In some embodiments, at least one R^4 is $-\text{CH}_2\text{NHR}^{16}$, R^{16} is $-\text{CH}_2\text{cyclobutyl}(\text{R}^{20})_j$, and j is 0.
- [0388] In some embodiments, R^4 is a $-\text{CH}_2\text{NHR}^{16}$, R^{16} is $-\text{CH}_2\text{cyclobutyl}(\text{R}^{20})_j$, and j is 0.
- [0389] In some embodiments, at least one R^4 is $-\text{CH}_2\text{NHR}^{16}$, R^{16} is $-\text{CH}_2\text{cyclopentyl}(\text{R}^{20})_j$, and j is 0.
- [0390] In some embodiments, R^4 is a $-\text{CH}_2\text{NHR}^{16}$, R^{16} is $-\text{CH}_2\text{cyclopentyl}(\text{R}^{20})_j$, and j is 0.
- [0391] In some embodiments, at least one R^4 is $-\text{CH}_2\text{NHR}^{16}$, R^{16} is $-\text{CH}_2\text{cyclohexyl}(\text{R}^{20})_j$, and j is 0.
- [0392] In some embodiments, R^4 is a $-\text{CH}_2\text{NHR}^{16}$, R^{16} is $-\text{CH}_2\text{cyclohexyl}(\text{R}^{20})_j$, and j is 0.
- [0393] In some embodiments, at least one R^4 is $-\text{OR}^{22}$.
- [0394] In some embodiments, at least one R^4 is $-\text{OH}$.
- [0395] In some embodiments, R^4 is a $-\text{OH}$.
- [0396] In some embodiments, at least one R^4 is $-\text{OR}^{22}$ and R^{22} is $-(\text{C}_{1-3} \text{ alkyl})$.
- [0397] In some embodiments, at least one R^4 is $-\text{OR}^{22}$ and R^{22} is $-(\text{C}_{1-2} \text{ alkyl})$.
- [0398] In some embodiments, at least one R^4 is $-\text{OMe}$.
- [0399] In some embodiments, R^4 is a $-\text{OMe}$.
- [0400] In some embodiments, at least one R^4 is $-\text{OR}^{22}$, R^{22} is $-\text{heterocyclyl}(\text{R}^{21})_h$, and h is 0.

- [0401] In some embodiments, R^4 is a $-OR^{22}$, R^{22} is $-\text{heterocyclyl}(R^{21})_h$, and h is 0.
- [0402] In some embodiments, at least one R^4 is $-OR^{22}$, R^{22} is $-\text{carbocyclyl}(R^{20})_j$, and j is 0.
- [0403] In some embodiments, R^4 is a $-OR^{22}$, R^{22} is $-\text{carbocyclyl}(R^{20})_j$, and j is 0.
- [0404] In some embodiments, at least one R^4 is $-OR^{22}$, R^{22} is $-(C_{1-4}$ alkylene) $\text{heterocyclyl}(R^{21})_h$, and h is 0.
- [0405] In some embodiments, at least one R^4 is $-OR^{22}$, R^{22} is $-(CH_2CH_2)\text{heterocyclyl}(R^{21})_h$, and h is 0.
- [0406] In some embodiments, R^4 is a $-OR^{22}$, R^{22} is $-(CH_2CH_2)\text{heterocyclyl}(R^{21})_h$, and h is 0.
- [0407] In some embodiments, at least one R^4 is $-OR^{22}$, R^{22} is $-(C_{1-4}$ alkylene) $NR^{23}R^{24}$ and R^{23} and R^{24} are independently a $-(C_{1-4}$ alkyl).
- [0408] In some embodiments, at least one R^4 is $-OR^{22}$, R^{22} is $-(CH_2CH_2)NR^{23}R^{24}$ and R^{23} and R^{24} are independently a $-(C_{1-2}$ alkyl).
- [0409] In some embodiments, at least one R^4 is $-OR^{22}$, and R^{22} is $-(CH_2CH_2)NMe_2$.
- [0410] In some embodiments, R^4 is a $-OR^{22}$, and R^{22} is $-(CH_2CH_2)NMe_2$.
- [0411] In some embodiments, at least one R^4 is $-OR^{22}$, R^{22} is $-(C_{1-4}$ alkylene)aryl(R^{19}) $_k$, k is 0 or 1 and R^{19} is halide.
- [0412] In some embodiments, at least one R^4 is $-OR^{22}$, R^{22} is $-(CH_2CH_2)\text{phenyl}(R^{19})_k$, k is 0 or 1 and R^{19} is a halide.
- [0413] In some embodiments, R^4 is a $-OR^{22}$, R^{22} is $-(CH_2CH_2)\text{phenyl}(R^{19})_k$, k is 0 or 1 and R^{19} is a halide.
- [0414] In some embodiments, at least one R^4 is $-OR^{22}$, R^{22} is $-(CH_2)\text{phenyl}(R^{19})_k$, k is 0 or 1 and R^{19} is a halide.
- [0415] In some embodiments, R^4 is a $-OR^{22}$, R^{22} is $-(CH_2)\text{phenyl}(R^{19})_k$, k is 0 or 1 and R^{19} is a halide.
- [0416] In some embodiments, h is 0.
- [0417] In some embodiments, at least one R^5 is a halide.
- [0418] In some embodiments, at least one R^5 is a F.
- [0419] In some embodiments, at least one R^5 is $-(C_{1-6}$ alkyl).
- [0420] In some embodiments, at least one R^5 is $-(C_{1-5}$ alkyl).
- [0421] In some embodiments, at least one R^5 is $-(C_{1-4}$ alkyl).
- [0422] In some embodiments, at least one R^5 is $-(C_{1-3}$ alkyl).
- [0423] In some embodiments, at least one R^5 is $-(C_{1-2}$ alkyl).

- [0424]** In some embodiments, at least one R⁵ is methyl.
- [0425]** In some embodiments, at least one R⁶ is a halide.
- [0426]** In some embodiments, at least one R⁶ is a F.
- [0427]** In some embodiments, at least one R⁶ is -(C₁₋₄ alkyl).
- [0428]** In some embodiments, at least one R⁶ is -(C₁₋₃ alkyl).
- [0429]** In some embodiments, at least one R⁶ is -(C₁₋₂ alkyl).
- [0430]** In some embodiments, at least one R⁶ is methyl.
- [0431]** In some embodiments, R⁶ is a methyl.
- [0432]** In some embodiments, at least one R⁶ is -C(=O)(C₁₋₃ alkyl).
- [0433]** In some embodiments, at least one R⁶ is -C(=O)Me.
- [0434]** In some embodiments, R⁶ is a -C(=O)Me.
- [0435]** In some embodiments, R² is H; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is -NHC(=O)R¹²; R¹² is -(C₂₋₅ alkyl); R³ is -phenyl(R⁸)_k; k is 1 or 2; and R⁸ is F.
- [0436]** In some embodiments, R² is H; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is -NHC(=O)R¹²; R¹² is -(C₂₋₅ alkyl); R³ is -phenyl(R⁸)_k; k is 2; one R⁸ is F and the other R⁸ is -(C₁₋₂ alkylene)_pNHSO₂R¹⁷; p is 1; and R¹⁷ is -(C₁₋₃ alkyl).
- [0437]** In some embodiments, R² is H; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is -NHC(=O)R¹²; R¹² is -(C₂₋₅ alkyl); R³ is -phenyl(R⁸)_k; k is 2; one R⁸ is F and the other R⁸ is -NH(C₁₋₆ alkylene)NR¹³R¹⁴; and R¹³ and R¹⁴ are independently selected from -(C₁₋₃ alkyl).
- [0438]** In some embodiments, R² is H; R¹ is -pyridin-3-yl(R⁴)_q, wherein q is 1; R⁴ is -NHC(=O)R¹²; R¹² is -(C₂₋₅ alkyl); R³ is -heteroaryl(R⁶)_q, wherein q is 1; R⁶ is selected from the group consisting of halide, -(C₁₋₂ alkyl), and -C(=O)R¹⁷; R¹⁷ is -(C₁₋₃ alkyl); and the heteroaryl is selected from the group consisting of pyridine, furan, thiophene, and imidazole.
- [0439]** In some embodiments, R² is H; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is -NHC(=O)R¹²; R¹² is -(C₂₋₅ alkyl); R³ is -heterocyclyl(R⁷)_h; h is 1 or 2; and R⁷ is selected from the group consisting of halide and -(C₁₋₂ alkyl).
- [0440]** In some embodiments, R² is H; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is -NHC(=O)R¹²; R¹² is -carbocyclyl(R²⁰)_j; j is 0; R³ is -phenyl(R⁸)_k; k is 1 or 2; R⁸ is F; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.
- [0441]** In some embodiments, R² is H; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is -NHC(=O)R¹²; R¹² is -carbocyclyl(R²⁰)_j; j is 0; R³ is -phenyl(R⁸)_k; k is 2; one R⁸ is F and the other R⁸ is -(C₁₋₂ alkylene)_pNHSO₂R¹⁷; p is 1; R¹⁷ is -(C₁₋₃ alkyl); and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0442] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-\text{carbocyclyl}(R^{20})_j$; j is 0; R^3 is $-\text{phenyl}(R^8)_k$; k is 2; one R^8 is F and the other R^8 is $-\text{NH}(\text{C}_{1-6} \text{ alkylene})\text{NR}^{13}\text{R}^{14}$; R^{13} and R^{14} are independently selected from $-(\text{C}_{1-3} \text{ alkyl})$; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0443] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$, wherein q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-\text{carbocyclyl}(R^{20})_j$; j is 0; R^3 is $-\text{heteroaryl}(R^6)_q$, wherein q is 1; R^6 is selected from the group consisting of halide, $-(\text{C}_{1-2} \text{ alkyl})$, and $-\text{C}(=\text{O})R^{17}$; R^{17} is C_{1-3} alkyl; the heteroaryl is selected from the group consisting of pyridine, furan, thiophene, and imidazole; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0444] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-\text{carbocyclyl}(R^{20})_j$; j is 0; R^3 is $-\text{heterocyclyl}(R^7)_h$; h is 1 or 2; R^7 is selected from the group consisting of halide and $-(\text{C}_{1-2} \text{ alkyl})$; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0445] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is selected from the group consisting of $-\text{NR}^{13}\text{R}^{14}$ and $-\text{CH}_2\text{NR}^{15}\text{R}^{16}$; R^{13} and R^{15} are independently selected from the group consisting of H and $-(\text{C}_{1-3} \text{ alkyl})$; R^{14} and R^{16} are independently selected from the group consisting of H, $-(\text{C}_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$, wherein j and k are 0; R^3 is $-\text{phenyl}(R^8)_k$, wherein k is 1 or 2; R^8 is F; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0446] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is selected from the group consisting of $-\text{NR}^{13}\text{R}^{14}$ and $-\text{CH}_2\text{NR}^{15}\text{R}^{16}$; R^{13} and R^{15} are independently selected from the group consisting of H and $-(\text{C}_{1-3} \text{ alkyl})$; R^{14} and R^{16} are independently selected from the group consisting of H, $-(\text{C}_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$, wherein j and k are 0; R^3 is $-\text{phenyl}(R^8)_k$, wherein k is 2; one R^8 is F and the other R^8 is $-(\text{C}_{1-2} \text{ alkylene})_p\text{NHSO}_2\text{R}^{17}$; p is 1; R^{17} is $-(\text{C}_{1-3} \text{ alkyl})$; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0447] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is selected from the group consisting of $-\text{NR}^{13}\text{R}^{14}$ and $-\text{CH}_2\text{NR}^{15}\text{R}^{16}$, wherein R^{13} and R^{15} are independently selected from the group consisting of H and $-(\text{C}_{1-3} \text{ alkyl})$, and R^{14} and R^{16} are independently selected from the group consisting of H, $-(\text{C}_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$, wherein k and j are 0; R^3 is $-\text{phenyl}(R^8)_k$, wherein k is 2; one R^8 is F and the other R^8 is $-\text{NH}(\text{C}_{1-6} \text{ alkylene})\text{NR}^{13}\text{R}^{14}$, wherein R^{13} and R^{14} are independently selected from $-(\text{C}_{1-3} \text{ alkyl})$; and the

carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0448] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$, wherein q is 1; R^4 is selected from the group consisting of $-\text{NR}^{13}\text{R}^{14}$ and $-\text{CH}_2\text{NR}^{15}\text{R}^{16}$; R^{13} and R^{15} are independently selected from the group consisting of H and $-(\text{C}_{1-3} \text{ alkyl})$; R^{14} and R^{16} are independently selected from the group consisting of H, $-(\text{C}_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$; k and j are 0; R^3 is $-\text{heteroaryl}(R^6)_q$, wherein q is 1; R^6 is selected from the group consisting of halide, $-(\text{C}_{1-2} \text{ alkyl})$, and $-\text{C}(=\text{O})\text{R}^{17}$; R^{17} is $-(\text{C}_{1-3} \text{ alkyl})$; the heteroaryl is selected from the group consisting of pyridine, furan, thiophene, and imidazole; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0449] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is selected from the group consisting of $-\text{NR}^{13}\text{R}^{14}$ and $-\text{CH}_2\text{NR}^{15}\text{R}^{16}$; R^{13} and R^{15} are independently selected from the group consisting of H and $-(\text{C}_{1-3} \text{ alkyl})$; R^{14} and R^{16} are independently selected from the group consisting of H, $-(\text{C}_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$; k and j are 0; R^3 is $-\text{heterocyclyl}(R^7)_h$; h is 1 or 2; R^7 is selected from the group consisting of halide and $-(\text{C}_{1-2} \text{ alkyl})$; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0450] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{CH}_2\text{heterocyclyl}(R^9)_h$; h is 0–2; R^9 is F; R^3 is $-\text{phenyl}(R^8)_k$; k is 1 or 2; R^8 is F; and the heterocyclyl is selected from the group consisting of pyrrolidine and piperidine.

[0451] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{CH}_2\text{heterocyclyl}(R^9)_h$; h is 0–2; R^9 is F; R^3 is $-\text{phenyl}(R^8)_k$; k is 2; one R^8 is F and the other R^8 is $-(\text{C}_{1-2} \text{ alkylene})_p\text{NH}\text{SO}_2\text{R}^{17}$; p is 1; R^{17} is $-(\text{C}_{1-3} \text{ alkyl})$; and the heterocyclyl is selected from the group consisting of pyrrolidine and piperidine.

[0452] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{CH}_2\text{heterocyclyl}(R^9)_h$; h is 0–2; R^9 is F; R^3 is $-\text{phenyl}(R^8)_k$; k is 2; and R^8 is one F and the other R^8 is $-\text{NH}(\text{C}_{1-6} \text{ alkylene})\text{NR}^{13}\text{R}^{14}$; R^{13} and R^{14} are independently selected from $-(\text{C}_{1-3} \text{ alkyl})$; and the heterocyclyl is selected from the group consisting of pyrrolidine and piperidine.

[0453] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$, wherein q is 1; R^4 is $-\text{CH}_2\text{heterocyclyl}(R^9)_h$; h is 0–2; R^9 is F; R^3 is $-\text{heteroaryl}(R^6)_q$, wherein q is 1; R^6 is selected from the group consisting of halide, $-(\text{C}_{1-2} \text{ alkyl})$, and $-\text{C}(=\text{O})\text{R}^{17}$; R^{17} is $-(\text{C}_{1-3} \text{ alkyl})$; the heteroaryl is selected from the group consisting of pyridine, furan, thiophene, and imidazole; and the heterocyclyl is selected from the group consisting of pyrrolidine and piperidine.

[0454] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{CH}_2\text{heterocyclyl}(R^9)_h$, wherein h is 0–2; R^9 is F; R^3 is $-\text{heterocyclyl}(R^7)_h$, wherein h is 1 or 2; R^7 is selected from the group consisting of halide and $-(C_{1-2} \text{ alkyl})$; and the heterocyclyl is selected from the group consisting of pyrrolidine and piperidine.

[0455] In some embodiments, R^2 is H; R^1 is $-\text{pyrimidinyl}(R^4)_q$; q is 0; R^3 is $-\text{phenyl}(R^8)_k$; k is 1 or 2; and R^8 is F.

[0456] In some embodiments, R^2 is H; R^1 is $-\text{pyrimidinyl}(R^4)_q$; q is 0; R^3 is $-\text{phenyl}(R^8)_k$; k is 2; one R^8 is F and the other R^8 is $-(C_{1-2} \text{ alkylene})_p\text{NHSO}_2R^{17}$; p is 1; and R^{17} is $-(C_{1-3} \text{ alkyl})$.

[0457] In some embodiments, R^2 is H; R^1 is $-\text{pyrimidinyl}(R^4)_q$; q is 0; R^3 is $-\text{phenyl}(R^8)_k$; k is 2; one R^8 is F and the other R^8 is $-\text{NH}(C_{1-6} \text{ alkylene})\text{NR}^{13}R^{14}$; and R^{13} and R^{14} are independently selected from $-(C_{1-3} \text{ alkyl})$.

[0458] In some embodiments, R^2 is H; R^1 is $-\text{pyrimidinyl}(R^4)_q$, wherein q is 0; R^3 is $-\text{heteroaryl}(R^6)_q$, wherein q is 1; R^6 is selected from the group consisting of halide, $-(C_{1-2} \text{ alkyl})$, and $-\text{C}(=\text{O})R^{17}$; R^{17} is $-(C_{1-3} \text{ alkyl})$; and the heteroaryl is selected from the group consisting of pyridine, furan, thiophene, and imidazole.

[0459] In some embodiments, R^2 is H; R^1 is $-\text{pyrimidinyl}(R^4)_q$; q is 0; R^3 is $-\text{heterocyclyl}(R^7)_h$; h is 1 or 2; R^7 is selected from the group consisting of halide and $-(C_{1-2} \text{ alkyl})$.

[0460] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-(C_{2-5} \text{ alkyl})$; R^3 is $-\text{phenyl}(R^8)_k$; k is 1 or 2; and R^8 is F.

[0461] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-(C_{2-5} \text{ alkyl})$; R^3 is $-\text{phenyl}(R^8)_k$; k is 2; one R^8 is F and the other R^8 is $-(C_{1-2} \text{ alkylene})_p\text{NHSO}_2R^{17}$; p is 1; and R^{17} is $-(C_{1-3} \text{ alkyl})$.

[0462] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-(C_{2-5} \text{ alkyl})$; R^3 is $-\text{phenyl}(R^8)_k$; k is 2; one R^8 is F and the other R^8 is $-\text{NH}(C_{1-6} \text{ alkylene})\text{NR}^{13}R^{14}$; and R^{13} and R^{14} are independently selected from $-(C_{1-3} \text{ alkyl})$.

[0463] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$, wherein q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-(C_{2-5} \text{ alkyl})$; R^3 is $-\text{heteroaryl}(R^6)_q$, wherein q is 1; R^6 is selected from the group consisting of halide, $-(C_{1-2} \text{ alkyl})$, and $-\text{C}(=\text{O})R^{17}$; R^{17} is $-(C_{1-3} \text{ alkyl})$; and the heteroaryl is selected from the group consisting of pyridine, furan, thiophene, and imidazole.

[0464] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-(C_{2-5} \text{ alkyl})$; R^3 is $-\text{heterocyclyl}(R^7)_h$; h is 1 or 2; and R^7 is selected from the group consisting of halide and $-(C_{1-2} \text{ alkyl})$.

[0465] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-\text{carbocyclyl}(R^{20})_j$; j is 0; R^3 is $-\text{phenyl}(R^8)_k$; k is 1 or 2; R^8 is F; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0466] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-\text{carbocyclyl}(R^{20})_j$; j is 0; R^3 is $-\text{phenyl}(R^8)_k$; k is 2; one R^8 is F and the other R^8 is $-(C_{1-2} \text{ alkylene})_p\text{NHSO}_2R^{17}$; p is 1; R^{17} is $-(C_{1-3} \text{ alkyl})$; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0467] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-\text{carbocyclyl}(R^{20})_j$; j is 0; R^3 is $-\text{phenyl}(R^8)_k$; k is 2; one R^8 is F and the other R^8 is $-\text{NH}(C_{1-6} \text{ alkylene})\text{NR}^{13}\text{R}^{14}$; R^{13} and R^{14} are independently selected from $-(C_{1-3} \text{ alkyl})$; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0468] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$, wherein q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-\text{carbocyclyl}(R^{20})_j$; j is 0; R^3 is $-\text{heteroaryl}(R^6)_q$, wherein q is 1; R^6 is selected from the group consisting of halide, $-(C_{1-2} \text{ alkyl})$, and $-\text{C}(=\text{O})R^{17}$; R^{17} is $C_{1-3} \text{ alkyl}$; the heteroaryl is selected from the group consisting of pyridine, furan, thiophene, and imidazole; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0469] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is $-\text{NHC}(=\text{O})R^{12}$; R^{12} is $-\text{carbocyclyl}(R^{20})_j$; j is 0; R^3 is $-\text{heterocyclyl}(R^7)_h$; h is 1 or 2; R^7 is selected from the group consisting of halide and $-(C_{1-2} \text{ alkyl})$; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0470] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is selected from the group consisting of $-\text{NR}^{13}\text{R}^{14}$ and $-\text{CH}_2\text{NR}^{15}\text{R}^{16}$; R^{13} and R^{15} are independently selected from the group consisting of H and $-(C_{1-3} \text{ alkyl})$; R^{14} and R^{16} are independently selected from the group consisting of H, $-(C_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$, wherein k and j are 0; R^3 is $-\text{phenyl}(R^8)_k$, wherein k is 1 or 2; R^8 is F; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0471] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(R^4)_q$; q is 1; R^4 is selected from the group consisting of $-\text{NR}^{13}\text{R}^{14}$ and $-\text{CH}_2\text{NR}^{15}\text{R}^{16}$; R^{13} and R^{15} are independently selected from the group consisting of H and $-(C_{1-3} \text{ alkyl})$; R^{14} and R^{16} are independently selected from the group consisting of H, $-(C_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$, wherein k and j are 0; R^3 is $-\text{phenyl}(R^8)_k$, wherein k is 2; one R^8 is F and the other R^8 is $-(C_{1-2} \text{ alkylene})_p\text{NHSO}_2R^{17}$;

p is 1; R¹⁷ is -(C₁₋₃ alkyl); and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0472] In some embodiments, R² is F; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is selected from the group consisting of -NR¹³R¹⁴ and -CH₂NR¹⁵R¹⁶, wherein R¹³ and R¹⁵ are independently selected from the group consisting of H and -(C₁₋₃ alkyl); R¹⁴ and R¹⁶ are independently selected from the group consisting of H, -(C₁₋₃ alkyl), -CH₂phenyl(R¹⁹)_k, and -CH₂carbocyclyl(R²⁰)_j, wherein k and j are 0; R³ is -phenyl(R⁸)_k, wherein k is 2; one R⁸ is F and the other R⁸ is -NH(C₁₋₆ alkylene)NR¹³R¹⁴, wherein R¹³ and R¹⁴ are independently selected from -(C₁₋₃ alkyl); and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0473] In some embodiments, R² is F; R¹ is -pyridin-3-yl(R⁴)_q, wherein q is 1; R⁴ is selected from the group consisting of -NR¹³R¹⁴ and -CH₂NR¹⁵R¹⁶; R¹³ and R¹⁵ are independently selected from the group consisting of H and -(C₁₋₃ alkyl); R¹⁴ and R¹⁶ are independently selected from the group consisting of H, -(C₁₋₃ alkyl), -CH₂phenyl(R¹⁹)_k, and -CH₂carbocyclyl(R²⁰)_j; k and j are 0; R³ is -heteroaryl(R⁶)_q, wherein q is 1; R⁶ is selected from the group consisting of halide, -(C₁₋₂ alkyl), and -C(=O)R¹⁷; R¹⁷ is -(C₁₋₃ alkyl); the heteroaryl is selected from the group consisting of pyridine, furan, thiophene, and imidazole; and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0474] In some embodiments, R² is F; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is selected from the group consisting of -NR¹³R¹⁴ and -CH₂NR¹⁵R¹⁶; R¹³ and R¹⁵ are independently selected from the group consisting of H and -(C₁₋₃ alkyl); R¹⁴ and R¹⁶ are independently selected from the group consisting of H, -(C₁₋₃ alkyl), -CH₂phenyl(R¹⁹)_k, and -CH₂carbocyclyl(R²⁰)_j; k and j are 0; R³ is -heterocyclyl(R⁷)_h; h is 1 or 2; R⁷ is selected from the group consisting of halide and -(C₁₋₂ alkyl); and the carbocyclyl is selected from the group consisting of cyclopropyl, cyclobutyl, cyclopentyl, and cyclohexyl.

[0475] In some embodiments, R² is F; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is -CH₂heterocyclyl(R⁹)_h; h is 0-2; R⁹ is F; R³ is -phenyl(R⁸)_k; k is 1 or 2; R⁸ is F; and the heterocyclyl is selected from the group consisting of pyrrolidine and piperidine.

[0476] In some embodiments, R² is F; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is -CH₂heterocyclyl(R⁹)_h; h is 0-2; R⁹ is F; R³ is -phenyl(R⁸)_k; k is 2; one R⁸ is F and the other R⁸ is -(C₁₋₂ alkylene)_pNHSO₂R¹⁷; p is 1; R¹⁷ is -(C₁₋₃ alkyl); and the heterocyclyl is selected from the group consisting of pyrrolidine and piperidine.

[0477] In some embodiments, R² is F; R¹ is -pyridin-3-yl(R⁴)_q; q is 1; R⁴ is -CH₂heterocyclyl(R⁹)_h; h is 0-2; R⁹ is F; R³ is -phenyl(R⁸)_k; k is 2; and R⁸ is one F and the other R⁸

is $-\text{NH}(\text{C}_{1-6} \text{ alkylene})\text{NR}^{13}\text{R}^{14}$; R^{13} and R^{14} are independently selected from $-(\text{C}_{1-3} \text{ alkyl})$; and the heterocyclyl is selected from the group consisting of pyrrolidine and piperidine.

[0478] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(\text{R}^4)_q$, wherein q is 1; R^4 is $-\text{CH}_2\text{heterocyclyl}(\text{R}^9)_h$; h is 0–2; R^9 is F; R^3 is $-\text{heteroaryl}(\text{R}^6)_q$, wherein q is 1; R^6 is selected from the group consisting of halide, $-(\text{C}_{1-2} \text{ alkyl})$, and $-\text{C}(=\text{O})\text{R}^{17}$; R^{17} is $-(\text{C}_{1-3} \text{ alkyl})$; the heteroaryl is selected from the group consisting of pyridine, furan, thiophene, and imidazole; and the heterocyclyl is selected from the group consisting of pyrrolidine and piperidine.

[0479] In some embodiments, R^2 is F; R^1 is $-\text{pyridin-3-yl}(\text{R}^4)_q$; q is 1; R^4 is $-\text{CH}_2\text{heterocyclyl}(\text{R}^9)_h$, wherein h is 0–2; R^9 is F; R^3 is $-\text{heterocyclyl}(\text{R}^7)_h$, wherein h is 1 or 2; R^7 is selected from the group consisting of halide and $-(\text{C}_{1-2} \text{ alkyl})$; and the heterocyclyl is selected from the group consisting of pyrrolidine and piperidine.

[0480] In some embodiments, R^2 is F; R^1 is $-\text{pyrimidinyl}(\text{R}^4)_q$; q is 0; R^3 is $-\text{phenyl}(\text{R}^8)_k$; k is 1 or 2; and R^8 is F.

[0481] In some embodiments, R^2 is F; R^1 is $-\text{pyrimidinyl}(\text{R}^4)_q$; q is 0; R^3 is $-\text{phenyl}(\text{R}^8)_k$; k is 2; one R^8 is F and the other R^8 is $-(\text{C}_{1-2} \text{ alkylene})_p\text{NHSO}_2\text{R}^{17}$; p is 1; and R^{17} is $-(\text{C}_{1-3} \text{ alkyl})$.

[0482] In some embodiments, R^2 is F; R^1 is $-\text{pyrimidinyl}(\text{R}^4)_q$; q is 0; R^3 is $-\text{phenyl}(\text{R}^8)_k$; k is 2; one R^8 is F and the other R^8 is $-\text{NH}(\text{C}_{1-6} \text{ alkylene})\text{NR}^{13}\text{R}^{14}$; and R^{13} and R^{14} are independently selected from $-(\text{C}_{1-3} \text{ alkyl})$.

[0483] In some embodiments, R^2 is F; R^1 is $-\text{pyrimidinyl}(\text{R}^4)_q$, wherein q is 0; R^3 is $-\text{heteroaryl}(\text{R}^6)_q$, wherein q is 1; R^6 is selected from the group consisting of halide, $-(\text{C}_{1-2} \text{ alkyl})$, and $-\text{C}(=\text{O})\text{R}^{17}$; R^{17} is $-(\text{C}_{1-3} \text{ alkyl})$; and the heteroaryl is selected from the group consisting of pyridine, furan, thiophene, and imidazole.

[0484] In some embodiments, R^2 is F; R^1 is $-\text{pyrimidinyl}(\text{R}^4)_q$; q is 0; R^3 is $-\text{heterocyclyl}(\text{R}^7)_h$; h is 1 or 2; R^7 is selected from the group consisting of halide and $-(\text{C}_{1-2} \text{ alkyl})$.

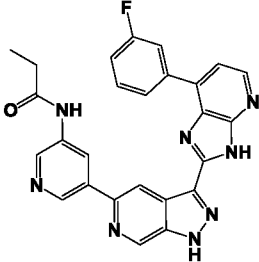
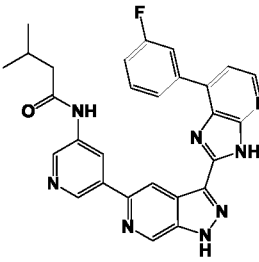
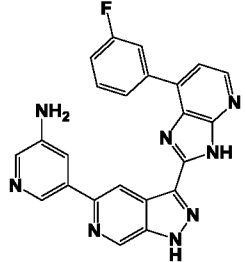
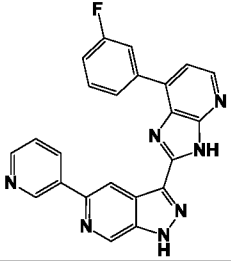
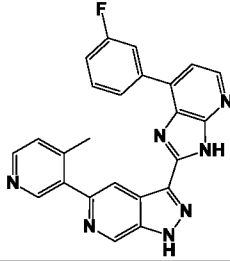
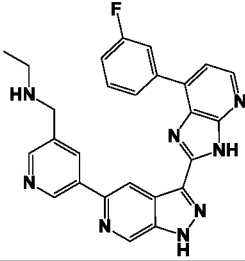
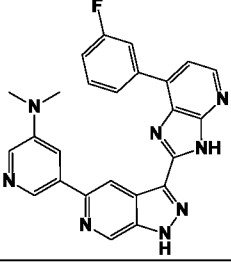
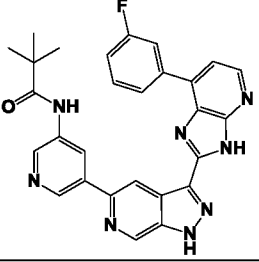
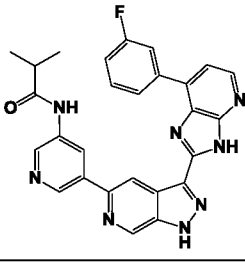
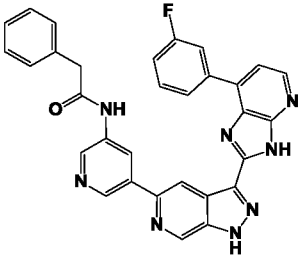
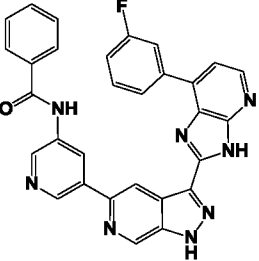
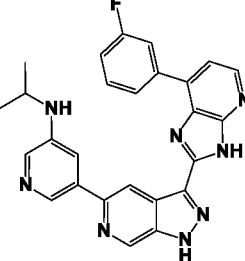
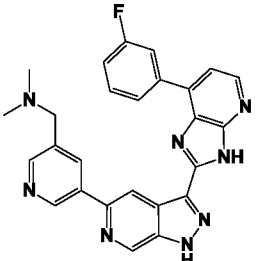
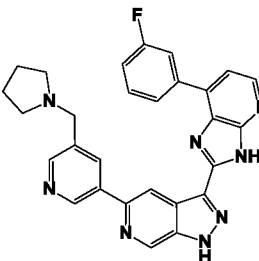
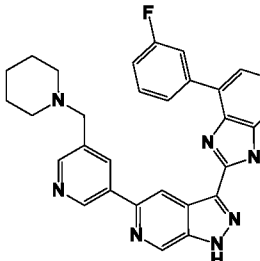
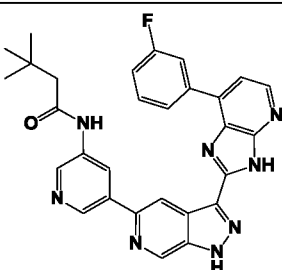
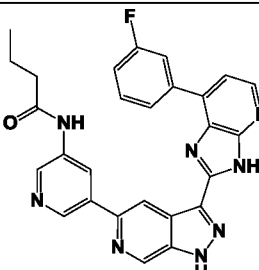
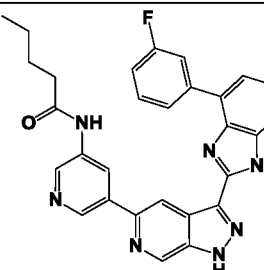
[0485] In some embodiments, R^2 is H; R^1 is $-\text{pyrazol-4-yl}(\text{R}^4)_q$; q is 0 or 1; R^4 is $-(\text{C}_{1-3} \text{ alkyl})$; R^3 is $-\text{phenyl}(\text{R}^8)_k$; k is 1 or 2; and R^8 is F.

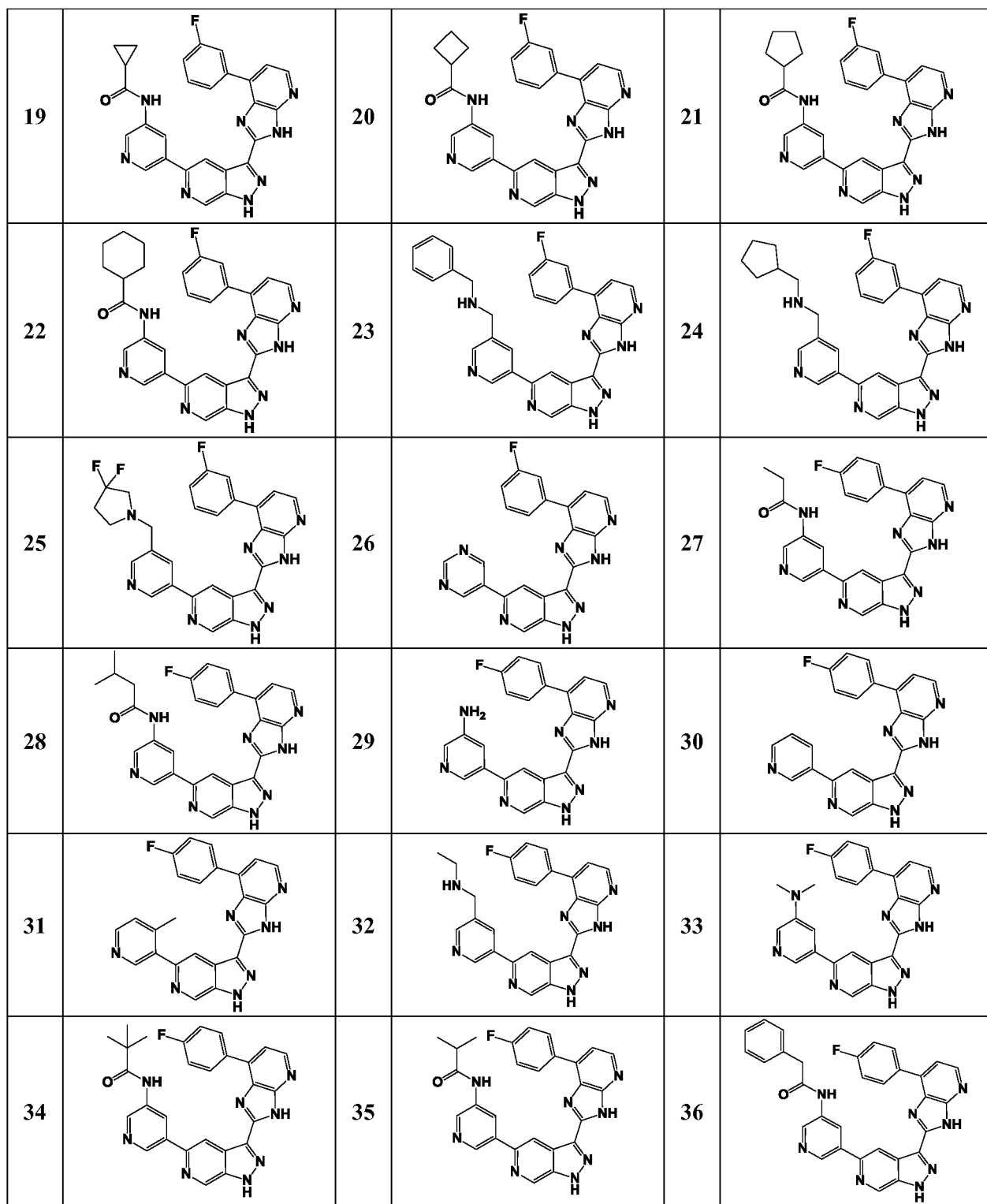
[0486] In some embodiments, R^2 is H; R^1 is $-\text{imidazol-5-yl}(\text{R}^4)_q$; q is 1 or 2; each R^4 is independently selected from $-(\text{C}_{1-3} \text{ alkyl})$; R^3 is $-\text{phenyl}(\text{R}^8)_k$; k is 1 or 2; and R^8 is F.

[0487] In some embodiments, R^2 is H; R^1 is $-\text{pyridin-3-yl}(\text{R}^4)_q$; q is 1; R^4 is $-\text{OR}^{22}$; R^{22} is selected from the group consisting of H and $-(\text{C}_{1-3} \text{ alkyl})$; R^3 is $-\text{phenyl}(\text{R}^8)_k$; k is 1 or 2; and R^8 is F.

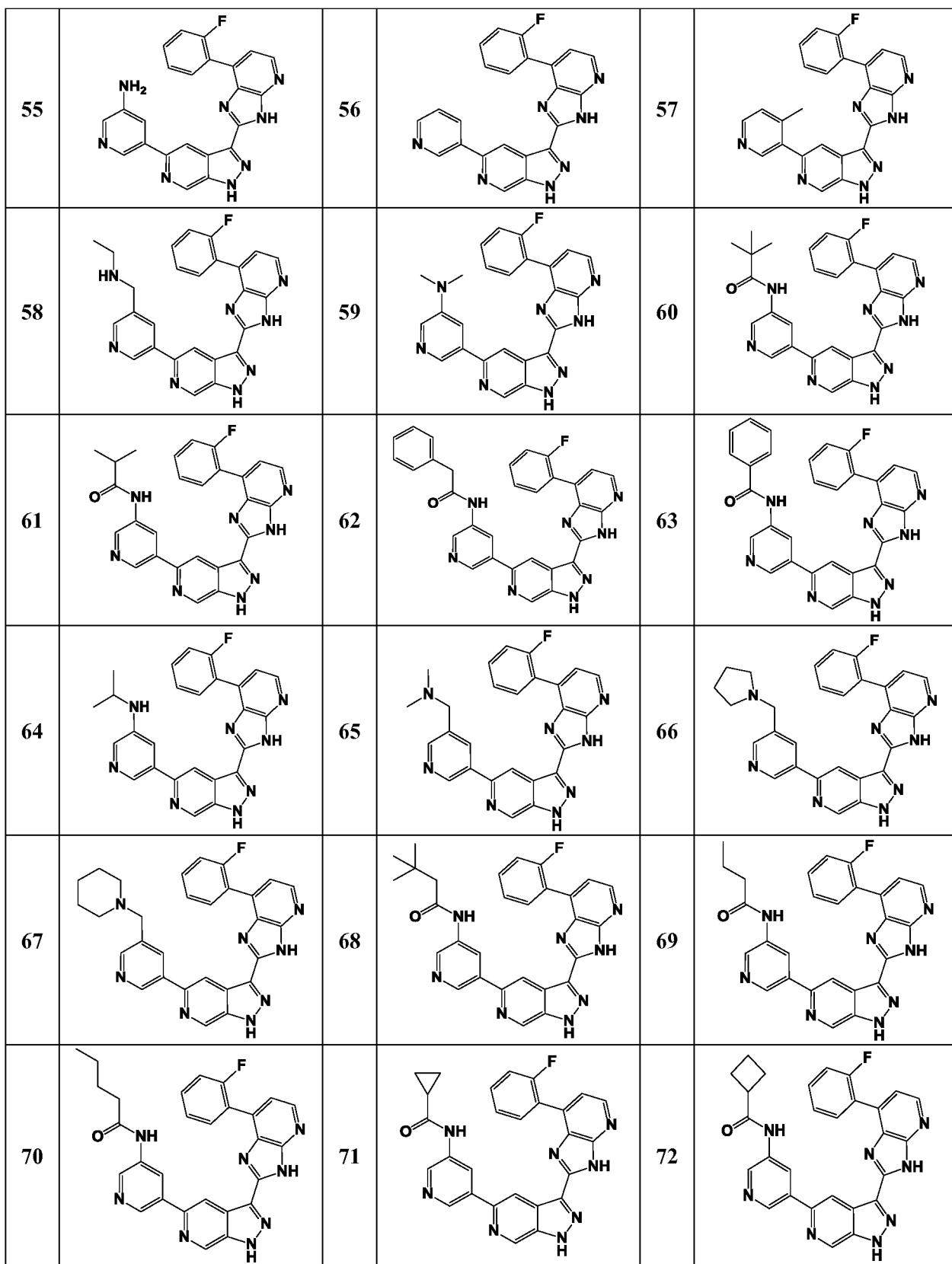
[0488] Illustrative compounds of Formula (I) are shown in Table 1.

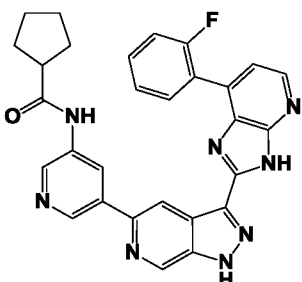
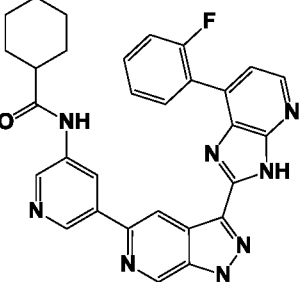
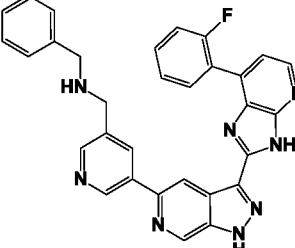
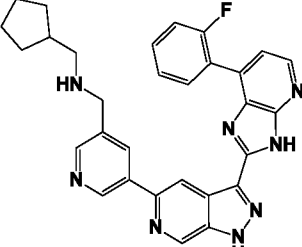
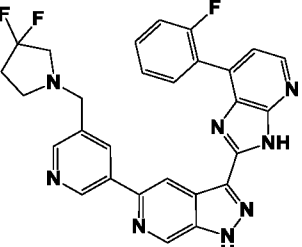
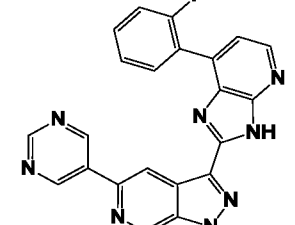
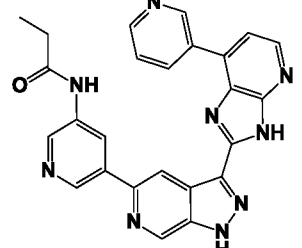
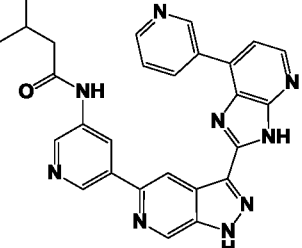
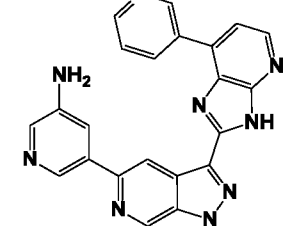
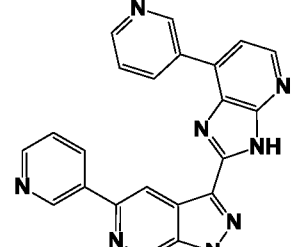
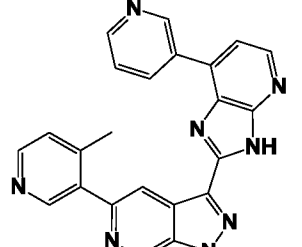
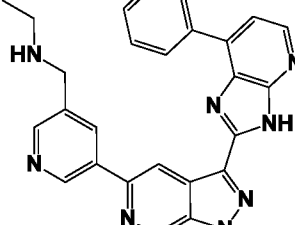
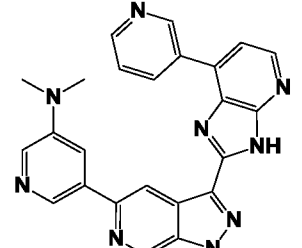
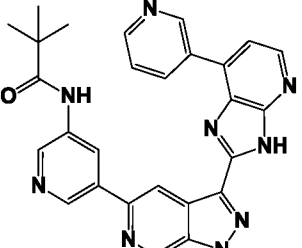
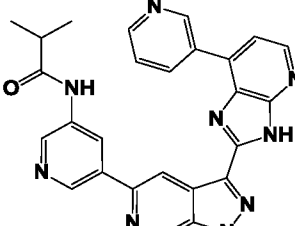
Table 1.

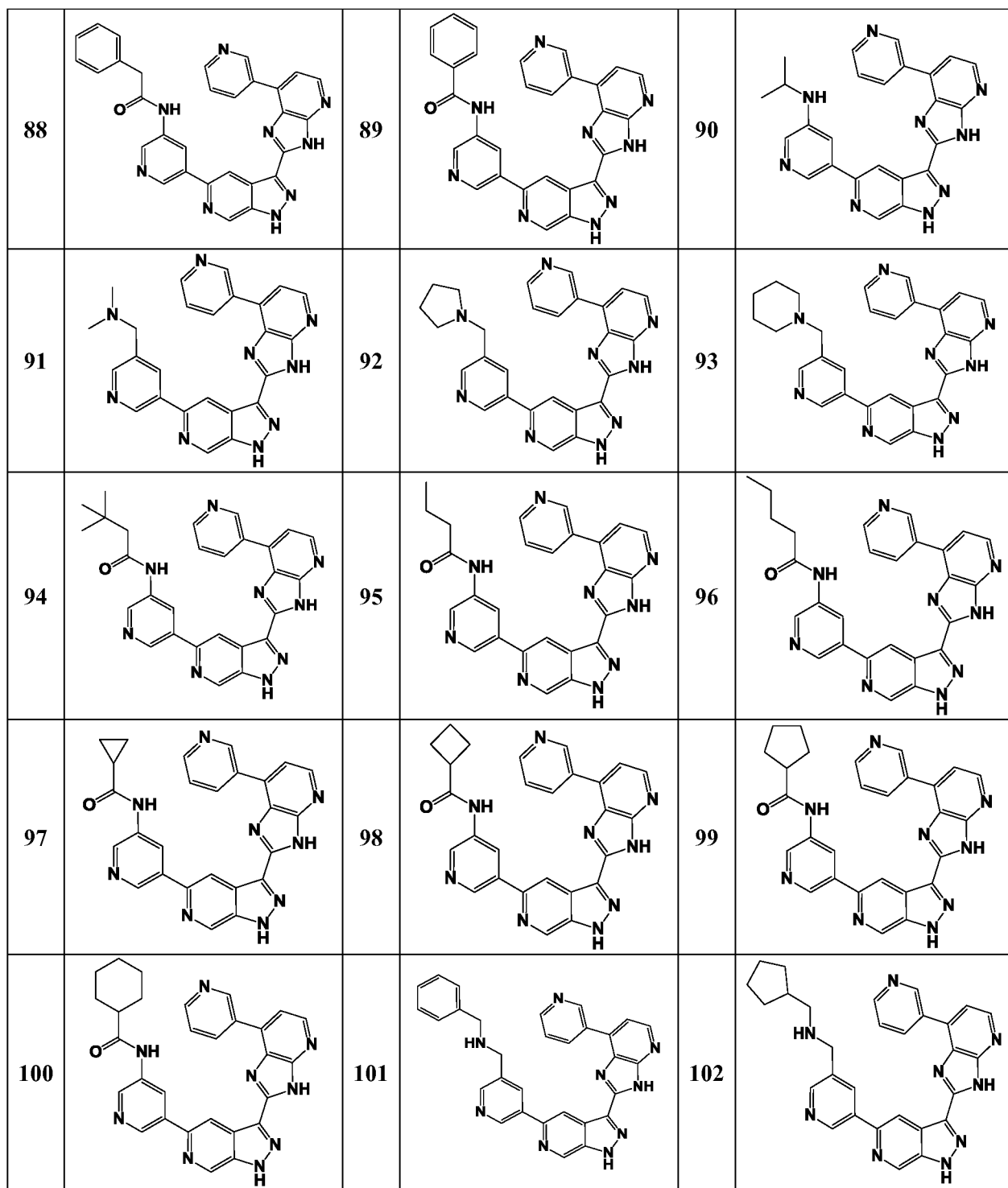
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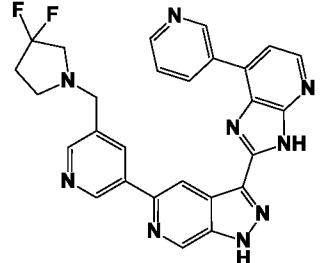
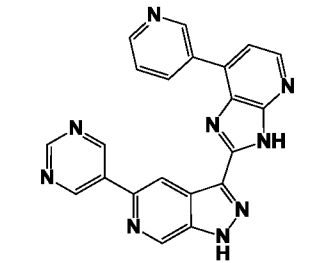
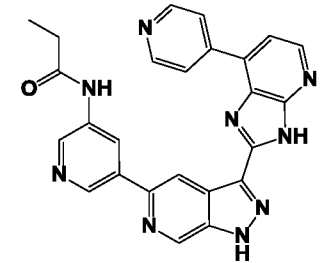
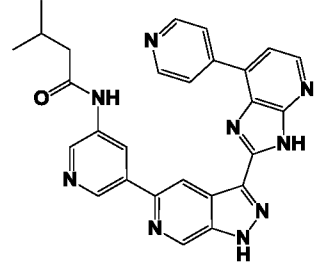
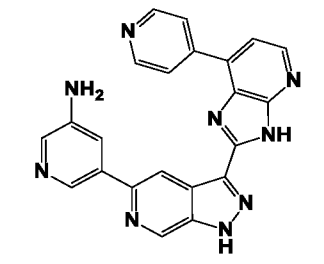
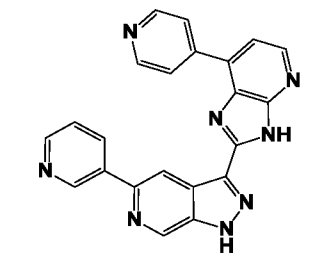
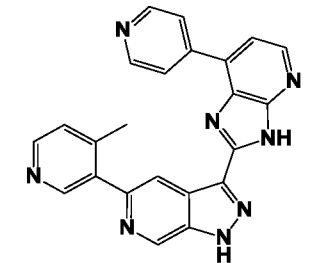
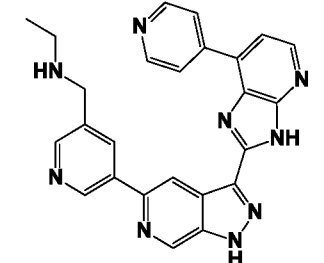
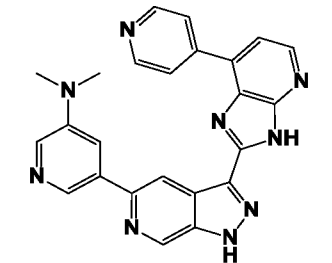
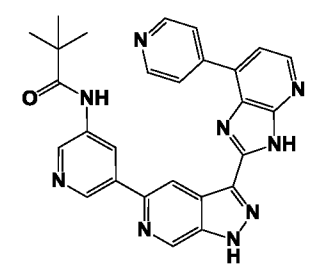
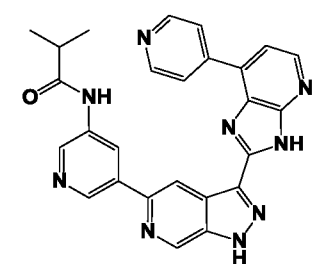
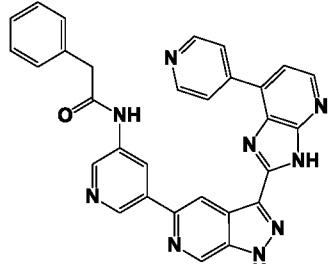
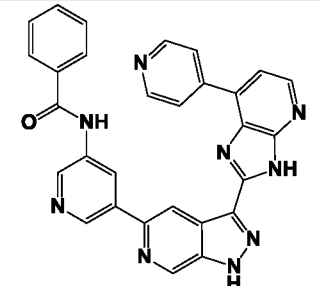
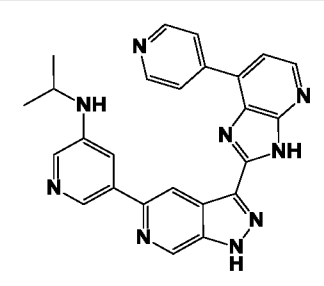
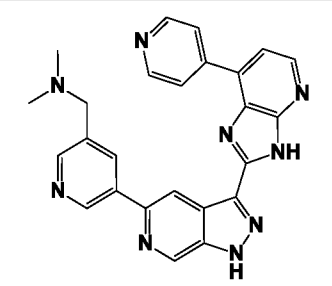
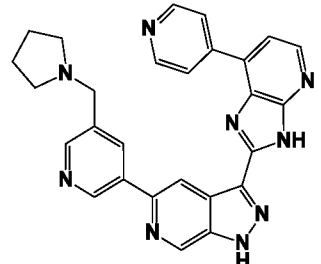
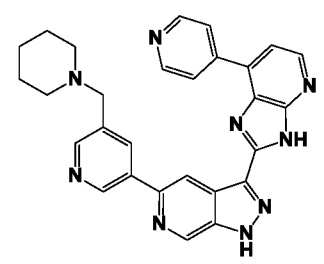
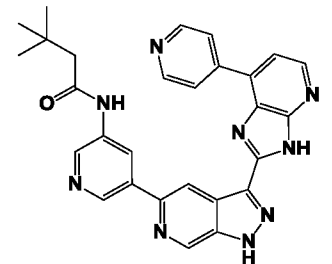


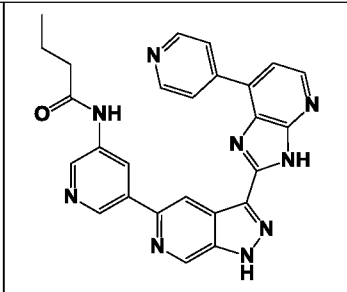
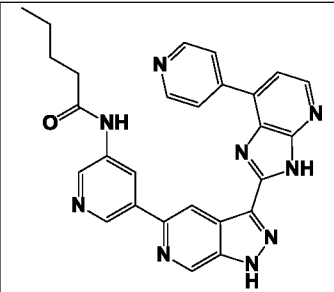
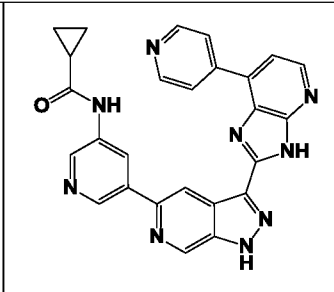
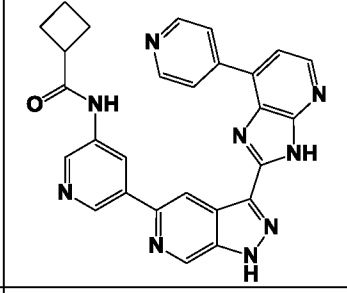
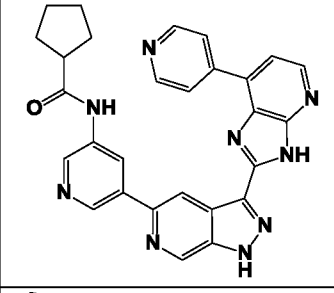
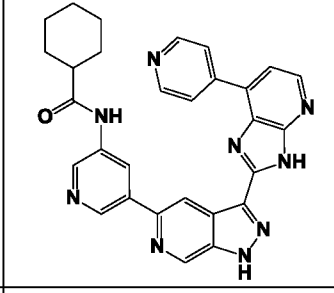
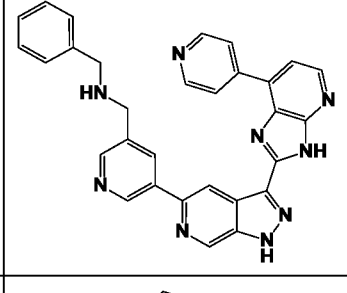
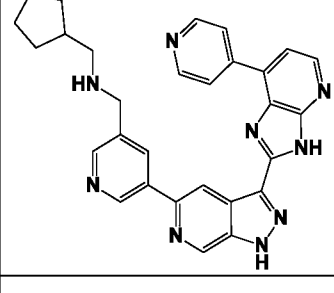
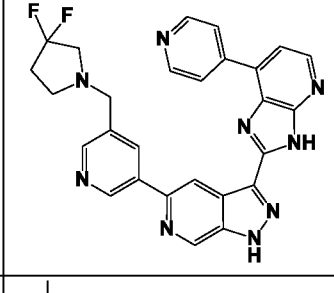
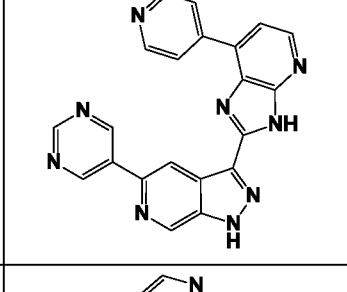
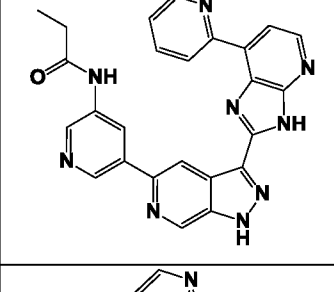
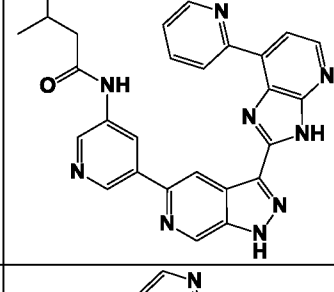
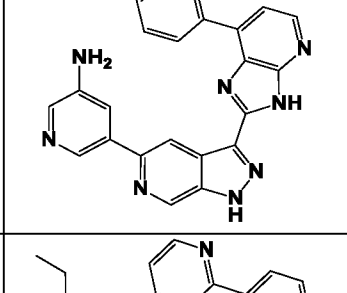
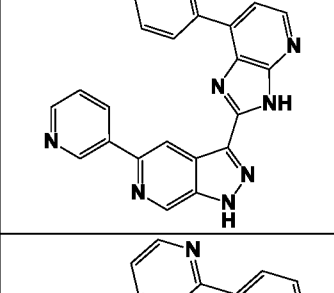
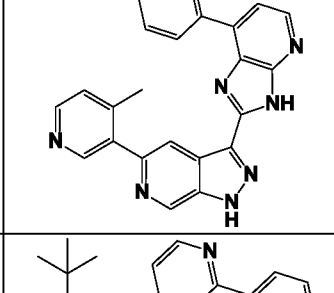
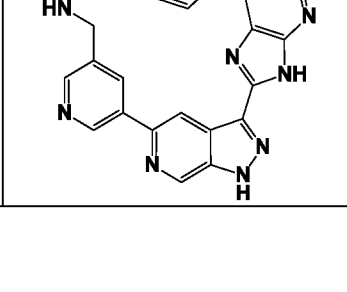
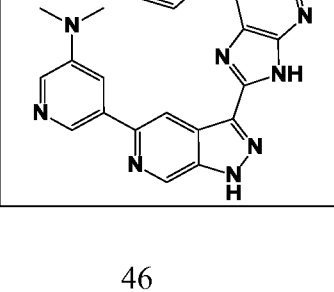
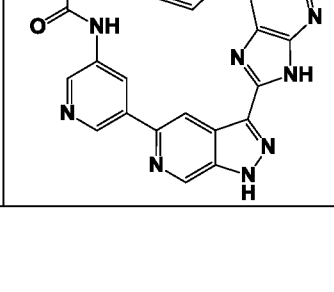
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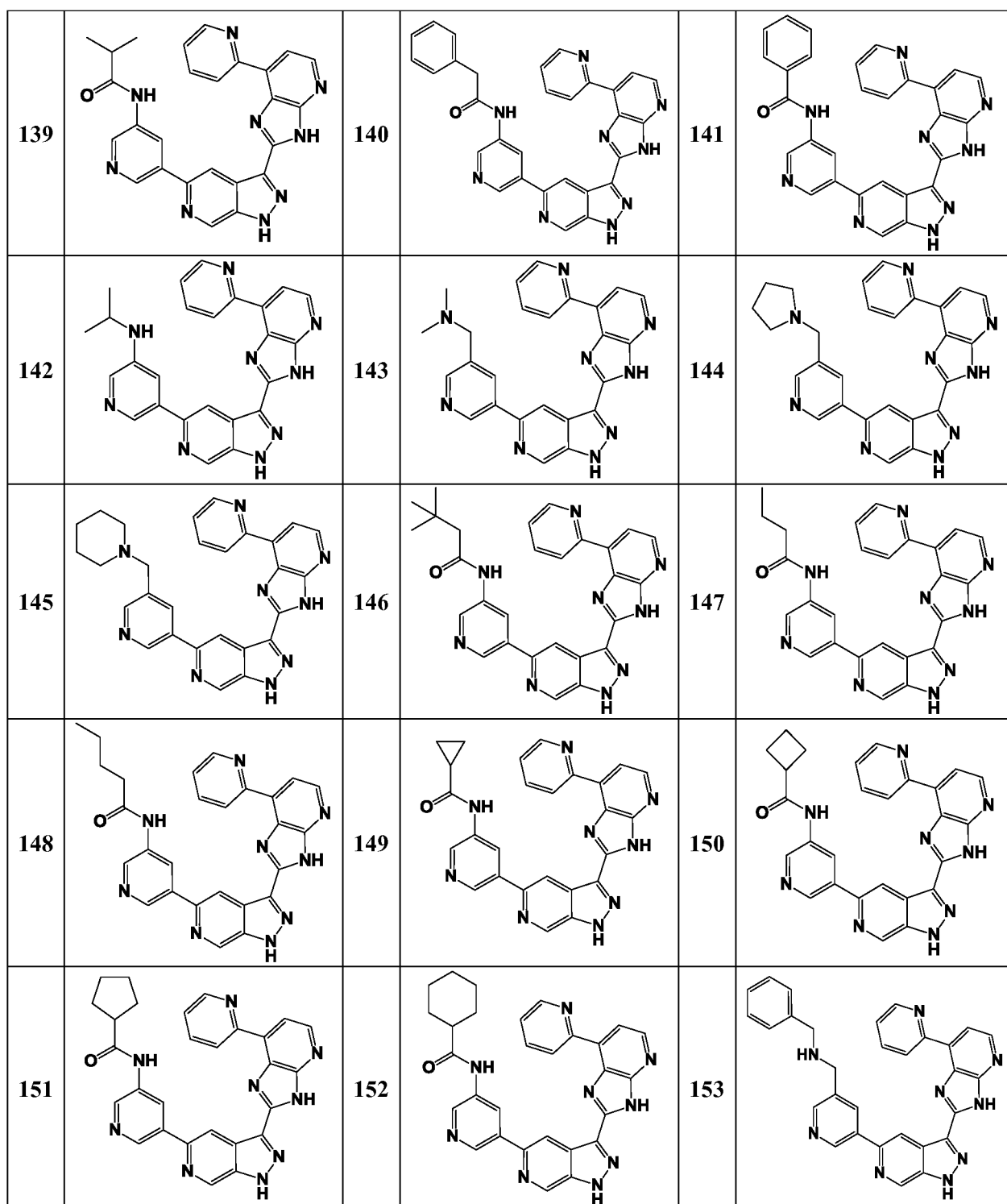


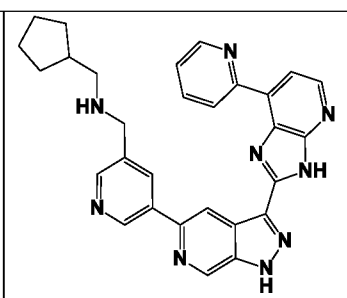
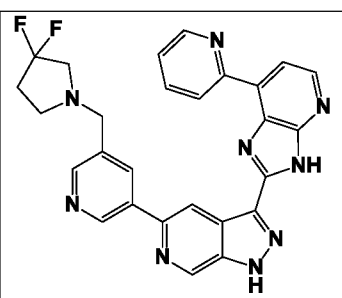
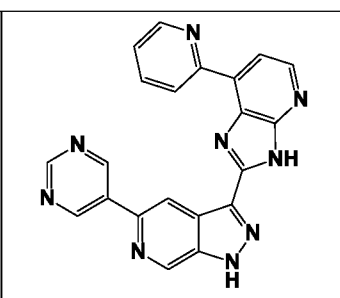
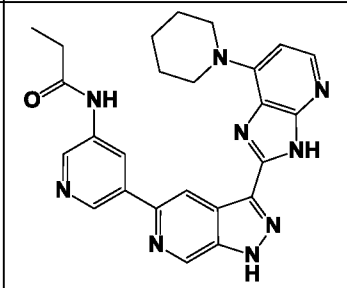
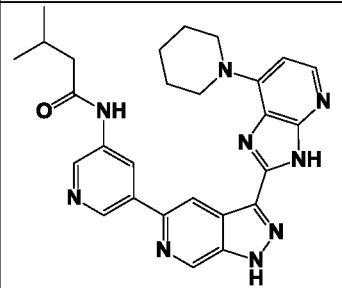
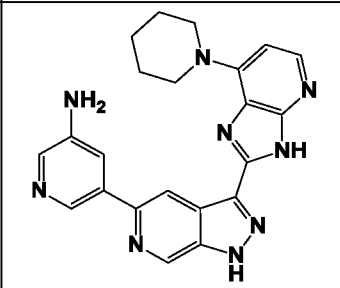
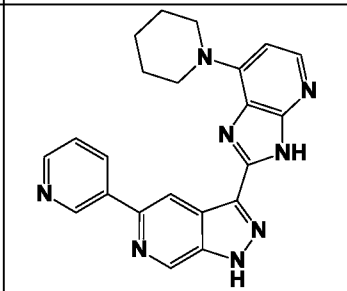
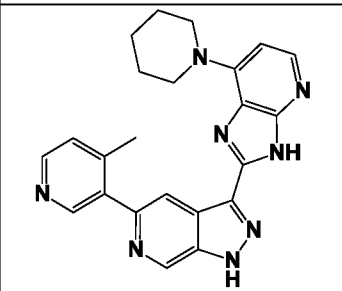
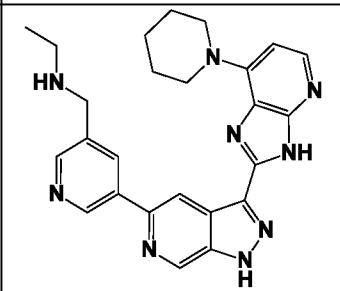
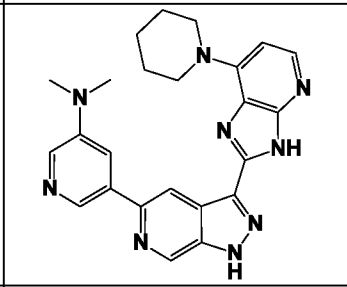
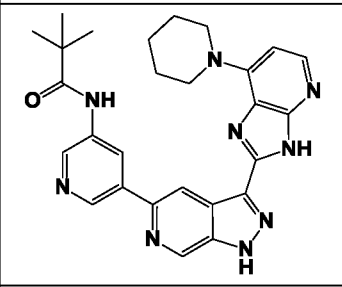
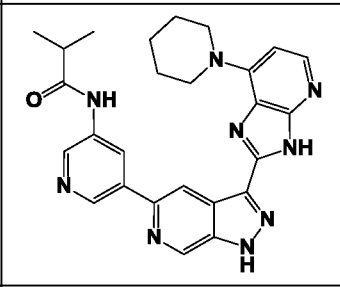
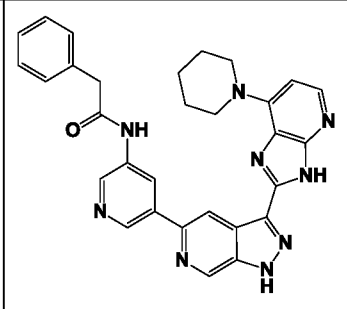
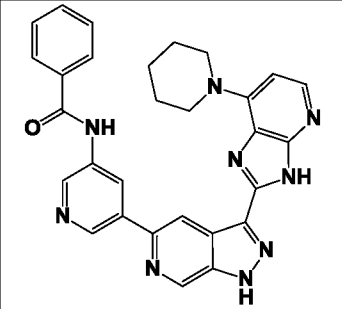
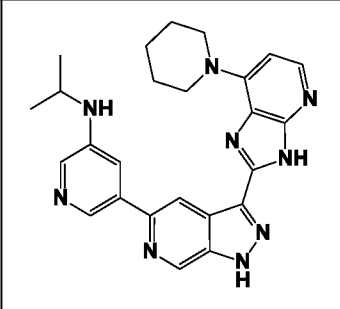
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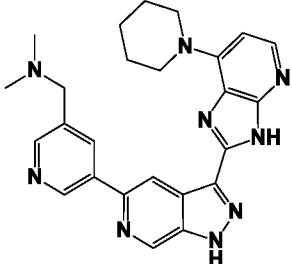
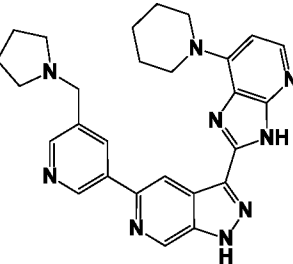
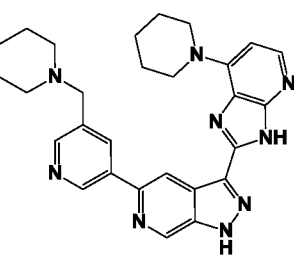
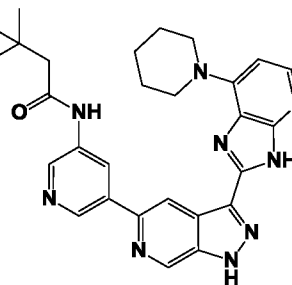
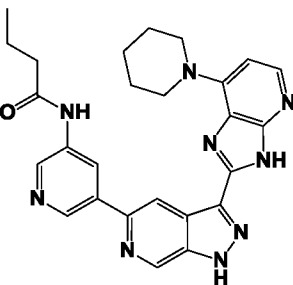
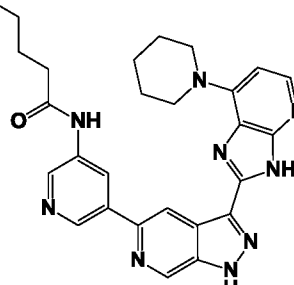
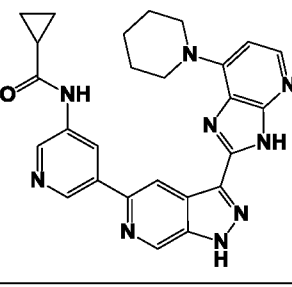
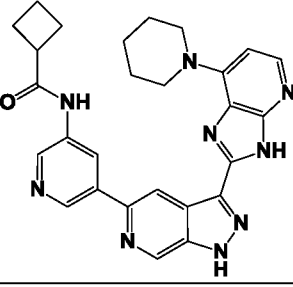
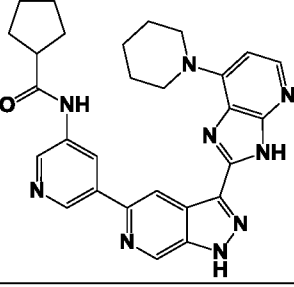
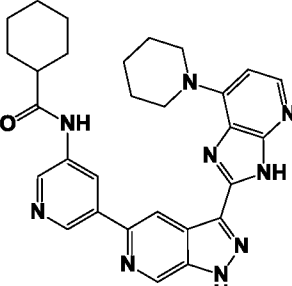
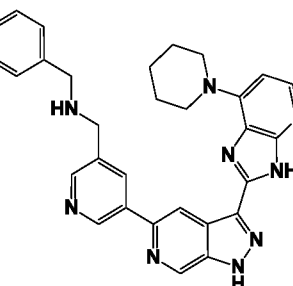
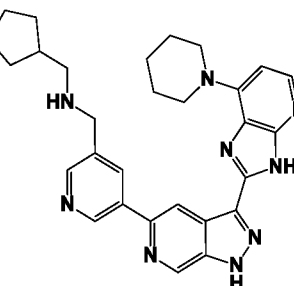
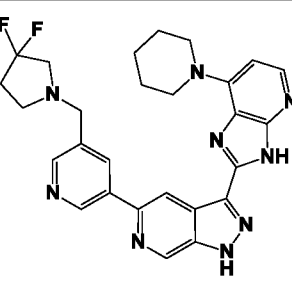
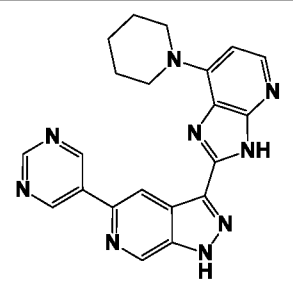
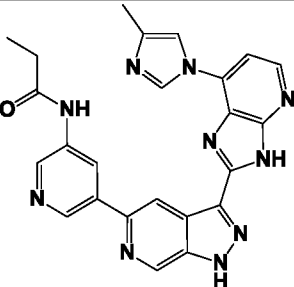


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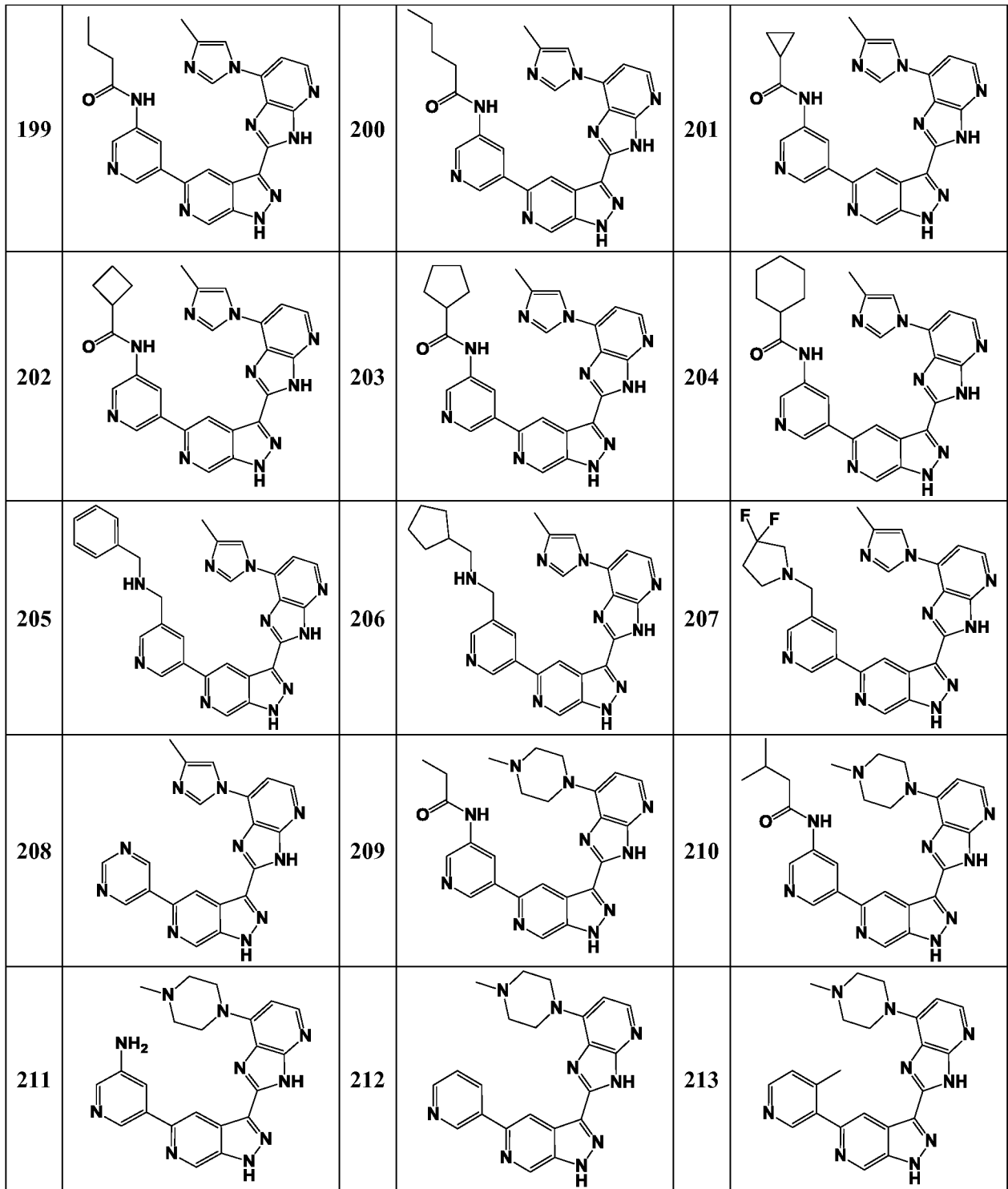
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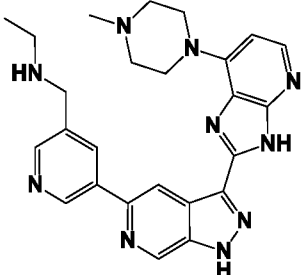
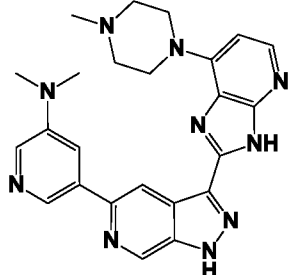
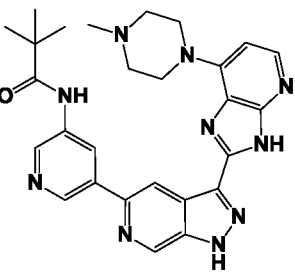
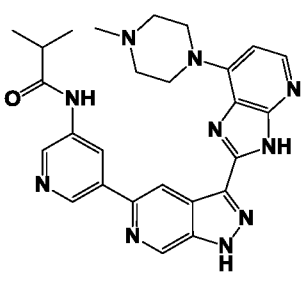
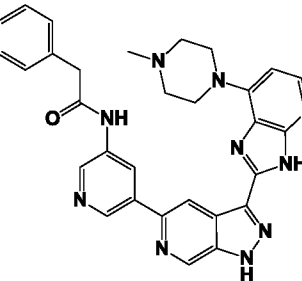
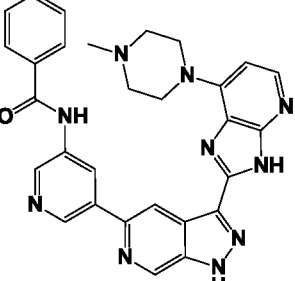
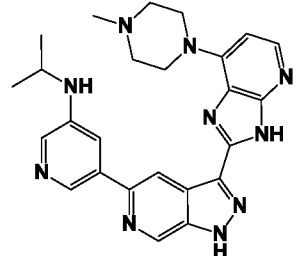
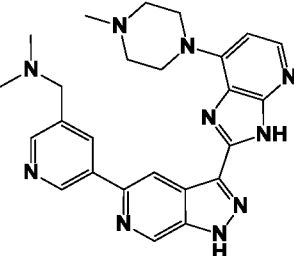
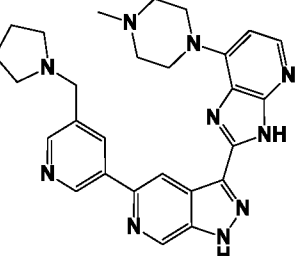
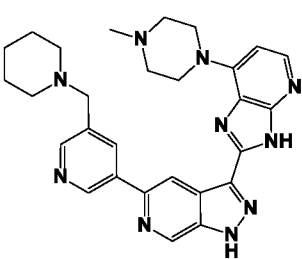
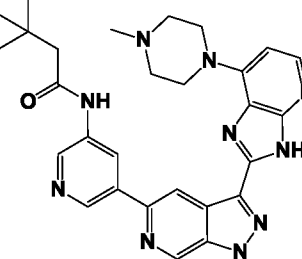
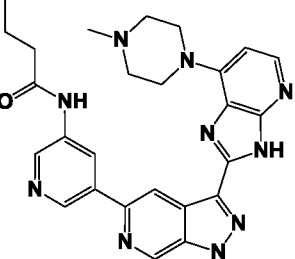
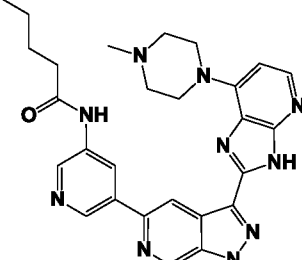
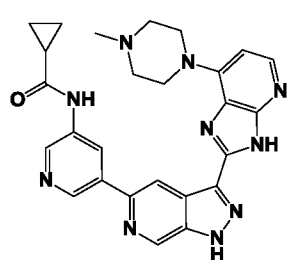
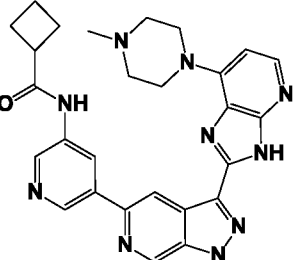


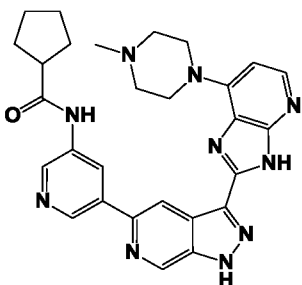
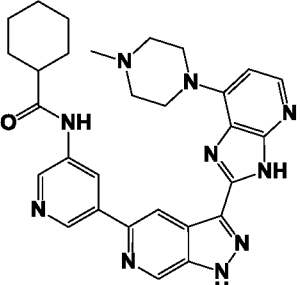
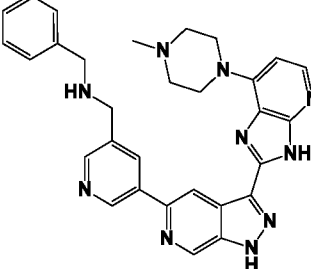
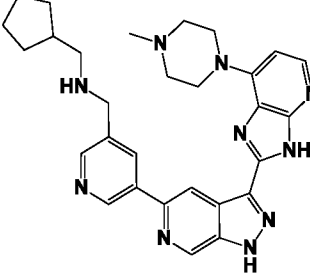
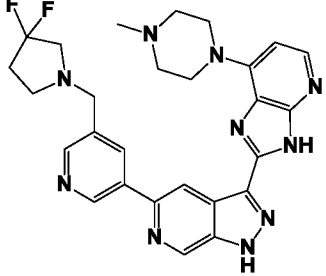
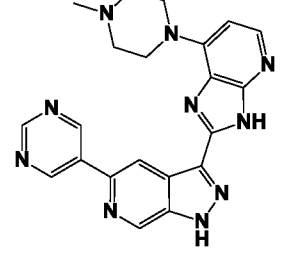
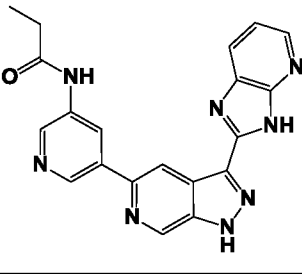
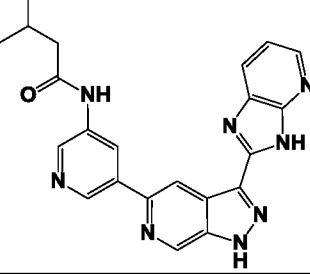
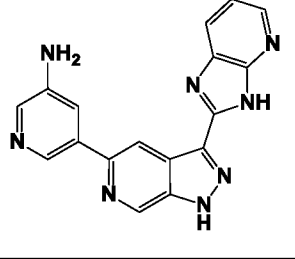
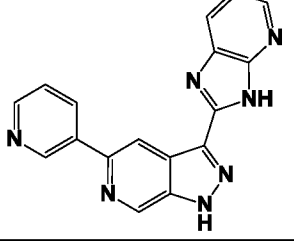
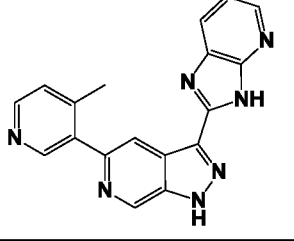
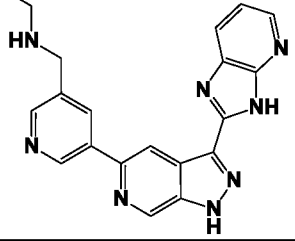
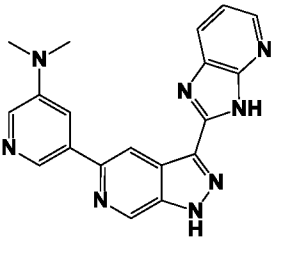
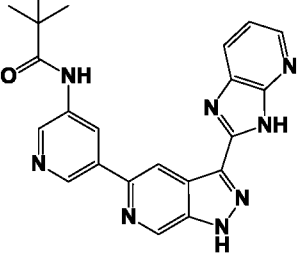
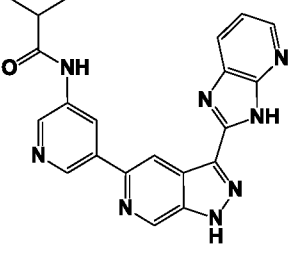
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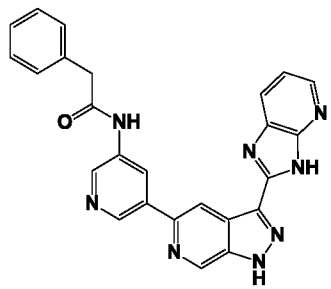
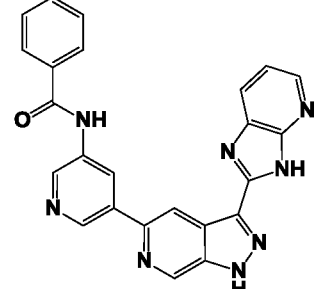
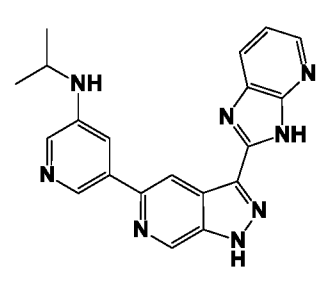
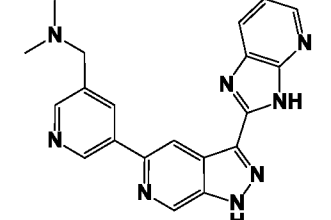
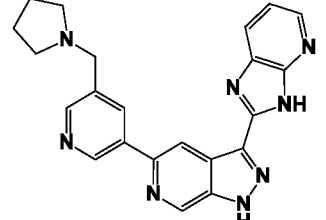
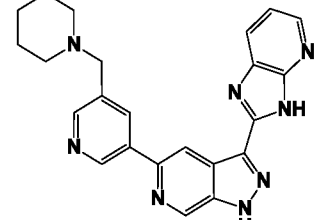
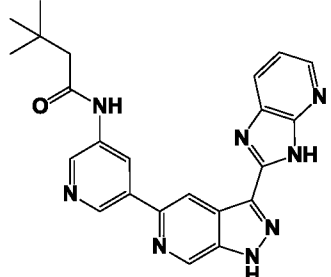
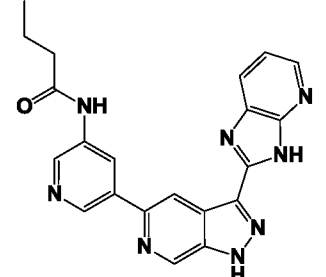
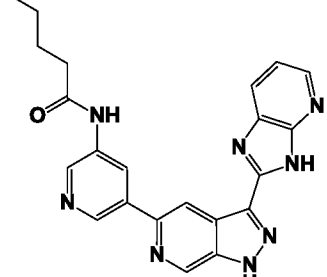
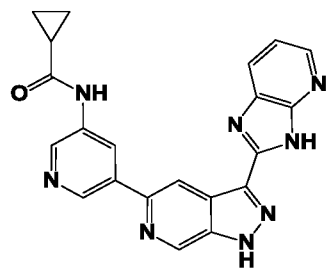
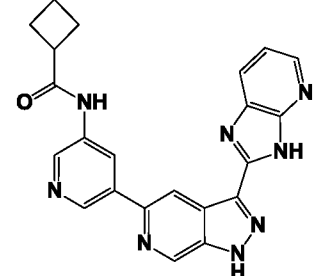
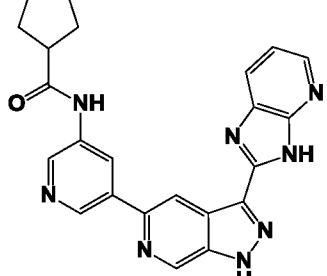
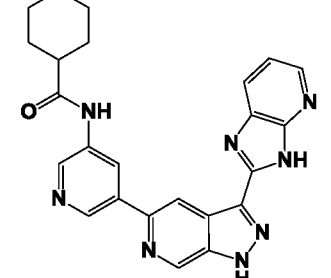
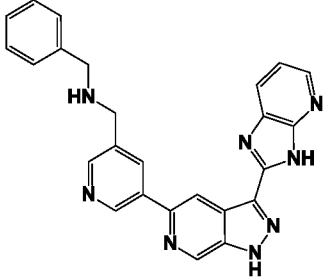
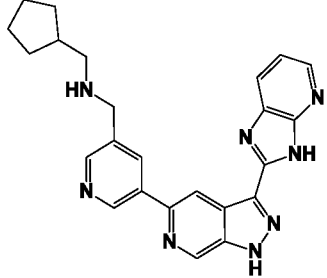
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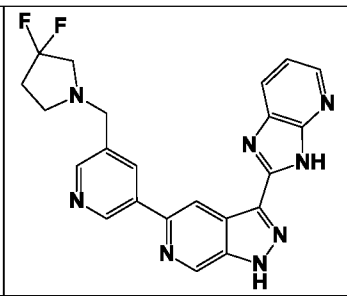
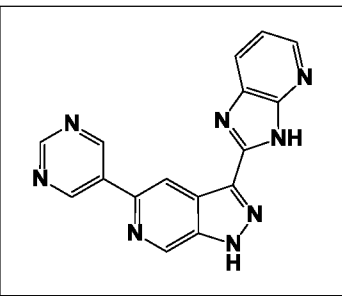
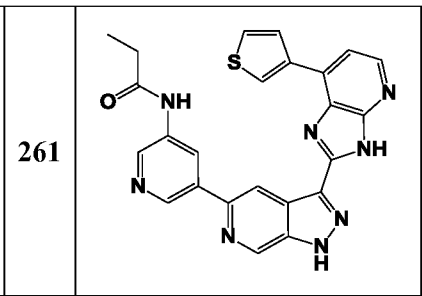
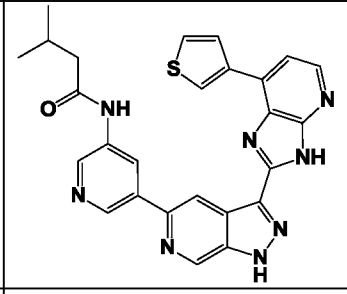
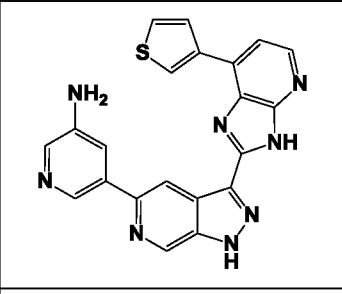
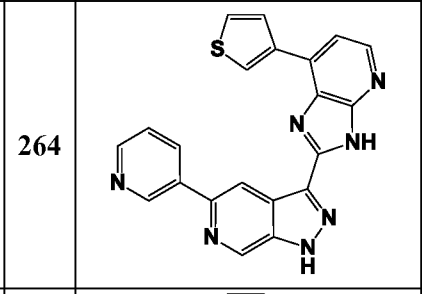
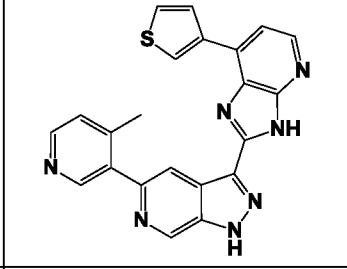
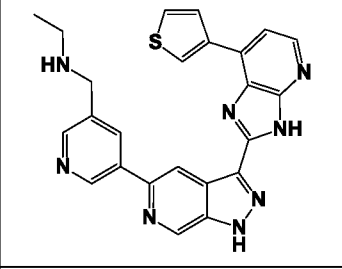
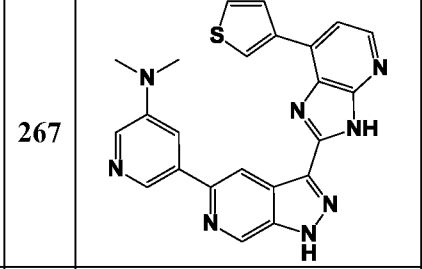
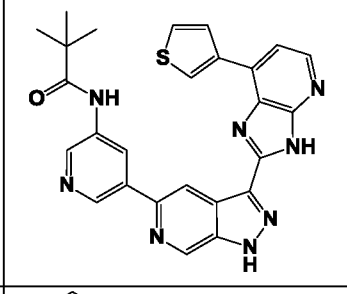
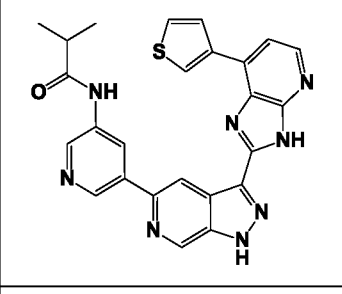
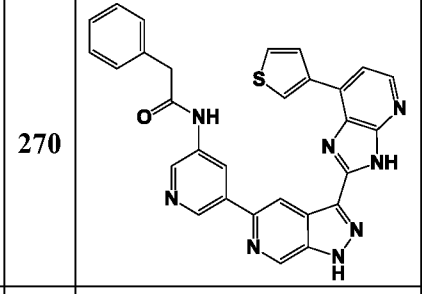
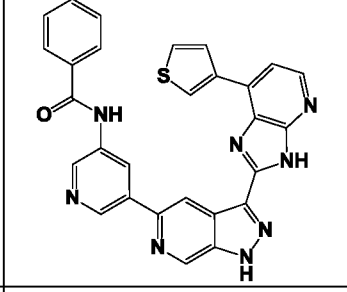
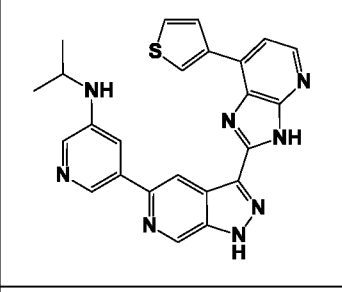
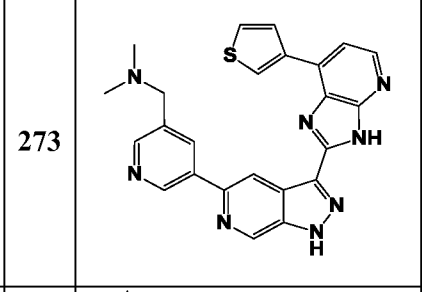
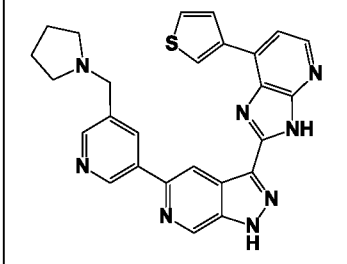
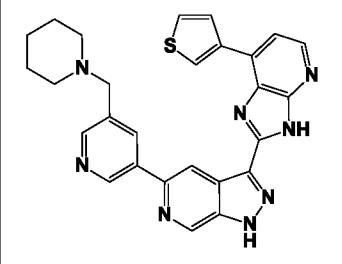
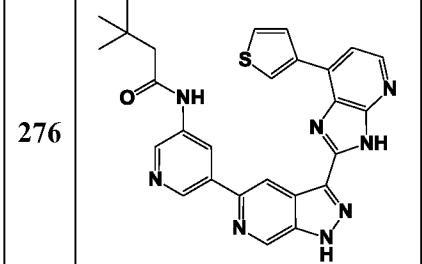
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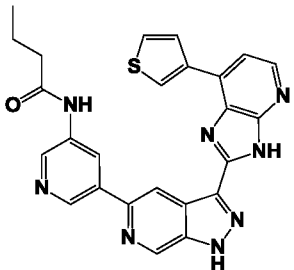
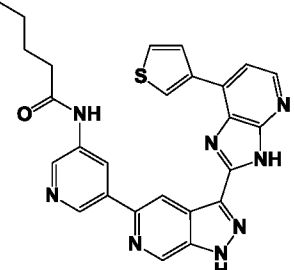
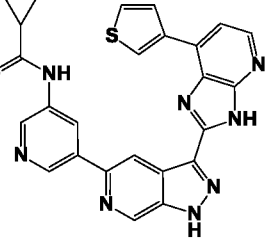
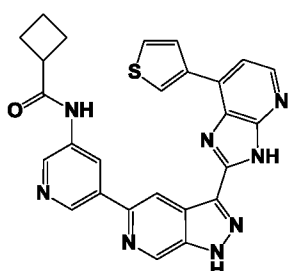
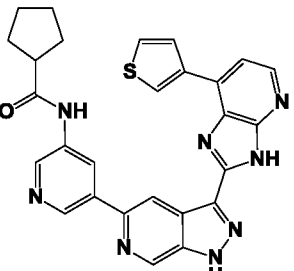
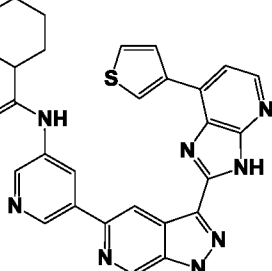
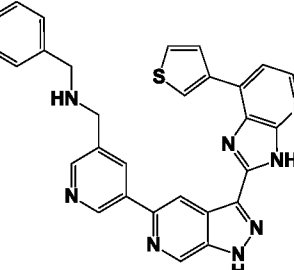
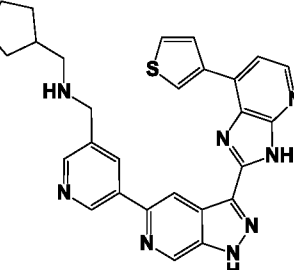
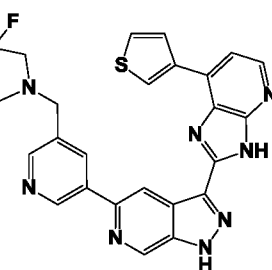
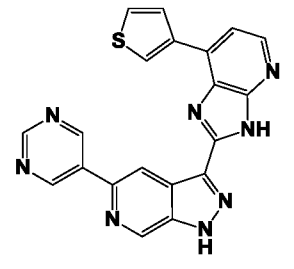
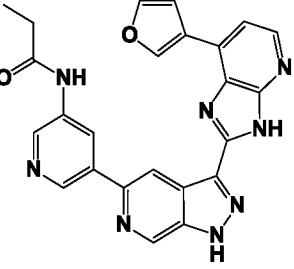
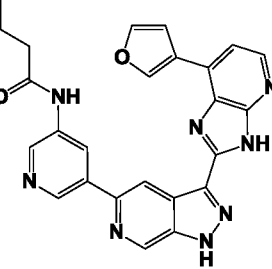
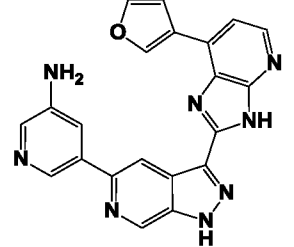
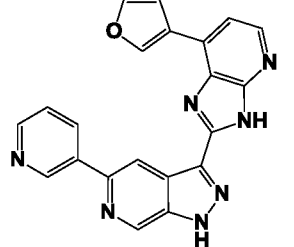
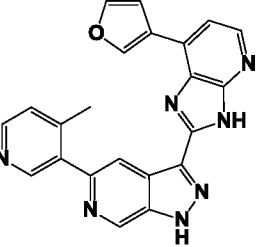
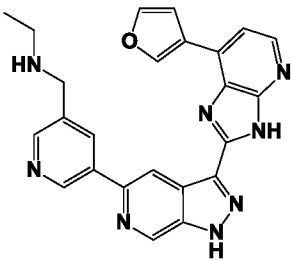
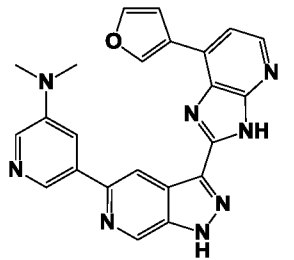
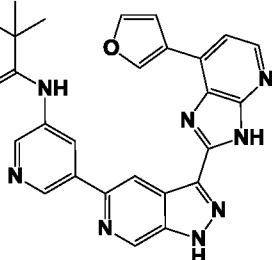


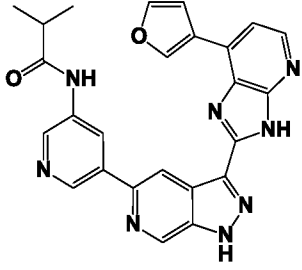
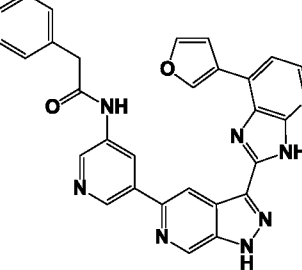
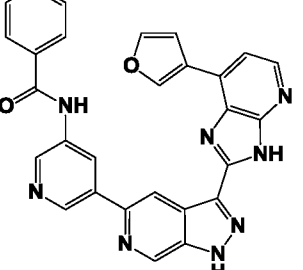
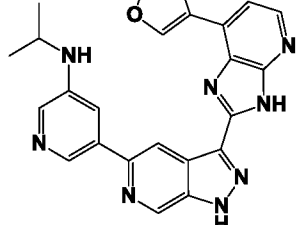
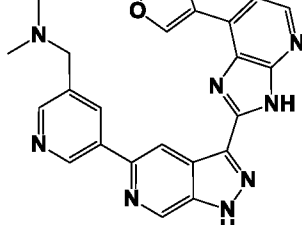
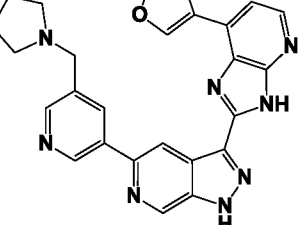
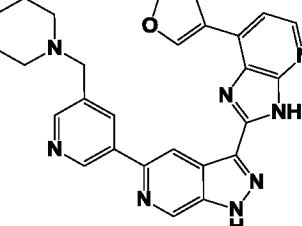
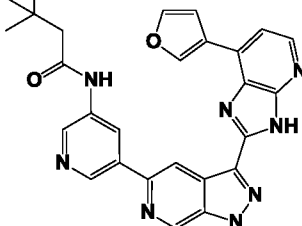
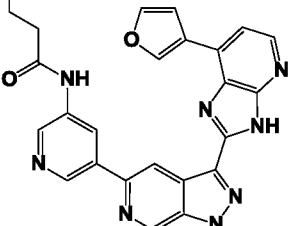
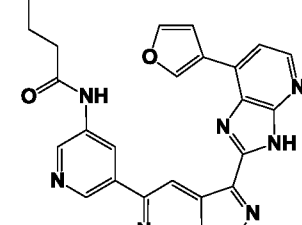
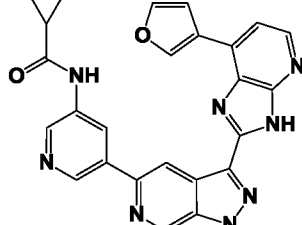
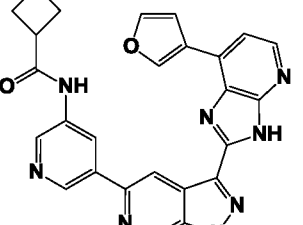
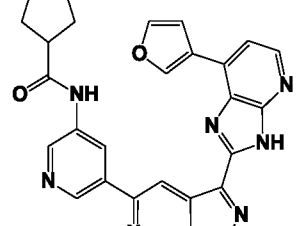
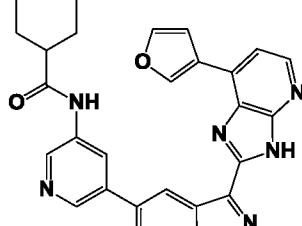
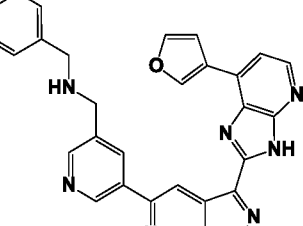
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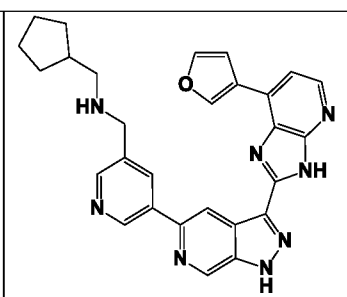
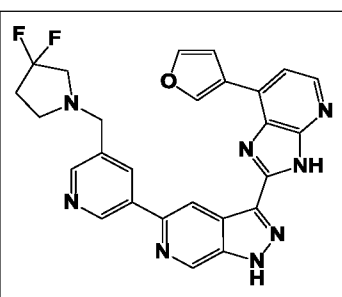
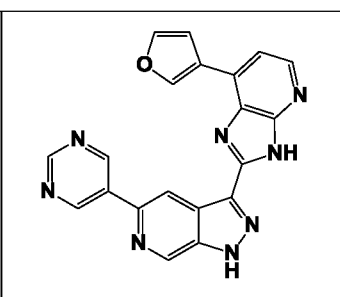
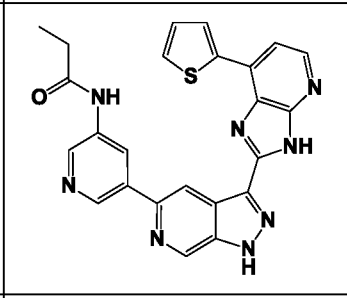
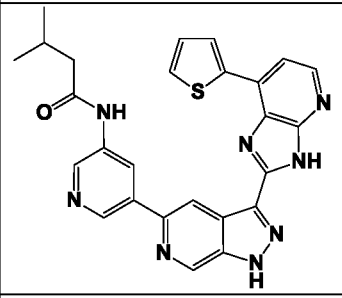
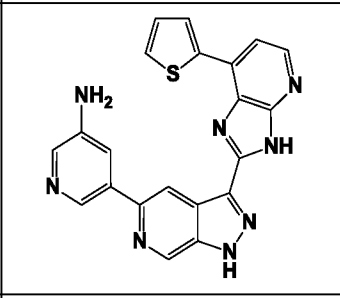
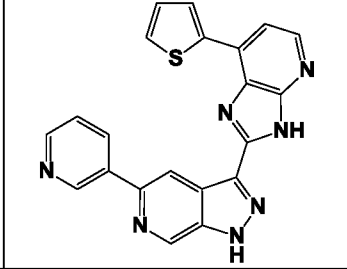
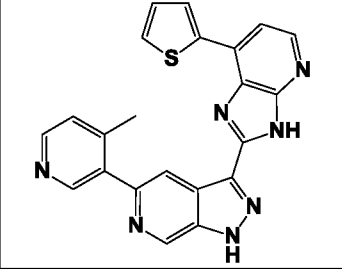
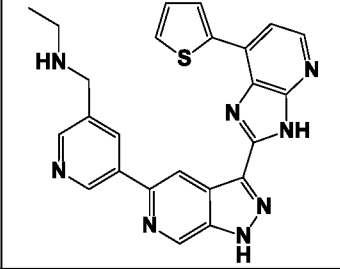
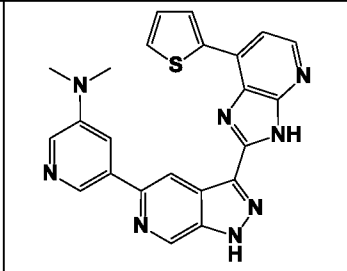
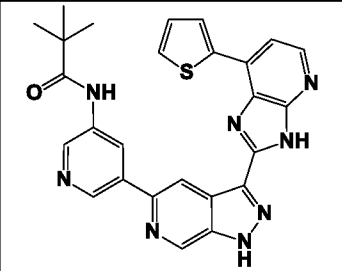
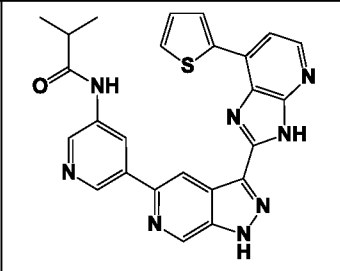
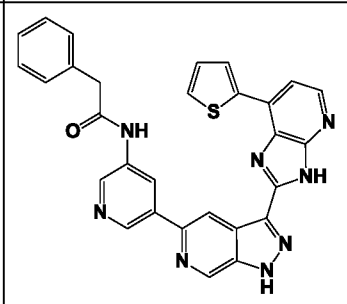
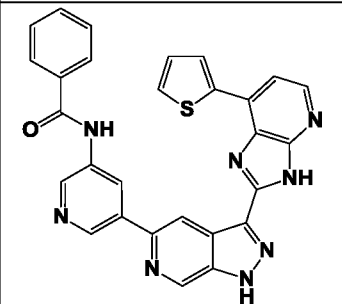
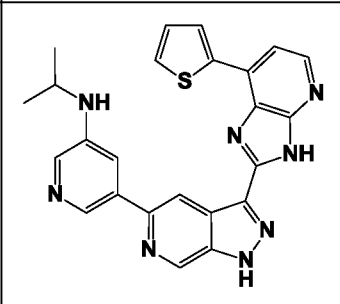
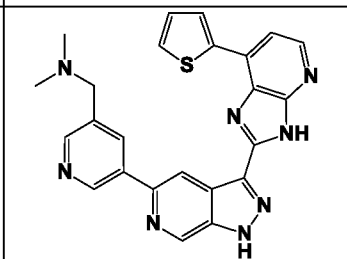
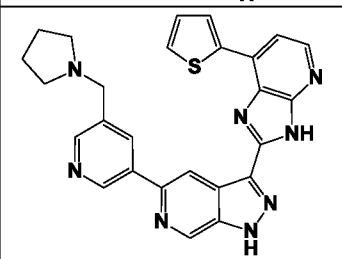
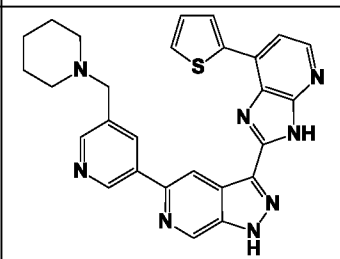
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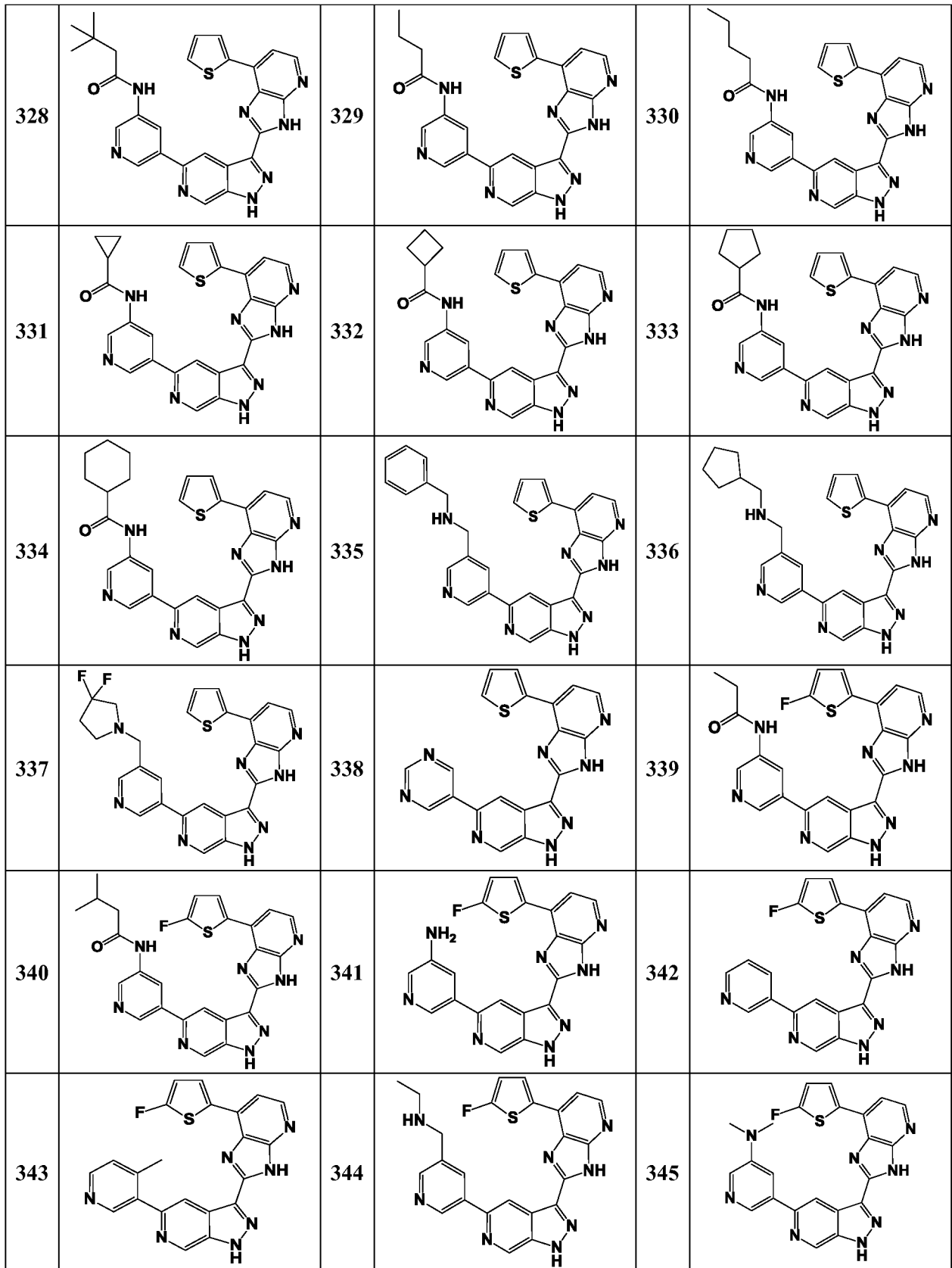
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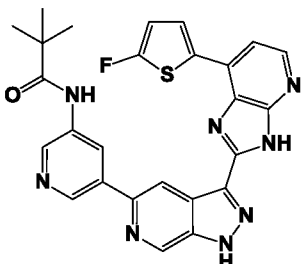
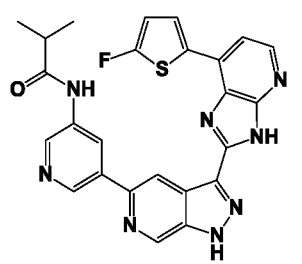
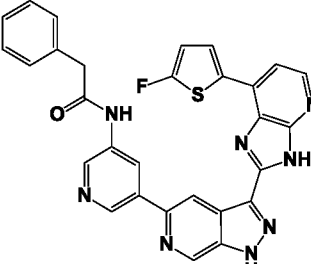
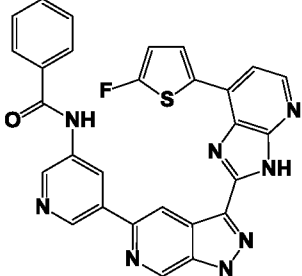
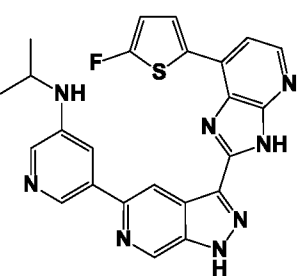
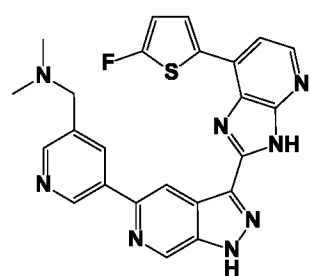
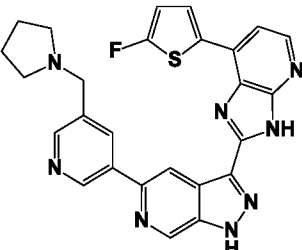
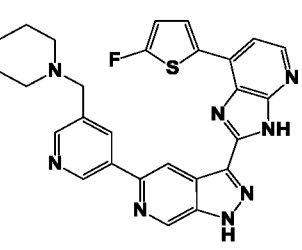
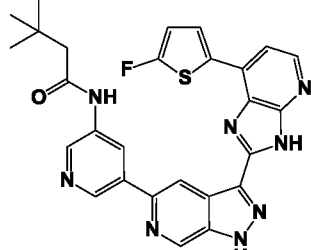
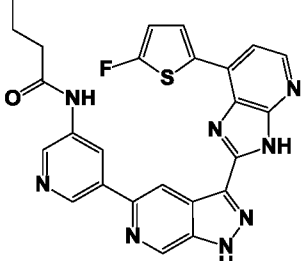
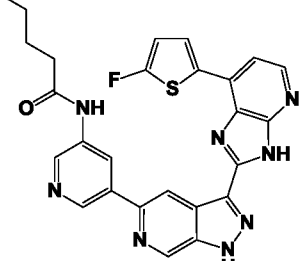
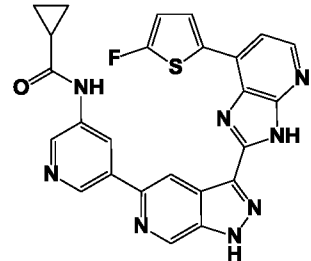
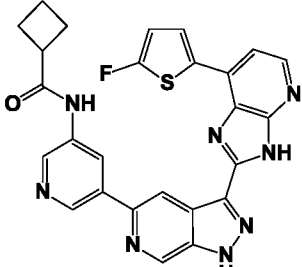
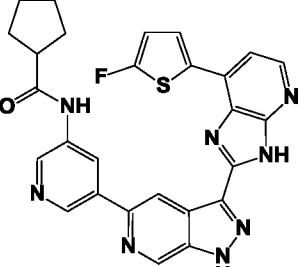
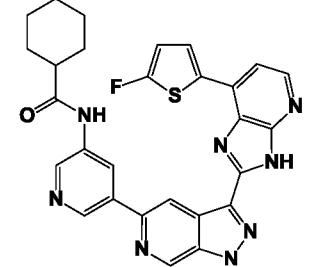
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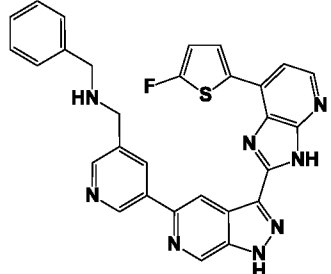
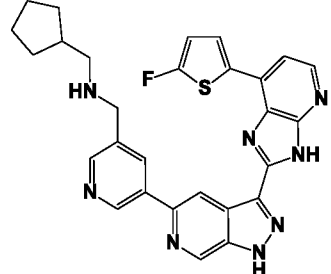
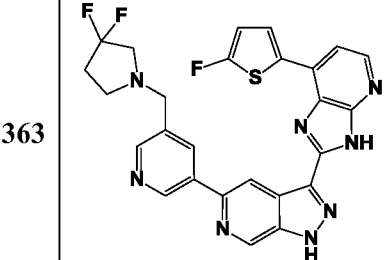
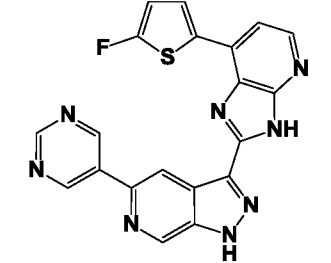
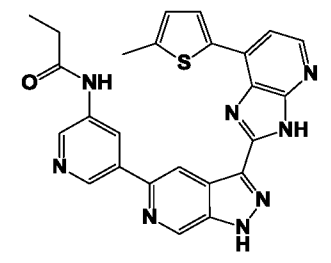
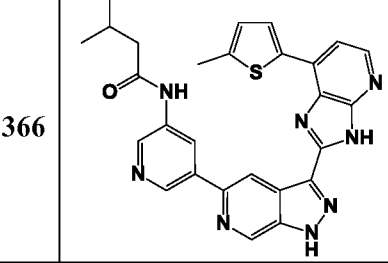
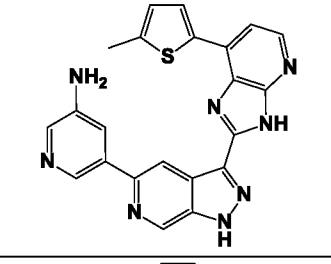
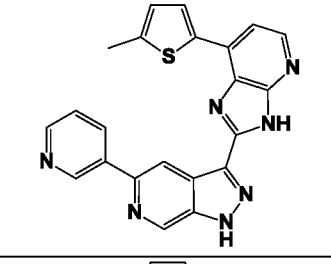
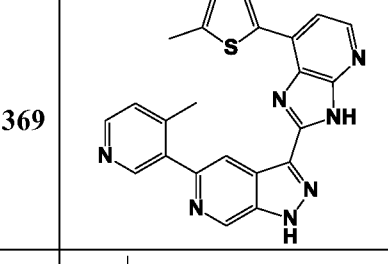
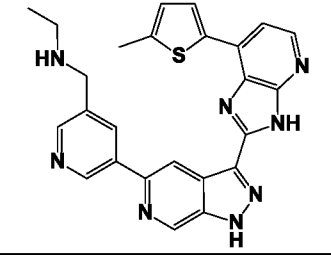
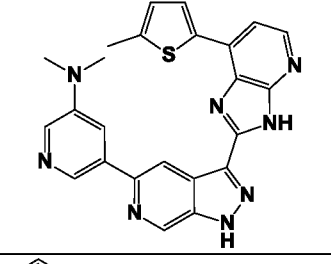
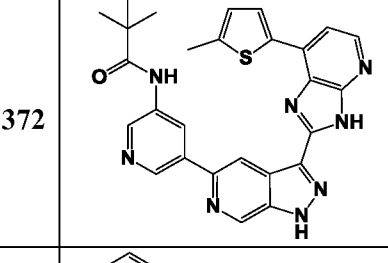
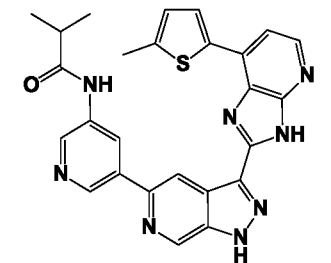
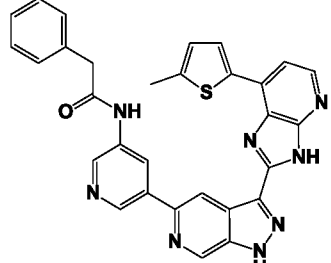
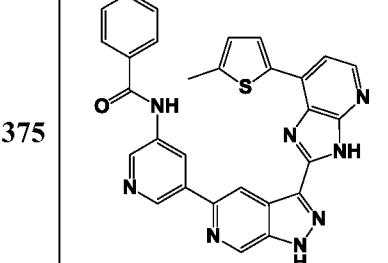
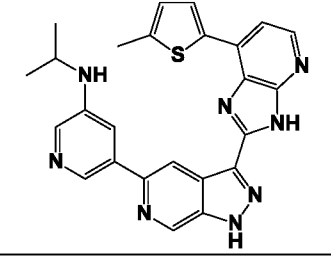
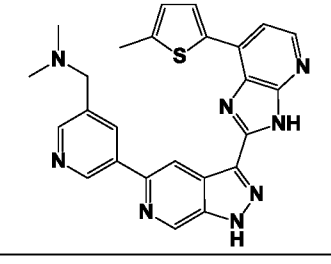
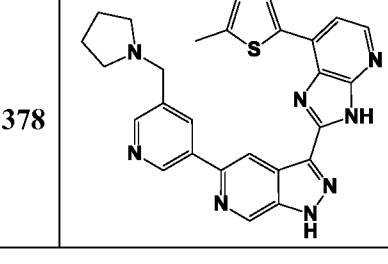
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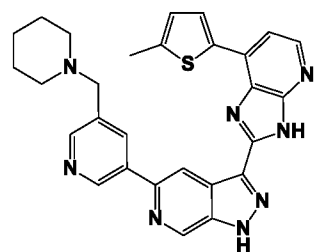
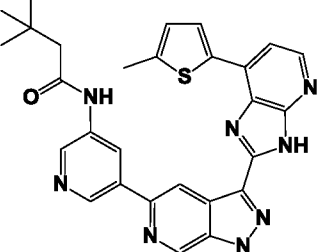
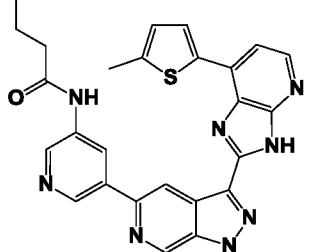
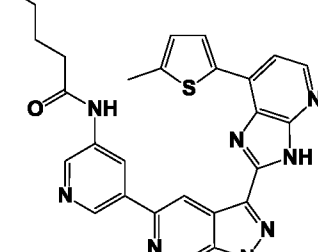
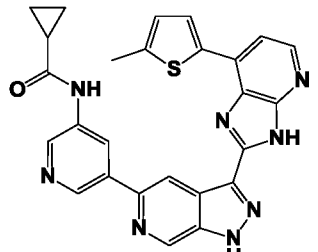
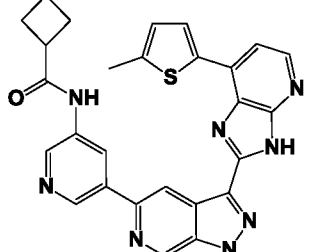
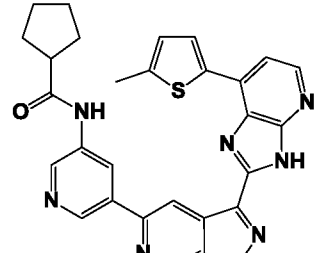
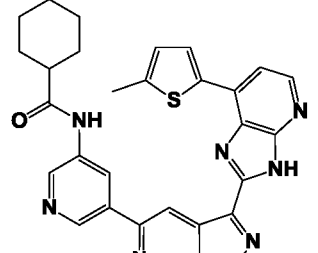
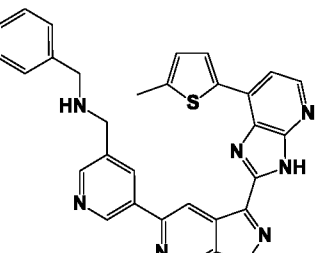
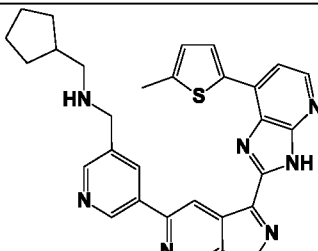
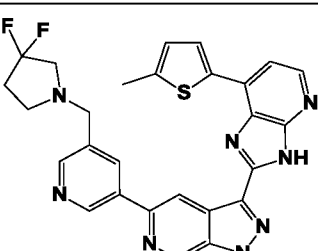
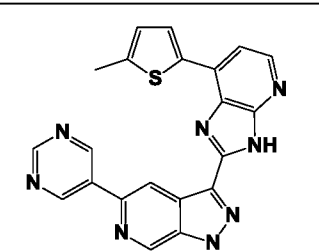
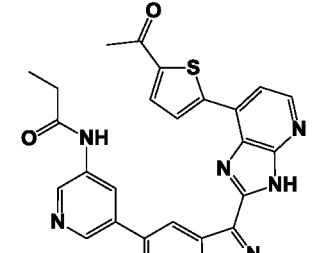
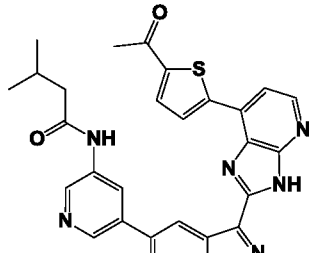
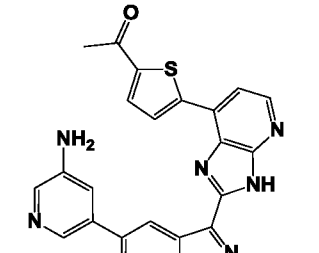
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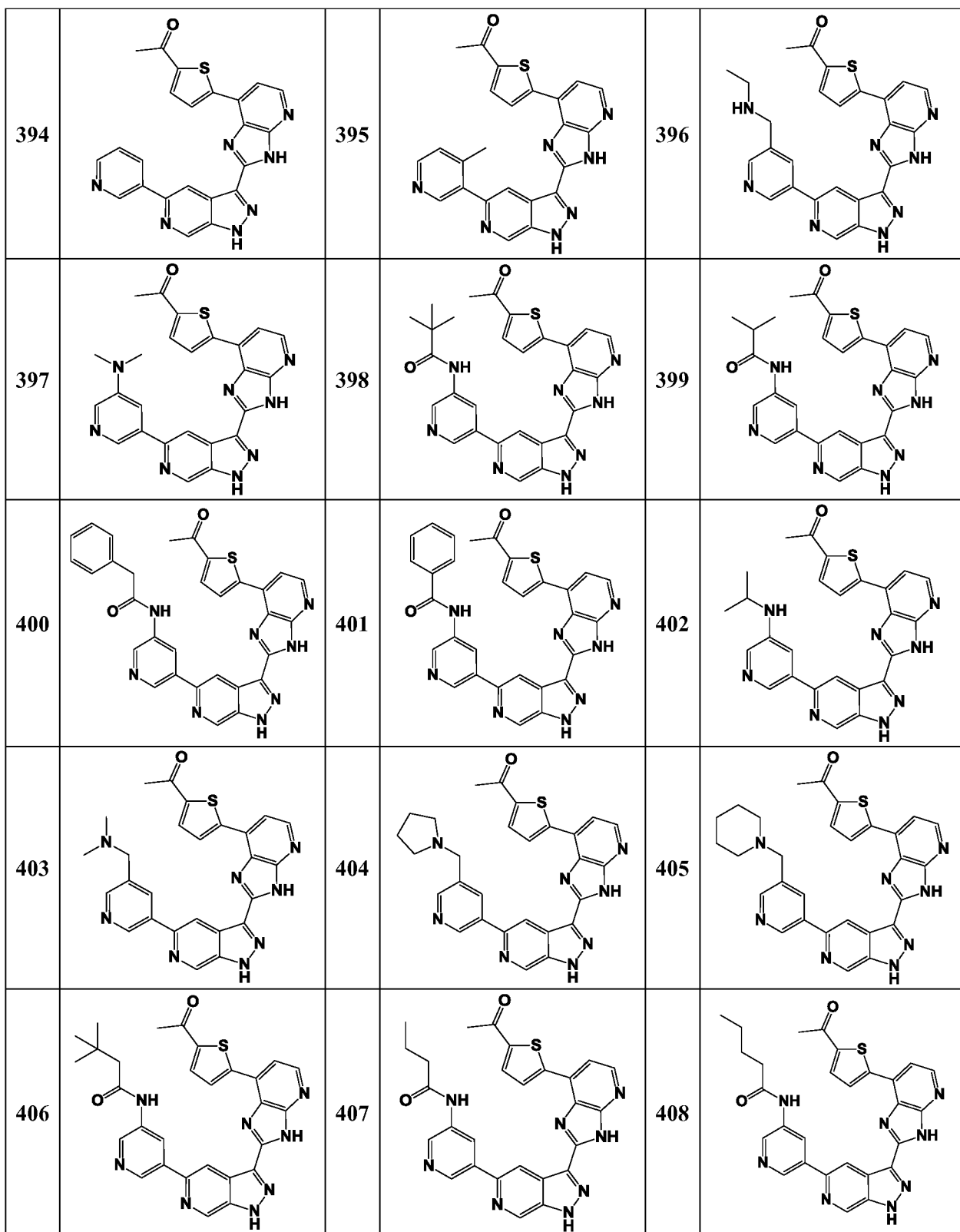
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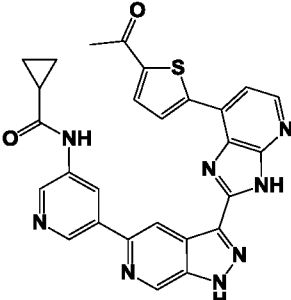
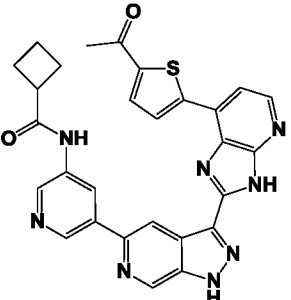
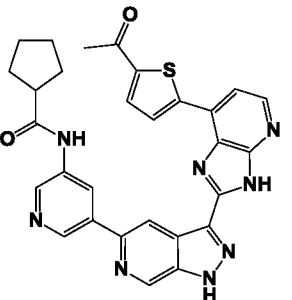
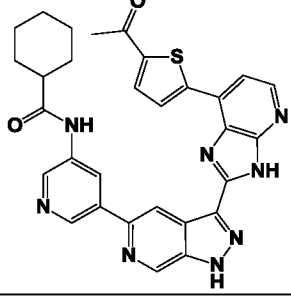
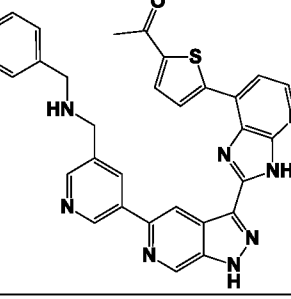
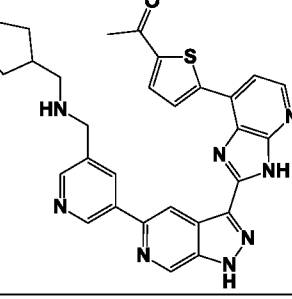
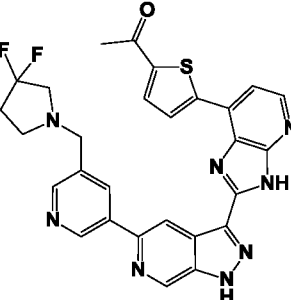
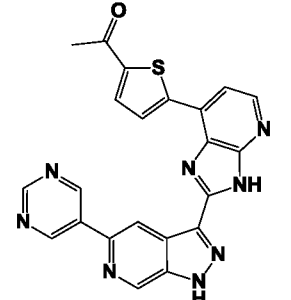
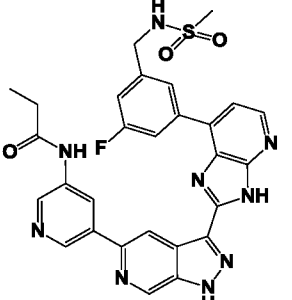
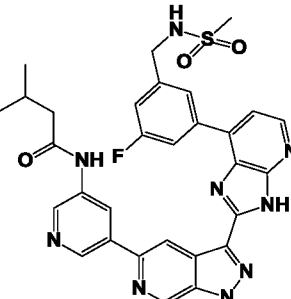
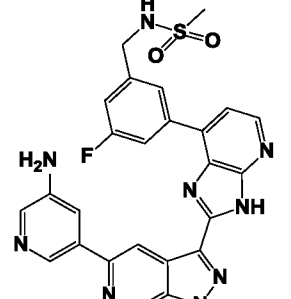
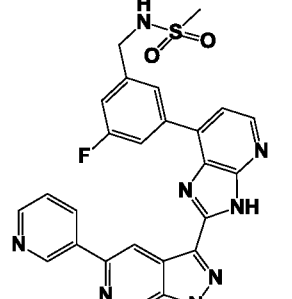
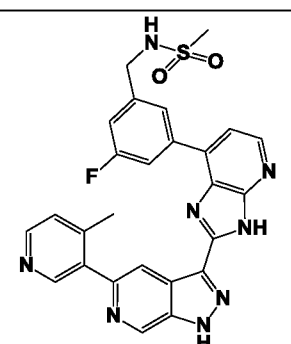
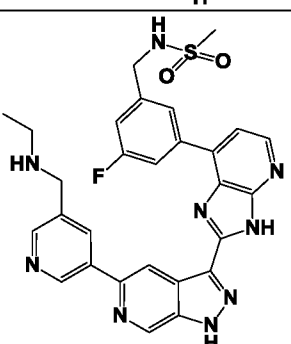
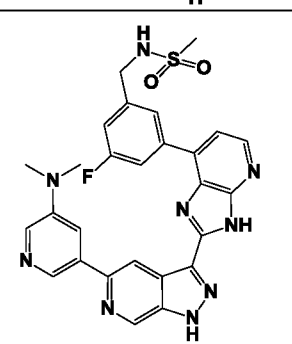


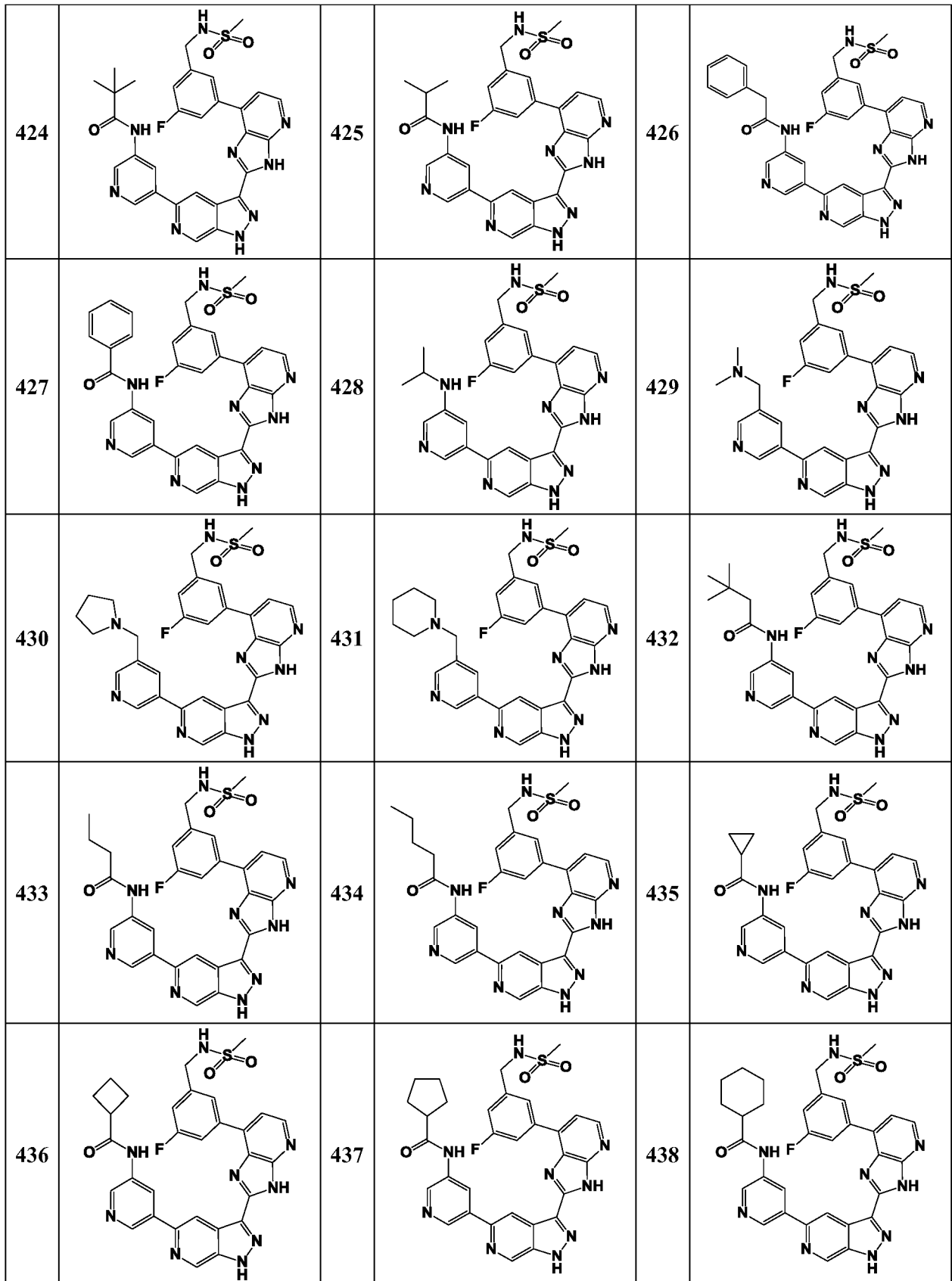
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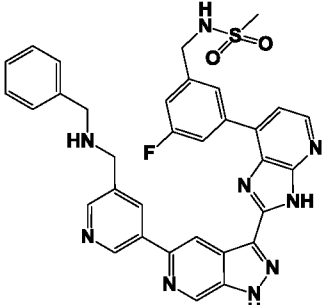
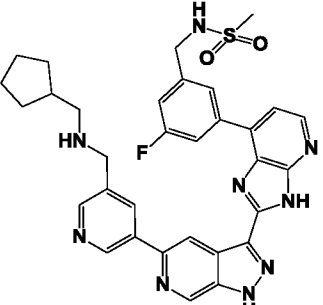
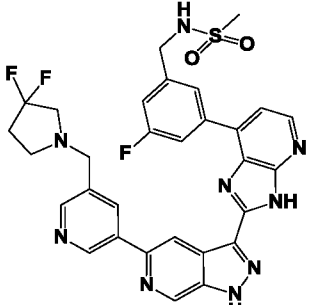
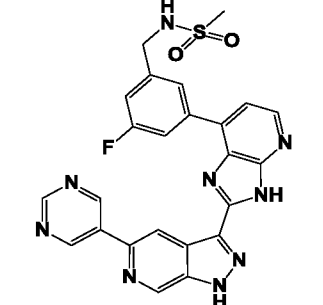
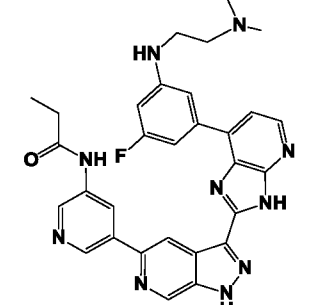
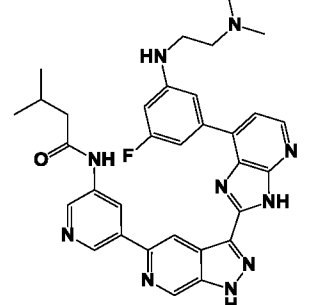
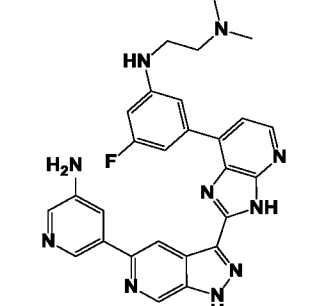
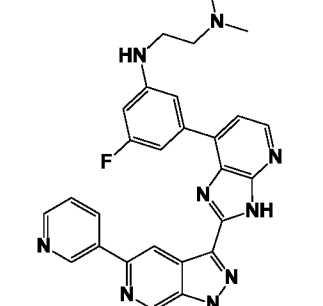
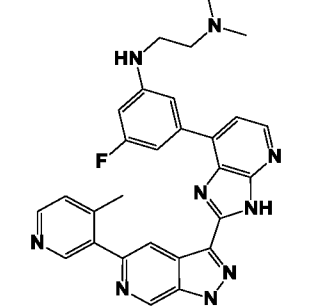
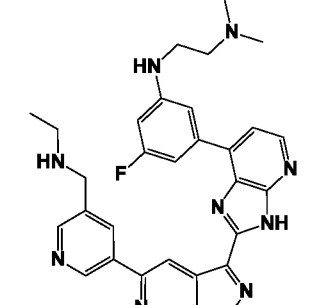
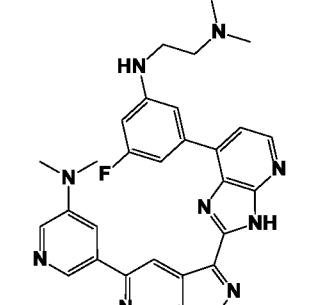
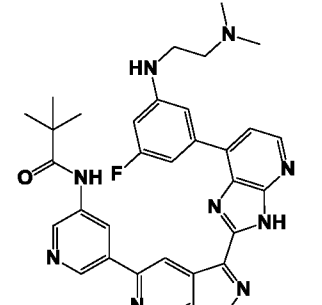
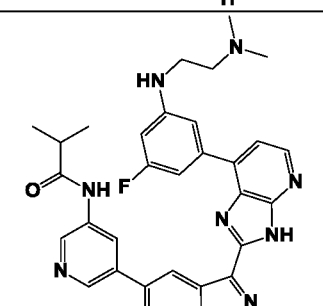
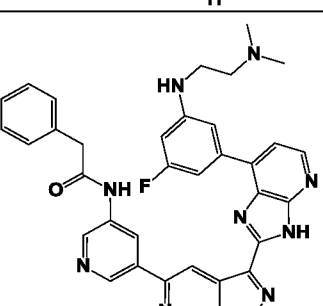
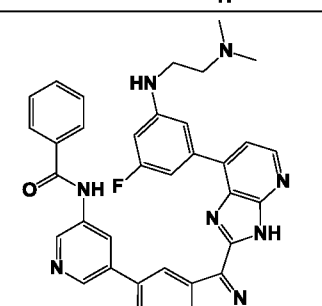
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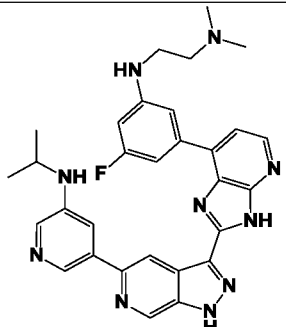
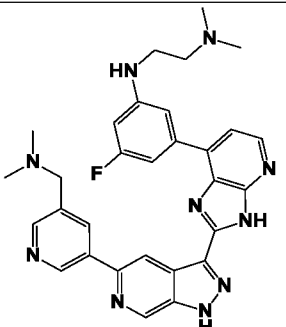
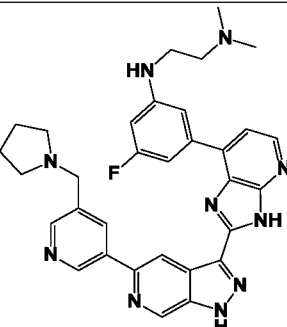
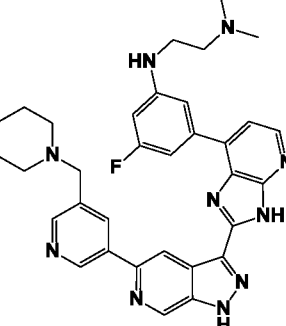
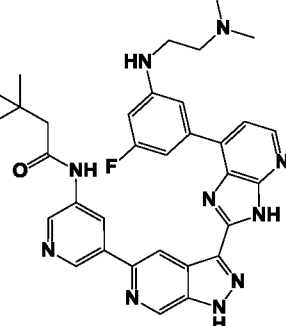
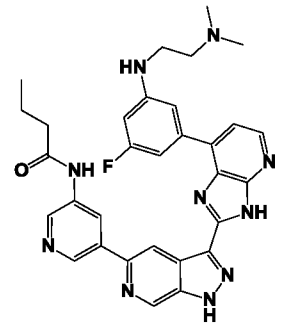
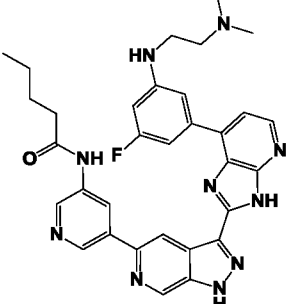
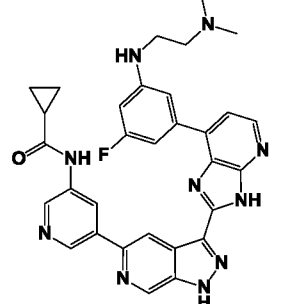
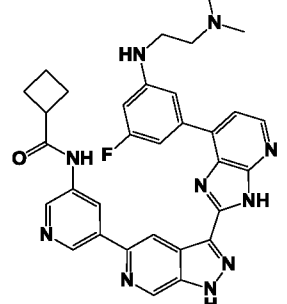
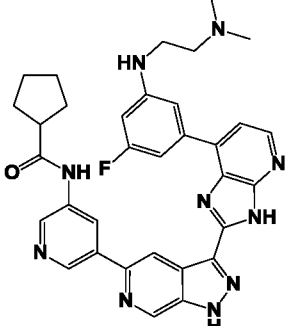
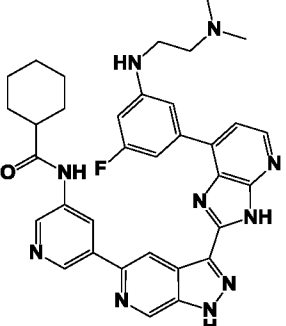
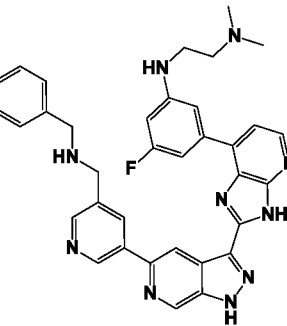
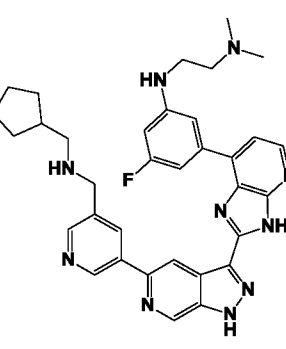
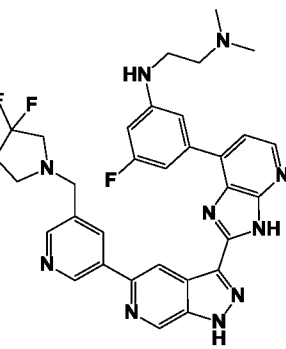
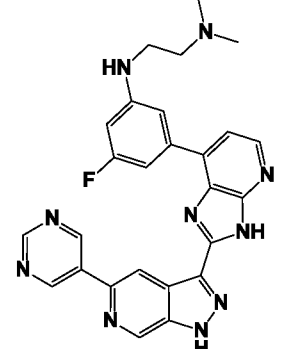
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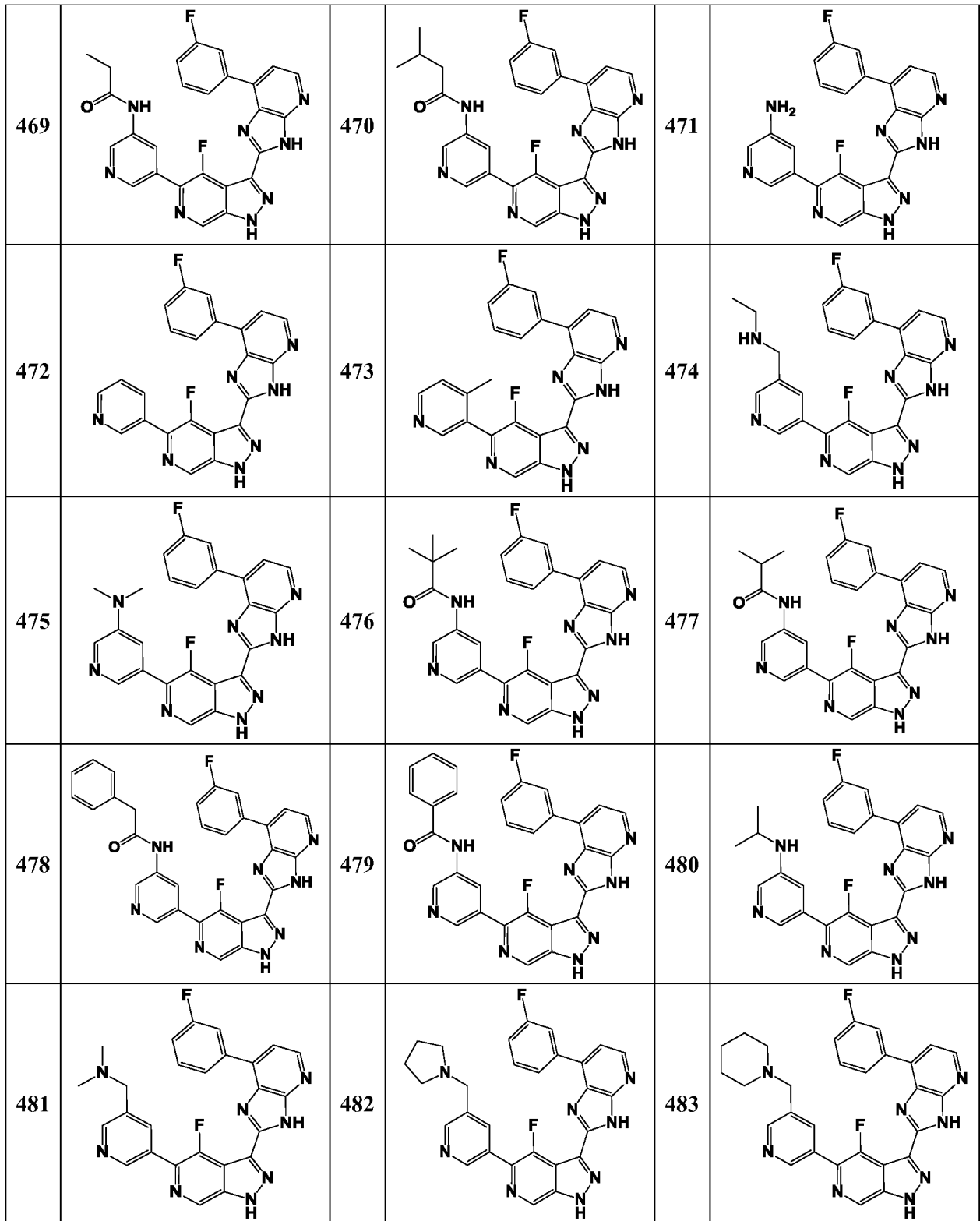


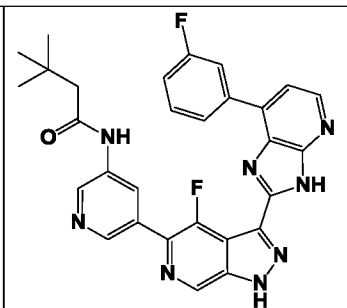
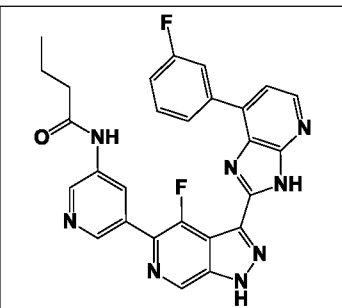
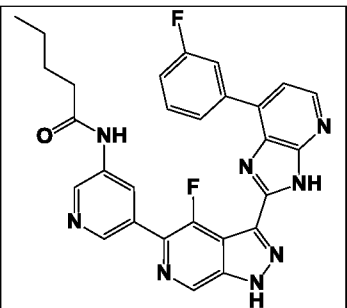
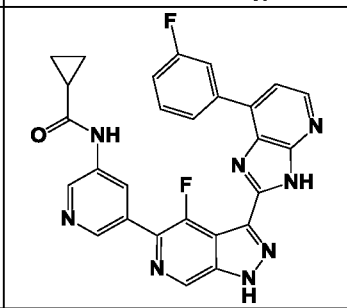
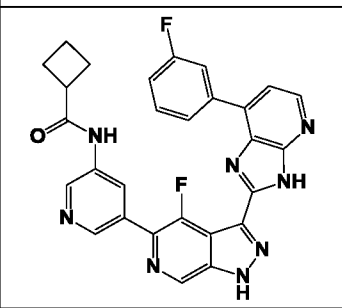
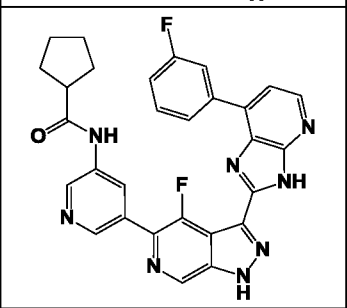
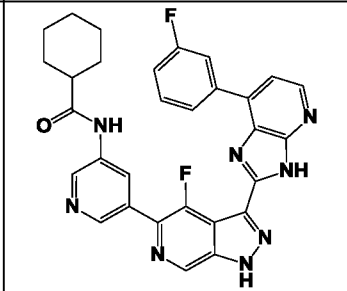
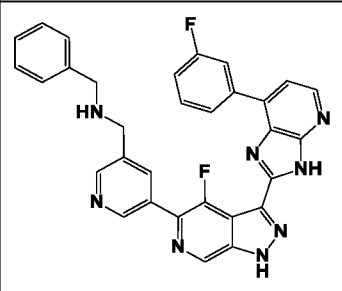
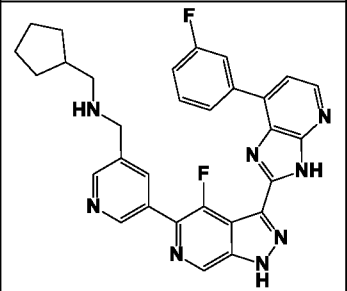
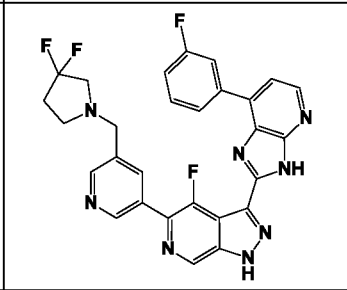
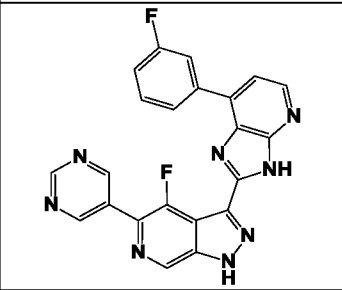
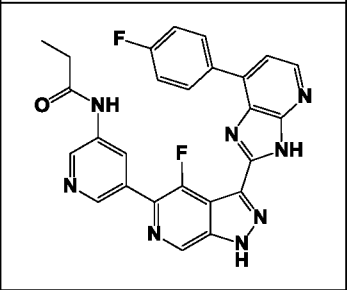
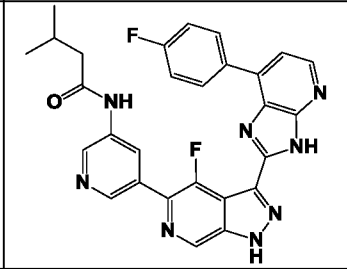
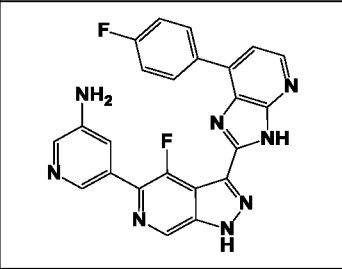
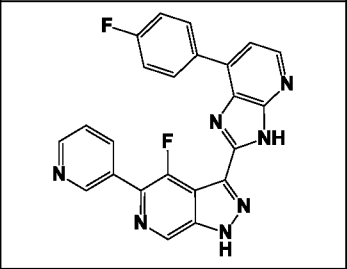
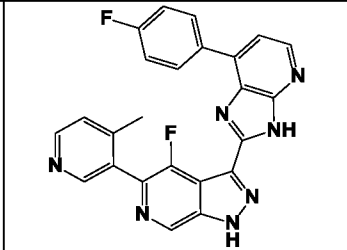
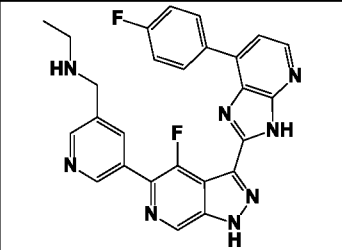
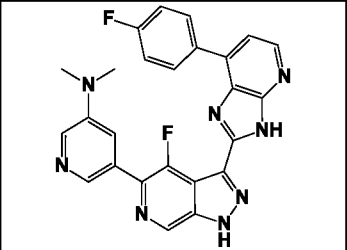
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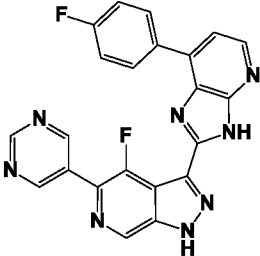
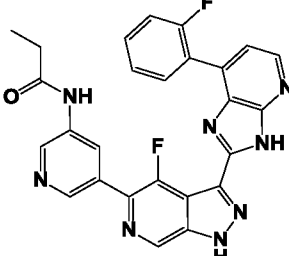
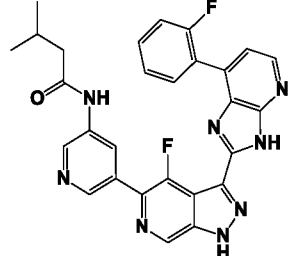
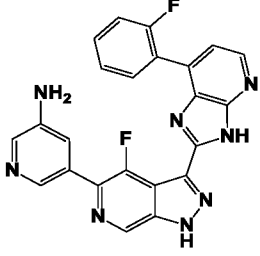
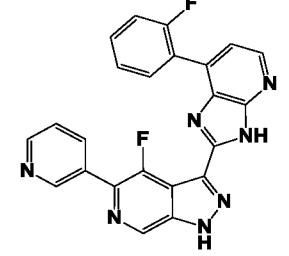
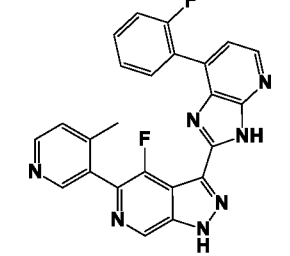
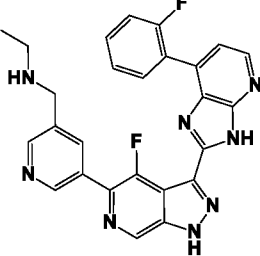
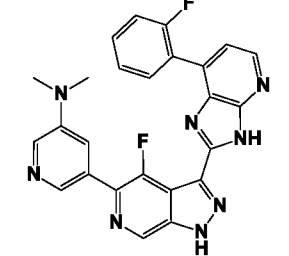
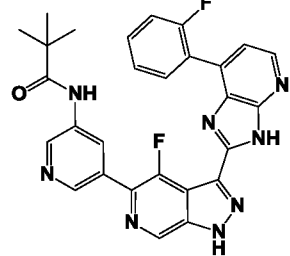
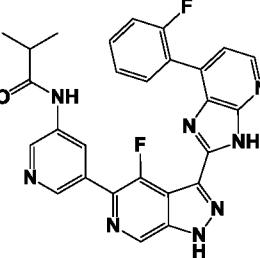
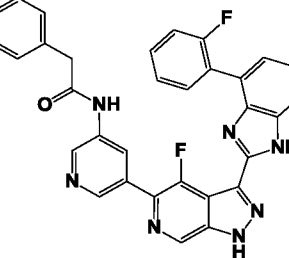
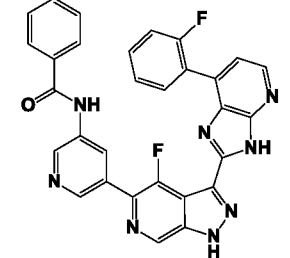
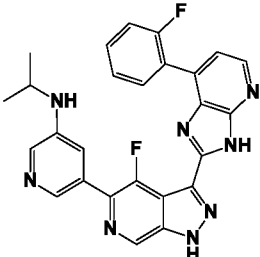
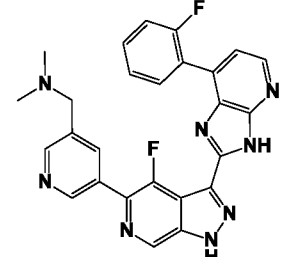
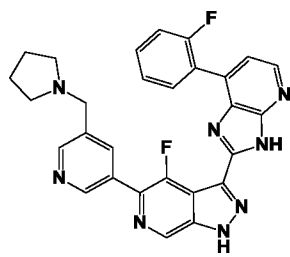
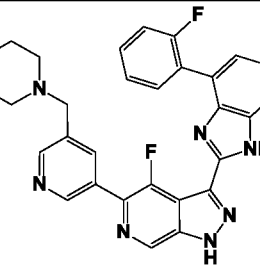
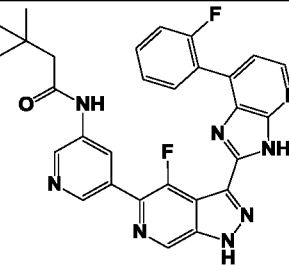
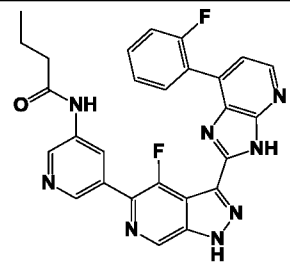
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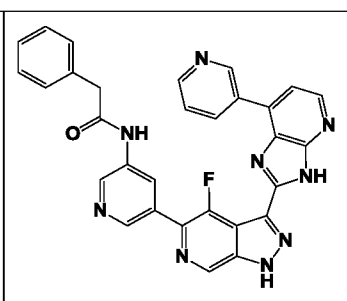
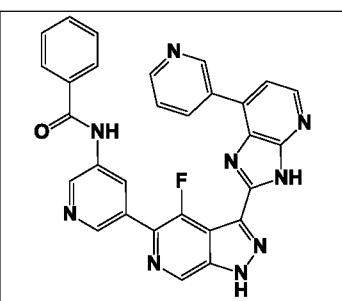
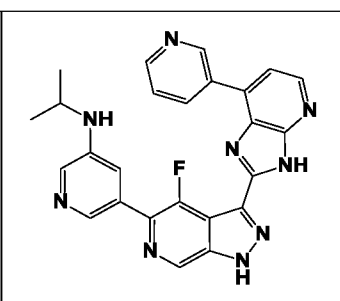
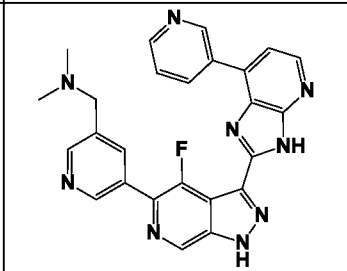
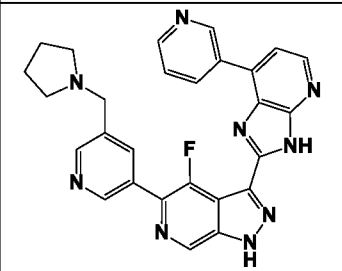
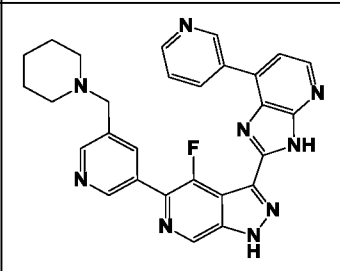
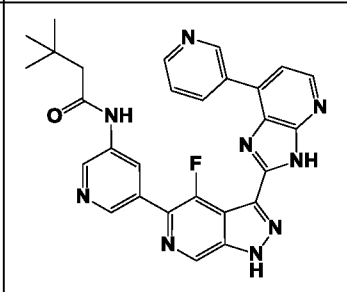
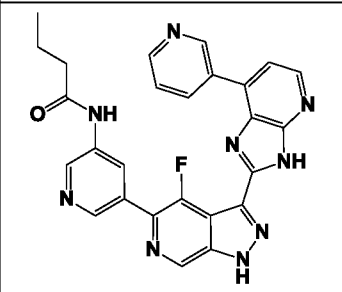
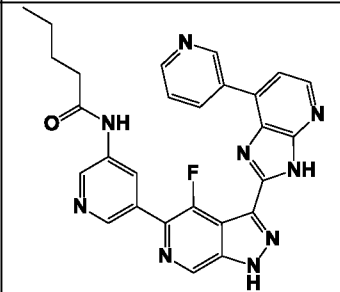
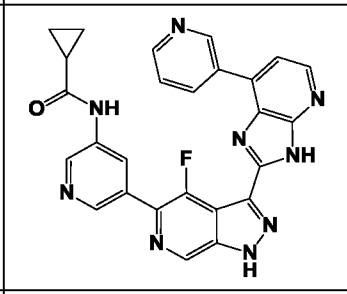
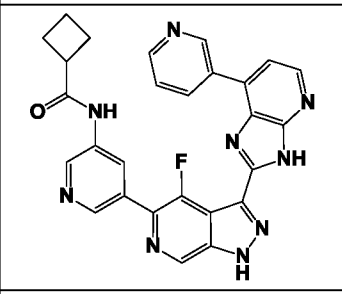
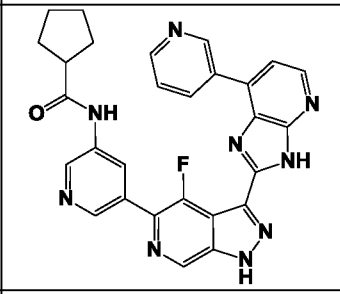
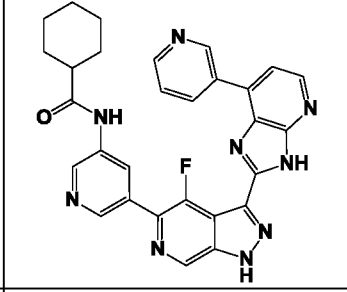
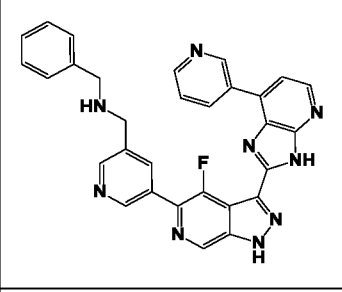
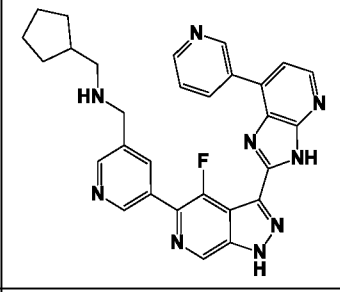
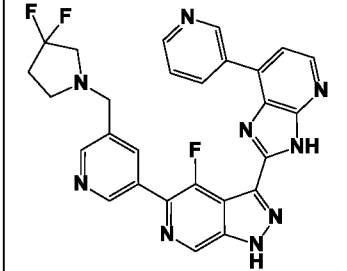
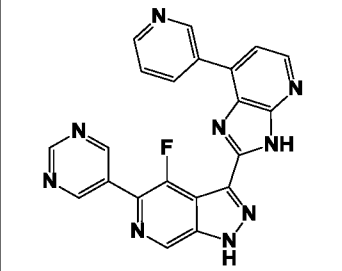
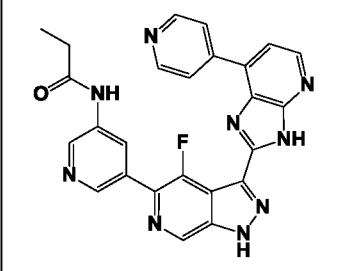
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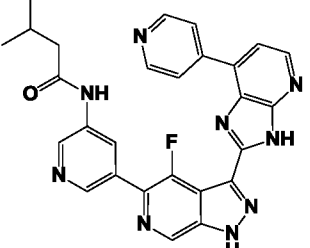
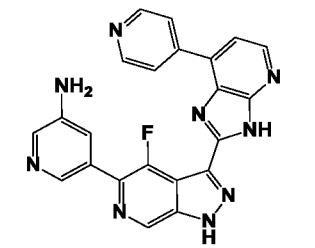
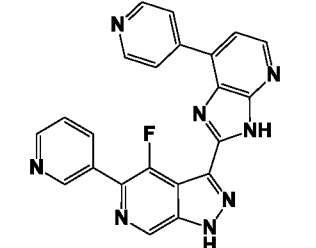
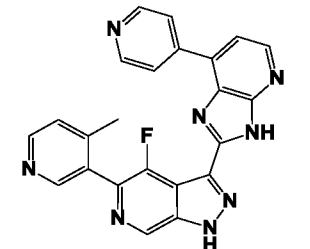
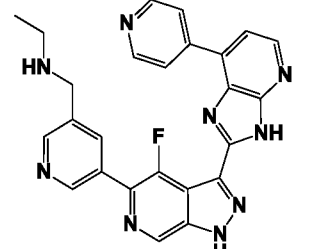
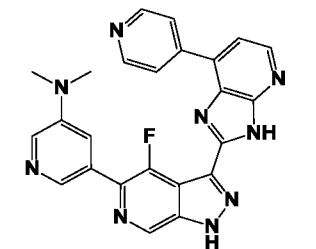
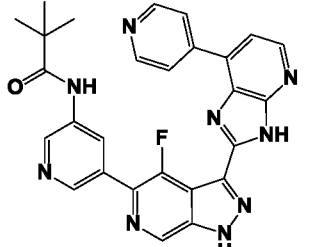
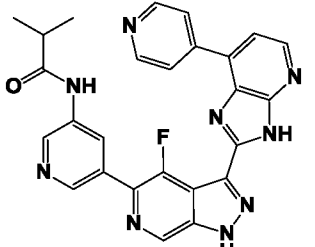
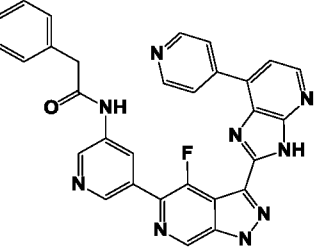
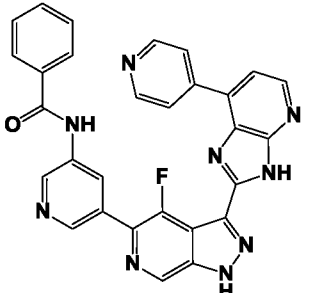
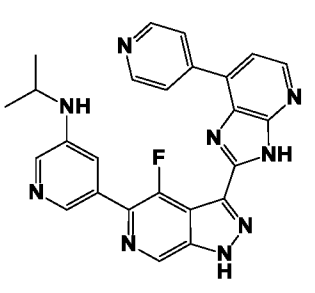
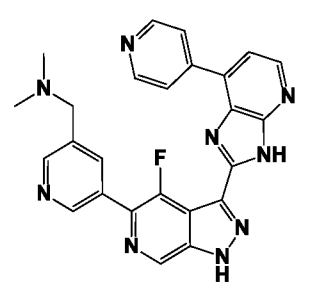
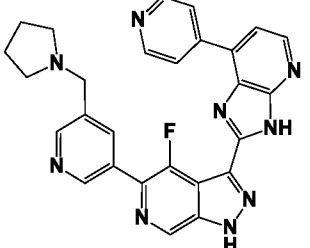
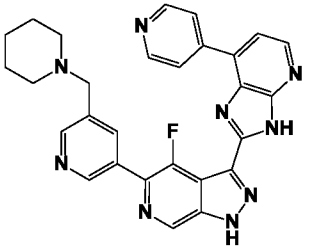
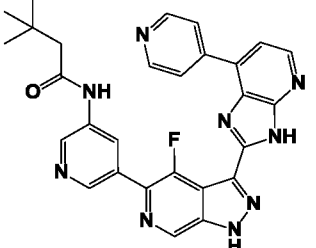
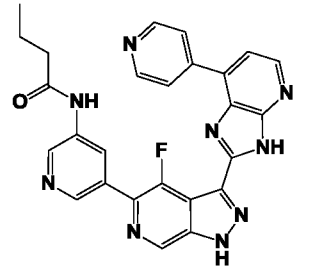
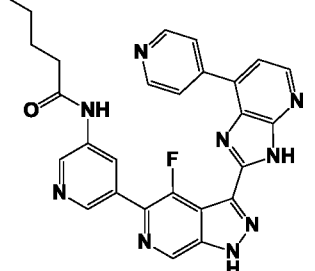
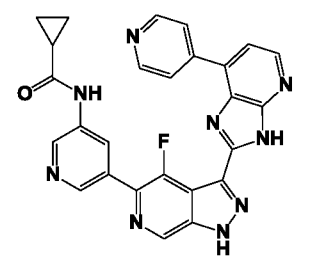
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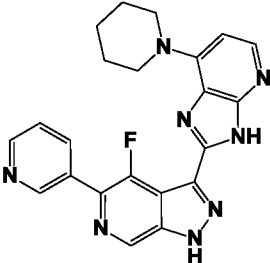
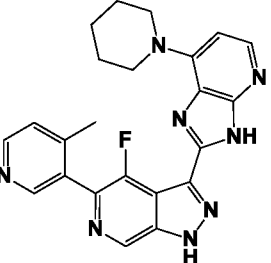
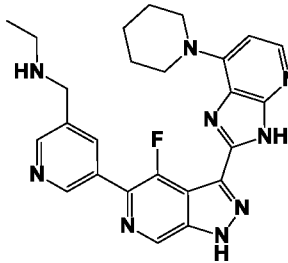
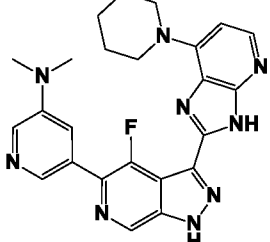
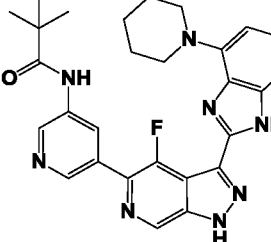
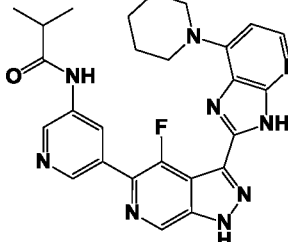
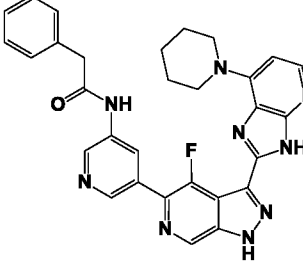
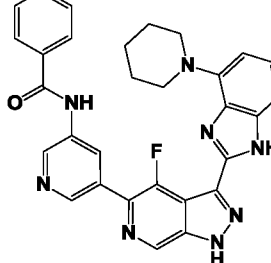
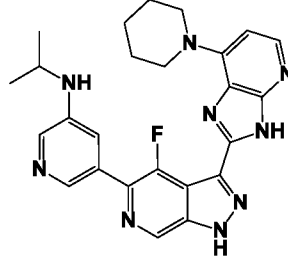
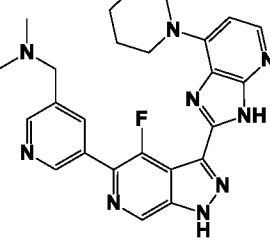
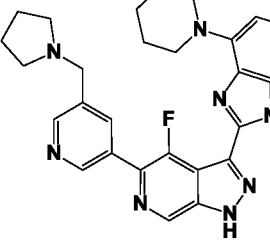
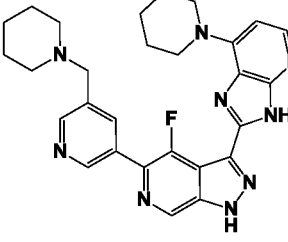
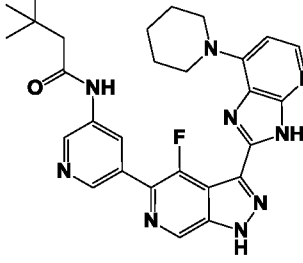
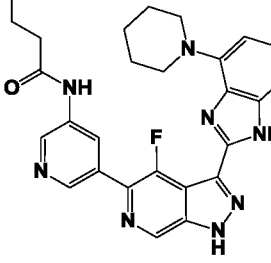
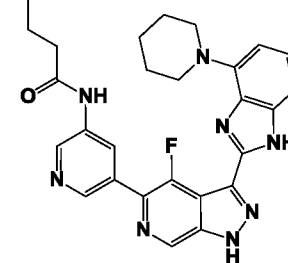
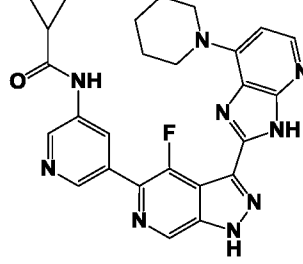
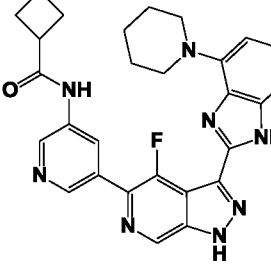
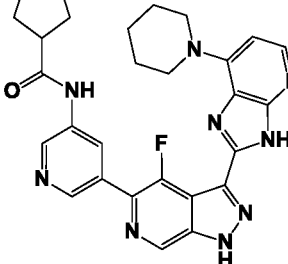
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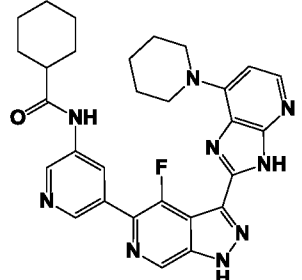
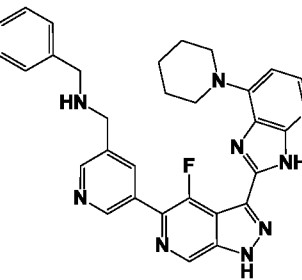
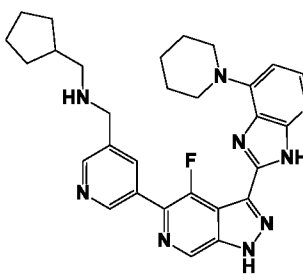
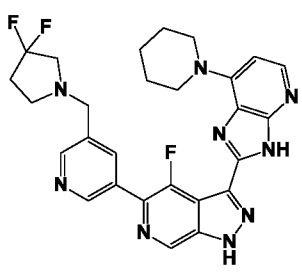
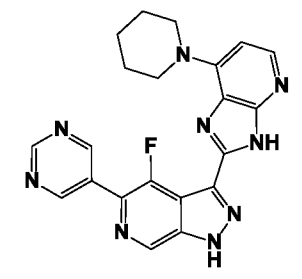
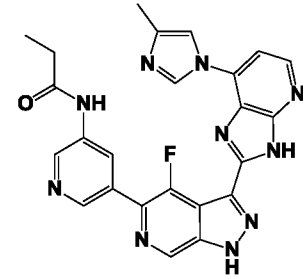
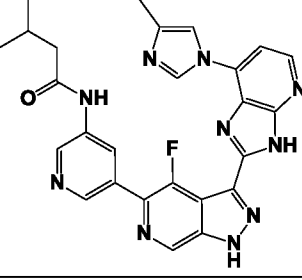
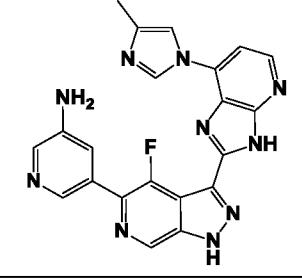
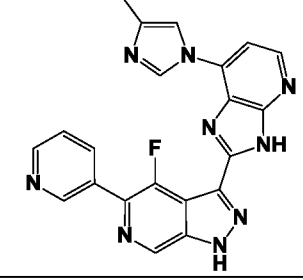
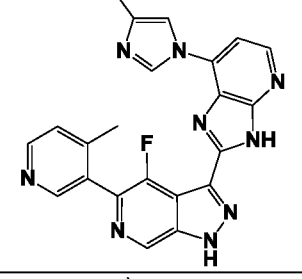
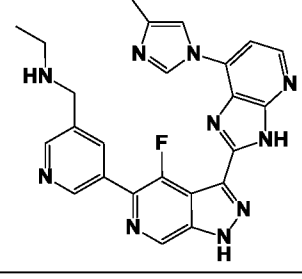
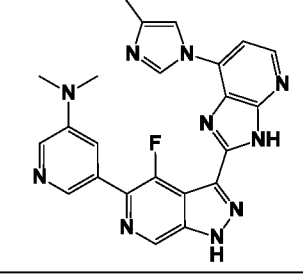
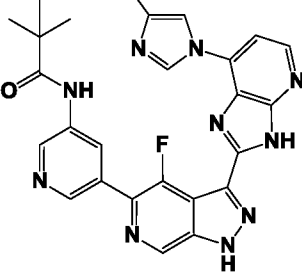
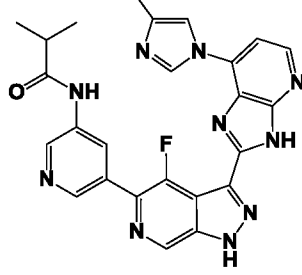
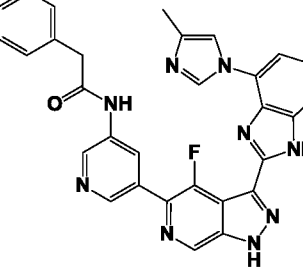
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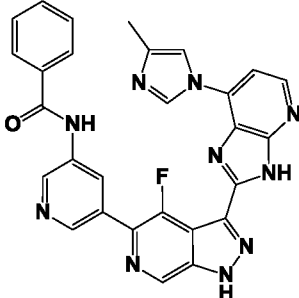
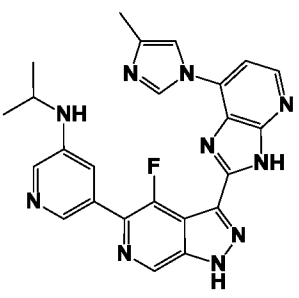
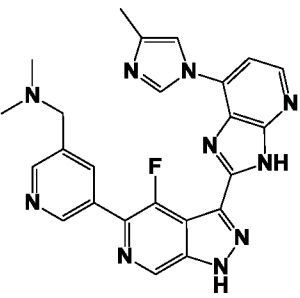
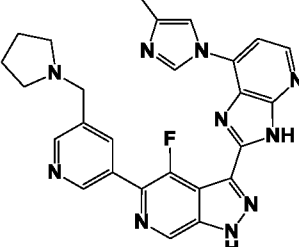
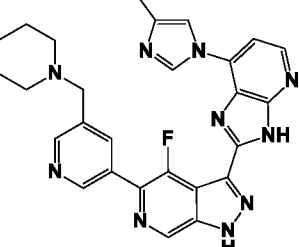
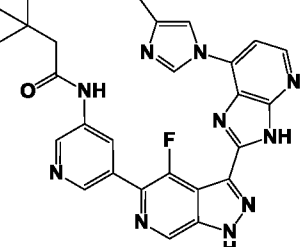
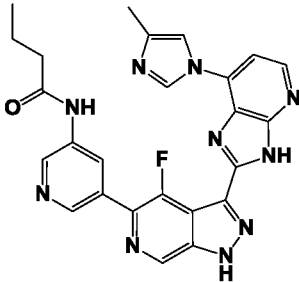
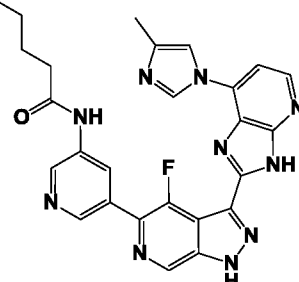
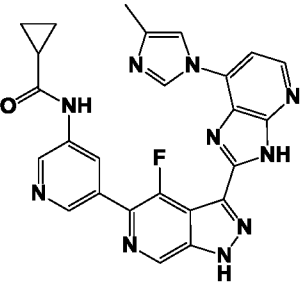
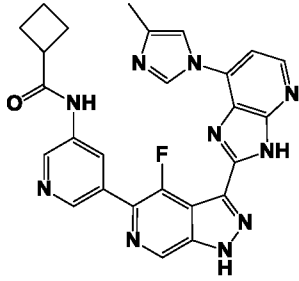
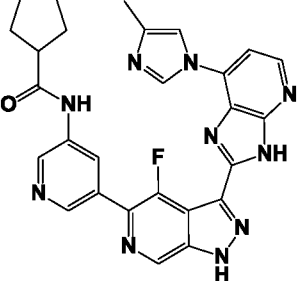
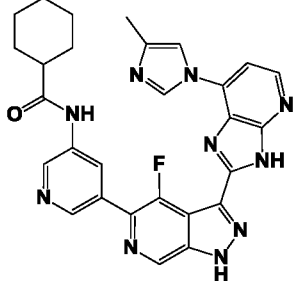
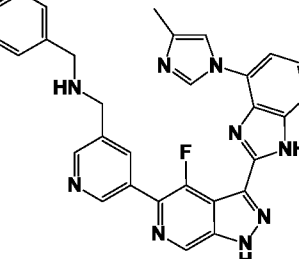
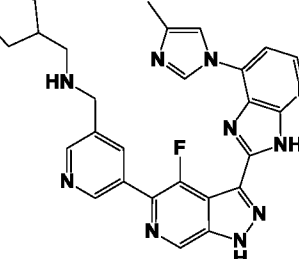
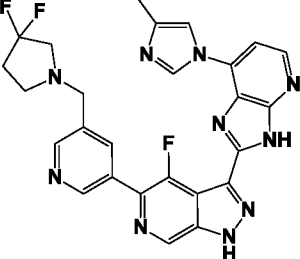
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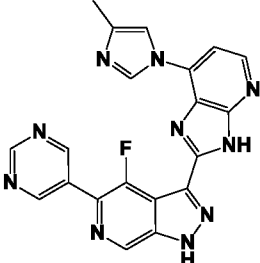
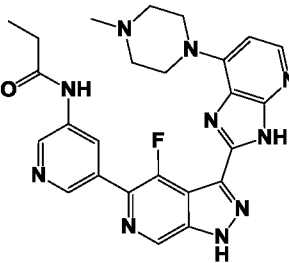
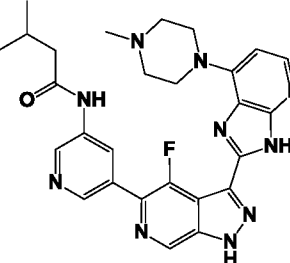
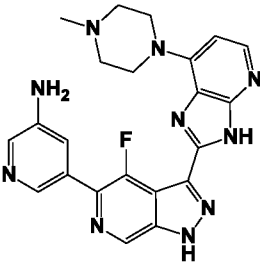
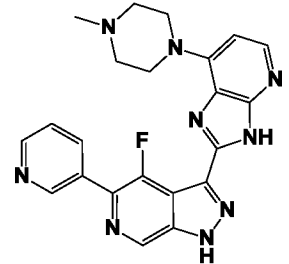
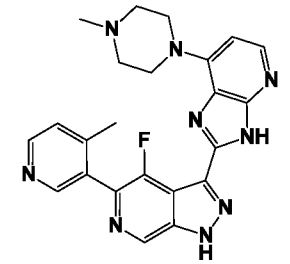
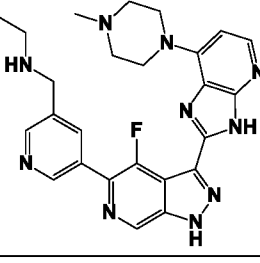
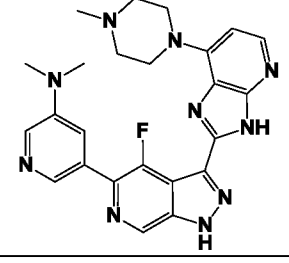
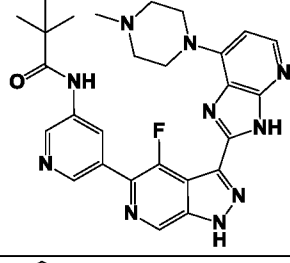
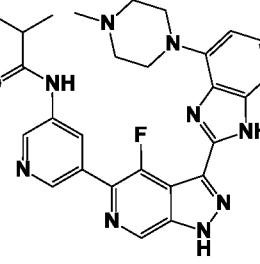
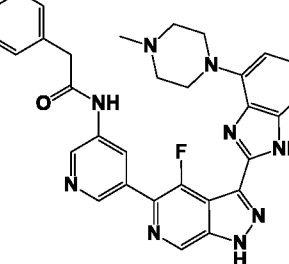
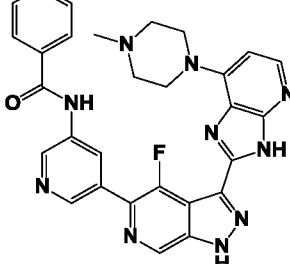
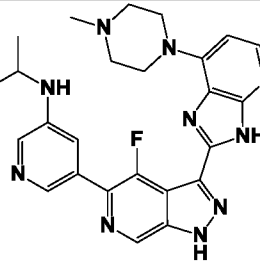
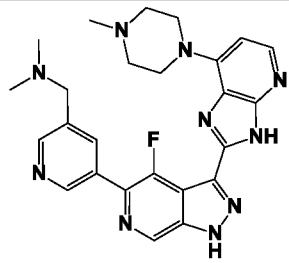
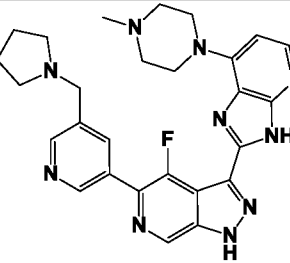
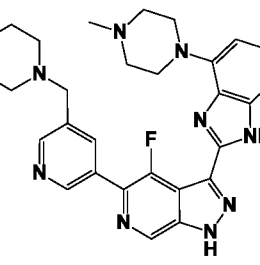
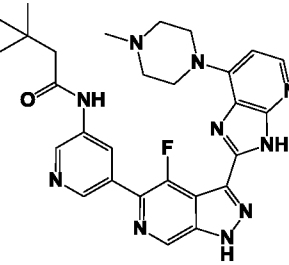
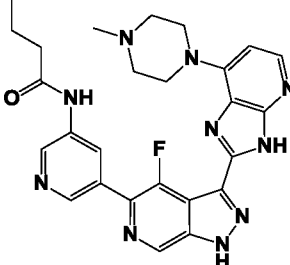
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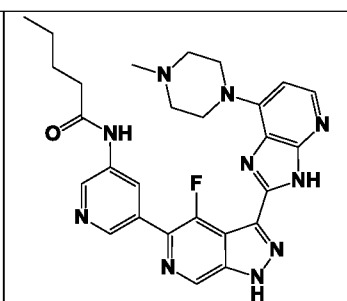
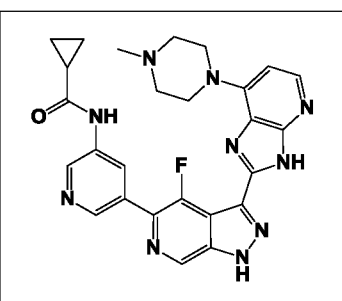
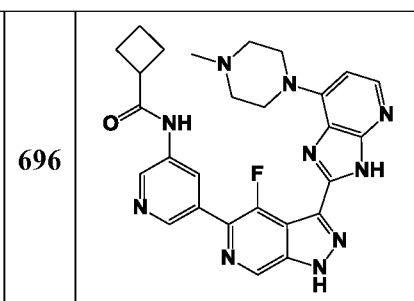
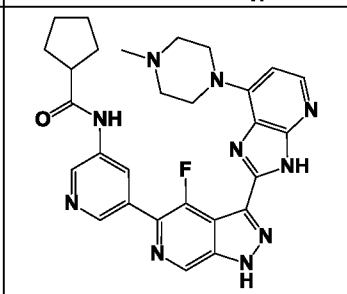
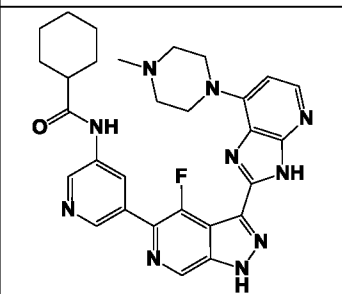
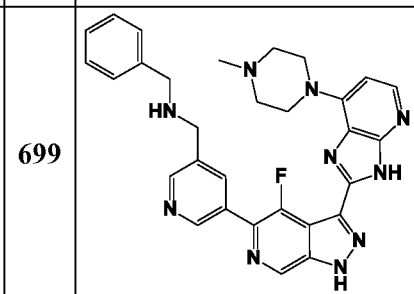
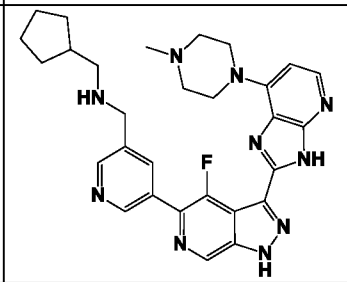
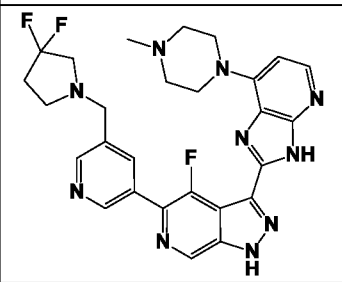
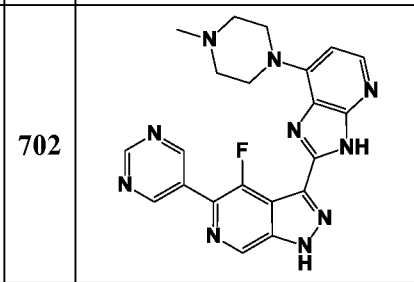
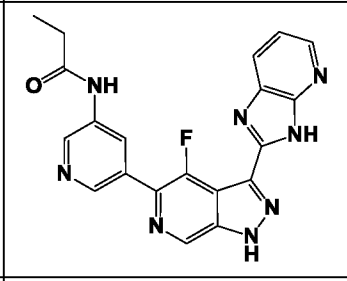
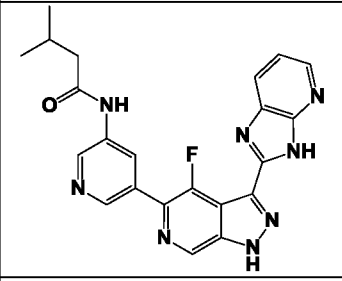
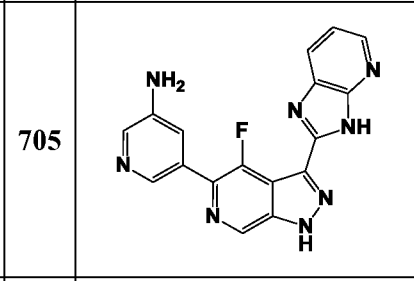
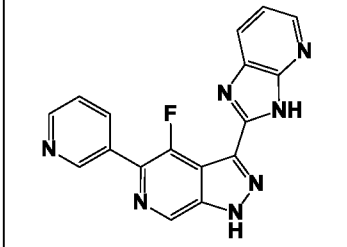
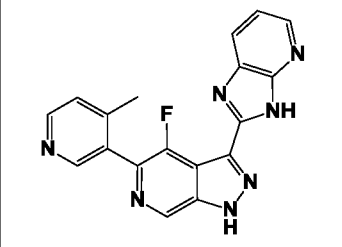
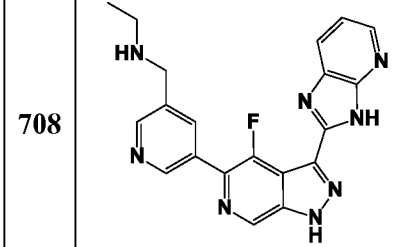
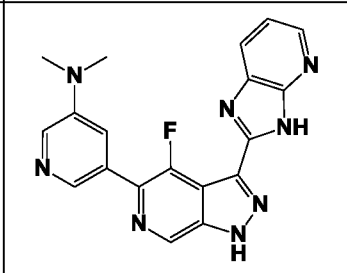
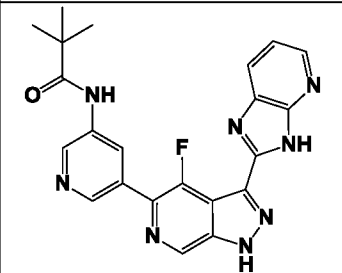
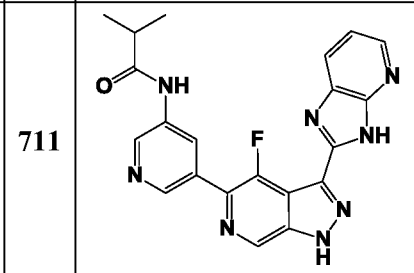
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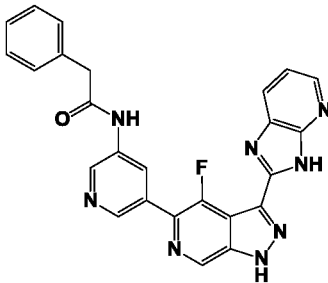
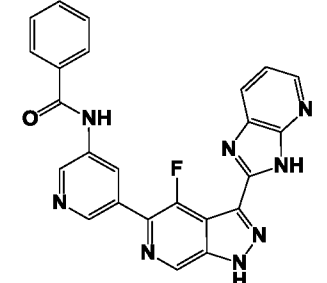
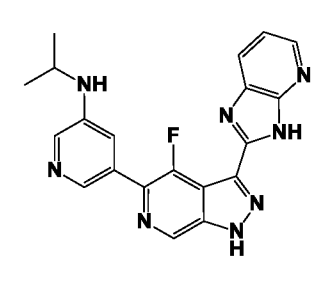
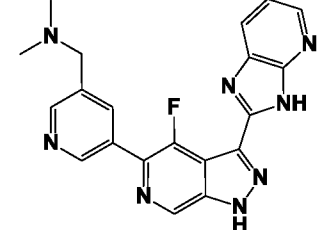
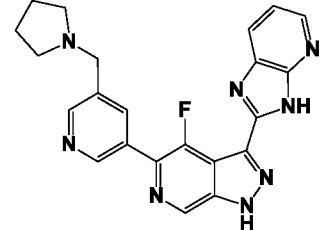
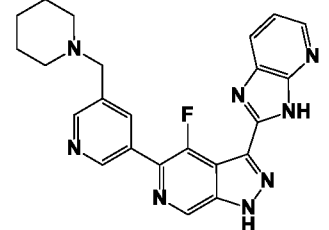
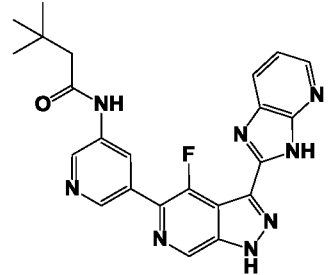
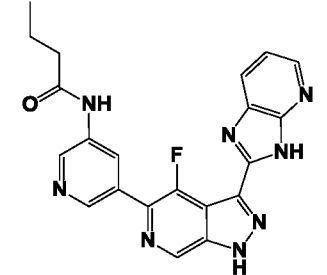
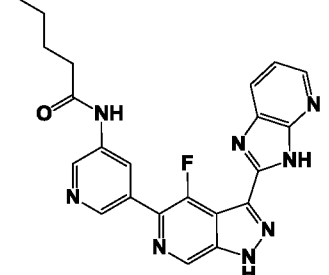
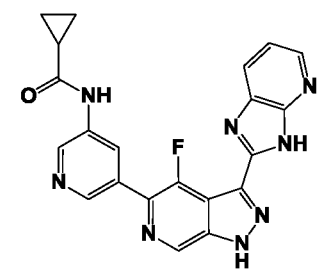
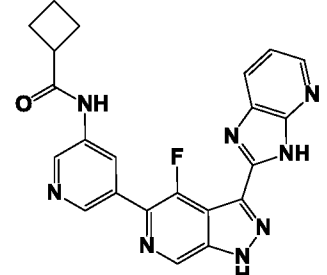
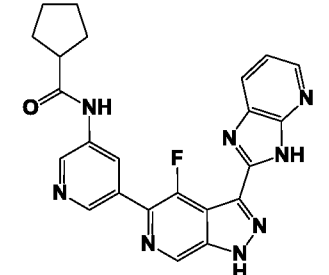
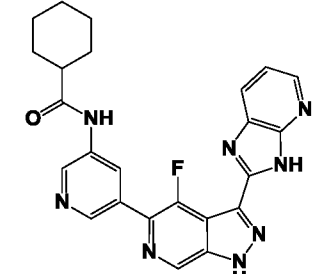
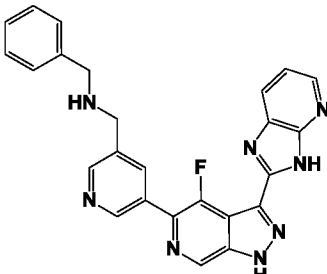
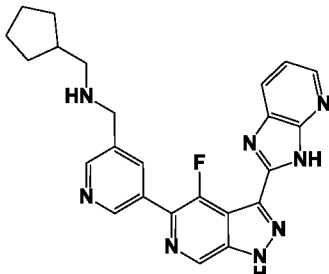
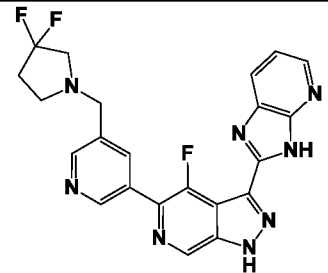
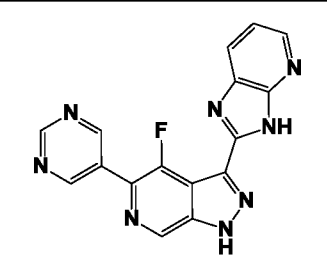
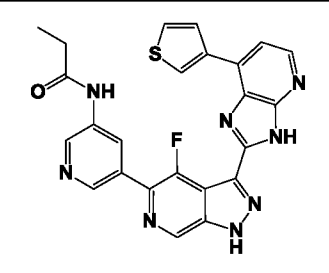
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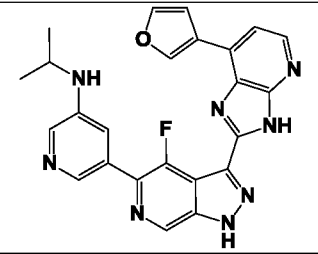
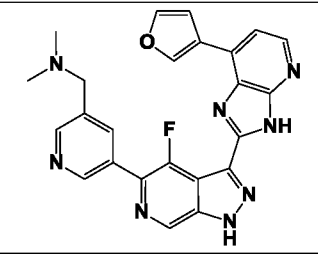
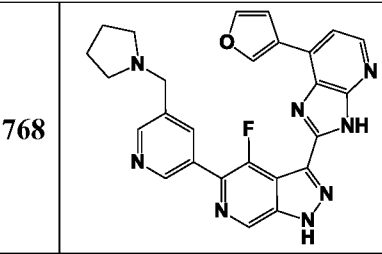
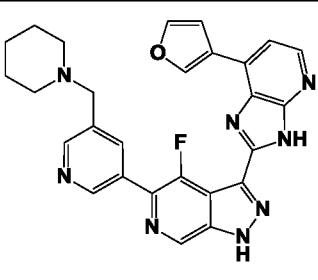
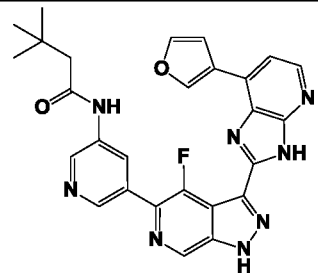
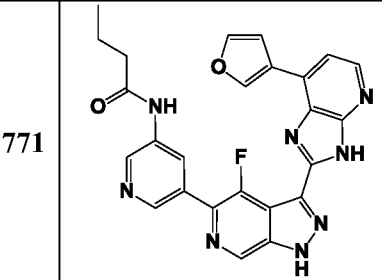
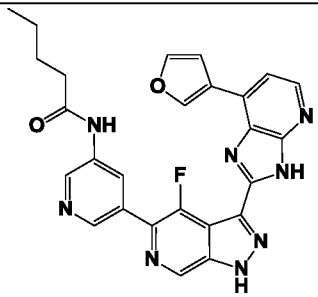
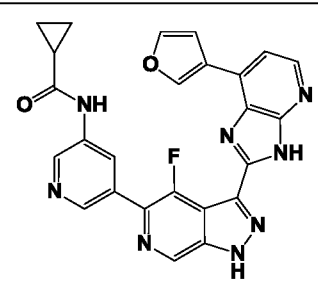
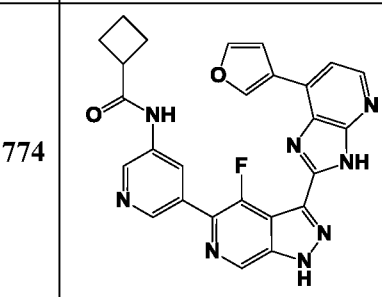
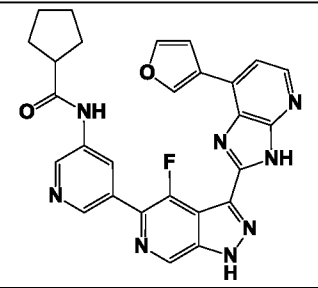
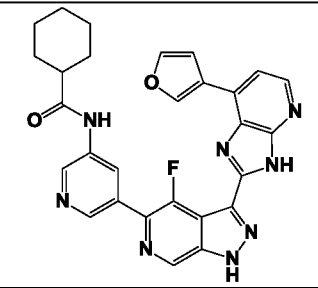
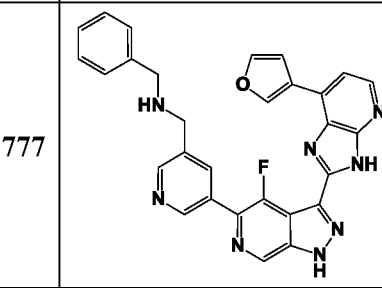
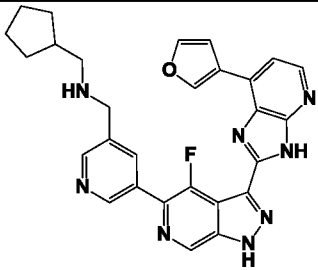
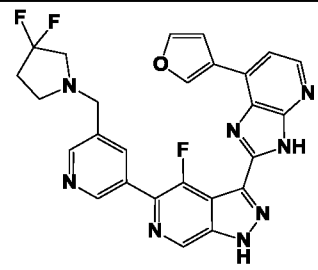
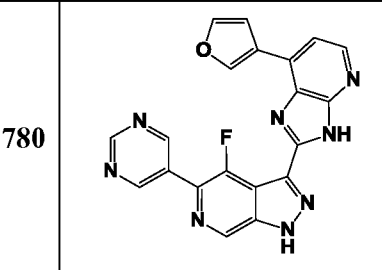
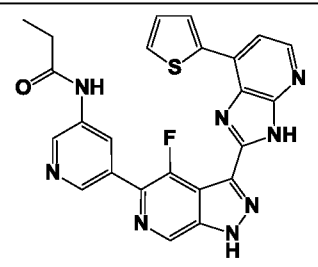
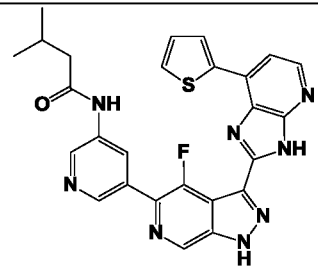
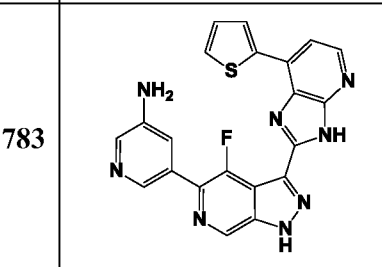
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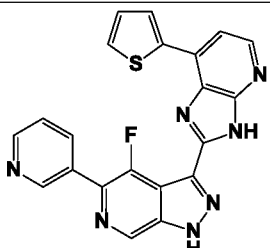
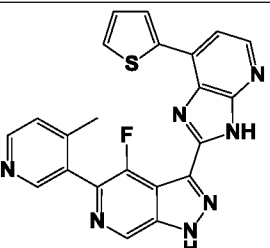
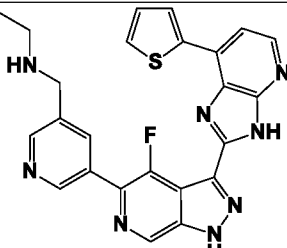
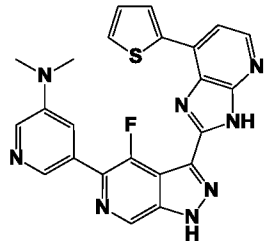
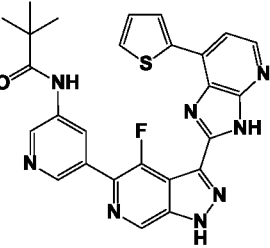
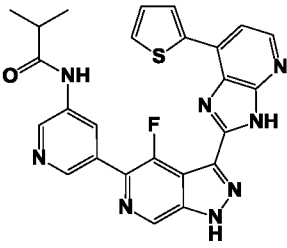
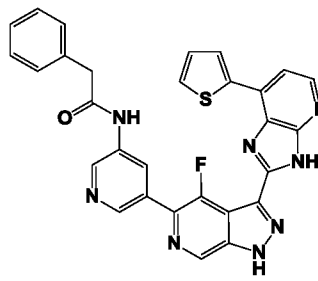
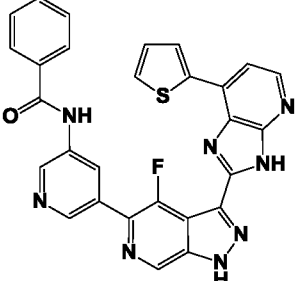
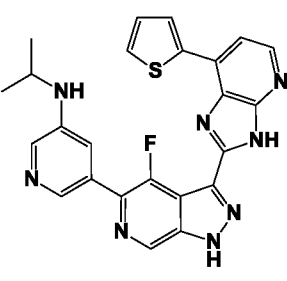
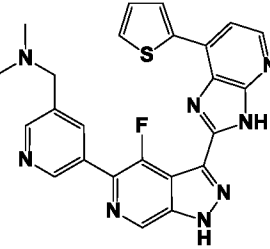
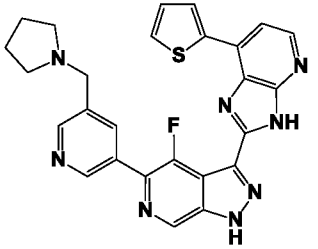
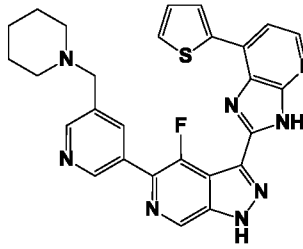
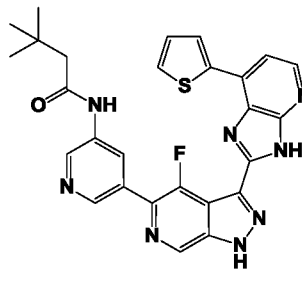
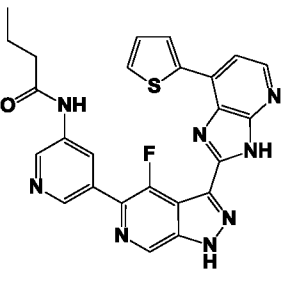
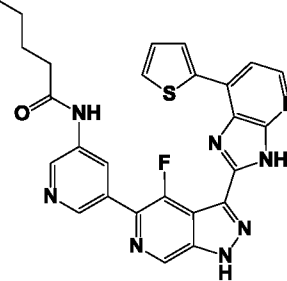
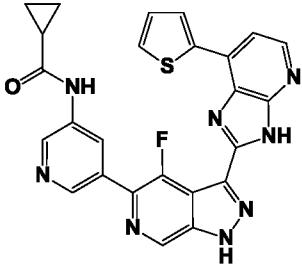
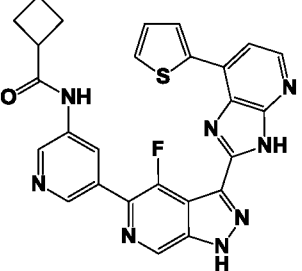
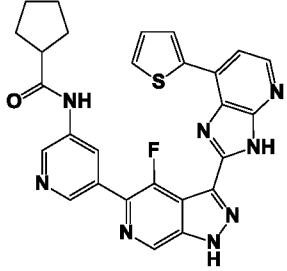
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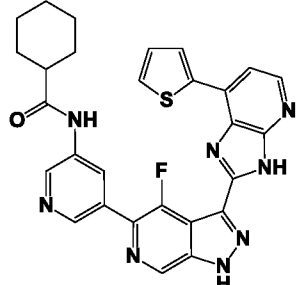
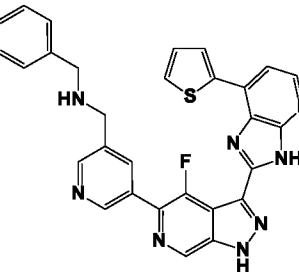
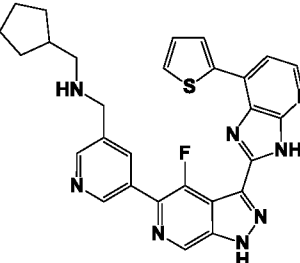
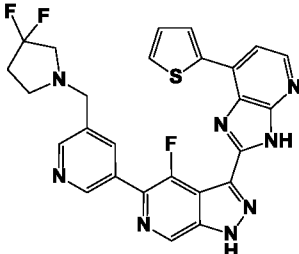
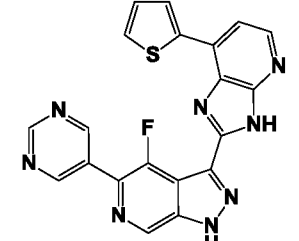
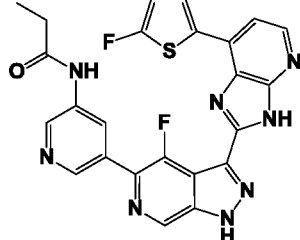
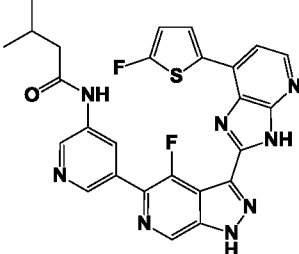
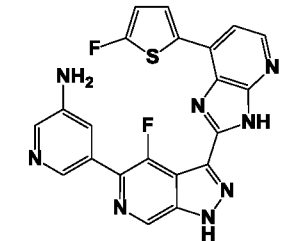
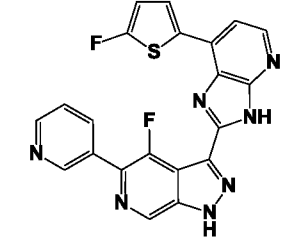
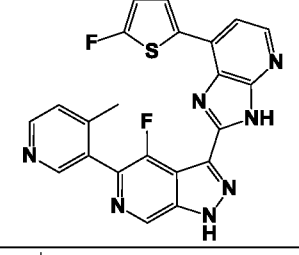
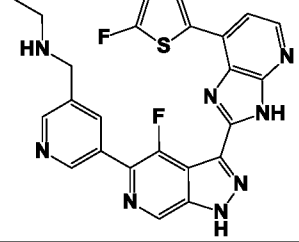
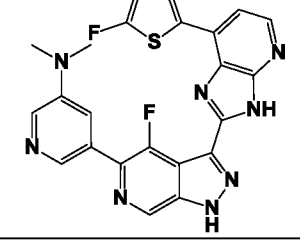
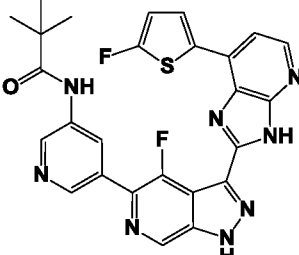
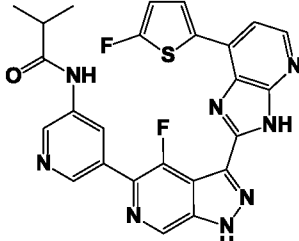
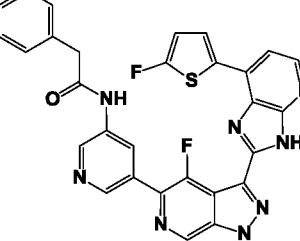
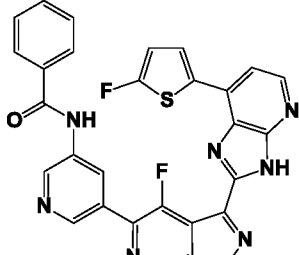
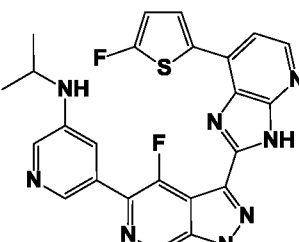
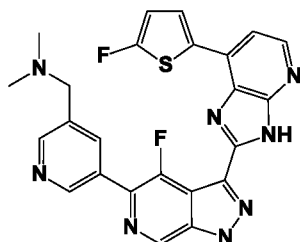
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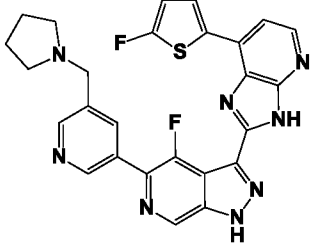
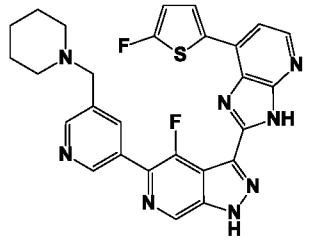
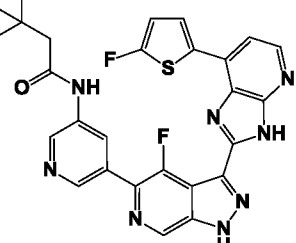
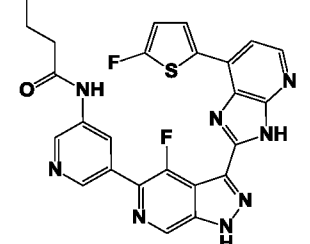
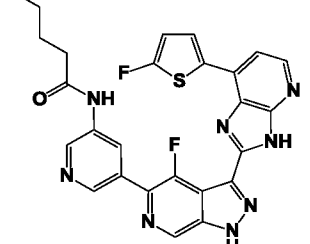
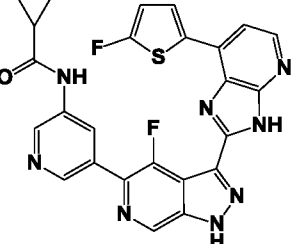
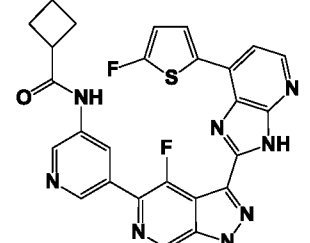
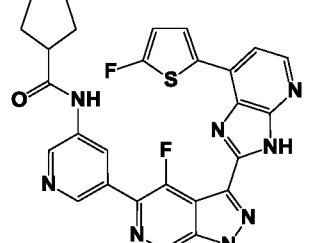
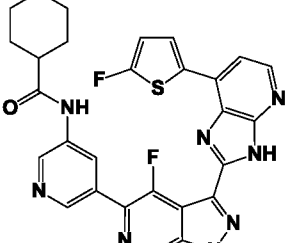
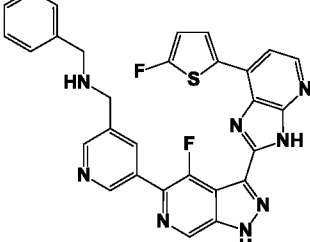
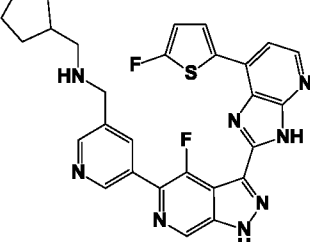
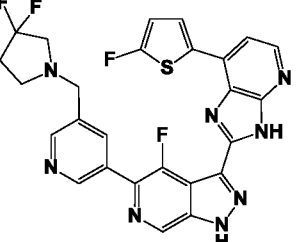
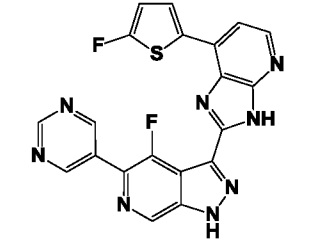
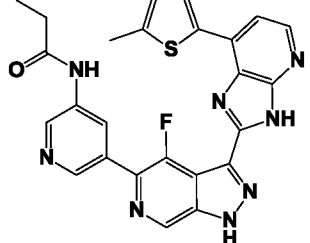
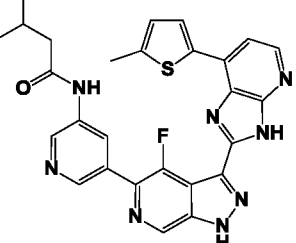
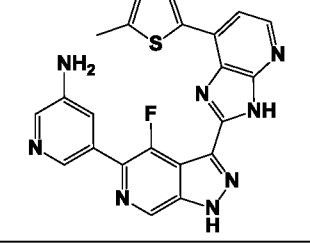
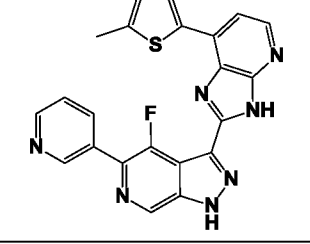
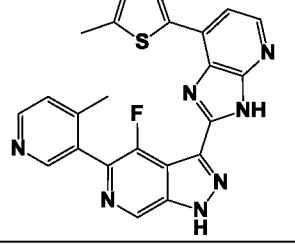
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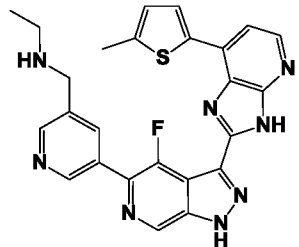
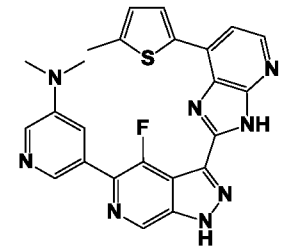
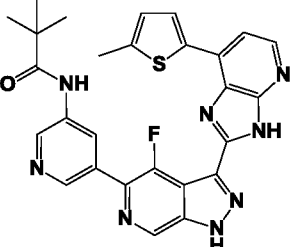
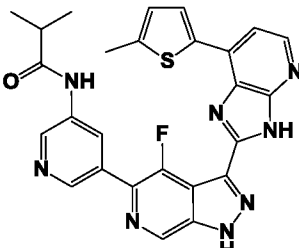
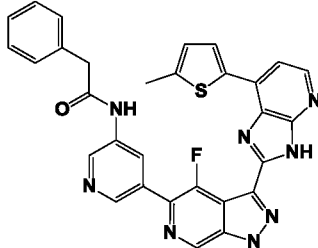
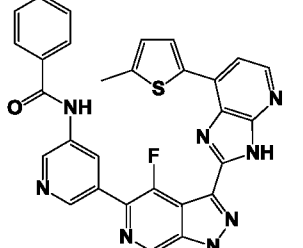
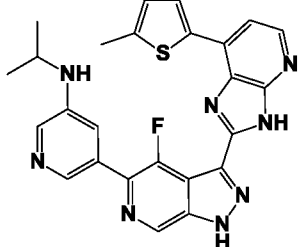
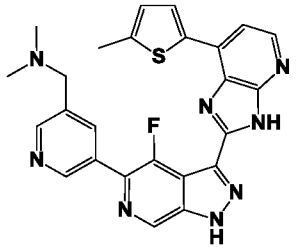
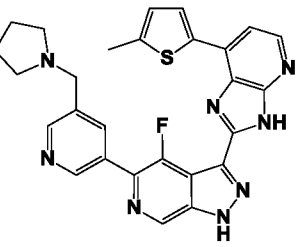
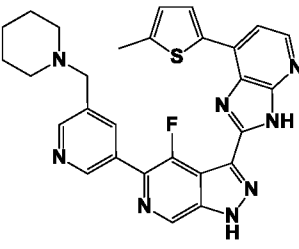
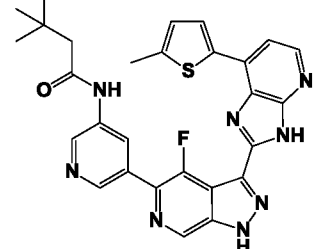
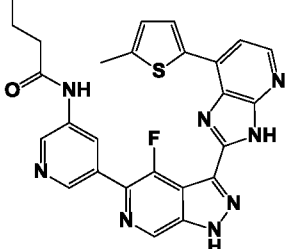
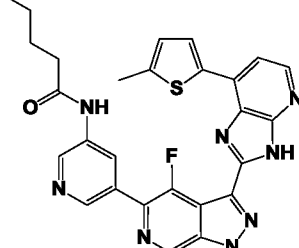
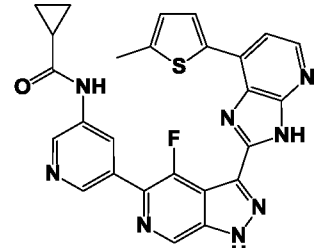
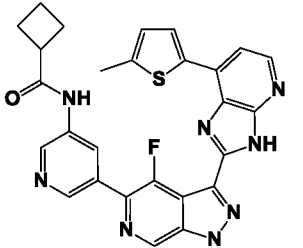
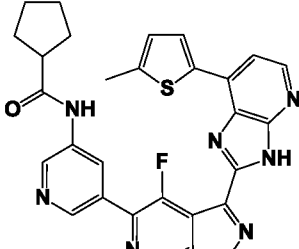
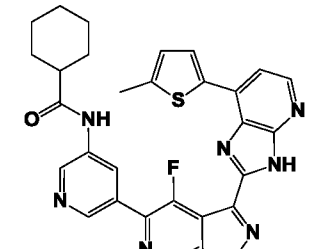
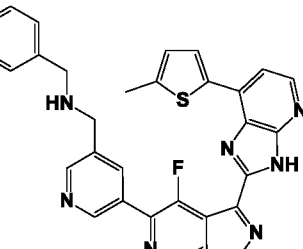
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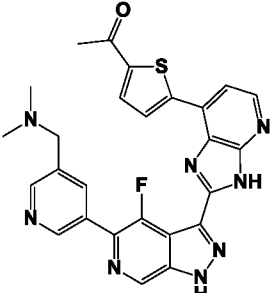
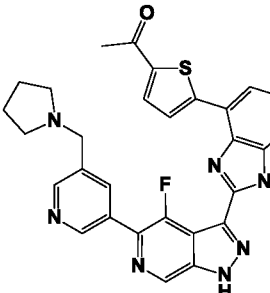
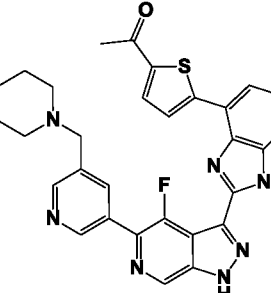
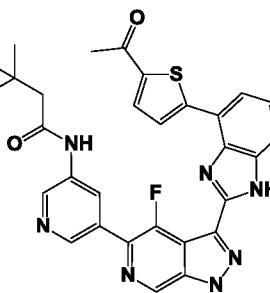
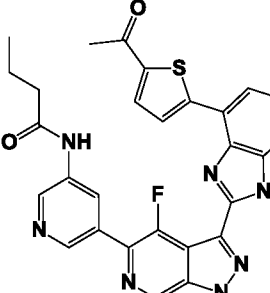
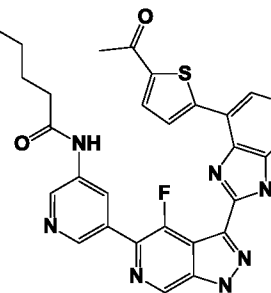
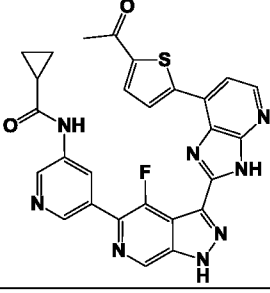
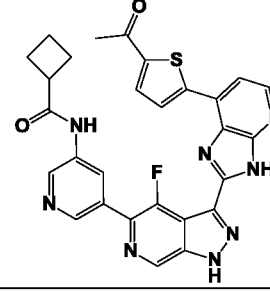
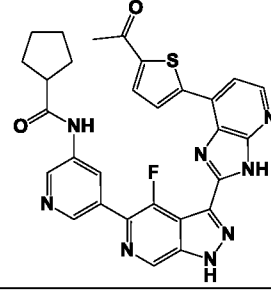
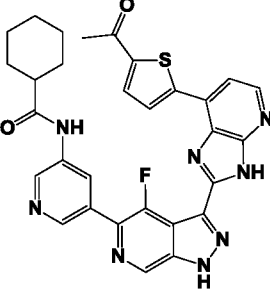
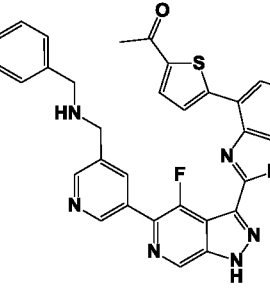
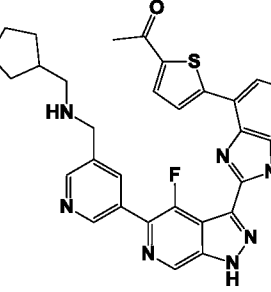
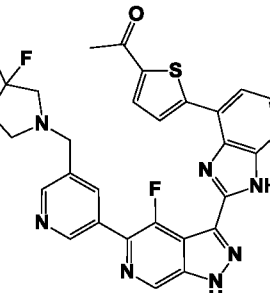
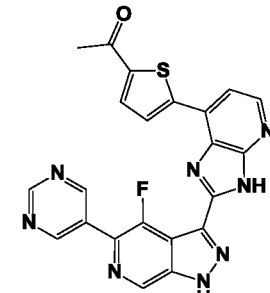
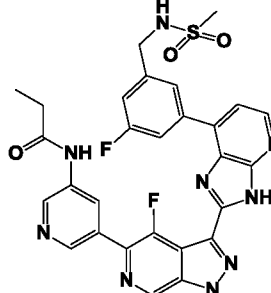
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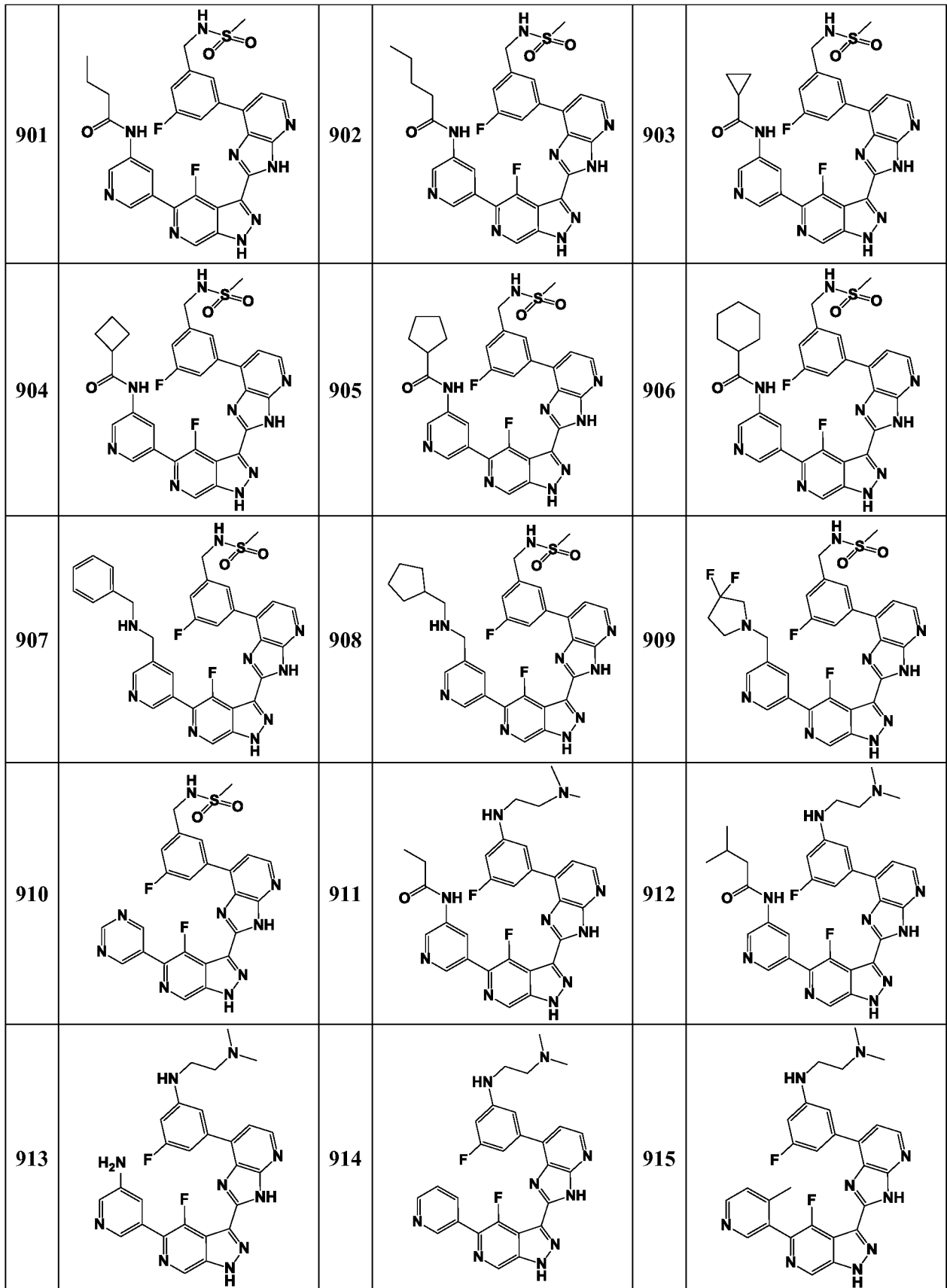
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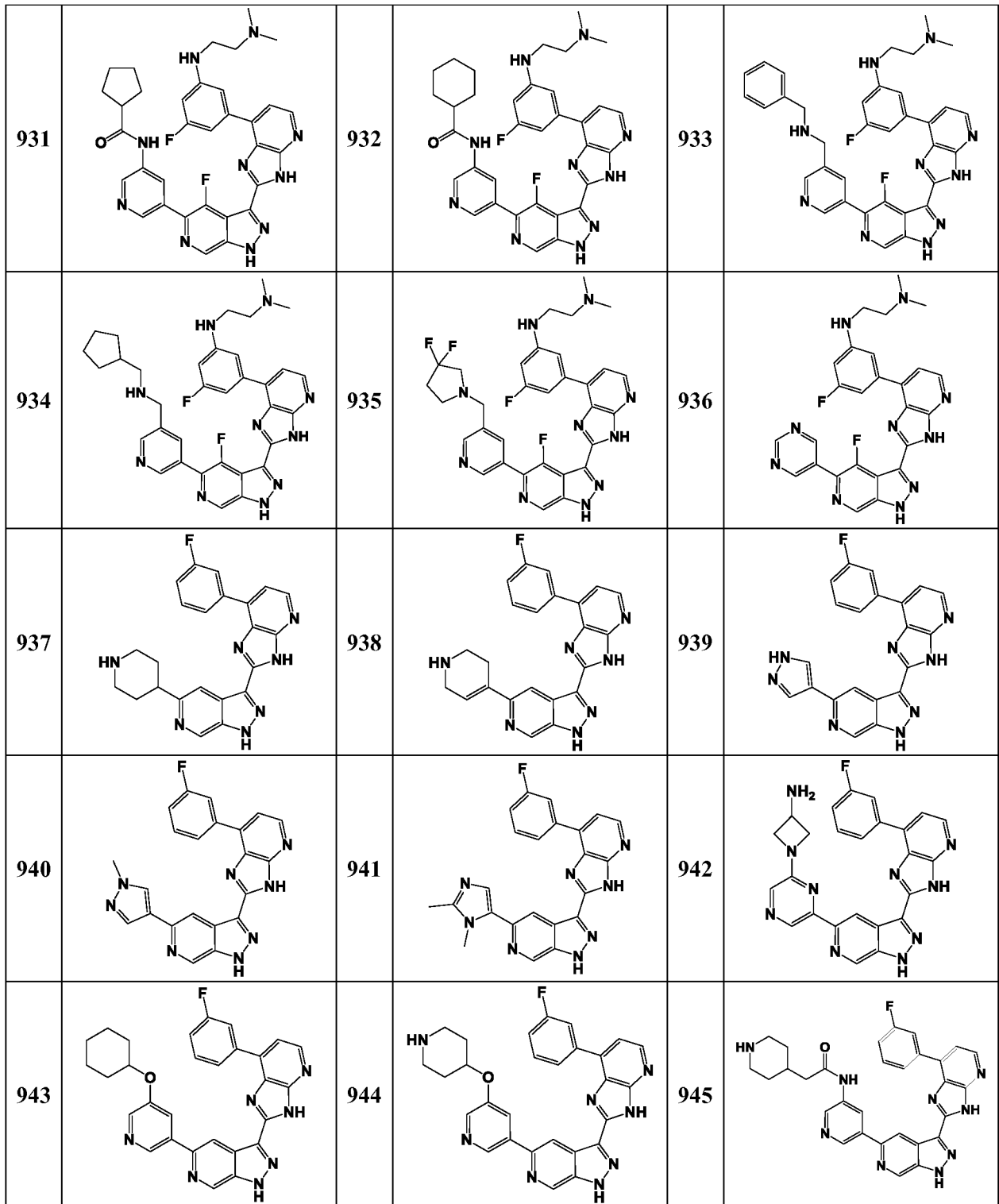
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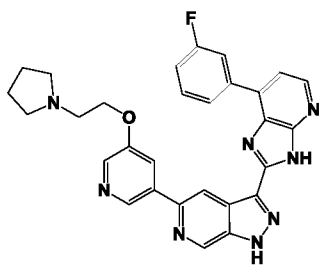
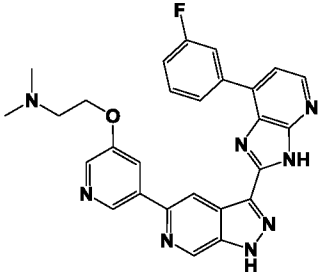
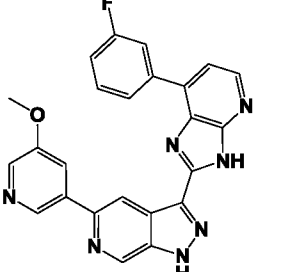
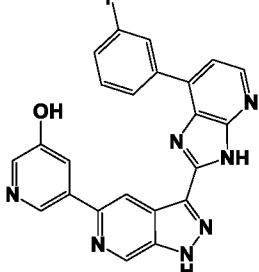
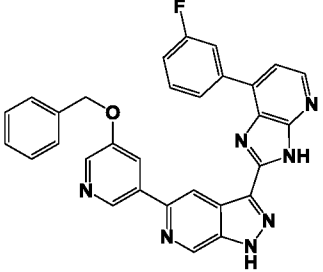
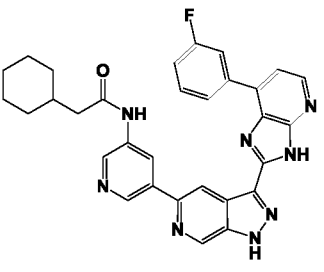
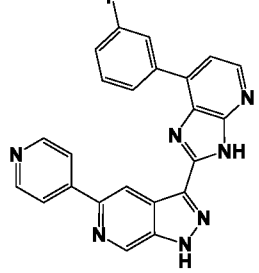
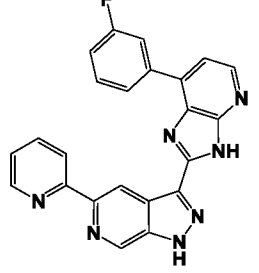
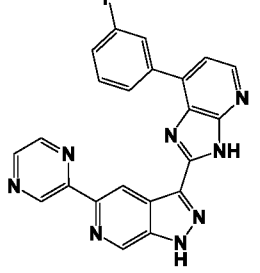
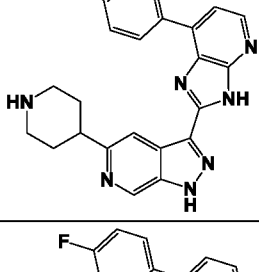
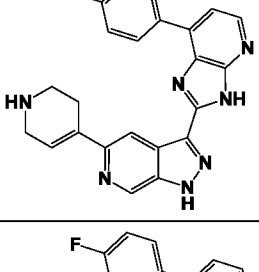
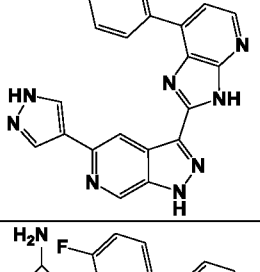
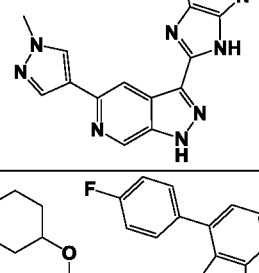
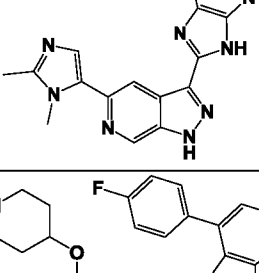
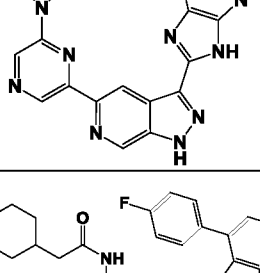
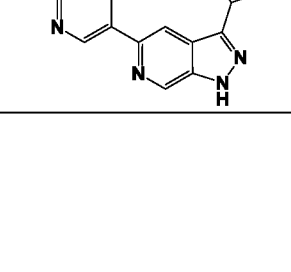
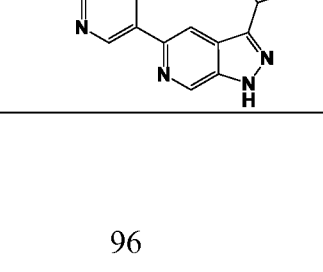
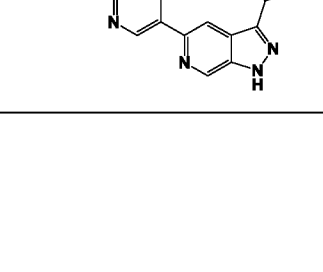
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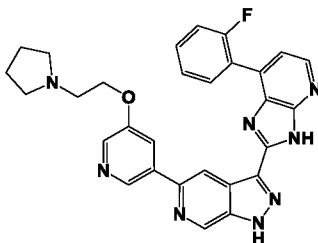
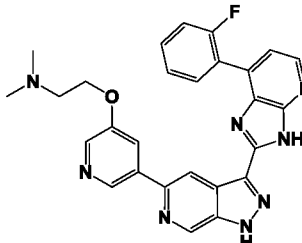
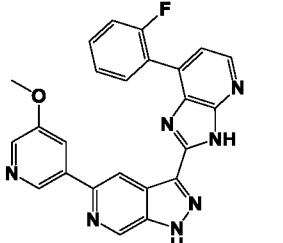
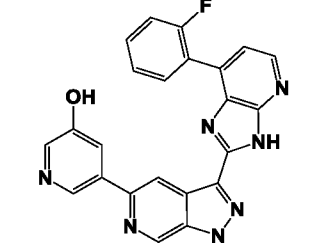
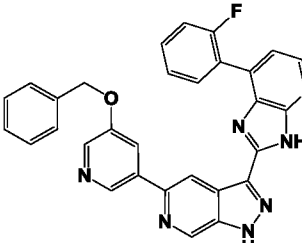
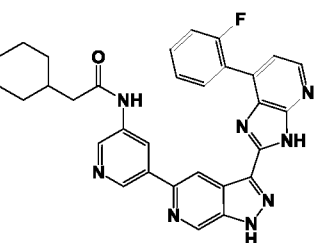
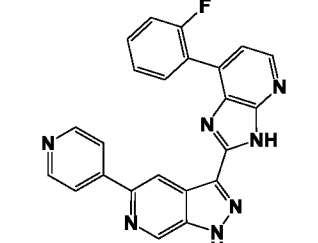
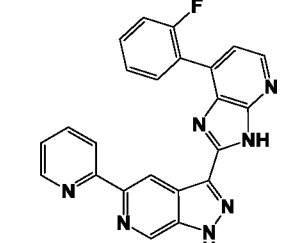
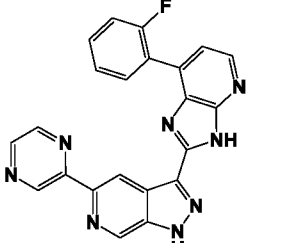
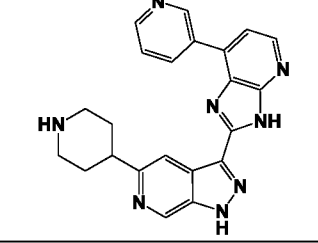
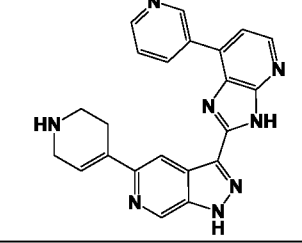
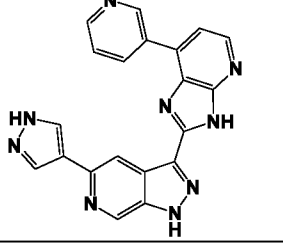
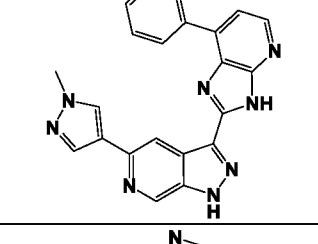
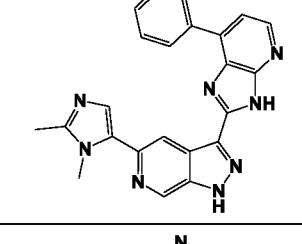
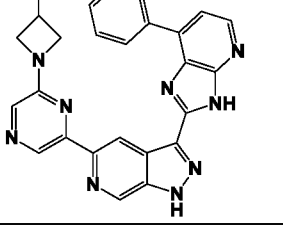
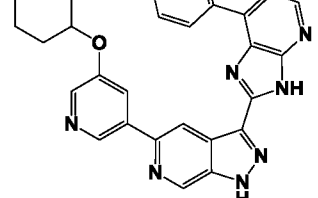
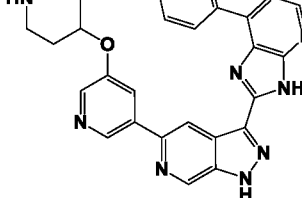
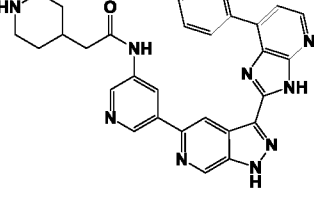


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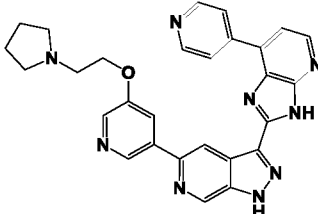
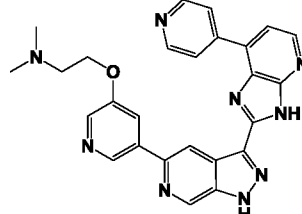
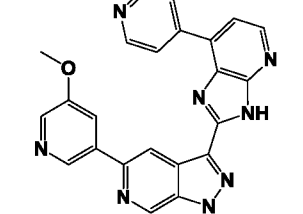
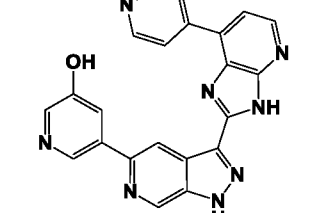
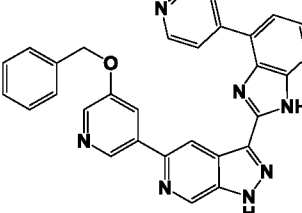
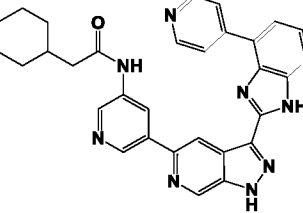
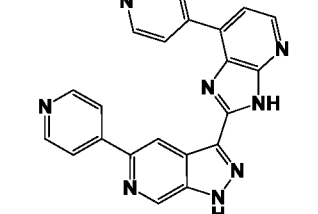
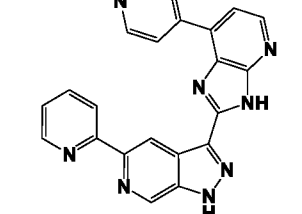
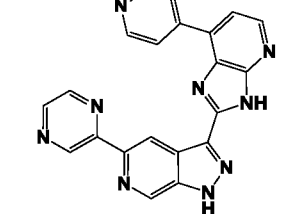
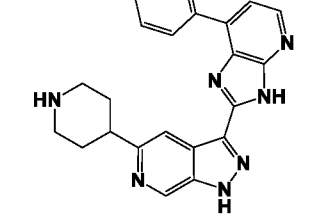
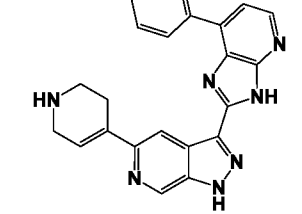
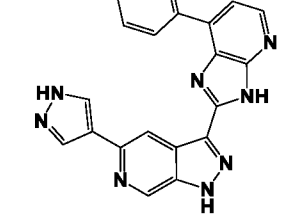
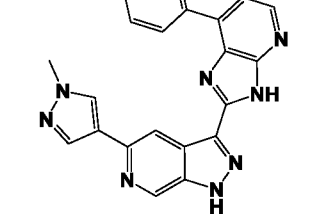
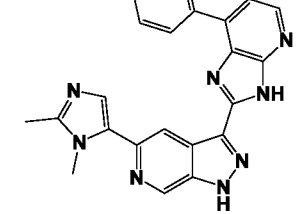
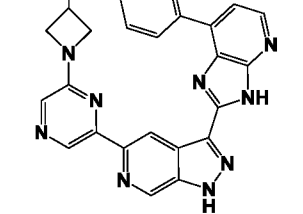
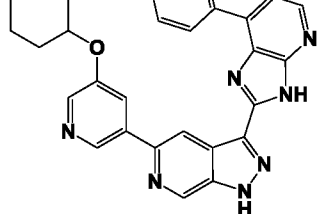
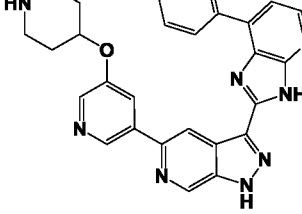
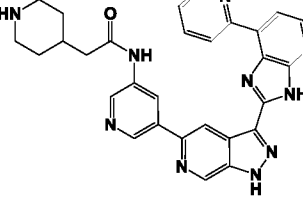


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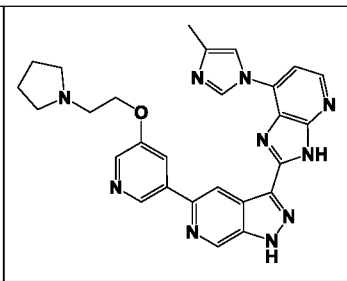
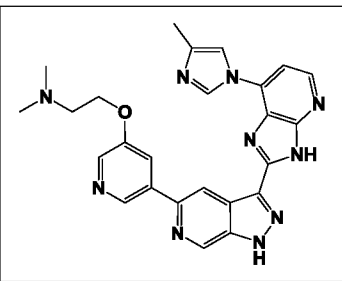
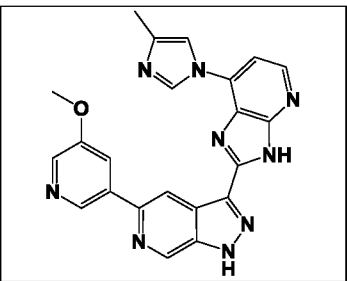
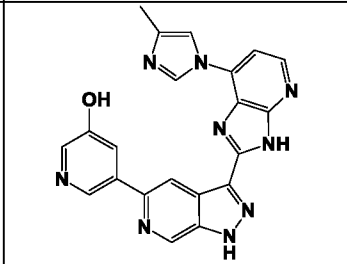
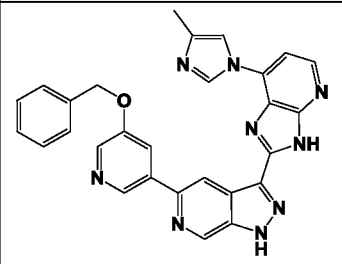
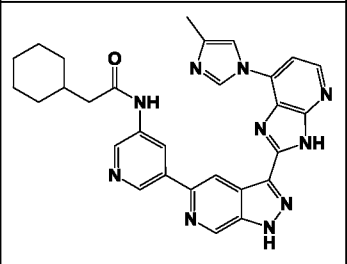
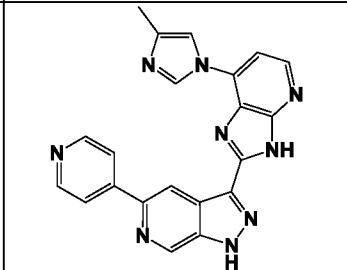
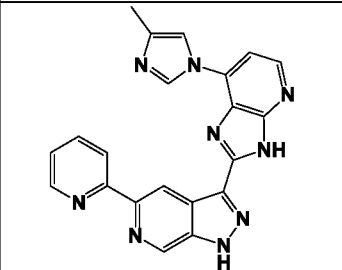
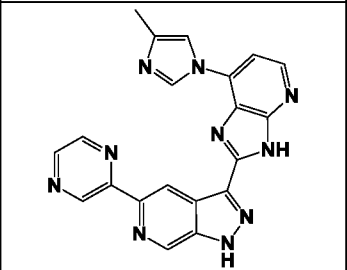
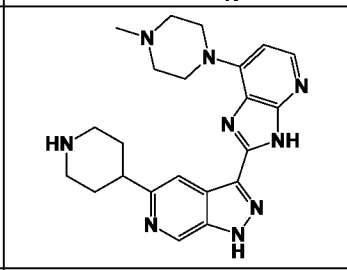
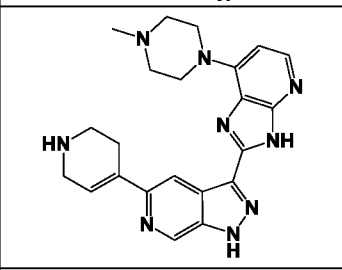
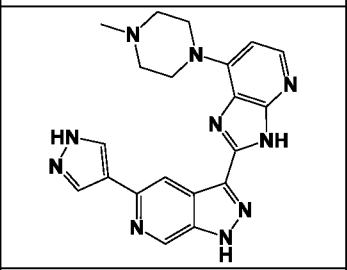
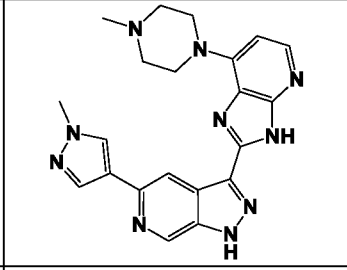
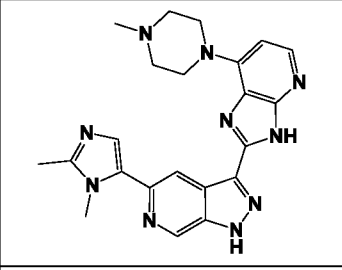
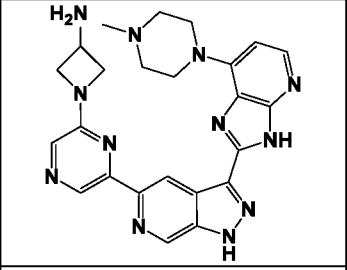
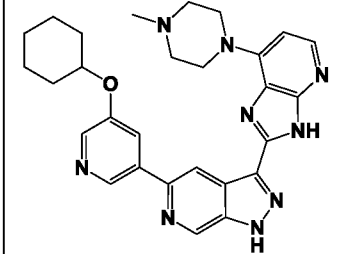
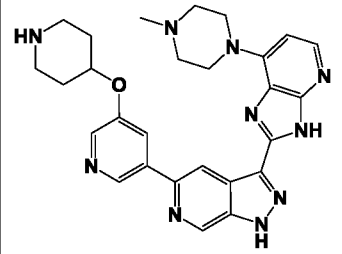
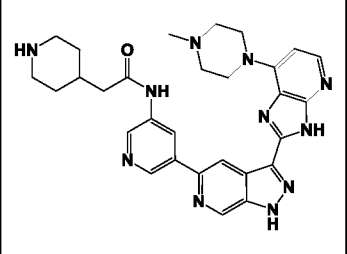
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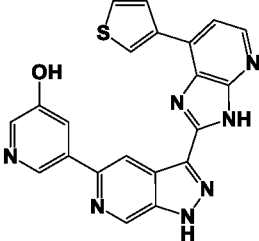
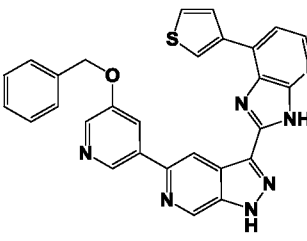
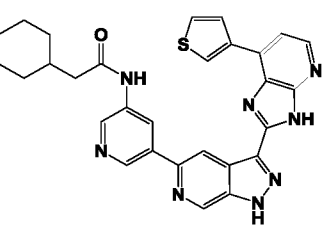
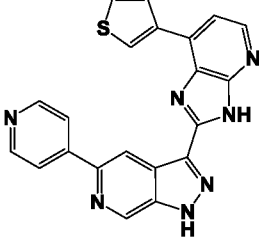
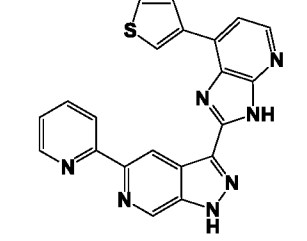
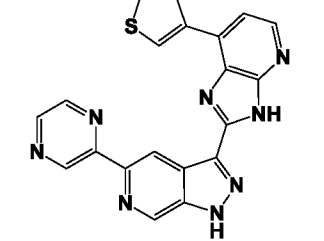
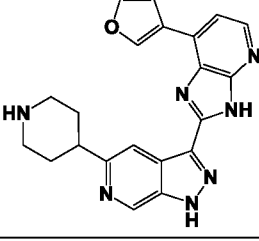
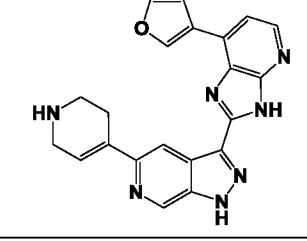
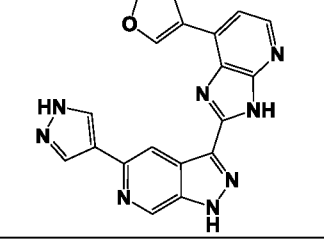
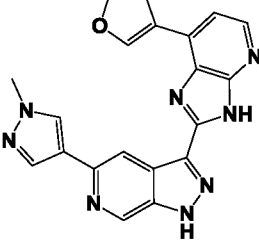
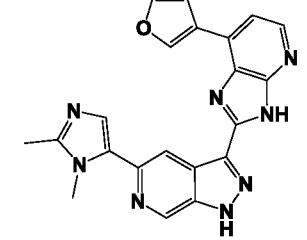
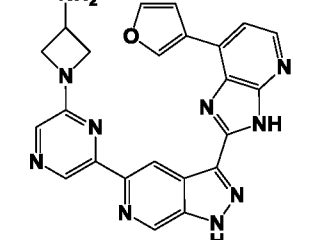
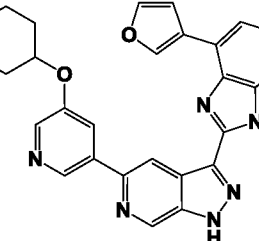
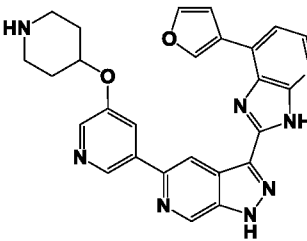
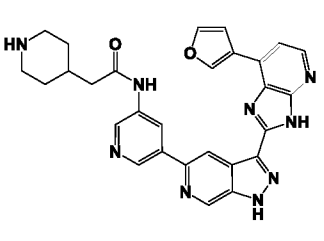
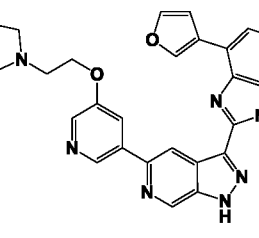
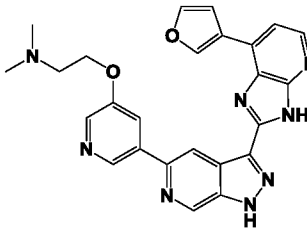
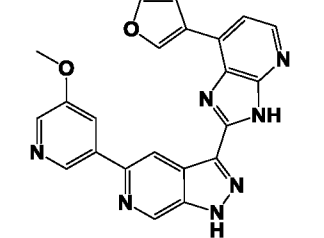
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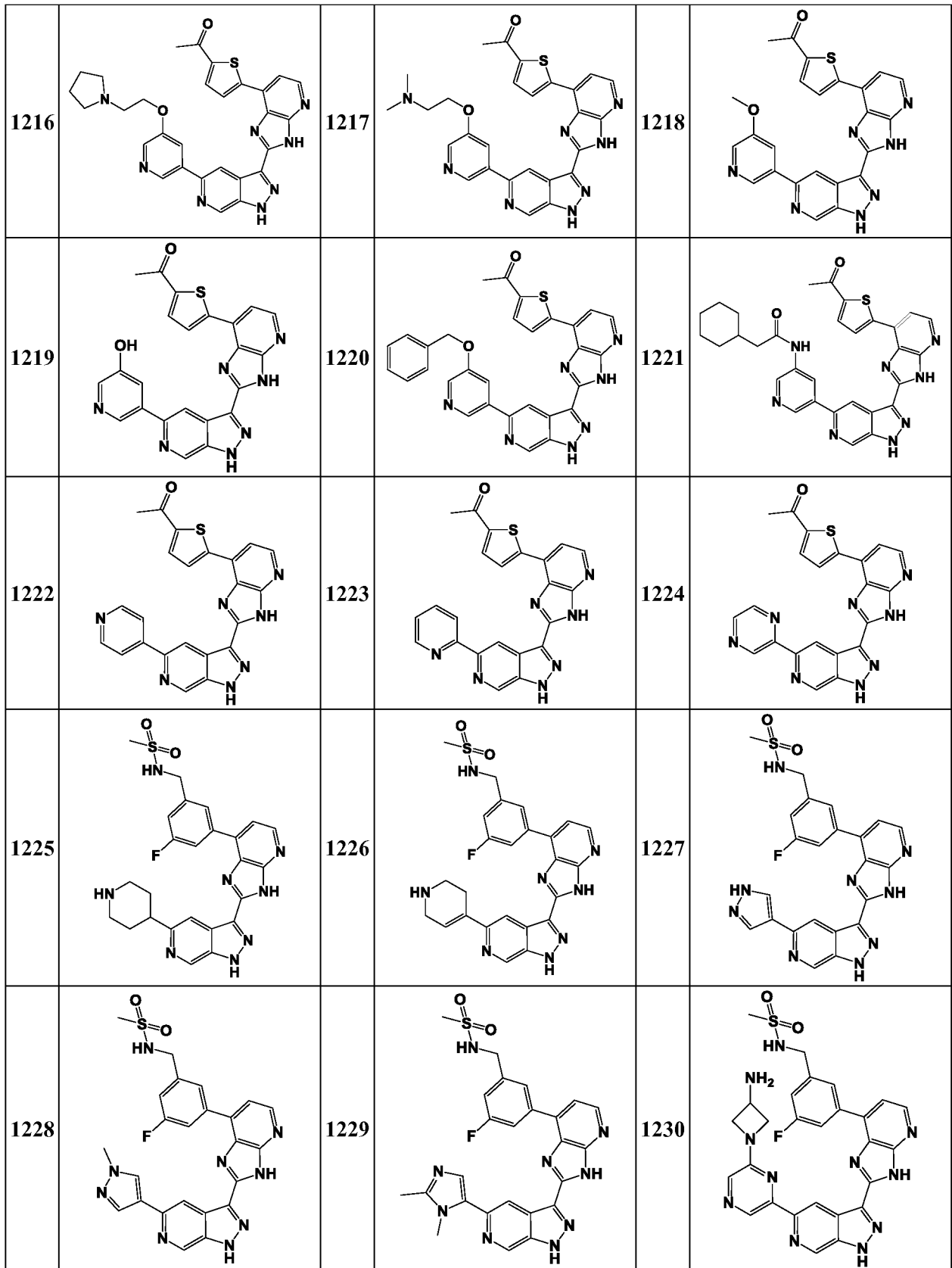
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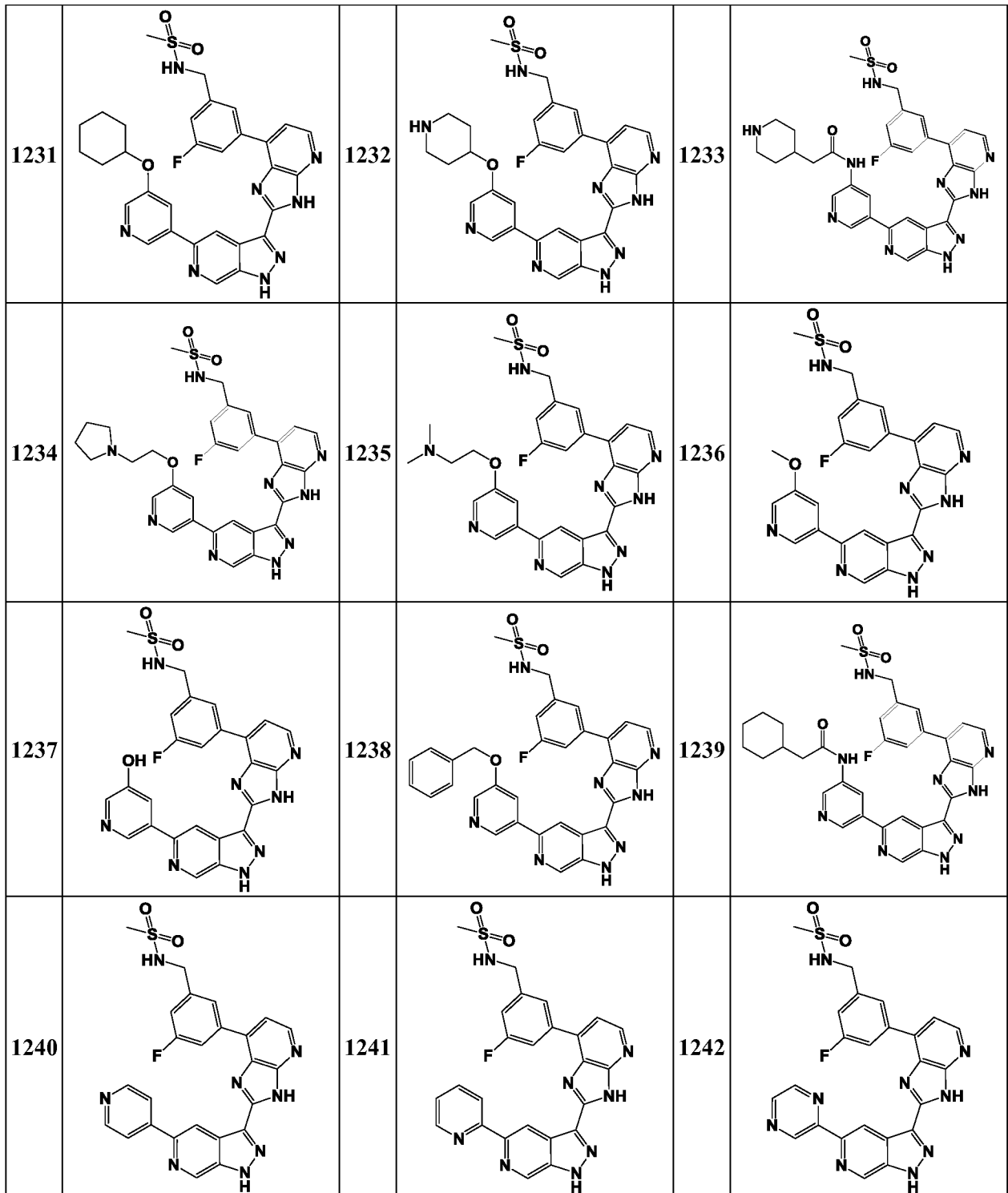
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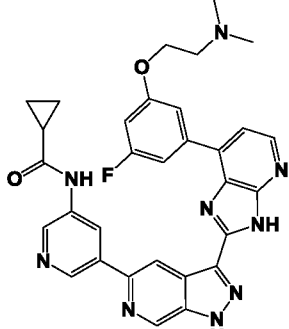
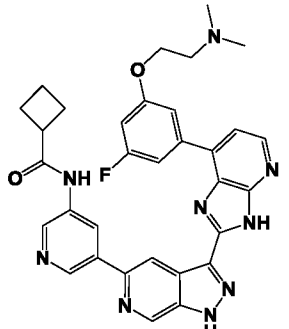
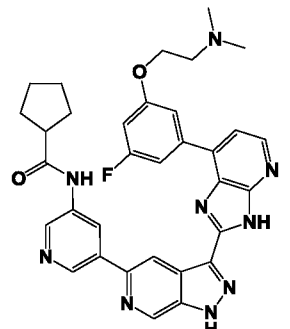
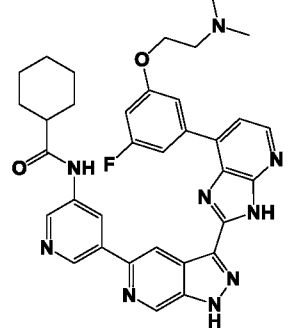
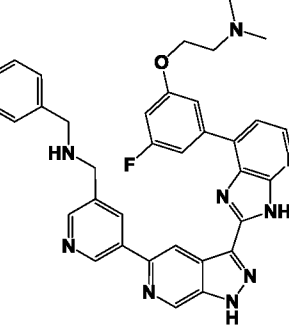
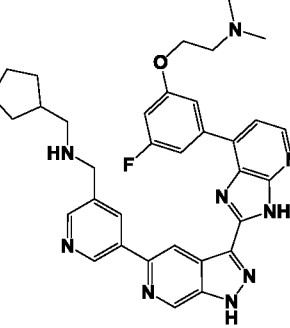
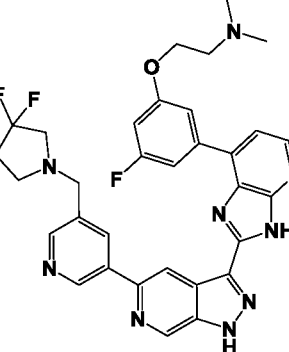
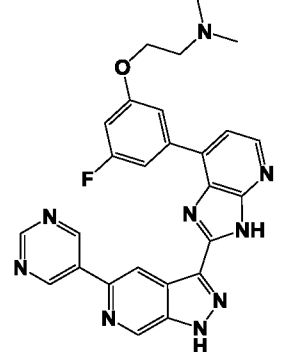
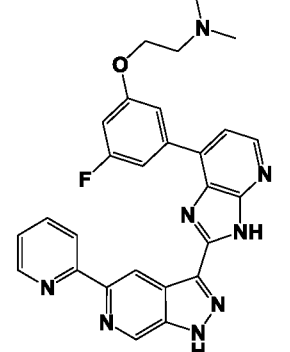
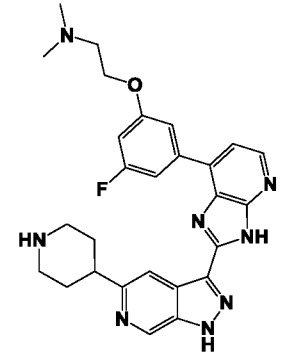
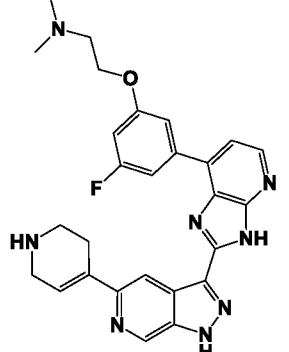
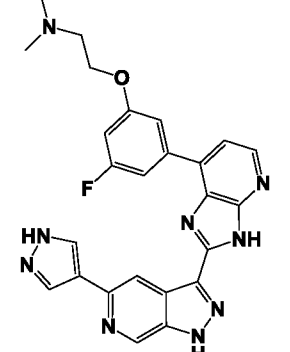
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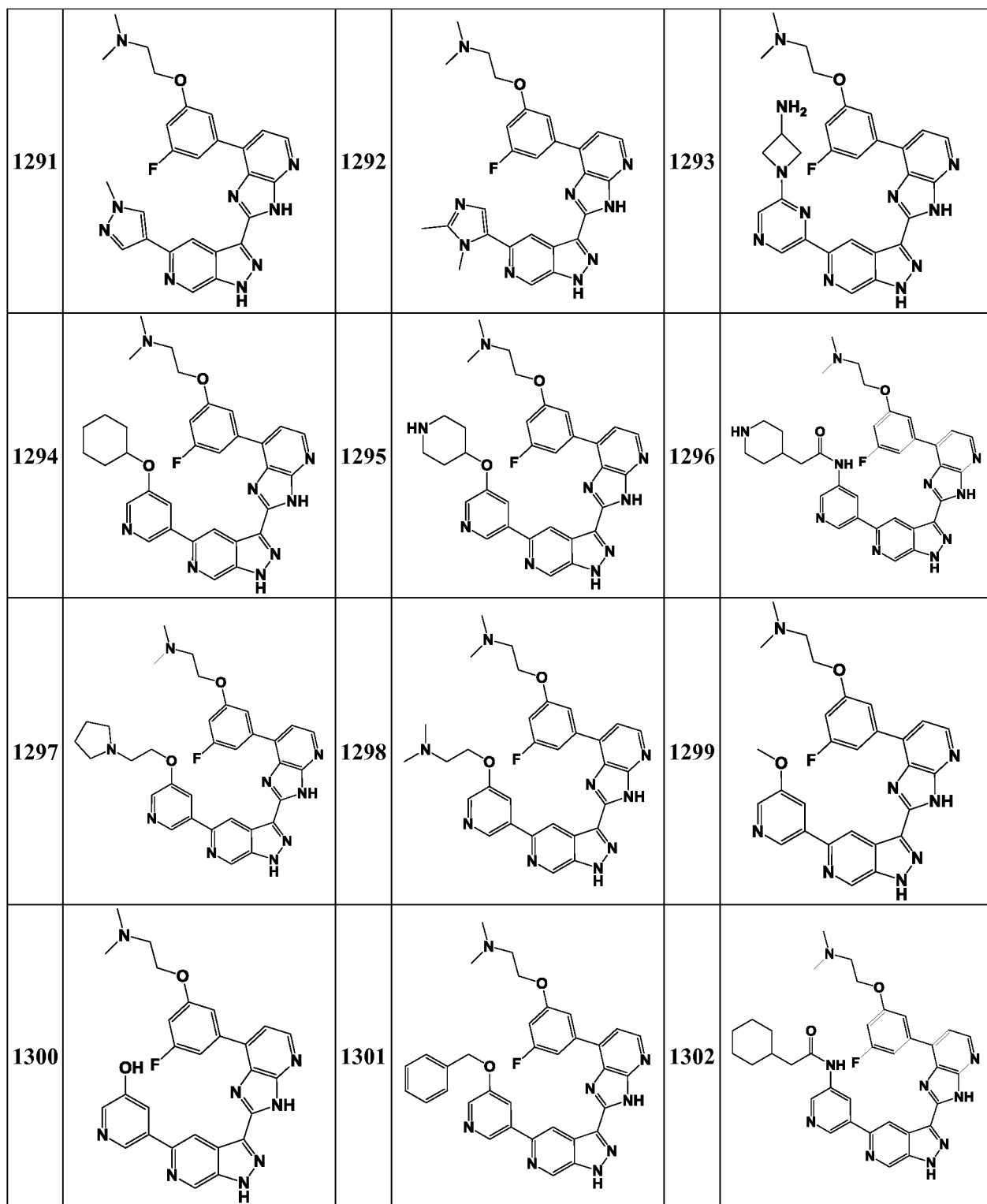


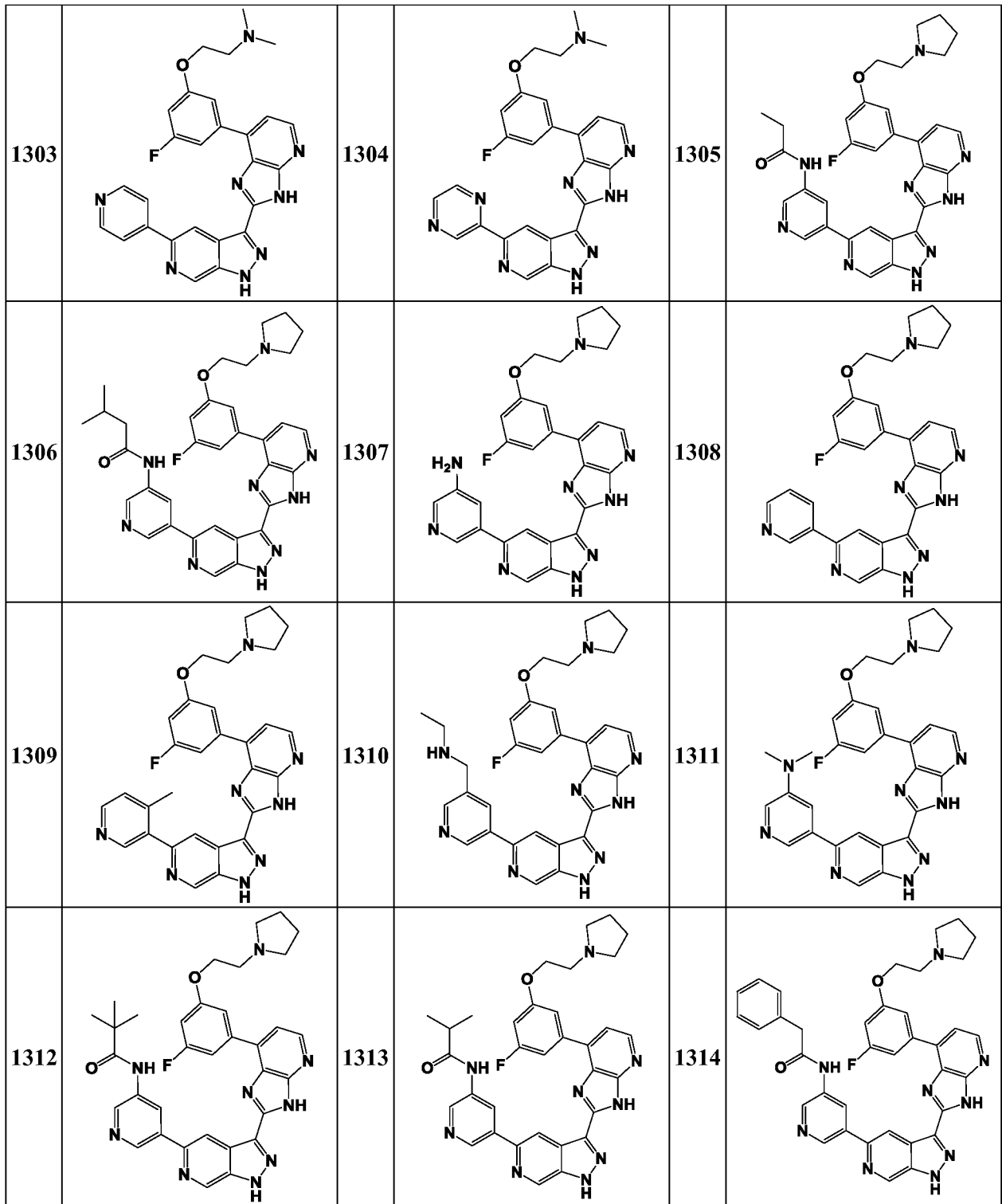


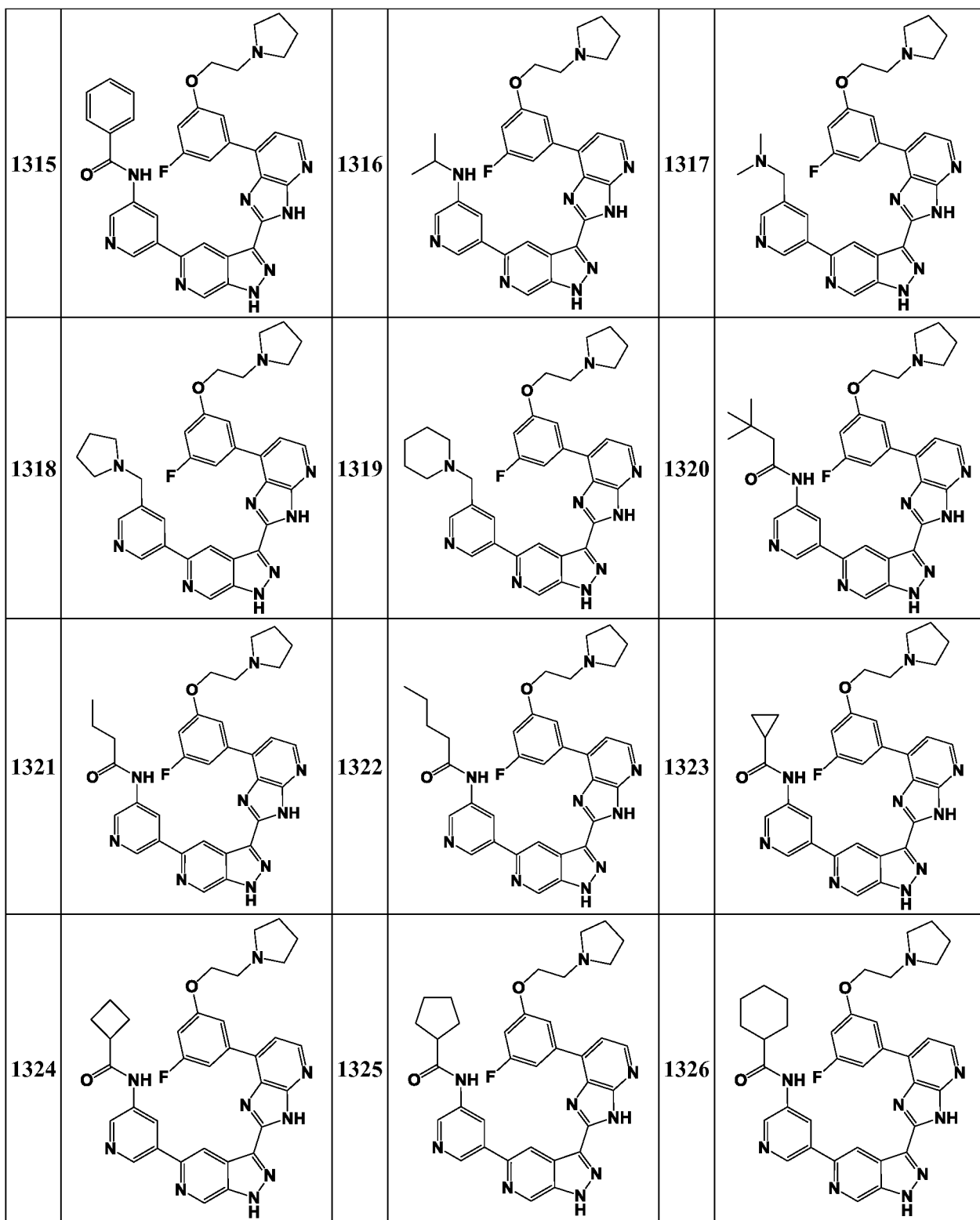
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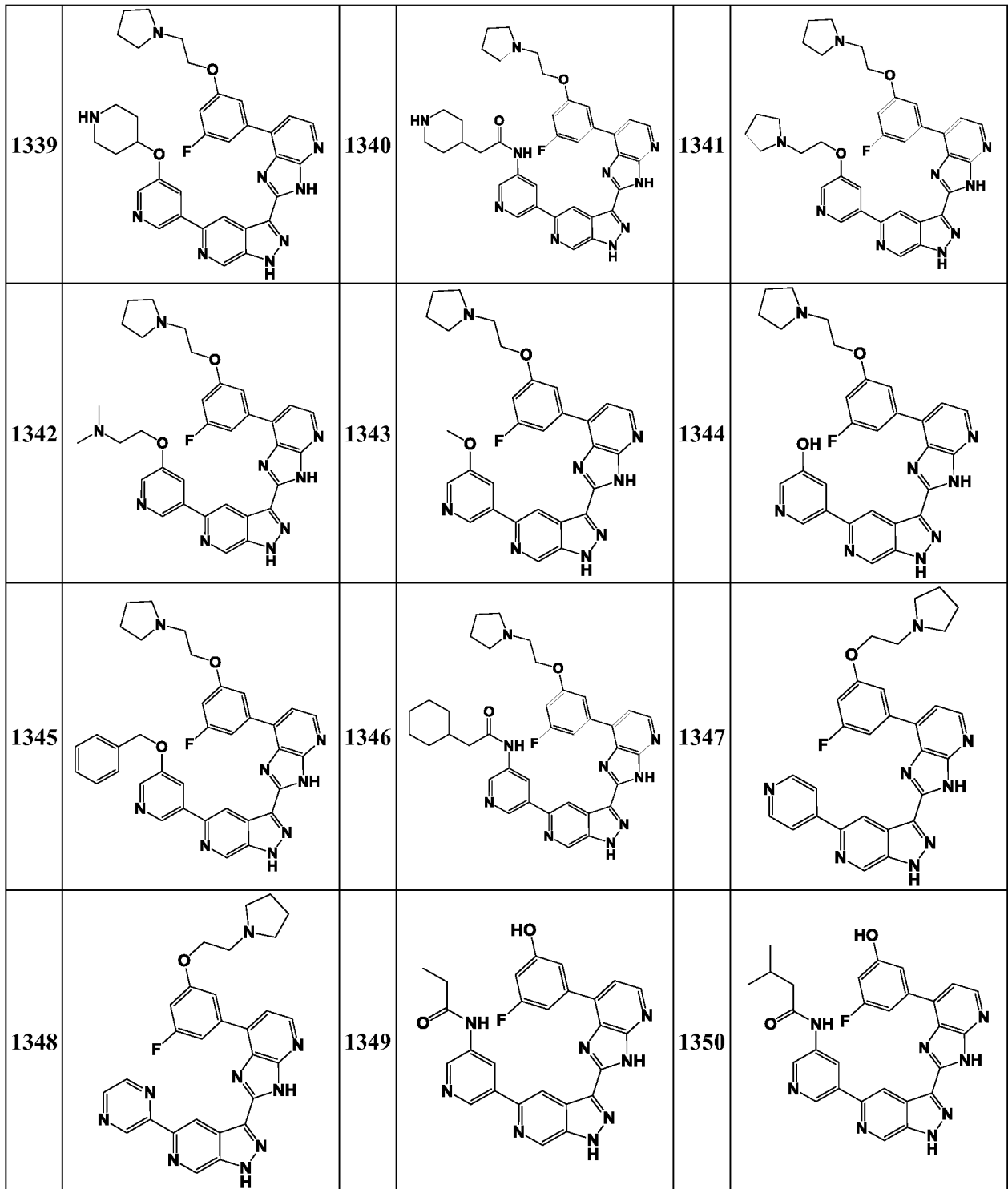
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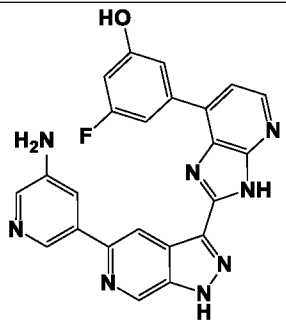
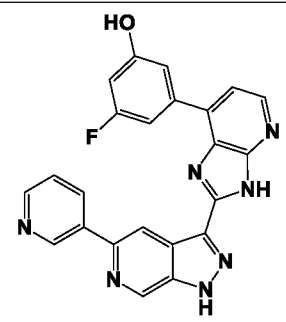
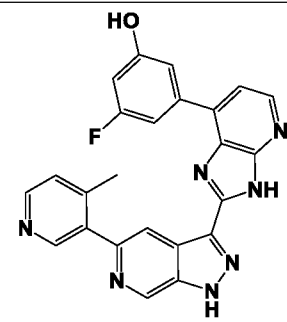
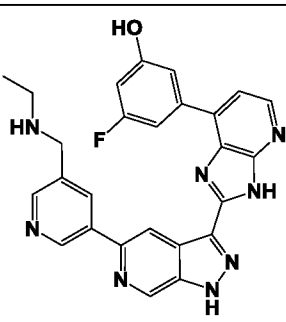
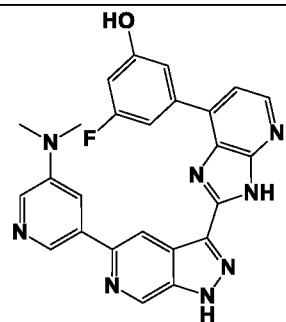
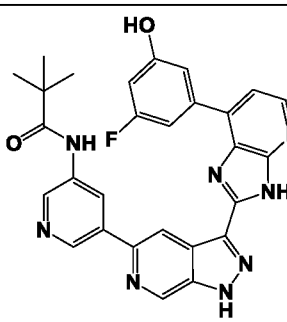
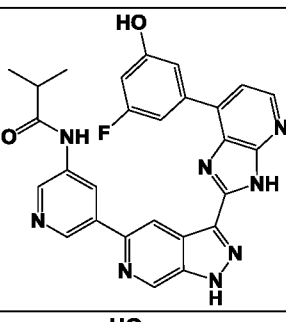
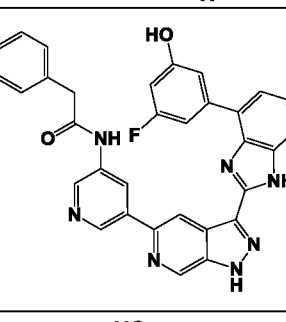
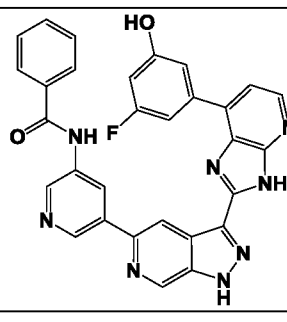
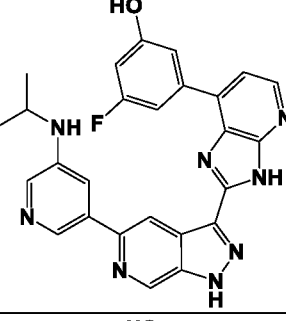
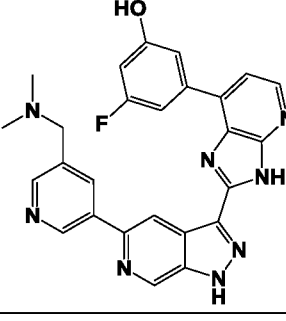
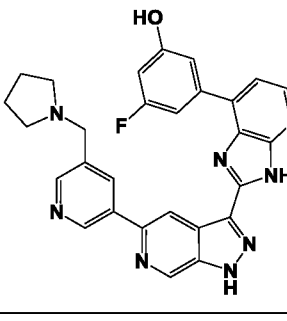
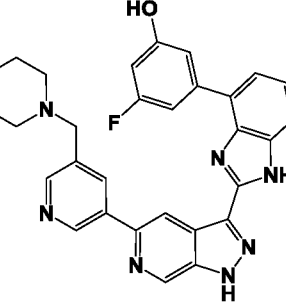
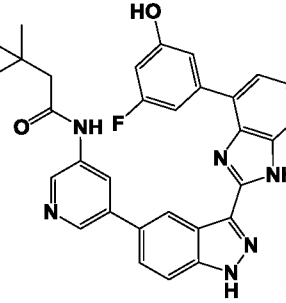
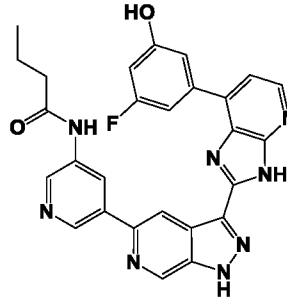
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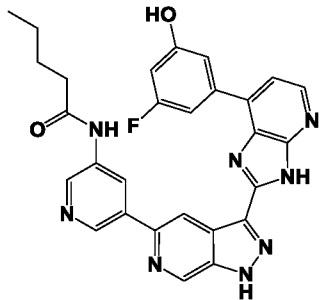
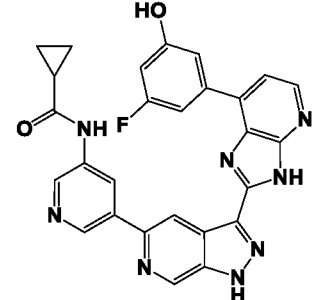
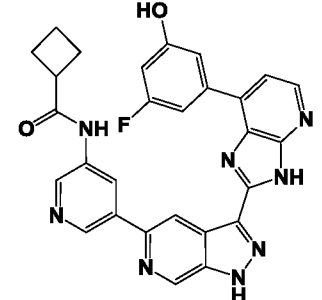
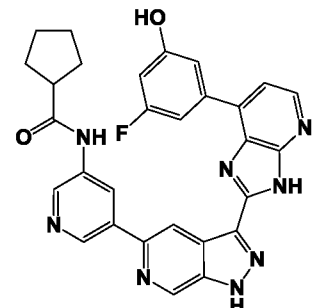
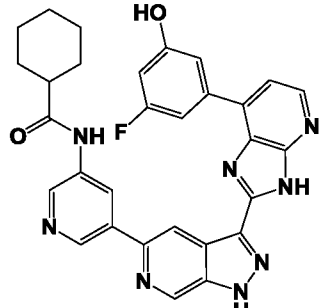
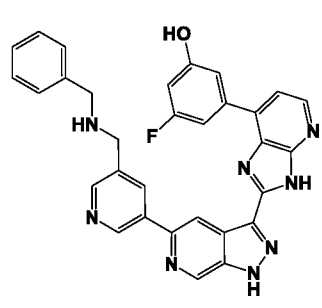
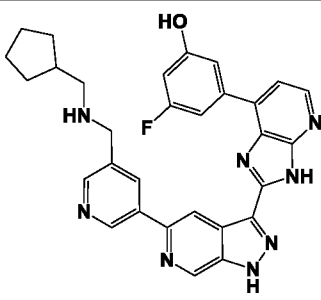
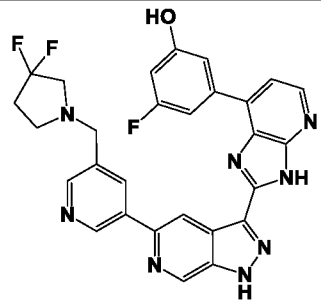
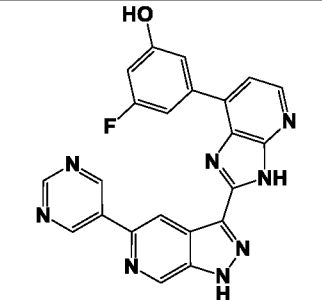
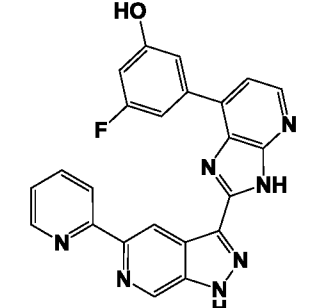
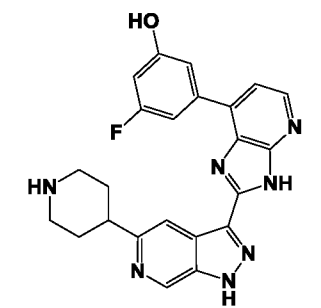
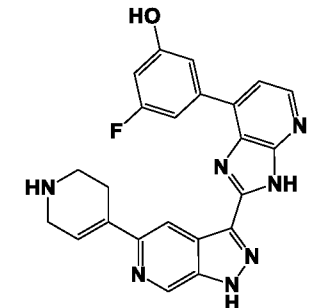
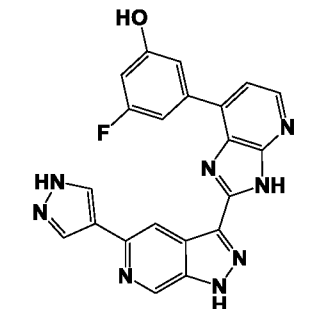
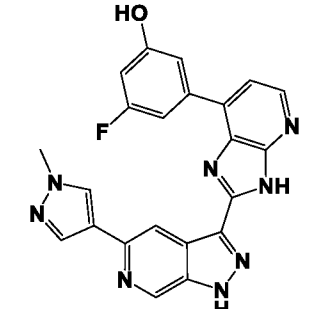
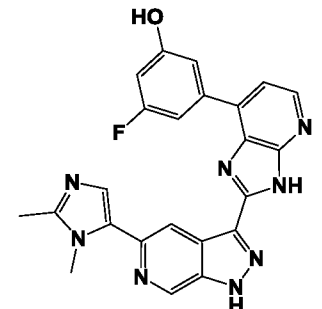




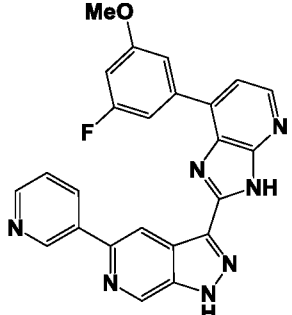
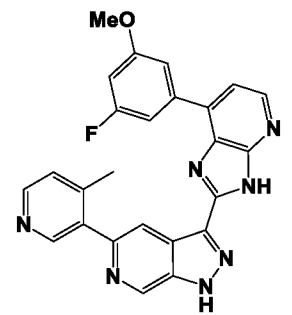
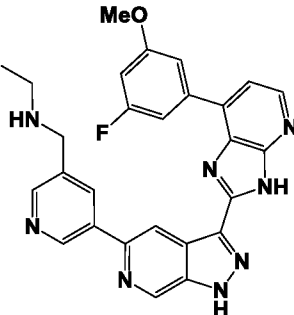
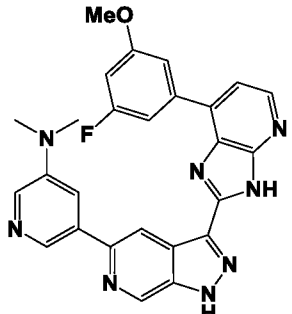
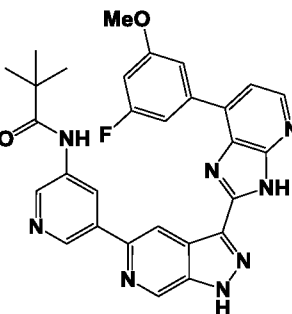
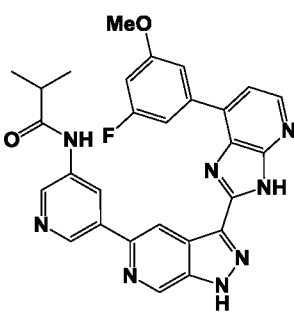
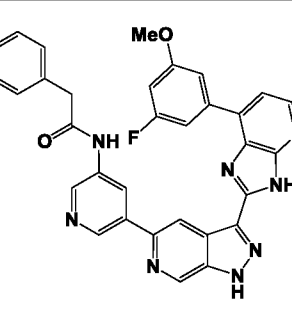
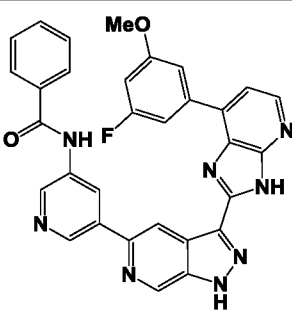
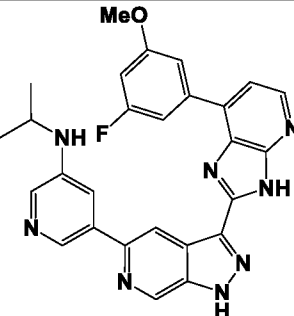
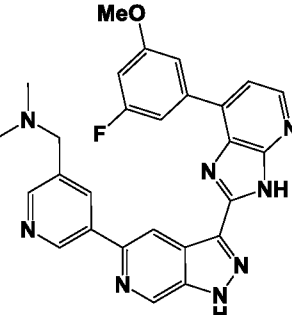
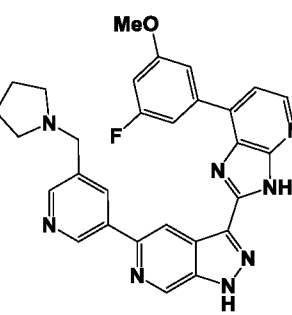
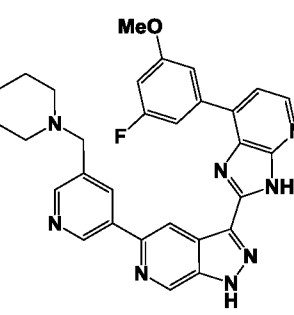
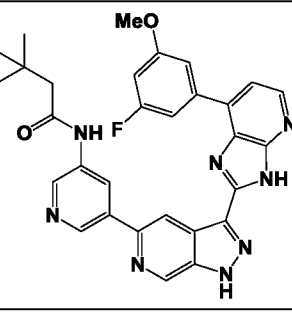
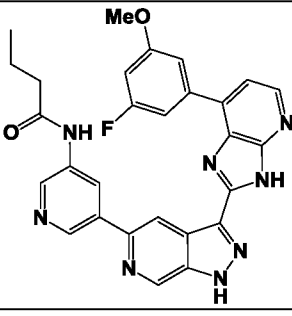
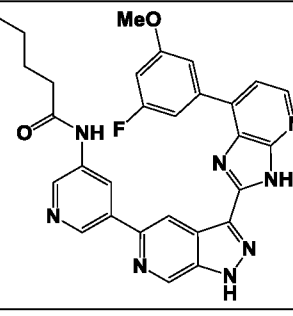


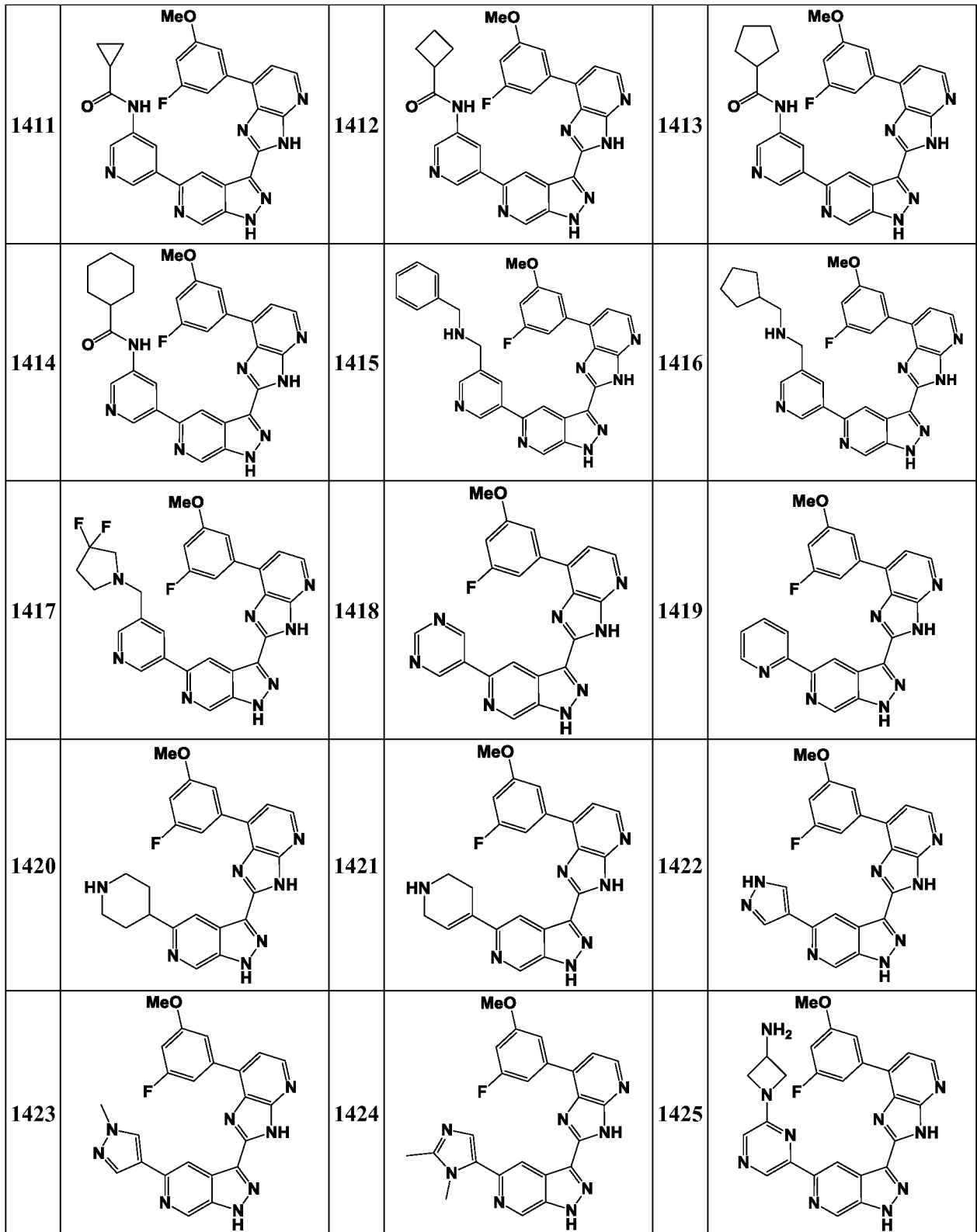


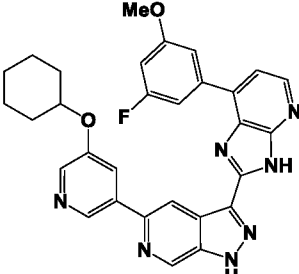
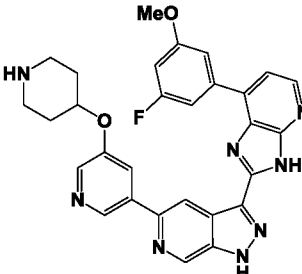
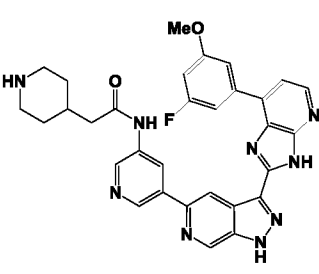
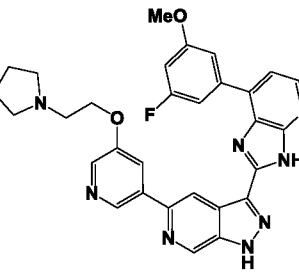
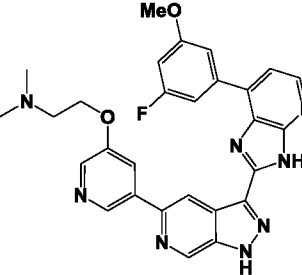
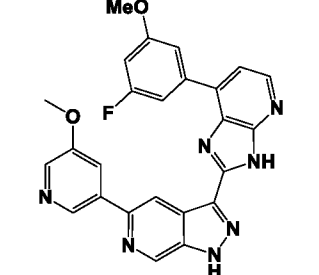
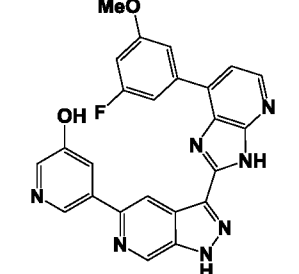
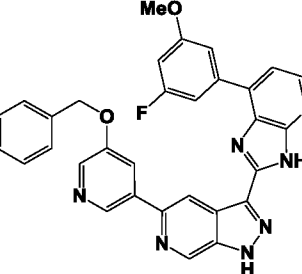
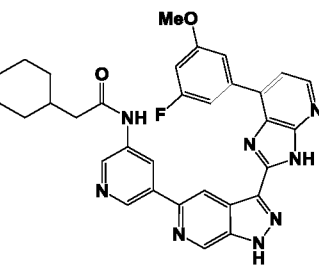
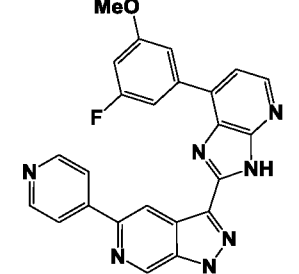
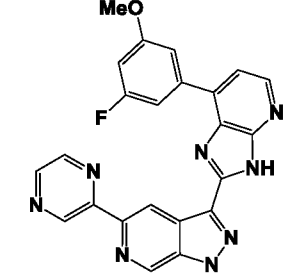
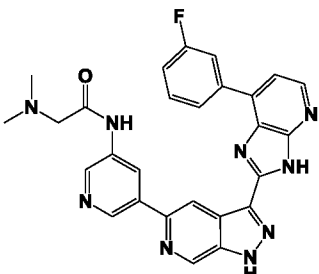
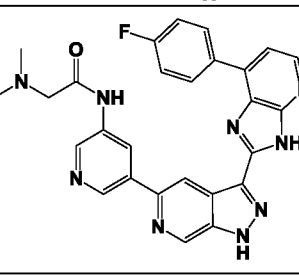
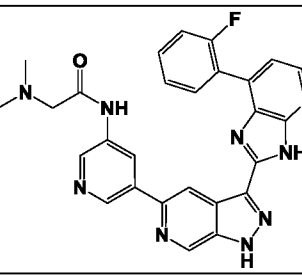
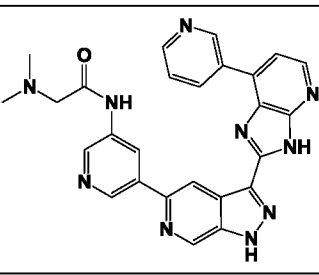
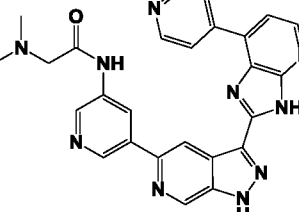
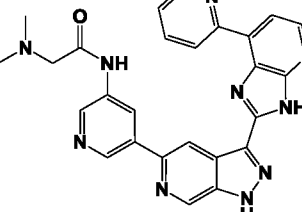
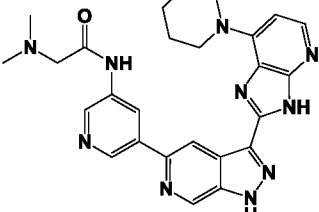
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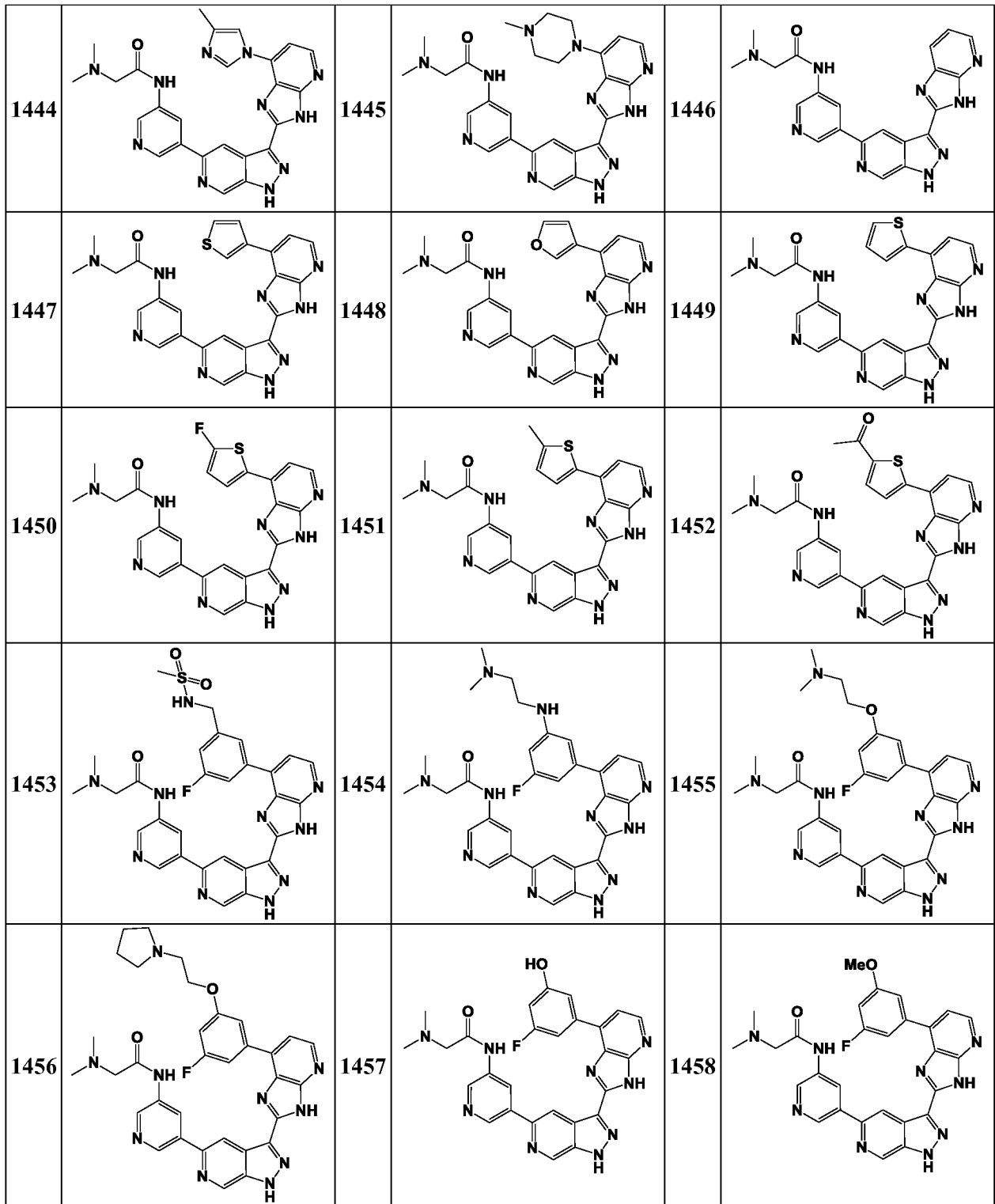
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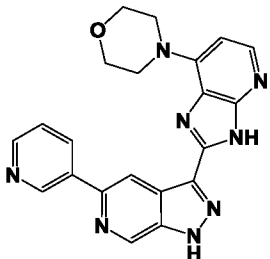
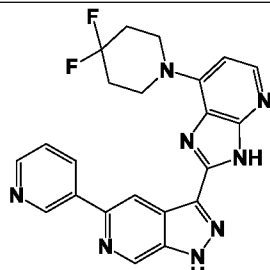
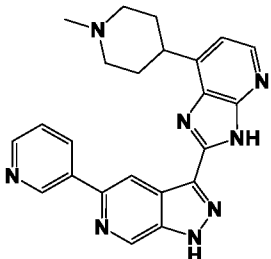
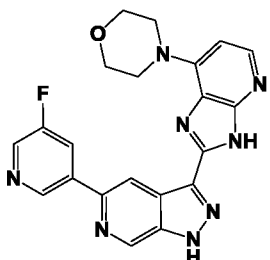
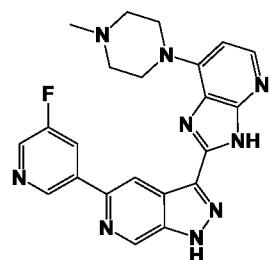
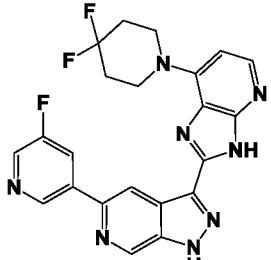
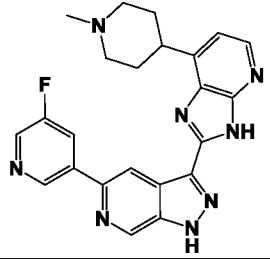
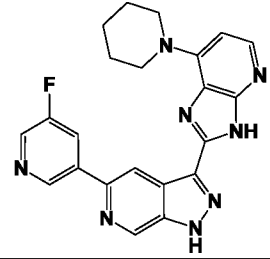
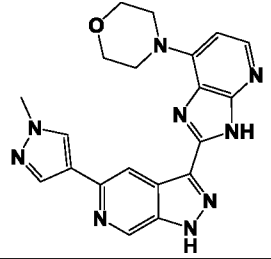
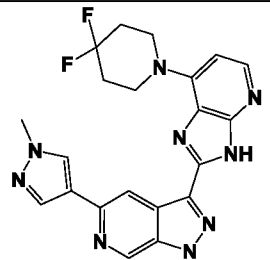
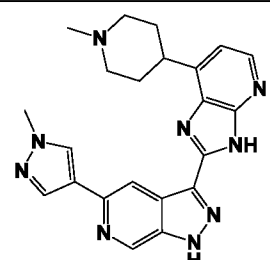
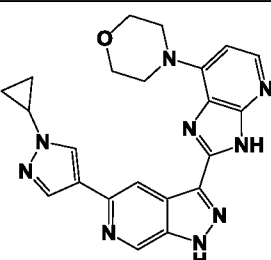
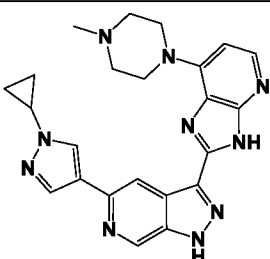
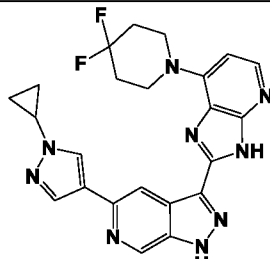
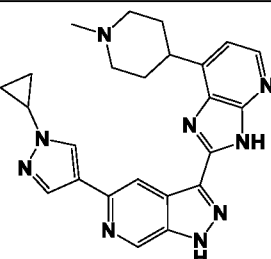
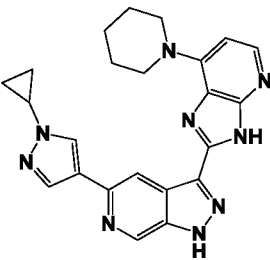
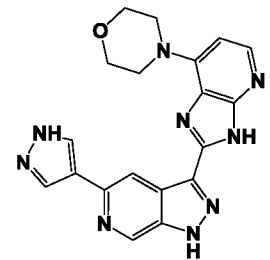
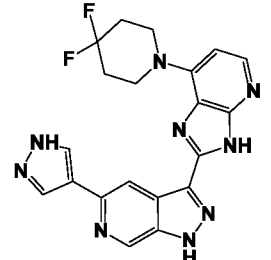
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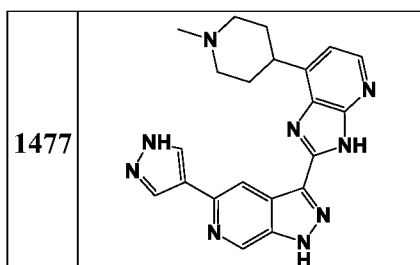
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Administration and Pharmaceutical Compositions

[0489] Some embodiments include pharmaceutical compositions comprising: (a) a therapeutically effective amount of a compound provided herein, or its corresponding enantiomer, diastereoisomer or tautomer, or pharmaceutically acceptable salt; and (b) a pharmaceutically acceptable carrier.

[0490] The compounds provided herein may also be useful in combination (administered together or sequentially) with other known agents.

[0491] Non-limiting examples of diseases which can be treated with a combination of a compound of Formula (I) and other known agents are colorectal cancer, ovarian cancer, retinitis pigmentosa, macular degeneration, diabetic retinopathy, idiopathic pulmonary fibrosis/pulmonary fibrosis, and osteoarthritis.

[0492] In some embodiments, colorectal cancer can be treated with a combination of a compound of Formula (I) and one or more of the following drugs: 5-Fluorouracil (5-FU), which can be administered with the vitamin-like drug leucovorin (also called folinic acid); capecitabine (XELODA[®]), irinotecan (CAMPOSTAR[®]), oxaliplatin (ELOXATIN[®]). Examples of combinations of these drugs which could be further combined with a compound of Formula (I) are FOLFOX (5-FU, leucovorin, and oxaliplatin), FOLFIRI (5-FU, leucovorin, and irinotecan), FOLFOXIRI (leucovorin, 5-FU, oxaliplatin, and irinotecan) and CapeOx (Capecitabine and oxaliplatin). For rectal cancer, chemo with 5-FU or capecitabine combined with radiation may be given before surgery (neoadjuvant treatment).

[0493] In some embodiments, ovarian cancer can be treated with a combination of a compound of Formula (I) and one or more of the following drugs: Topotecan, Liposomal doxorubicin (DOXIL[®]), Gemcitabine (GEMZAR[®]), Cyclophosphamide (CYTOXAN[®]), Vinorelbine (NAVELBINE[®]), Ifosfamide (IFEX[®]), Etoposide (VP-16), Altretamine (HEXALEN[®]), Capecitabine (XELODA[®]), Irinotecan (CPT-11, CAMPTOSAR[®]), Melphalan, Pemetrexed (ALIMTA[®]) and Albumin bound paclitaxel (nab-paclitaxel, ABRAXANE[®]). Examples of combinations of these drugs which could be further combined with a compound of

Formula (I) are TIP (paclitaxel [Taxol], ifosfamide, and cisplatin), VeIP (vinblastine, ifosfamide, and cisplatin) and VIP (etoposide [VP-16], ifosfamide, and cisplatin).

[0494] In some embodiments, a compound of Formula (I) can be used to treat cancer in combination with any of the following methods: (a) Hormone therapy such as aromatase inhibitors, LHRH [luteinizing hormone-releasing hormone] analogs and inhibitors, and others; (b) Ablation or embolization procedures such as radiofrequency ablation (RFA), ethanol (alcohol) ablation, microwave thermotherapy and cryosurgery (cryotherapy); (c) Chemotherapy using alkylating agents such as cisplatin and carboplatin, oxaliplatin, mechlorethamine, cyclophosphamide, chlorambucil and ifosfamide; (d) Chemotherapy using anti-metabolites such as azathioprine and mercaptopurine; (e) Chemotherapy using plant alkaloids and terpenoids such as vinca alkaloids (i.e. Vincristine, Vinblastine, Vinorelbine and Vindesine) and taxanes; (f) Chemotherapy using podophyllotoxin, etoposide, teniposide and docetaxel; (g) Chemotherapy using topoisomerase inhibitors such as irinotecan, topotecan, amsacrine, etoposide, etoposide phosphate, and teniposide; (h) Chemotherapy using cytotoxic antibiotics such as actinomycin, anthracyclines, doxorubicin, daunorubicin, valrubicin, idarubicin, epirubicin, bleomycin, plicamycin and mitomycin; (i) Chemotherapy using tyrosine-kinase inhibitors such as Imatinib mesylate (GLEEVEC[®], also known as STI-571), Gefitinib (Iressa, also known as ZD1839), Erlotinib (marketed as TARCEVA[®]), Bortezomib (VELCADE[®]), tamoxifen, tofacitinib, crizotinib, Bcl-2 inhibitors (e.g. obatoclax in clinical trials, ABT-263, and Gossypol), PARP inhibitors (e.g. Iniparib, Olaparib in clinical trials), PI3K inhibitors (eg. perifosine in a phase III trial), VEGF Receptor 2 inhibitors (e.g. Apatinib), AN-152, (AEZS-108), Braf inhibitors (e.g. vemurafenib, dabrafenib and LGX818), MEK inhibitors (e.g. trametinib and MEK162), CDK inhibitors, (e.g. PD-0332991), salinomycin and Sorafenib; (j) Chemotherapy using monoclonal antibodies such as Rituximab (marketed as MABTHERA[®] or RITUXAN[®]), Trastuzumab (Herceptin also known as ErbB2), Cetuximab (marketed as ERBITUX[®]), and Bevacizumab (marketed as AVASTIN[®]); and (k) radiation therapy.

[0495] In some embodiments, diabetic retinopathy can be treated with a combination of a compound of Formula (I) and one or more of the following natural supplements: Bilberry, Butcher's broom, Ginkgo, Grape seed extract, and Pycnogenol (Pine bark).

[0496] In some embodiments, idiopathic pulmonary fibrosis/pulmonary fibrosis can be treated with a combination of a compound of Formula (I) and one or more of the following drugs: pirfenidone (pirfenidone was approved for use in 2011 in Europe under the brand name Esbriet[®]), prednisone, azathioprine, N-acetylcysteine, interferon- γ 1b, bosentan (bosentan is currently being studied in patients with IPF, [*The American Journal of Respiratory and Critical*

Care Medicine (2011), 184(1), 92-9]), Nintedanib (BIBF 1120 and Vargatef), QAX576 [*British Journal of Pharmacology* (2011), 163(1), 141–172], and anti-inflammatory agents such as corticosteroids.

[0497] In some embodiments, a compound of Formula (I) can be used to treat idiopathic pulmonary fibrosis/pulmonary fibrosis in combination with any of the following methods: oxygen therapy, pulmonary rehabilitation and surgery.

[0498] In some embodiments, a compound of Formula (I) can be used to treat osteoarthritis in combination with any of the following methods: (a) Nonsteroidal anti-inflammatory drugs (NSAIDs) such as ibuprofen, naproxen, aspirin and acetaminophen; (b) physical therapy; (c) injections of corticosteroid medications; (d) injections of hyaluronic acid derivatives (e.g. Hyalgan, Synvisc); (e) narcotics, like codeine; (f) in combination with braces and/or shoe inserts or any device that can immobilize or support your joint to help you keep pressure off it (e.g., splints, braces, shoe inserts or other medical devices); (g) realigning bones (osteotomy); (h) joint replacement (arthroplasty); and (i) in combination with a chronic pain class.

[0499] In some embodiments, macular degeneration can be treated with a combination of a compound of Formula (I) and one or more of the following drugs: Bevacizumab (Avastin[®]), Ranibizumab (Lucentis[®]), Pegaptanib (Macugen), Aflibercept (Eylea[®]), verteporfin (Visudyne[®]) in combination with photodynamic therapy (PDT) or with any of the following methods: (a) in combination with laser to destroy abnormal blood vessels (photocoagulation); and (b) in combination with increased vitamin intake of antioxidant vitamins and zinc.

[0500] In some embodiments, retinitis pigmentosa can be treated with a combination of a compound of Formula (I) and one or more of the following drugs: UF-021 (Ocuseva[™]), vitamin A palmitate and pikachurin or with any of the following methods: (a) with the Argus[®] II retinal implant; and (b) with stem cell and/or gene therapy.

[0501] Administration of the compounds disclosed herein or the pharmaceutically acceptable salts thereof can be via any of the accepted modes of administration, including, but not limited to, orally, subcutaneously, intravenously, intranasally, topically, transdermally, intraperitoneally, intramuscularly, intrapulmonarily, vaginally, rectally, ontologically, neuro-otologically, intraocularly, subconjunctivally, via anterior eye chamber injection, intravitreally, intraperitoneally, intrathecally, intracystically, intrapleurally, via wound irrigation, intrabuccally, intra-abdominally, intra-articularly, intra-aurally, intrabronchially, intracapsularly, intrameningeally, via inhalation, via endotracheal or endobronchial instillation, via direct instillation into pulmonary cavities, intraspinally, intrasynovially, intrathoracically, via thoracostomy irrigation, epidurally, intratympanically, intracisternally, intravascularly,

intraventricularly, intraosseously, via irrigation of infected bone, or via application as part of any admixture with a prosthetic devices. In some embodiments, the administration method includes oral or parenteral administration.

[0502] Compounds provided herein intended for pharmaceutical use may be administered as crystalline or amorphous products. Pharmaceutically acceptable compositions may include solid, semi-solid, liquid, solutions, colloidal, liposomes, emulsions, suspensions, complexes, coacervates and aerosols. Dosage forms, such as, *e.g.*, tablets, capsules, powders, liquids, suspensions, suppositories, aerosols, implants, controlled release or the like. They may be obtained, for example, as solid plugs, powders, or films by methods such as precipitation, crystallization, milling, grinding, supercritical fluid processing, coacervation, complex coacervation, encapsulation, emulsification, complexation, freeze drying, spray drying, or evaporative drying. Microwave or radio frequency drying may be used for this purpose. The compounds can also be administered in sustained or controlled release dosage forms, including depot injections, osmotic pumps, pills (tablets and or capsules), transdermal (including electrotransport) patches, implants and the like, for prolonged and/or timed, pulsed administration at a predetermined rate.

[0503] The compounds can be administered either alone or in combination with a conventional pharmaceutical carrier, excipient or the like. Pharmaceutically acceptable excipients include, but are not limited to, ion exchangers, alumina, aluminum stearate, lecithin, self-emulsifying drug delivery systems (SEDDS) such as *d*- α -tocopherol polyethylene glycol 1000 succinate, surfactants used in pharmaceutical dosage forms such as Tweens, poloxamers or other similar polymeric delivery matrices, serum proteins, such as human serum albumin, buffer substances such as phosphates, tris, glycine, sorbic acid, potassium sorbate, partial glyceride mixtures of saturated vegetable fatty acids, water, salts or electrolytes, such as protamine sulfate, disodium hydrogen phosphate, potassium hydrogen phosphate, sodium-chloride, zinc salts, colloidal silica, magnesium trisilicate, polyvinyl pyrrolidone, cellulose-based substances, polyethylene glycol, sodium carboxymethyl cellulose, polyacrylates, waxes, polyethylene-polyoxypropylene-block polymers, and wool fat. Cyclodextrins such as α -, β -, and γ -cyclodextrin, or chemically modified derivatives such as hydroxyalkylcyclodextrins, including 2- and 3-hydroxypropyl- β -cyclodextrins, or other solubilized derivatives can also be used to enhance delivery of compounds described herein. Dosage forms or compositions containing a compound as described herein in the range of 0.005% to 100% with the balance made up from non-toxic carrier may be prepared. The contemplated compositions may contain 0.001%-100% of a compound provided herein, in one embodiment 0.1-95%, in another embodiment 75-85%, in a further

embodiment 20-80%. Actual methods of preparing such dosage forms are known, or will be apparent, to those skilled in this art; for example, see *Remington: The Science and Practice of Pharmacy*, 22nd Edition (Pharmaceutical Press, London, UK, 2012).

[0504] In one embodiment, the compositions will take the form of a unit dosage form such as a pill or tablet and thus the composition may contain, along with a compound provided herein, a diluent such as lactose, sucrose, dicalcium phosphate, or the like; a lubricant such as magnesium stearate or the like; and a binder such as starch, gum acacia, polyvinylpyrrolidone, gelatin, cellulose, cellulose derivatives or the like. In another solid dosage form, a powder, marume, solution or suspension (*e.g.*, in propylene carbonate, vegetable oils, PEG's, poloxamer 124 or triglycerides) is encapsulated in a capsule (gelatin or cellulose base capsule). Unit dosage forms in which one or more compounds provided herein or additional active agents are physically separated are also contemplated; *e.g.*, capsules with granules (or tablets in a capsule) of each drug; two-layer tablets; two-compartment gel caps, etc. Enteric coated or delayed release oral dosage forms are also contemplated.

[0505] Liquid pharmaceutically administrable compositions can, for example, be prepared by dissolving, dispersing, etc. a compound provided herein and optional pharmaceutical adjuvants in a carrier (*e.g.*, water, saline, aqueous dextrose, glycerol, glycols, ethanol or the like) to form a solution, colloid, liposome, emulsion, complexes, coacervate or suspension. If desired, the pharmaceutical composition can also contain minor amounts of nontoxic auxiliary substances such as wetting agents, emulsifying agents, co-solvents, solubilizing agents, pH buffering agents and the like (*e.g.*, sodium acetate, sodium citrate, cyclodextrin derivatives, sorbitan monolaurate, triethanolamine acetate, triethanolamine oleate, and the like).

[0506] In some embodiments, the unit dosage of compounds of Formula (I) is about 0.25 mg/Kg to about 50 mg/Kg in humans.

[0507] In some embodiments, the unit dosage of compounds of Formula (I) is about 0.25 mg/Kg to about 20 mg/Kg in humans.

[0508] In some embodiments, the unit dosage of compounds of Formula (I) is about 0.50 mg/Kg to about 19 mg/Kg in humans.

[0509] In some embodiments, the unit dosage of compounds of Formula (I) is about 0.75 mg/Kg to about 18 mg/Kg in humans.

[0510] In some embodiments, the unit dosage of compounds of Formula (I) is about 1.0 mg/Kg to about 17 mg/Kg in humans.

[0511] In some embodiments, the unit dosage of compounds of Formula (I) is about 1.25 mg/Kg to about 16 mg/Kg in humans.

[0512] In some embodiments, the unit dosage of compounds of Formula (I) is about 1.50 mg/Kg to about 15 mg/Kg in humans.

[0513] In some embodiments, the unit dosage of compounds of Formula (I) is about 1.75 mg/Kg to about 14 mg/Kg in humans.

[0514] In some embodiments, the unit dosage of compounds of Formula (I) is about 2.0 mg/Kg to about 13 mg/Kg in humans.

[0515] In some embodiments, the unit dosage of compounds of Formula (I) is about 3.0 mg/Kg to about 12 mg/Kg in humans.

[0516] In some embodiments, the unit dosage of compounds of Formula (I) is about 4.0 mg/Kg to about 11 mg/Kg in humans.

[0517] In some embodiments, the unit dosage of compounds of Formula (I) is about 5.0 mg/Kg to about 10 mg/Kg in humans.

[0518] In some embodiments, the compositions are provided in unit dosage forms suitable for single administration.

[0519] In some embodiments, the compositions are provided in unit dosage forms suitable for twice a day administration.

[0520] In some embodiments, the compositions are provided in unit dosage forms suitable for three times a day administration.

[0521] Injectables can be prepared in conventional forms, either as liquid solutions, colloid, liposomes, complexes, coacervate or suspensions, as emulsions, or in solid forms suitable for reconstitution in liquid prior to injection. The percentage of a compound provided herein contained in such parenteral compositions is highly dependent on the specific nature thereof, as well as the activity of the compound and the needs of the patient. However, percentages of active ingredient of 0.01% to 10% in solution are employable, and could be higher if the composition is a solid or suspension, which could be subsequently diluted to the above percentages.

[0522] In some embodiments, the composition will comprise about 0.1-10% of the active agent in solution.

[0523] In some embodiments, the composition will comprise about 0.1-5% of the active agent in solution.

[0524] In some embodiments, the composition will comprise about 0.1-4% of the active agent in solution.

[0525] In some embodiments, the composition will comprise about 0.15-3% of the active agent in solution.

[0526] In some embodiments, the composition will comprise about 0.2-2% of the active agent in solution.

[0527] In some embodiments, the compositions are provided in dosage forms suitable for continuous dosage by intravenous infusion over a period of about 1-96 hours.

[0528] In some embodiments, the compositions are provided in dosage forms suitable for continuous dosage by intravenous infusion over a period of about 1-72 hours.

[0529] In some embodiments, the compositions are provided in dosage forms suitable for continuous dosage by intravenous infusion over a period of about 1-48 hours.

[0530] In some embodiments, the compositions are provided in dosage forms suitable for continuous dosage by intravenous infusion over a period of about 1-24 hours.

[0531] In some embodiments, the compositions are provided in dosage forms suitable for continuous dosage by intravenous infusion over a period of about 1-12 hours.

[0532] In some embodiments, the compositions are provided in dosage forms suitable for continuous dosage by intravenous infusion over a period of about 1-6 hours.

[0533] In some embodiments, these compositions can be administered by intravenous infusion to humans at doses of about 5 mg/m² to about 300 mg/m².

[0534] In some embodiments, these compositions can be administered by intravenous infusion to humans at doses of about 5 mg/m² to about 200 mg/m².

[0535] In some embodiments, these compositions can be administered by intravenous infusion to humans at doses of about 5 mg/m² to about 100 mg/m².

[0536] In some embodiments, these compositions can be administered by intravenous infusion to humans at doses of about 10 mg/m² to about 50 mg/m².

[0537] In some embodiments, these compositions can be administered by intravenous infusion to humans at doses of about 50 mg/m² to about 200 mg/m².

[0538] In some embodiments, these compositions can be administered by intravenous infusion to humans at doses of about 75 mg/m² to about 175 mg/m².

[0539] In some embodiments, these compositions can be administered by intravenous infusion to humans at doses of about 100 mg/m² to about 150 mg/m².

[0540] It is to be noted that concentrations and dosage values may also vary depending on the specific compound and the severity of the condition to be alleviated. It is to be further understood that for any particular patient, specific dosage regimens should be adjusted over time according to the individual need and the professional judgment of the person administering or supervising the administration of the compositions, and that the concentration ranges set forth

herein are exemplary only and are not intended to limit the scope or practice of the claimed compositions.

[0541] In one embodiment, the compositions can be administered to the respiratory tract (including nasal and pulmonary) e.g., through a nebulizer, metered-dose inhalers, atomizer, mister, aerosol, dry powder inhaler, insufflator, liquid instillation or other suitable device or technique.

[0542] In some embodiments, aerosols intended for delivery to the nasal mucosa are provided for inhalation through the nose. For optimal delivery to the nasal cavities, inhaled particle sizes of about 5 to about 100 microns are useful, with particle sizes of about 10 to about 60 microns being preferred. For nasal delivery, a larger inhaled particle size may be desired to maximize impaction on the nasal mucosa and to minimize or prevent pulmonary deposition of the administered formulation. In some embodiments, aerosols intended for delivery to the lung are provided for inhalation through the nose or the mouth. For delivery to the lung, inhaled aerodynamic particle sizes of about less than 10 μm are useful (e.g., about 1 to about 10 microns). Inhaled particles may be defined as liquid droplets containing dissolved drug, liquid droplets containing suspended drug particles (in cases where the drug is insoluble in the suspending medium), dry particles of pure drug substance, drug substance incorporated with excipients, liposomes, emulsions, colloidal systems, coacervates, aggregates of drug nanoparticles, or dry particles of a diluent which contain embedded drug nanoparticles.

[0543] In some embodiments, compounds of Formula (I) disclosed herein intended for respiratory delivery (either systemic or local) can be administered as aqueous formulations, as non-aqueous solutions or suspensions, as suspensions or solutions in halogenated hydrocarbon propellants with or without alcohol, as a colloidal system, as emulsions, coacervates, or as dry powders. Aqueous formulations may be aerosolized by liquid nebulizers employing either hydraulic or ultrasonic atomization or by modified micropump systems (like the soft mist inhalers, the Aerodose[®] or the AERx[®] systems). Propellant-based systems may use suitable pressurized metered-dose inhalers (pMDIs). Dry powders may use dry powder inhaler devices (DPIs), which are capable of dispersing the drug substance effectively. A desired particle size and distribution may be obtained by choosing an appropriate device.

[0544] In some embodiments, the compositions of Formula (I) disclosed herein can be administered to the ear by various methods. For example, a round window catheter (e.g., U.S. Pat. Nos. 6,440,102 and 6,648,873) can be used.

[0545] Alternatively, formulations can be incorporated into a wick for use between the outer and middle ear (e.g., U.S. Pat. No. 6,120,484) or absorbed to collagen sponge or other solid support (e.g., U.S. Pat. No. 4,164,559).

[0546] If desired, formulations of the invention can be incorporated into a gel formulation (e.g., U.S. Pat. Nos. 4,474,752 and 6,911,211).

[0547] In some embodiments, compounds of Formula (I) disclosed herein intended for delivery to the ear can be administered via an implanted pump and delivery system through a needle directly into the middle or inner ear (cochlea) or through a cochlear implant stylet electrode channel or alternative prepared drug delivery channel such as but not limited to a needle through temporal bone into the cochlea.

[0548] Other options include delivery via a pump through a thin film coated onto a multichannel electrode or electrode with a specially imbedded drug delivery channel (pathways) carved into the thin film for this purpose. In other embodiments the acidic or basic solid compound of Formula (I) can be delivered from the reservoir of an external or internal implanted pumping system.

[0549] Formulations of the invention also can be administered to the ear by intratympanic injection into the middle ear, inner ear, or cochlea (e.g., U.S. Pat. No. 6,377,849 and Ser. No. 11/337,815).

[0550] Intratympanic injection of therapeutic agents is the technique of injecting a therapeutic agent behind the tympanic membrane into the middle and/or inner ear. In one embodiment, the formulations described herein are administered directly onto the round window membrane via transtympanic injection. In another embodiment, the ion channel modulating agent auris-acceptable formulations described herein are administered onto the round window membrane via a non-transtympanic approach to the inner ear. In additional embodiments, the formulation described herein is administered onto the round window membrane via a surgical approach to the round window membrane comprising modification of the crista fenestrae cochleae.

[0551] In some embodiments, the compounds of Formula (I) are formulated in rectal compositions such as enemas, rectal gels, rectal foams, rectal aerosols, suppositories, jelly suppositories, or retention enemas, containing conventional suppository bases such as cocoa butter or other glycerides, as well as synthetic polymers such as polyvinylpyrrolidone, PEG (like PEG ointments), and the like.

[0552] Suppositories for rectal administration of the drug (either as a solution, colloid, suspension or a complex) can be prepared by mixing a compound provided herein with a suitable non-irritating excipient that is solid at ordinary temperatures but liquid at the rectal temperature

and will therefore melt or erode/dissolve in the rectum and release the compound. Such materials include cocoa butter, glycerinated gelatin, hydrogenated vegetable oils, poloxamers, mixtures of polyethylene glycols of various molecular weights and fatty acid esters of polyethylene glycol. In suppository forms of the compositions, a low-melting wax such as, but not limited to, a mixture of fatty acid glycerides, optionally in combination with cocoa butter, is first melted.

[0553] Solid compositions can be provided in various different types of dosage forms, depending on the physicochemical properties of the compound provided herein, the desired dissolution rate, cost considerations, and other criteria. In one of the embodiments, the solid composition is a single unit. This implies that one unit dose of the compound is comprised in a single, physically shaped solid form or article. In other words, the solid composition is coherent, which is in contrast to a multiple unit dosage form, in which the units are incoherent.

[0554] Examples of single units which may be used as dosage forms for the solid composition include tablets, such as compressed tablets, film-like units, foil-like units, wafers, lyophilized matrix units, and the like. In one embodiment, the solid composition is a highly porous lyophilized form. Such lyophilizates, sometimes also called wafers or lyophilized tablets, are particularly useful for their rapid disintegration, which also enables the rapid dissolution of the compound.

[0555] On the other hand, for some applications the solid composition may also be formed as a multiple unit dosage form as defined above. Examples of multiple units are powders, granules, microparticles, pellets, mini-tablets, beads, lyophilized powders, and the like. In one embodiment, the solid composition is a lyophilized powder. Such a dispersed lyophilized system comprises a multitude of powder particles, and due to the lyophilization process used in the formation of the powder, each particle has an irregular, porous microstructure through which the powder is capable of absorbing water very rapidly, resulting in quick dissolution. Effervescent compositions are also contemplated to aid the quick dispersion and absorption of the compound.

[0556] Another type of multiparticulate system which is also capable of achieving rapid drug dissolution is that of powders, granules, or pellets from water-soluble excipients which are coated with a compound provided herein so that the compound is located at the outer surface of the individual particles. In this type of system, the water-soluble low molecular weight excipient may be useful for preparing the cores of such coated particles, which can be subsequently coated with a coating composition comprising the compound and, for example, one or more additional excipients, such as a binder, a pore former, a saccharide, a sugar alcohol, a film-forming polymer, a plasticizer, or other excipients used in pharmaceutical coating compositions.

[0557] Also provided herein are kits. Typically, a kit includes one or more compounds or compositions as described herein. In certain embodiments, a kit can include one or more delivery systems, e.g., for delivering or administering a compound as provided herein, and directions for use of the kit (e.g., instructions for treating a patient). In another embodiment, the kit can include a compound or composition as described herein and a label that indicates that the contents are to be administered to a patient with cancer. In another embodiment, the kit can include a compound or composition as described herein and a label that indicates that the contents are to be administered to a patient with one or more of hepatocellular carcinoma, colon cancer, leukemia, lymphoma, sarcoma, ovarian cancer, diabetic retinopathy, pulmonary fibrosis, rheumatoid arthritis, sepsis, ankylosing spondylitis, psoriasis, scleroderma, mycotic and viral infections, bone and cartilage diseases, Alzheimer's disease, lung disease, bone/osteoporotic (wrist, spine, shoulder and hip) fractures, articular cartilage (chondral) defects, degenerative disc disease (or intervertebral disc degeneration), polyposis coli, bone density and vascular defects in the eye (Osteoporosis-pseudoglioma Syndrome, OPPG) and other eye diseases or syndromes associated with defects and/or damaged photoreceptors, familial exudative vitreoretinopathy, retinal angiogenesis, early coronary disease, tetra-amelia, Müllerian-duct regression and virilization, SERKAL syndrome, type II diabetes, Fuhrmann syndrome, Al-Awadi/Raas-Rothschild/Schinzel phocomelia syndrome, odonto-onycho-dermal dysplasia, obesity, split-hand/foot malformation, caudal duplication, tooth agenesis, Wilms tumor, skeletal dysplasia, focal dermal hypoplasia, autosomal recessive anonychia, neural tube defects, alpha-thalassemia (ATRX) syndrome, fragile X syndrome, ICF syndrome, Angelman's syndrome, Prader-Willi syndrome, Beckwith-Wiedemann Syndrome, Norrie disease, and Rett syndrome.

Methods of Treatment

[0558] The compounds and compositions provided herein can be used as inhibitors and/or modulators of one or more components of the Wnt pathway, which may include one or more Wnt proteins, and thus can be used to treat a variety of disorders and diseases in which aberrant Wnt signaling is implicated, such as cancer and other diseases associated with abnormal angiogenesis, cellular proliferation, and cell cycling. Accordingly, the compounds and compositions provided herein can be used to treat cancer, to reduce or inhibit angiogenesis, to reduce or inhibit cellular proliferation, to correct a genetic disorder, and/or to treat a neurological condition/disorder/disease due to mutations or dysregulation of the Wnt pathway and/or of one or more of Wnt signaling components. Non-limiting examples of diseases which can be treated with the compounds and compositions provided herein include a variety of cancers, diabetic retinopathy,

pulmonary fibrosis, rheumatoid arthritis, scleroderma, mycotic and viral infections, bone and cartilage diseases, neurological conditions/diseases such as Alzheimer's disease, amyotrophic lateral sclerosis (ALS), motor neuron disease, multiple sclerosis or autism, lung disease, bone/osteoporotic (wrist, spine, shoulder and hip) fractures, polyposis coli, bone density and vascular defects in the eye (Osteoporosis-pseudoglioma Syndrome, OPPG) and other eye diseases or syndromes associated with defects and/or damaged photoreceptors, familial exudative vitreoretinopathy, retinal angiogenesis, early coronary disease, tetra-amelia, Müllerian-duct regression and virilization, SERKAL syndrome, type II diabetes, Fuhrmann syndrome, Al-Awadi/Raas-Rothschild/Schinzel phocomelia syndrome, odonto-onycho-dermal dysplasia, obesity, split-hand/foot malformation, caudal duplication, tooth agenesis, Wilms tumor, skeletal dysplasia, focal dermal hypoplasia, autosomal recessive anonychia, neural tube defects, alpha-thalassemia (ATRX) syndrome, fragile X syndrome, ICF syndrome, Angelman's syndrome, Prader-Willi syndrome, Beckwith-Wiedemann Syndrome, Norrie disease and Rett syndrome.

[0559] In some embodiments, non-limiting examples of eye diseases which can be treated with the compounds and compositions provided herein include age-related macular degeneration (AMD or ARMD), rod cone dystrophy, retinitis pigmentosa (RP), acute idiopathic blind spot enlargement (AIBSE), acute zonal occult outer retinopathy (AZOOR), acute macular neuroretinopathy (AMN), multiple evanescent white dot syndrome (MEWDS), multifocal choroiditis, opticneuropathy. Further causes of photoreceptor damage that can be treated with the compounds and compositions provided herein include retinal detachment, vascular disturbance, eye tumors or extreme light damage.

[0560] With respect to cancer, the Wnt pathway is known to be constitutively activated in a variety of cancers including, for example, colon cancer, hepatocellular carcinoma, lung cancer, ovarian cancer, prostate cancer, pancreatic cancer and leukemias such as CML, CLL and T-ALL. Accordingly, the compounds and compositions described herein may be used to treat these cancers in which the Wnt pathway is constitutively activated. In certain embodiments, the cancer is chosen from hepatocellular carcinoma, colon cancer, leukemia, lymphoma, sarcoma and ovarian cancer.

[0561] Other cancers can also be treated with the compounds and compositions described herein.

[0562] More particularly, cancers that may be treated by the compounds, compositions and methods described herein include, but are not limited to, the following:

[0563] 1) Breast cancers, including, for example ER⁺ breast cancer, ER⁻ breast cancer, her2⁻ breast cancer, her2⁺ breast cancer, stromal tumors such as fibroadenomas, phyllodes tumors,

and sarcomas, and epithelial tumors such as large duct papillomas; carcinomas of the breast including *in situ* (noninvasive) carcinoma that includes ductal carcinoma *in situ* (including Paget's disease) and lobular carcinoma *in situ*, and invasive (infiltrating) carcinoma including, but not limited to, invasive ductal carcinoma, invasive lobular carcinoma, medullary carcinoma, colloid (mucinous) carcinoma, tubular carcinoma, and invasive papillary carcinoma; and miscellaneous malignant neoplasms. Further examples of breast cancers can include luminal A, luminal B, basal A, basal B, and triple negative breast cancer, which is estrogen receptor negative (ER⁻), progesterone receptor negative, and her2 negative (her2⁻). In some embodiments, the breast cancer may have a high risk Oncotype score.

[0564] 2) Cardiac cancers, including, for example sarcoma, e.g., angiosarcoma, fibrosarcoma, rhabdomyosarcoma, and liposarcoma; myxoma; rhabdomyoma; fibroma; lipoma and teratoma.

[0565] 3) Lung cancers, including, for example, bronchogenic carcinoma, e.g., squamous cell, undifferentiated small cell, undifferentiated large cell, and adenocarcinoma; alveolar and bronchiolar carcinoma; bronchial adenoma; sarcoma; lymphoma; chondromatous hamartoma; and mesothelioma.

[0566] 4) Gastrointestinal cancer, including, for example, cancers of the esophagus, e.g., squamous cell carcinoma, adenocarcinoma, leiomyosarcoma, and lymphoma; cancers of the stomach, e.g., carcinoma, lymphoma, and leiomyosarcoma; cancers of the pancreas, e.g., ductal adenocarcinoma, insulinoma, glucagonoma, gastrinoma, carcinoid tumors, and vipoma; cancers of the small bowel, e.g., adenocarcinoma, lymphoma, carcinoid tumors, Kaposi's sarcoma, leiomyoma, hemangioma, lipoma, neurofibroma, and fibroma; cancers of the large bowel, e.g., adenocarcinoma, tubular adenoma, villous adenoma, hamartoma, and leiomyoma.

[0567] 5) Genitourinary tract cancers, including, for example, cancers of the kidney, e.g., adenocarcinoma, Wilm's tumor (nephroblastoma), lymphoma, and leukemia; cancers of the bladder and urethra, e.g., squamous cell carcinoma, transitional cell carcinoma, and adenocarcinoma; cancers of the prostate, e.g., adenocarcinoma, and sarcoma; cancer of the testis, e.g., seminoma, teratoma, embryonal carcinoma, teratocarcinoma, choriocarcinoma, sarcoma, interstitial cell carcinoma, fibroma, fibroadenoma, adenomatoid tumors, and lipoma.

[0568] 6) Liver cancers, including, for example, hepatoma, e.g., hepatocellular carcinoma; cholangiocarcinoma; hepatoblastoma; angiosarcoma; hepatocellular adenoma; and hemangioma.

[0569] 7) Bone cancers, including, for example, osteogenic sarcoma (osteosarcoma), fibrosarcoma, malignant fibrous histiocytoma, chondrosarcoma, Ewing's sarcoma, malignant

lymphoma (reticulum cell sarcoma), multiple myeloma, malignant giant cell tumor chordoma, osteochondroma (osteocartilaginous exostoses), benign chondroma, chondroblastoma, chondromyxofibroma, osteoid osteoma and giant cell tumors.

[0570] 8) Nervous system cancers, including, for example, cancers of the skull, e.g., osteoma, hemangioma, granuloma, xanthoma, and osteitis deformans; cancers of the meninges, e.g., meningioma, meningiosarcoma, and gliomatosis; cancers of the brain, e.g., astrocytoma, medulloblastoma, glioma, ependymoma, germinoma (pinealoma), glioblastoma multiform, oligodendroglioma, schwannoma, retinoblastoma, and congenital tumors; and cancers of the spinal cord, e.g., neurofibroma, meningioma, glioma, and sarcoma.

[0571] 9) Gynecological cancers, including, for example, cancers of the uterus, e.g., endometrial carcinoma; cancers of the cervix, e.g., cervical carcinoma, and pre tumor cervical dysplasia; cancers of the ovaries, e.g., ovarian carcinoma, including serous cystadenocarcinoma, mucinous cystadenocarcinoma, unclassified carcinoma, granulosa theca cell tumors, Sertoli Leydig cell tumors, dysgerminoma, and malignant teratoma; cancers of the vulva, e.g., squamous cell carcinoma, intraepithelial carcinoma, adenocarcinoma, fibrosarcoma, and melanoma; cancers of the vagina, e.g., clear cell carcinoma, squamous cell carcinoma, botryoid sarcoma, and embryonal rhabdomyosarcoma; and cancers of the fallopian tubes, e.g., carcinoma.

[0572] 10) Hematologic cancers, including, for example, cancers of the blood, e.g., acute myeloid leukemia, chronic myeloid leukemia, acute lymphoblastic leukemia, chronic lymphocytic leukemia, myeloproliferative diseases, multiple myeloma, and myelodysplastic syndrome, Hodgkin's lymphoma, non-Hodgkin's lymphoma (malignant lymphoma) and Waldenström's macroglobulinemia.

[0573] 11) Skin cancers and skin disorders, including, for example, malignant melanoma and metastatic melanoma, basal cell carcinoma, squamous cell carcinoma, Kaposi's sarcoma, moles dysplastic nevi, lipoma, angioma, dermatofibroma, keloids, and scleroderma.

[0574] 12) Adrenal gland cancers, including, for example, neuroblastoma.

[0575] Cancers may be solid tumors that may or may not be metastatic. Cancers may also occur, as in leukemia, as a diffuse tissue. Thus, the term "tumor cell," as provided herein, includes a cell afflicted by any one of the above identified disorders.

[0576] A method of treating cancer using a compound or composition as described herein may be combined with existing methods of treating cancers, for example by chemotherapy, irradiation, or surgery (e.g., oophorectomy). In some embodiments, a compound or composition can be administered before, during, or after another anticancer agent or treatment.

[0577] The compounds and compositions described herein can be used as anti-angiogenesis agents and as agents for modulating and/or inhibiting the activity of protein kinases, thus providing treatments for cancer and other diseases associated with cellular proliferation mediated by protein kinases. For example, the compounds described herein can inhibit the activity of one or more kinases. Accordingly, provided herein is a method of treating cancer or preventing or reducing angiogenesis through kinase inhibition.

[0578] In addition, and including treatment of cancer, the compounds and compositions described herein can function as cell-cycle control agents for treating proliferative disorders in a patient. Disorders associated with excessive proliferation include, for example, cancers, scleroderma, immunological disorders involving undesired proliferation of leukocytes, and restenosis and other smooth muscle disorders. Furthermore, such compounds may be used to prevent de-differentiation of post-mitotic tissue and/or cells.

[0579] Diseases or disorders associated with uncontrolled or abnormal cellular proliferation include, but are not limited to, the following:

- a variety of cancers, including, but not limited to, carcinoma, hematopoietic tumors of lymphoid lineage, hematopoietic tumors of myeloid lineage, tumors of mesenchymal origin, tumors of the central and peripheral nervous system and other tumors including melanoma, seminoma and Kaposi's sarcoma.
- a disease process which features abnormal cellular proliferation, e.g., benign prostatic hyperplasia, familial adenomatous polyposis, neurofibromatosis, atherosclerosis, arthritis, glomerulonephritis, restenosis following angioplasty or vascular surgery, inflammatory bowel disease, transplantation rejection, endotoxic shock, and fungal infections. Fibrotic disorders such as skin fibrosis; scleroderma; progressive systemic fibrosis; lung fibrosis; muscle fibrosis; kidney fibrosis; glomerulosclerosis; glomerulonephritis; hypertrophic scar formation; uterine fibrosis; renal fibrosis; cirrhosis of the liver, liver fibrosis; fatty liver disease (FLD); adhesions, such as those occurring in the abdomen, pelvis, spine or tendons; chronic obstructive pulmonary disease; fibrosis following myocardial infarction; pulmonary fibrosis; fibrosis and scarring associated with diffuse/interstitial lung disease; central nervous system fibrosis, such as fibrosis following stroke; fibrosis associated with neuro-degenerative disorders such as Alzheimer's Disease or multiple sclerosis; fibrosis associated with proliferative vitreoretinopathy (PVR); restenosis; endometriosis; ischemic disease and radiation fibrosis.

- defective apoptosis-associated conditions, such as cancers (including but not limited to those types mentioned herein), viral infections (including but not limited to herpesvirus, poxvirus, Epstein-Barr virus, Sindbis virus and adenovirus), prevention of AIDS development in HIV-infected individuals, autoimmune diseases (including but not limited to systemic lupus erythematosus, rheumatoid arthritis, sepsis, ankylosing spondylitis, psoriasis, scleroderma, autoimmune mediated glomerulonephritis, inflammatory bowel disease and autoimmune diabetes mellitus), neuro-degenerative disorders (including but not limited to Alzheimer's disease, lung disease, amyotrophic lateral sclerosis, retinitis pigmentosa, Parkinson's disease, AIDS-related dementia, spinal muscular atrophy and cerebellar degeneration), myelodysplastic syndromes, aplastic anemia, ischemic injury associated with myocardial infarctions, stroke and reperfusion injury, arrhythmia, atherosclerosis, toxin-induced or alcohol related liver diseases, hematological diseases (including but not limited to chronic anemia and aplastic anemia), degenerative diseases of the musculoskeletal system (including but not limited to osteoporosis and arthritis), tendinopathies such as tendinitis and tendinosis, aspirin-sensitive rhinosinusitis, cystic fibrosis, multiple sclerosis, kidney diseases and cancer pain.
- genetic diseases due to mutations in Wnt signaling components, such as polyposis coli, bone density and vascular defects in the eye (Osteoporosis-pseudoglioma Syndrome, OPPG) and other eye diseases or syndromes associated with defects and/or damaged photoreceptors, familial exudative vitreoretinopathy, retinal angiogenesis, early coronary disease, tetra-amelia, Müllerian-duct regression and virilization, SERKAL syndrome, type II diabetes, Fuhrmann syndrome, Al-Awadi/Raas-Rothschild/Schinzle phocomelia syndrome, odonto-onycho-dermal dysplasia, obesity, split-hand/foot malformation, caudal duplication, tooth agenesis, Wilms tumor, skeletal dysplasia, focal dermal hypoplasia, autosomal recessive anonychia, neural tube defects, alpha-thalassemia (ATRX) syndrome, fragile X syndrome, ICF syndrome, Angelman's syndrome, Prader-Willi syndrome, Beckwith-Wiedemann Syndrome, Norrie disease and Rett syndrome.

[0580] The compounds and compositions described herein can be used to treat neurological conditions, disorders and/or diseases caused by dysfunction in the Wnt signaling pathway. Non-limiting examples of neurological conditions/disorders/diseases which can be treated with the compounds and compositions provided herein include Alzheimer's disease,

aphasia, apraxia, arachnoiditis, ataxia telangiectasia, attention deficit hyperactivity disorder, auditory processing disorder, autism, alcoholism, Bell's palsy, bipolar disorder, brachial plexus injury, Canavan disease, carpal tunnel syndrome, causalgia, central pain syndrome, central pontine myelinolysis, centronuclear myopathy, cephalic disorder, cerebral aneurysm, cerebral arteriosclerosis, cerebral atrophy, cerebral gigantism, cerebral palsy, cerebral vasculitis, cervical spinal stenosis, Charcot-Marie-Tooth disease, Chiari malformation, chronic fatigue syndrome, chronic inflammatory demyelinating polyneuropathy (CIDP), chronic pain, Coffin–Lowry syndrome, complex regional pain syndrome, compression neuropathy, congenital facial diplegia, corticobasal degeneration, cranial arteritis, craniosynostosis, Creutzfeldt-Jakob disease, cumulative trauma disorder, Cushing's syndrome, cytomegalic inclusion body disease (CIBD), Dandy-Walker syndrome, Dawson disease, de Morsier's syndrome, Dejerine-Klumpke palsy, Dejerine-Sottas disease, delayed sleep phase syndrome, dementia, dermatomyositis, developmental dyspraxia, diabetic neuropathy, diffuse sclerosis, Dravet syndrome, dysautonomia, dyscalculia, dysgraphia, dyslexia, dystonia, empty sella syndrome, encephalitis, encephalocele, encephalotrigeminal angiomatosis, encopresis, epilepsy, Erb's palsy, erythromelalgia, essential tremor, Fabry's disease, Fahr's syndrome, familial spastic paralysis, febrile seizure, Fisher syndrome, Friedreich's ataxia, fibromyalgia, Foville's syndrome, Gaucher's disease, Gerstmann's syndrome, giant cell arteritis, giant cell inclusion disease, globoid cell leukodystrophy, gray matter heterotopia, Guillain-Barré syndrome, HTLV-1 associated myelopathy, Hallervorden-Spatz disease, hemifacial spasm, hereditary spastic paraplegia, heredopathia atactica polyneuritiformis, herpes zoster oticus, herpes zoster, Hirayama syndrome, holoprosencephaly, Huntington's disease, hydranencephaly, hydrocephalus, hypercortisolism, hypoxia, immune-mediated encephalomyelitis, inclusion body myositis, incontinentia pigmenti, infantile phytanic acid storage disease, infantile Refsum disease, infantile spasms, inflammatory myopathy, intracranial cyst, intracranial hypertension, Joubert syndrome, Karak syndrome, Kearns-Sayre syndrome, Kennedy disease, Kinsbourne syndrome, Klippel Feil syndrome, Krabbe disease, Kugelberg-Welander disease, kuru, Lafora disease, Lambert-Eaton myasthenic syndrome, Landau-Kleffner syndrome, lateral medullary (Wallenberg) syndrome, Leigh's disease, Lennox-Gastaut syndrome, Lesch-Nyhan syndrome, leukodystrophy, Lewy body dementia, lissencephaly, locked-in syndrome, Lou Gehrig's disease, lumbar disc disease, lumbar spinal stenosis, Lyme disease, Machado-Joseph disease (Spinocerebellar ataxia type 3), macrencephaly, macropsia, megalencephaly, Melkersson-Rosenthal syndrome, Meniere's disease, meningitis, Menkes disease, metachromatic leukodystrophy, microcephaly, micropsia, Miller Fisher syndrome, misophonia, mitochondrial myopathy, Mobius syndrome, monomelic amyotrophy, motor neuron disease, motor skills disorder, Moyamoya disease,

mucopolysaccharidoses, multi-infarct dementia, multifocal motor neuropathy, multiple sclerosis, multiple system atrophy, muscular dystrophy, myalgic encephalomyelitis, myasthenia gravis, myelinoclastic diffuse sclerosis, myoclonic Encephalopathy of infants, myoclonus, myopathy, myotubular myopathy, myotonia congenital, narcolepsy, neurofibromatosis, neuroleptic malignant syndrome, lupus erythematosus, neuromyotonia, neuronal ceroid lipofuscinosis, Niemann-Pick disease, O'Sullivan-McLeod syndrome, occipital Neuralgia, occult Spinal Dysraphism Sequence, Ohtahara syndrome, olivopontocerebellar atrophy, opsoclonus myoclonus syndrome, optic neuritis, orthostatic hypotension, palinopsia, paresthesia, Parkinson's disease, paramyotonia Congenita, paraneoplastic diseases, paroxysmal attacks, Parry-Romberg syndrome, Pelizaeus-Merzbacher disease, periodic paralyses, peripheral neuropathy, photic sneeze reflex, phytanic acid storage disease, Pick's disease, polymicrogyria (PMG), polymyositis, porencephaly, post-polio syndrome, postherpetic neuralgia (PHN), postural hypotension, Prader-Willi syndrome, primary lateral sclerosis, prion diseases, progressive hemifacial atrophy, progressive multifocal leukoencephalopathy, progressive supranuclear palsy, pseudotumor cerebri, Ramsay Hunt syndrome type I, Ramsay Hunt syndrome type II, Ramsay Hunt syndrome type III, Rasmussen's encephalitis, reflex neurovascular dystrophy, Refsum disease, restless legs syndrome, retrovirus-associated myelopathy, Rett syndrome, Reye's syndrome, rhythmic movement disorder, Romberg syndrome, Saint Vitus dance, Sandhoff disease, schizophrenia, Schilder's disease, schizencephaly, sensory integration dysfunction, septo-optic dysplasia, Shy-Drager syndrome, Sjögren's syndrome, snatiation, Sotos syndrome, spasticity, spina bifida, spinal cord tumors, spinal muscular atrophy, spinocerebellar ataxia, Steele-Richardson-Olszewski syndrome, Stiff-person syndrome, stroke, Sturge-Weber syndrome, subacute sclerosing panencephalitis, subcortical arteriosclerotic encephalopathy, superficial siderosis, Sydenham's chorea, syncope, synesthesia, syringomyelia, tarsal tunnel syndrome, tardive dyskinesia, tardive dysphrenia, Tarlov cyst, Tay-Sachs disease, temporal arteritis, tetanus, tethered spinal cord syndrome, Thomsen disease, thoracic outlet syndrome, tic douloureux, Todd's paralysis, Tourette syndrome, toxic encephalopathy, transient ischemic attack, transmissible spongiform encephalopathies, transverse myelitis, tremor, trigeminal neuralgia, tropical spastic paraparesis, trypanosomiasis, tuberous sclerosis, ubisiosis, Von Hippel-Lindau disease (VHL), Viliuisk Encephalomyelitis (VE), Wallenberg's syndrome, Werdnig, Hoffman disease, west syndrome, Williams syndrome, Wilson's disease and Zellweger syndrome.

[0581] The compounds and compositions may also be useful in the inhibition of the development of invasive cancer, tumor angiogenesis and metastasis.

[0582] In some embodiments, the disclosure provides a method for treating a disease or disorder associated with aberrant cellular proliferation by administering to a patient in need of

such treatment an effective amount of one or more of the compounds of Formula (I), in combination (simultaneously or sequentially) with at least one other agent.

[0583] In some embodiments, the disclosure provides a method of treating or ameliorating in a patient a disorder or disease selected from the group consisting of: cancer, pulmonary fibrosis, idiopathic pulmonary fibrosis (IPF), degenerative disc disease, bone/osteoporotic fractures, bone or cartilage disease, and osteoarthritis, the method comprising administering to the patient a therapeutically effective amount of a compound according to claim 1, or a pharmaceutically acceptable salt thereof.

[0584] In some embodiments, the pharmaceutical composition comprises a therapeutically effective amount of a compound of Formula (I), or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable excipient.

[0585] In some embodiments, the method of treats a disorder or disease in which aberrant Wnt signaling is implicated in a patient, the method comprises administering to the patient a therapeutically effective amount of a compound of Formula (I), or a pharmaceutically acceptable salt thereof.

[0586] In some embodiments, the disorder or disease is cancer.

[0587] In some embodiments, the disorder or disease is systemic inflammation.

[0588] In some embodiments, the disorder or disease is metastatic melanoma.

[0589] In some embodiments, the disorder or disease is fatty liver disease.

[0590] In some embodiments, the disorder or disease is liver fibrosis.

[0591] In some embodiments, the disorder or disease is tendonitis.

[0592] In some embodiments, the disorder or disease is damage to a tendon which would benefit from tendon regeneration.

[0593] In some embodiments, the disorder or disease is diabetes.

[0594] In some embodiments, the disorder or disease is degenerative disc disease.

[0595] In some embodiments, the disorder or disease is osteoarthritis.

[0596] In some embodiments, the disorder or disease is diabetic retinopathy.

[0597] In some embodiments, the disorder or disease is pulmonary fibrosis.

[0598] In some embodiments, the disorder or disease is idiopathic pulmonary fibrosis (IPF).

[0599] In some embodiments, the disorder or disease is degenerative disc disease.

[0600] In some embodiments, the disorder or disease is rheumatoid arthritis.

[0601] In some embodiments, the disorder or disease is scleroderma.

[0602] In some embodiments, the disorder or disease is a mycotic or viral infection.

- [0603] In some embodiments, the disorder or disease is a bone or cartilage disease.
- [0604] In some embodiments, the disorder or disease is Alzheimer's disease.
- [0605] In some embodiments, the disorder or disease is osteoarthritis.
- [0606] In some embodiments, the disorder or disease is lung disease
- [0607] In some embodiments, the disorder or disease is a genetic disease caused by mutations in Wnt signaling components, wherein the genetic disease is selected from: polyposis coli, osteoporosis-pseudoglioma syndrome, familial exudative vitreoretinopathy, retinal angiogenesis, early coronary disease, tetra-amelia syndrome, Müllerian-duct regression and virilization, SERKAL syndrome, diabetes mellitus type 2, Fuhrmann syndrome, Al-Awadi/Raas-Rothschild/Schinzel phocomelia syndrome, odonto-onycho-dermal dysplasia, obesity, split-hand/foot malformation, caudal duplication syndrome, tooth agenesis, Wilms tumor, skeletal dysplasia, focal dermal hypoplasia, autosomal recessive anonychia, neural tube defects, alpha-thalassemia (ATRX) syndrome, fragile X syndrome, ICF syndrome, Angelman syndrome, Prader-Willi syndrome, Beckwith-Wiedemann Syndrome, Norrie disease and Rett syndrome.
- [0608] In some embodiments, the patient is a human.
- [0609] In some embodiments, the cancer is chosen from: hepatocellular carcinoma, colon cancer, breast cancer, pancreatic cancer, chronic myeloid leukemia (CML), chronic myelomonocytic leukemia, chronic lymphocytic leukemia (CLL), acute myeloid leukemia, acute lymphocytic leukemia, Hodgkin lymphoma, lymphoma, sarcoma and ovarian cancer.
- [0610] In some embodiments, the cancer is chosen from: lung cancer - non-small cell, lung cancer - small cell, multiple myeloma, nasopharyngeal cancer, neuroblastoma, osteosarcoma, penile cancer, pituitary tumors, prostate cancer, retinoblastoma, rhabdomyosarcoma, salivary gland cancer, skin cancer - basal and squamous cell, skin cancer – melanoma, small intestine cancer, stomach (gastric) cancers, testicular cancer, thymus cancer, thyroid cancer, uterine sarcoma, vaginal cancer, vulvar cancer, laryngeal or hypopharyngeal cancer, kidney cancer, Kaposi sarcoma, gestational trophoblastic disease, gastrointestinal stromal tumor, gastrointestinal carcinoid tumor, gallbladder cancer, eye cancer (melanoma and lymphoma), Ewing tumor, esophagus cancer, endometrial cancer, colorectal cancer, cervical cancer, brain or spinal cord tumor, bone metastasis, bone cancer, bladder cancer, bile duct cancer, anal cancer and adrenal cortical cancer.
- [0611] In some embodiments, the cancer is hepatocellular carcinoma.
- [0612] In some embodiments, the cancer is colon cancer.
- [0613] In some embodiments, the cancer is colorectal cancer.
- [0614] In some embodiments, the cancer is breast cancer.
- [0615] In some embodiments, the cancer is pancreatic cancer.

- [0616] In some embodiments, the cancer is chronic myeloid leukemia (CML).
- [0617] In some embodiments, the cancer is chronic myelomonocytic leukemia.
- [0618] In some embodiments, the cancer is chronic lymphocytic leukemia (CLL).
- [0619] In some embodiments, the cancer is acute myeloid leukemia.
- [0620] In some embodiments, the cancer is acute lymphocytic leukemia.
- [0621] In some embodiments, the cancer is Hodgkin lymphoma.
- [0622] In some embodiments, the cancer is lymphoma.
- [0623] In some embodiments, the cancer is sarcoma.
- [0624] In some embodiments, the cancer is ovarian cancer.
- [0625] In some embodiments, the cancer is lung cancer - non-small cell.
- [0626] In some embodiments, the cancer is lung cancer - small cell.
- [0627] In some embodiments, the cancer is multiple myeloma.
- [0628] In some embodiments, the cancer is nasopharyngeal cancer.
- [0629] In some embodiments, the cancer is neuroblastoma.
- [0630] In some embodiments, the cancer is osteosarcoma.
- [0631] In some embodiments, the cancer is penile cancer.
- [0632] In some embodiments, the cancer is pituitary tumors.
- [0633] In some embodiments, the cancer is prostate cancer.
- [0634] In some embodiments, the cancer is retinoblastoma.
- [0635] In some embodiments, the cancer is rhabdomyosarcoma.
- [0636] In some embodiments, the cancer is salivary gland cancer.
- [0637] In some embodiments, the cancer is skin cancer - basal and squamous cell.
- [0638] In some embodiments, the cancer is skin cancer – melanoma.
- [0639] In some embodiments, the cancer is small intestine cancer.
- [0640] In some embodiments, the cancer is stomach (gastric) cancers.
- [0641] In some embodiments, the cancer is testicular cancer.
- [0642] In some embodiments, the cancer is thymus cancer.
- [0643] In some embodiments, the cancer is thyroid cancer.
- [0644] In some embodiments, the cancer is uterine sarcoma.
- [0645] In some embodiments, the cancer is vaginal cancer.
- [0646] In some embodiments, the cancer is vulvar cancer.
- [0647] In some embodiments, the cancer is Wilms tumor.
- [0648] In some embodiments, the cancer is laryngeal or hypopharyngeal cancer.
- [0649] In some embodiments, the cancer is kidney cancer.

- [0650] In some embodiments, the cancer is Kaposi sarcoma.
- [0651] In some embodiments, the cancer is gestational trophoblastic disease.
- [0652] In some embodiments, the cancer is gastrointestinal stromal tumor.
- [0653] In some embodiments, the cancer is gastrointestinal carcinoid tumor.
- [0654] In some embodiments, the cancer is gallbladder cancer.
- [0655] In some embodiments, the cancer is eye cancer (melanoma and lymphoma).
- [0656] In some embodiments, the cancer is Ewing tumor.
- [0657] In some embodiments, the cancer is esophagus cancer.
- [0658] In some embodiments, the cancer is endometrial cancer.
- [0659] In some embodiments, the cancer is colorectal cancer.
- [0660] In some embodiments, the cancer is cervical cancer.
- [0661] In some embodiments, the cancer is brain or spinal cord tumor.
- [0662] In some embodiments, the cancer is bone metastasis.
- [0663] In some embodiments, the cancer is bone cancer.
- [0664] In some embodiments, the cancer is bladder cancer.
- [0665] In some embodiments, the cancer is bile duct cancer.
- [0666] In some embodiments, the cancer is anal cancer.
- [0667] In some embodiments, the cancer is adrenal cortical cancer.
- [0668] In some embodiments, the disorder or disease is a neurological condition, disorder or disease, wherein the neurological condition/disorder/disease is selected from: Alzheimer's disease, frontotemporal dementias, dementia with lewy bodies, prion diseases, Parkinson's disease, Huntington's disease, progressive supranuclear palsy, corticobasal degeneration, multiple system atrophy, amyotrophic lateral sclerosis (ALS), inclusion body myositis, autism, degenerative myopathies, diabetic neuropathy, other metabolic neuropathies, endocrine neuropathies, orthostatic hypotension, multiple sclerosis and Charcot-Marie-Tooth disease.
- [0669] In some embodiments, the compound of Formula (I) inhibits one or more proteins in the Wnt pathway.
- [0670] In some embodiments, the compound of Formula (I) inhibits signaling induced by one or more Wnt proteins.
- [0671] In some embodiments, the Wnt proteins are chosen from: WNT1, WNT2, WNT2B, WNT3, WNT3A, WNT4, WNT5A, WNT5B, WNT6, WNT7A, WNT7B, WNT8A, WNT8B, WNT9A, WNT9B, WNT10A, WNT10B, WNT11, and WNT16.

[0672] In some embodiments, the method inhibits one or more proteins in the Wnt pathway, the method comprises contacting a cell with an effective amount of a compound of Formula (I).

[0673] In some embodiments, the cell is a human cell.

[0674] In some embodiments, the human cell is a cancerous cell.

[0675] In some embodiments, the cancerous cell is a colon cancer cell.

[0676] In some embodiments, the contacting is *in vitro*.

[0677] In some embodiments, the compound of Formula (I) inhibits a kinase activity.

[0678] In some embodiments, the method treats a disease or disorder mediated by the Wnt pathway in a patient, the method comprises administering to the patient a therapeutically effective amount of a compound (or compounds) of Formula (I), or a pharmaceutically acceptable salt thereof.

[0679] In some embodiments, the compound of Formula (I) inhibits one or more Wnt proteins.

[0680] In some embodiments, the method treats a disease or disorder mediated by kinase activity in a patient, the method comprises administering to the patient a therapeutically effective amount of a compound (or compounds) of Formula (I), or a pharmaceutically acceptable salt thereof.

[0681] In some embodiments, the disease or disorder comprises tumor growth, cell proliferation, or angiogenesis.

[0682] In some embodiments, the method inhibits the activity of a protein kinase receptor, the method comprises contacting the receptor with an effective amount of a compound (or compounds) of Formula (I), or a pharmaceutically acceptable salt thereof.

[0683] In some embodiments, the method treats a disease or disorder associated with aberrant cellular proliferation in a patient; the method comprises administering to the patient a therapeutically effective amount of a compound (or compounds) of Formula (I), or a pharmaceutically acceptable salt thereof.

[0684] In some embodiments, the method prevents or reduces angiogenesis in a patient; the method comprises administering to the patient a therapeutically effective amount of a compound (or compounds) of Formula (I), or a pharmaceutically acceptable salt thereof.

[0685] In some embodiments, the method prevents or reduces abnormal cellular proliferation in a patient; the method comprises administering to the patient a therapeutically effective amount of a compound (or compounds) of Formula (I), or a pharmaceutically acceptable salt thereof.

[0686] In some embodiments, the method treats a disease or disorder associated with aberrant cellular proliferation in a patient, the method comprises administering to the patient a pharmaceutical composition comprising one or more of the compounds of claim 1 in combination with a pharmaceutically acceptable carrier and one or more other agents.

[0687] Moreover, the compounds and compositions, for example, as inhibitors of the cyclin-dependent kinases (CDKs), can modulate the level of cellular RNA and DNA synthesis and therefore are expected to be useful in the treatment of viral infections such as HIV, human papilloma virus, herpes virus, Epstein-Barr virus, adenovirus, Sindbis virus, pox virus and the like.

[0688] Compounds and compositions described herein can inhibit the kinase activity of, for example, CDK/cyclin complexes, such as those active in the G₀ or G₁ stage of the cell cycle, e.g., CDK2, CDK4, and/or CDK6 complexes.

Evaluation of Biological Activity

[0689] The biological activity of the compounds described herein can be tested using any suitable assay known to those of skill in the art, see, e.g., WO 2001/053268 and WO 2005/009997. For example, the activity of a compound may be tested using one or more of the test methods outlined below.

[0690] In one example, tumor cells may be screened for Wnt independent growth. In such a method, tumor cells of interest are contacted with a compound (i.e. inhibitor) of interest, and the proliferation of the cells, e.g. by uptake of tritiated thymidine, is monitored. In some embodiments, tumor cells may be isolated from a candidate patient who has been screened for the presence of a cancer that is associated with a mutation in the Wnt signaling pathway. Candidate cancers include, without limitation, those listed above.

[0691] In another example, one may utilize *in vitro* assays for Wnt biological activity, e.g. stabilization of β -catenin and promoting growth of stem cells. Assays for biological activity of Wnt include stabilization of β -catenin, which can be measured, for example, by serial dilutions of a candidate inhibitor composition. An exemplary assay for Wnt biological activity contacts a candidate inhibitor with cells containing constitutively active Wnt/ β -catenin signaling. The cells are cultured for a period of time sufficient to stabilize β -catenin, usually at least about 1 hour, and lysed. The cell lysate is resolved by SDS PAGE, then transferred to nitrocellulose and probed with antibodies specific for β -catenin.

[0692] In a further example, the activity of a candidate compound can be measured in a *Xenopus* secondary axis bioassay (Leyns, L. *et al. Cell* (1997), 88(6), 747-756).

[0693] To further illustrate this invention, the following examples are included. The examples should not, of course, be construed as specifically limiting the invention. Variations of these examples within the scope of the claims are within the purview of one skilled in the art and are considered to fall within the scope of the invention as described, and claimed herein. The reader will recognize that the skilled artisan, armed with the present disclosure, and skill in the art is able to prepare and use the invention without exhaustive examples.

EXAMPLES

Compound preparation

[0694] The starting materials used in preparing the compounds of the invention are known, made by known methods, or are commercially available. It will be apparent to the skilled artisan that methods for preparing precursors and functionality related to the compounds claimed herein are generally described in the literature. The skilled artisan given the literature and this disclosure is well equipped to prepare any of the compounds.

[0695] It is recognized that the skilled artisan in the art of organic chemistry can readily carry out manipulations without further direction, that is, it is well within the scope and practice of the skilled artisan to carry out these manipulations. These include reduction of carbonyl compounds to their corresponding alcohols, oxidations, acylations, aromatic substitutions, both electrophilic and nucleophilic, etherifications, esterification and saponification and the like. These manipulations are discussed in standard texts such as *March's Advanced Organic Chemistry: Reactions, Mechanisms, and Structure* 7th Ed., John Wiley & Sons (2013), Carey and Sundberg, *Advanced Organic Chemistry* 5th Ed., Springer (2007), *Comprehensive Organic Transformations: A Guide to Functional Group Transformations*, 2nd Ed., John Wiley & Sons (1999) (incorporated herein by reference in its entirety) and the like.

[0696] The skilled artisan will readily appreciate that certain reactions are best carried out when other functionality is masked or protected in the molecule, thus avoiding any undesirable side reactions and/or increasing the yield of the reaction. Often the skilled artisan utilizes protecting groups to accomplish such increased yields or to avoid the undesired reactions. These reactions are found in the literature and are also well within the scope of the skilled artisan. Examples of many of these manipulations can be found for example in T. Greene and P. Wuts *Protective Groups in Organic Synthesis*, 4th Ed., John Wiley & Sons (2007), incorporated herein by reference in its entirety.

[0697] Trademarks used herein are examples only and reflect illustrative materials used at the time of the invention. The skilled artisan will recognize that variations in lot,

manufacturing processes, and the like, are expected. Hence the examples, and the trademarks used in them are non-limiting, and they are not intended to be limiting, but are merely an illustration of how a skilled artisan may choose to perform one or more of the embodiments of the invention.

[0698] ^1H nuclear magnetic resonance spectra (NMR) were measured in the indicated solvents on a Bruker NMR spectrometer (Avance TM DRX300, 300 MHz for ^1H or Avance TM DRX500, 500 MHz for ^1H) or Varian NMR spectrometer (Mercury 400BB, 400 MHz for ^1H). Peak positions are expressed in parts per million (ppm) downfield from tetramethylsilane. The peak multiplicities are denoted as follows, s, singlet; d, doublet; t, triplet; q, quartet; ABq, AB quartet; quin, quintet; sex, sextet; sep, septet; non, nonet; dd, doublet of doublets; ddd, doublet of doublets of doublets; d/ABq, doublet of AB quartet; dt, doublet of triplets; td, triplet of doublets; dq, doublet of quartets; m, multiplet.

[0699] The following abbreviations have the indicated meanings:

$\text{BH}_3\text{-Me}_2\text{S}$ = borane dimethyl sulfide complex

$(\text{Boc})_2\text{O}$ = di-*tert*-butyl dicarbonate

brine = saturated aqueous sodium chloride

CDCl_3 = deuterated chloroform

CD_3OD = deuterated methanol

DCAD = di-(4-chlorobenzyl)azodicarboxylate

DCE = dichloroethane

DCM = dichloromethane

DEAD = diethyl azodicarboxylate

DHP = dihydropyran

DMAP = 4-dimethylaminopyridine

DMF = N,N-dimethylformamide

DMSO-d_6 = deuterated dimethylsulfoxide

ESIMS = electron spray mass spectrometry

EtOAc = ethyl acetate

EtOH = ethanol

HCl = hydrochloric acid

HOAc = acetic acid

K_2CO_3 = potassium carbonate

KOAc = potassium acetate

LC/MS = liquid chromatography–mass spectrometry

MeOH = methanol

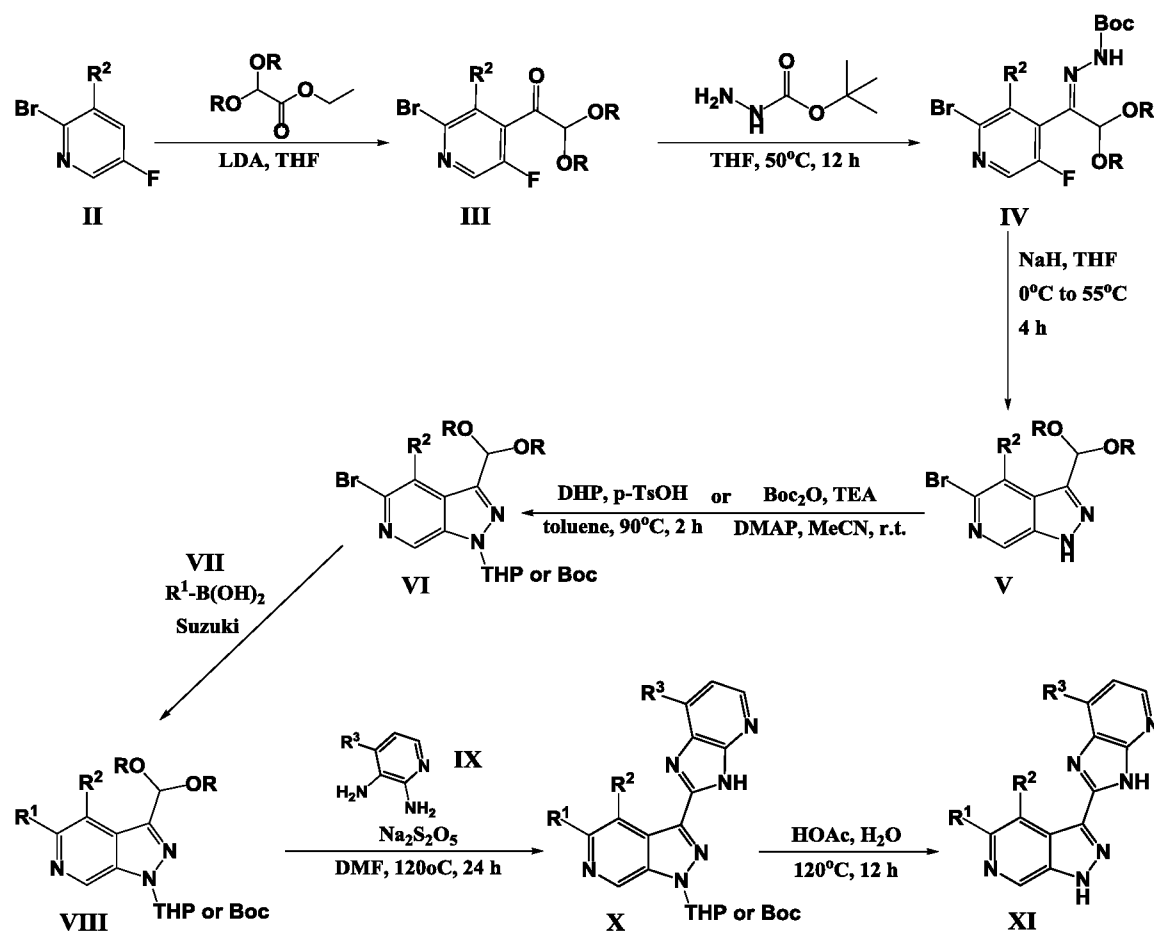
MgSO₄ = magnesium sulfate
MsCl = methanesulfonyl chloride or mesyl chloride
MW = microwave
NaBH₄ = sodium borohydride
NaBH(OAc)₃ = sodium triacetoxyborohydride
NaCNBH₃ = sodium cyanoborohydride
NaHCO₃ = sodium bicarbonate
NaOH = sodium hydroxide
Na₂S₂O₅ = sodium metabisulfite or sodium pyrosulfite
NH₄OH = ammonium hydroxide
NMR = nuclear magnetic resonance
ON = overnight
Pd/C = palladium(0) on carbon
Pd(dppf)Cl₂ = 1,1'-bis(diphenylphosphino)ferrocene]palladium(II) chloride
Pd(PPh₃)₄ = tetrakis(triphenylphosphine)palladium(0)
Pd(PPh₃)₂Cl₂ = bis(triphenylphosphine)palladium(II) dichloride
PE = petroleum ether
Pin₂B₂ = bis(pinacolato)diboron
PPh₃ = triphenylphosphine
PPTS = pyridinium p-toluenesulfonate
r.t. = room temperature
SEM-Cl = 2-(trimethylsilyl)ethoxymethyl chloride
TEA = triethylamine
TFA = trifluoroacetic acid
THF = tetrahydrofuran
THP = tetrahydropyran
TLC = thin layer chromatography
p-TsOH = p-toluenesulfonic acid

[0700] The following example schemes are provided for the guidance of the reader, and collectively represent an example method for making the compounds provided herein. Furthermore, other methods for preparing compounds of the invention will be readily apparent to the person of ordinary skill in the art in light of the following reaction schemes and examples. The skilled artisan is thoroughly equipped to prepare these compounds by those methods given the

literature and this disclosure. The compound numberings used in the synthetic schemes depicted below are meant for those specific schemes only, and should not be construed as or confused with same numberings in other sections of the application. Unless otherwise indicated, all variables are as defined above.

General procedure

[0701] Compounds of Formula (I) of the present invention can be prepared as depicted in Scheme 1.



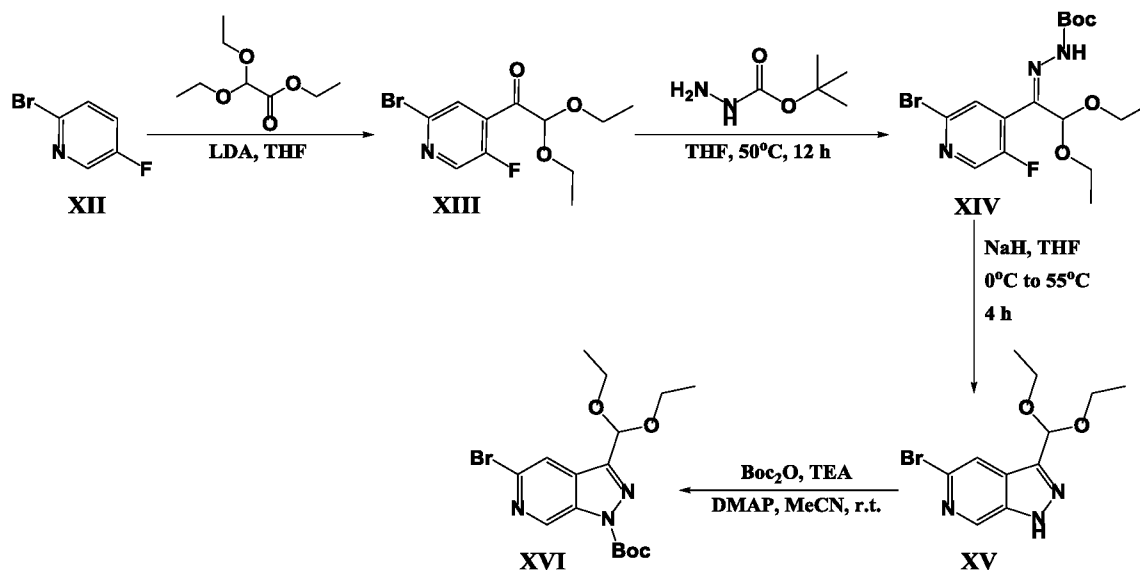
Scheme 1

[0702] Scheme 1 describes an alternative method for preparation of 1H-pyrazolo[3,4-c]pyridine derivatives (XI) by first acylating a 2-bromo-5-fluoropyridine (II) with an ethyl 2,2-dialkoxyacetate to produce the acetal protected oxoacetaldehyde (III). The keto group was then converted to the Boc-protected hydrazone (IV) followed by base cyclization to the 1H-pyrazolo[3,4-c]pyridine (V). The pyrazolopyridine (V) is then protected with either a Boc or THP

(VI) followed by Suzuki coupling with various boronic acids (VII). The pyrazolopyridine acetyl (VIII) is reacted with various 1,2-diamines (IX) to produce (X). Final deprotection of the pyrazole nitrogen yields the desired 1H-pyrazolo[3,4-*c*]pyridine derivatives (XI).

Illustrative Compound Examples

[0703] Preparation of Boc-protected intermediate (XVI) is depicted below in Scheme 2.



Scheme 2

Step 1

[0704] To a solution of 2-bromo-5-fluoropyridine (XII) (100.0 g, 568.21 mmol, 1.0 eq) in THF (1000 mL) was added a solution of LDA (66.95 g, 625.04 mmol, 1.10 eq) drop-wise at -78°C over a period of 1 h under N_2 . The reaction mixture was stirred at -78°C for 30 min. Then a solution of ethyl 2,2-diethoxyacetate (120.15 g, 681.86 mmol, 1.20 eq) was added at -78°C over a period of 1 h. The reaction mixture was stirred at -78°C for another 1 h. TLC (PE:EtOAc=10:1) showed that the starting material was consumed completely. The reaction was quenched by NH_4Cl slowly and then extracted with EtOAc (1000 mL x 2). The combined organic phase was washed with saturated brine (500 mL x 2), dried over anhydrous Na_2SO_4 , filtered and concentrated in vacuo. The residue was purified by chromatography on silica gel (PE:EtOAc=10:1) to give 1-(2-bromo-5-fluoropyridin-4-yl)-2,2-diethoxyethan-1-one (XIII) (160.0 g, 522.65 mmol, 92.0% yield) as a yellow solid. ^1H NMR (CDCl_3 , 400 MHz) δ ppm 1.26 (t, $J=6.8\text{Hz}$, 6H), 3.69 (q, $J=7.2\text{Hz}$, 2H), 3.79 (q, $J=7.2\text{Hz}$, 2H), 5.23 (d, $J=2\text{Hz}$, 1H), 7.85 (d, $J=4.8\text{Hz}$, 1H), 8.39 (d, $J=1.2\text{Hz}$, 1H); ESIMS found for $\text{C}_{11}\text{H}_{13}\text{BrFNO}_3$ m/z 306.1 (M+H).

Step 2

[0705] To a mixture of 1-(2-bromo-5-fluoropyridin-4-yl)-2,2-diethoxyethan-1-one (**XIII**) (142.0 g, 463.86 mmol, 1.0 eq) in THF (2 L) *tert*-butyl hydrazinecarboxylate (61.30 g, 463.86 mmol, 1.0 eq) in one portion at room temperature. The mixture was stirred at 55°C for 60 h. TLC (PE:EtOAc=2:1) showed that most of the starting material was consumed. The crude mixture of *tert*-butyl 2-(1-(2-bromo-5-fluoropyridin-4-yl)-2,2-diethoxyethylidene)hydrazine-1-carboxylate (**XIV**) was used in the next step without further purification.

Step 3

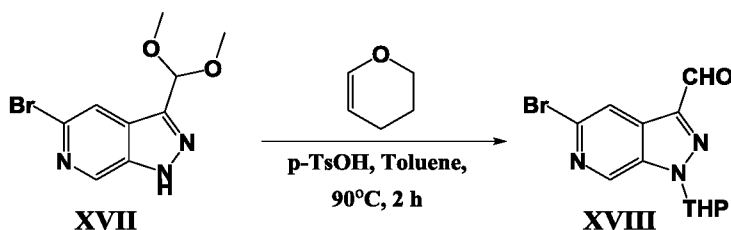
[0706] To a solution of *tert*-butyl 2-(1-(2-bromo-5-fluoropyridin-4-yl)-2,2-diethoxyethylidene)hydrazine-1-carboxylate (**XIV**) (190.0 g, 452.09 mmol, 1.0 eq) in THF (2 L) was added NaH (36.17 g, 904.18 mmol, 2.0 eq) in portions at 0°C over 0.5 h. The mixture was stirred at 55°C for 4 hours. TLC (PE:EtOAc=2:1) showed the material was consumed completely. The mixture was cooled to 0°C. The mixture was poured into 10% aqueous NH₄Cl (1000 mL). The aqueous phase was extracted with EtOAc (800 mL x 3). The combined organic phase was washed with saturated brine (800 mL), dried with anhydrous Na₂SO₄, filtered and concentrated in vacuum. The residue was purified by silica gel chromatography (PE/EtOAc=10:1→5:1→1:4) to afford 5-bromo-3-(diethoxymethyl)-1H-pyrazolo[3,4-*c*]pyridine (**XV**) (67.0 g, 223.22 mmol, 49.4% yield) as yellow solid. ¹H NMR (CDCl₃, 400 MHz) δ ppm 1.30 (t, *J*=7.03Hz, 6H), 3.60-3.86 (m, 4H), 5.98 (s, 1 H), 8.12 (d, *J*=1.13Hz, 1H), 8.96 (s, 1 H), 11.82 (brs, 1H); ESIMS found for C₁₁H₁₄BrN₃O₂ *m/z* 300.0 (M+H).

Step 4

[0707] To a solution of 5-bromo-3-(diethoxymethyl)-1H-pyrazolo[3,4-*c*]pyridine (**XV**) (20.0 g, 66.63 mmol, 1.0 eq) in CH₃CN (100 mL) was added Boc₂O (21.81 g, 99.95 mmol, 1.5 eq), DMAP (814.06 mg, 6.66 mmol, 0.10 eq) and TEA (13.49 g, 133.27 mmol, 2.0 eq) at room temperature. The mixture was stirred at 15°C for 1 hr. TLC (PE:EtOAc=5:1) showed that starting the material was consumed completely. The mixture was added water (50 mL) and extracted with EtOAc (40 mL x 2). The organic layers were washed with brine (60 mL) and concentrated under vacuum. The residue was purified by chromatography on silica gel (PE:EtOAc=10:1) to produce *tert*-butyl 5-bromo-3-(diethoxymethyl)-1H-pyrazolo[3,4-*c*]pyridine-1-carboxylate (**XVI**) (23.70 g, 59.21 mmol, 88.9% yield) as a light yellow oil. ¹H NMR (CDCl₃, 400 MHz) δ ppm 1.29 (t, *J*=6.8Hz,

6H), 1.75 (s, 9H), 3.60-3.71 (m, 2H), 3.75-3.87 (m, 2H), 5.79 (s, 1H), 8.14 (d, $J=1.00\text{Hz}$, 1H), 9.25 (s, 1H); ESIMS found for $\text{C}_{16}\text{H}_{22}\text{BrN}_3\text{O}_4$ m/z 400.0 (M+H).

[0708] Preparation of THP protected intermediate (XVIII) is depicted below in Scheme 3.

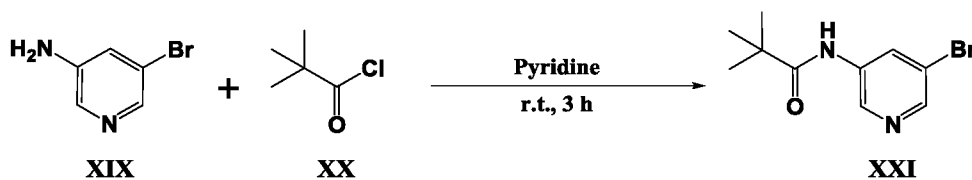


Scheme 3

Step 1

[0709] To a mixture of 5-bromo-3-(dimethoxymethyl)-1H-pyrazolo[3,4-c]pyridine (XVII) (26.0 g, 69.85 mmol, 1.0 eq) and 3,4-dihydro-2H-pyran (14.69 g, 174.63 mmol, 2.5 eq) in toluene (100 mL) was added 4-methylbenzenesulfonic acid (2.41 g, 13.97 mmol, 0.20 eq) in one portion at room temperature under N_2 . The mixture was heated to 90°C and stirred for 2 hr. LC/MS showed the reaction was completed. The mixture was extracted with EtOAc (100 mL x 3), washed with water (50 mL x 2) and brine (50 mL x 2). The organic layer was dried and concentrated to give a residue. The residue was purified by a column (PE:EtOAc=10:1 \rightarrow 8:1 \rightarrow 5:1) to give 5-bromo-1-(tetrahydro-2H-pyran-2-yl)-1H-pyrazolo[3,4-c]pyridine-3-carbaldehyde (XVIII) (8.50 g, 30.65 mmol, 43.9% yield). ^1H NMR (CDCl_3 , 500 MHz) δ ppm 1.68-1.80 (m, 2H), 1.80-1.91 (m, 1H), 2.08-2.27 (m, 2H), 2.41-2.56 (m, 1H), 3.77-3.88 (m, 1H), 3.90-4.00 (m, 1H), 5.93 (dd, $J=2.8\text{Hz}$, $J=8\text{Hz}$, 1H), 8.31 (s, 1H), 9.01 (s, 1H), 10.24 (s, 1H); ESIMS found $\text{C}_{12}\text{H}_{12}\text{BrN}_3\text{O}_2$ m/z 310.1 (M+H).

[0710] Preparation of intermediate N-(5-bromopyridin-3-yl)pivalamide (XXI) is depicted below in Scheme 4.

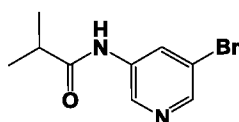


Scheme 4

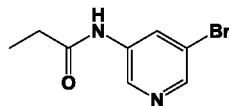
Step 1

[0711] To a solution of 3-amino-5-bromo pyridine (**XIX**) (1.0 g, 5.78 mmol) in dry pyridine (10 mL) was added pivaloyl chloride (**XX**) (769 mg, 6.38 mmol). The reaction mixture was stirred at room temperature for 3 h. The reaction was poured into an ice water/saturated aqueous NaHCO₃ mixture and stirred for 30 min. The precipitate was filtered, washed with cold water and dried at room temperature to yield N-(5-bromopyridin-3-yl)pivalamide (**XXI**) as an off-white solid (1.082 g, 4.22 mmol, 73.1% yield). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 1.23 (s, 9H), 8.37 (d, *J*=2Hz, 1H), 8.39 (t, *J*=2Hz, 1H), 8.80 (d, *J*=2Hz, 1H), 9.58 (brs, 1H); ESIMS found C₁₀H₁₃BrN₂O *m/z* 258.9 (Br⁸¹M+H).

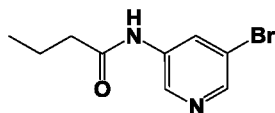
[0712] The following intermediates were prepared in accordance with the procedure described in the above Scheme 4.

**XXII**

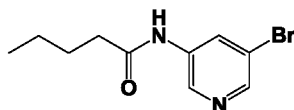
[0713] N-(5-Bromopyridin-3-yl)isobutyramide (**XXII**): Off-white solid, (71% yield). ¹H NMR (CDCl₃) δ ppm 8.55-8.35 (m, 3H), 7.32 (s, 1H), 2.59-2.48 (m, 1H), 1.28-1.27 (d, 6H); ESIMS found C₉H₁₁BrN₂O *m/z* 242.9 (Br⁷⁹M+H).

**XXIII**

[0714] N-(5-Bromopyridin-3-yl)propionamide (**XXIII**): Off white solid (92% yield). ¹H NMR (DMSO-*d*₆) δ ppm 1.09 (t, *J*=7.54 Hz, 3H), 2.36 (q, *J*=7.54 Hz, 2H), 8.36 (m, 2H), 8.65 (d, *J*=2.07 Hz, 1H), 10.26 (s, 1H); ESIMS found C₈H₉BrN₂O *m/z* 231.1 (Br⁸¹M+H).

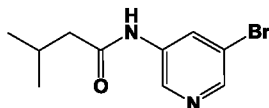
**XXIV**

[0715] N-(5-Bromopyridin-3-yl)butyramide (**XXIV**): Yellow solid (2.1 g, 8.64 mmol, 88.8% yield). ¹H NMR (CD₃OD, 400 MHz) δ ppm 1.02 (t, *J*=7.2Hz, 3H), 1.74 (sxt, *J*=7.2Hz, 2H), 2.40 (t, *J*=7.2Hz, 2H), 8.35 (d, *J*=2Hz, 1H), 8.46 (t, *J*=2Hz, 1H), 8.63 (d, *J*=2Hz, 1H); ESIMS found C₉H₁₁BrN₂O *m/z* 243.1 (Br⁷⁹M+H).



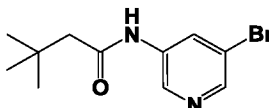
XXV

[0716] N-(5-Bromopyridin-3-yl)pentanamide (XXV): Yellow solid (2.0 g, 7.78 mmol, 85.3% yield). $^1\text{H NMR}$ (CD_3OD , 400 MHz) δ ppm 0.98 (t, $J=7.4\text{Hz}$, 3H), 1.43 (sxt, $J=7.4\text{Hz}$, 2H), 1.70 (quin, $J=7.4\text{Hz}$, 2H), 2.43 (t, $J=7.6\text{Hz}$, 2H), 8.35 (s, 1H), 8.45 (d, $J=2\text{Hz}$, 1H), 8.64 (d, $J=2\text{Hz}$, 1H); ESIMS found $\text{C}_{10}\text{H}_{13}\text{BrN}_2\text{O}$ m/z 256.9 ($\text{Br}^{79}\text{M}+\text{H}$).



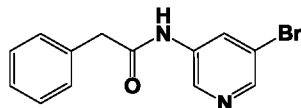
XXVI

[0717] N-(5-Bromopyridin-3-yl)-3-methylbutanamide (XXVI): Off white solid, (67% yield), $^1\text{H NMR}$ (CDCl_3 , 500 MHz) δ ppm 8.55-8.42 (m, 3H), 7.62 (s, 1H), 2.31-2.18 (m, 3H), 1.02-1.01 (d, $J=6\text{Hz}$, 6H); ESIMS found $\text{C}_{10}\text{H}_{13}\text{BrN}_2\text{O}$ m/z 258.9 ($\text{Br}^{81}\text{M}+\text{H}$).



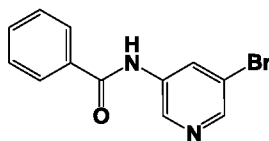
XXVII

[0718] N-(5-Bromopyridin-3-yl)-3,3-dimethylbutanamide (XXVII): Yellow solid (1.7 g, 6.27 mmol, 78.6% yield). $^1\text{H NMR}$ (CD_3OD , 400 MHz) δ ppm 1.10 (s, 9H), 2.29 (s, 2H), 8.36 (d, $J=1.6\text{Hz}$, 1H), 8.46 (d, $J=2.0\text{Hz}$, 1H), 8.64 (d, $J=2.0\text{Hz}$, 1H); ESIMS found $\text{C}_{11}\text{H}_{15}\text{BrN}_2\text{O}$ m/z 273.1 ($\text{Br}^{81}\text{M}+\text{H}$).



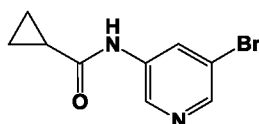
XXVIII

[0719] N-(5-Bromopyridin-3-yl)-2-phenylacetamide (XXVIII): White solid (2.5 g, 8.59 mmol, 77.9% yield). $^1\text{H NMR}$ (CDCl_3 , 400 MHz) δ ppm 3.76 (s, 2H), 7.26-7.45 (m, 5H), 7.57 (brs, 1H), 8.33 (s, 1H), 8.37 (s, 2H); ESIMS found $\text{C}_{13}\text{H}_{11}\text{BrN}_2\text{O}$ m/z 292.8 ($\text{Br}^{81}\text{M}+\text{H}$).



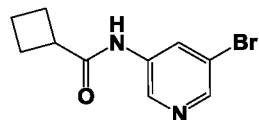
XXIX

[0720] N-(5-Bromopyridin-3-yl)benzamide (XXIX): White solid (2.7 g, 9.74 mmol, 60% yield). $^1\text{H NMR}$ (CDCl_3 , 400 MHz) δ ppm 7.40-7.52 (m, 2H), 7.52-7.62 (m, 1H), 7.86 (d, $J=7.2\text{Hz}$, 2H), 8.39 (d, $J=1.6\text{Hz}$, 1H), 8.46 (s, 1H), 8.55 (d, $J=1.6\text{Hz}$, 1H), 8.57 (d, $J=2.0\text{Hz}$, 1H); ESIMS found $\text{C}_{12}\text{H}_9\text{BrN}_2\text{O}$ m/z 278.8 ($\text{Br}^{81}\text{M}+\text{H}$).



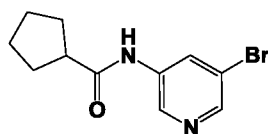
XXX

[0721] N-(5-Bromopyridin-3-yl)cyclopropanecarboxamide (**XXX**): Off-white solid, (83% yield), $^1\text{H NMR}$ (CDCl_3 , 500 MHz) δ ppm 8.46-8.39 (m, 3H), 7.54 (bs, 1H), 1.56-1.50 (m, 1H), 1.13-1.07 (m, 2H), 0.96-0.90 (m, 2H); ESIMS found for $\text{C}_9\text{H}_9\text{BrN}_2\text{O}$ m/z 240.9 ($\text{Br}^{79}\text{M}+\text{H}$).



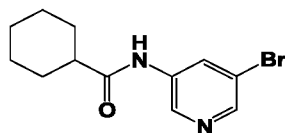
XXXI

[0722] N-(5-Bromopyridin-3-yl)cyclobutanecarboxamide (**XXXI**): Yellow solid (2.1 g, 6.27 mmol, 86.6% yield). $^1\text{H NMR}$ (CD_3OD , 400 MHz) δ ppm 1.80-1.99 (m, 1H), 1.99-2.15 (m, 1H), 2.16-2.30 (m, 2H), 2.30-2.45 (m, 2H), 3.25-3.35 (m, 1H), 8.34 (d, $J=2.0\text{Hz}$, 1H), 8.47 (s, 1H), 8.64 (d, $J=2.0\text{Hz}$, 1H); ESIMS found $\text{C}_{10}\text{H}_{11}\text{BrN}_2\text{O}$ m/z 257.1 ($\text{Br}^{81}\text{M}+\text{H}$).



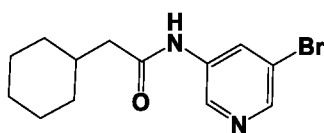
XXXII

[0723] N-(5-Bromopyridin-3-yl)cyclopentanecarboxamide (**XXXII**): Yellow solid (1.9 g, 7.06 mmol, 80.2% yield). $^1\text{H NMR}$ (CD_3OD , 400 MHz) δ ppm 1.57-1.74 (m, 2H), 1.74-1.91 (m, 4H), 1.91-2.07 (m, 2H), 2.77-2.92 (m, 1H), 8.34 (d, $J=1.6\text{Hz}$, 1H), 8.45 (s, 1H), 8.65 (d, $J=2.0\text{Hz}$, 1H); ESIMS found $\text{C}_{11}\text{H}_{13}\text{BrN}_2\text{O}$ m/z 271.1 ($\text{Br}^{81}\text{M}+\text{H}$).



XXXIII

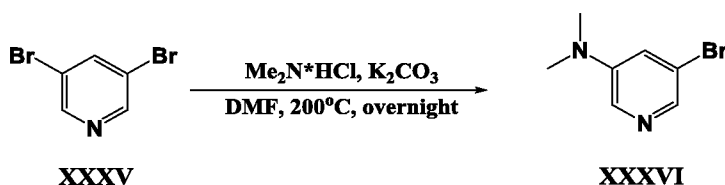
[0724] N-(5-bromopyridin-3-yl)cyclohexanecarboxamide (**XXXIII**): Yellow solid (2.0 g, 7.06 mmol, 84.3% yield). $^1\text{H NMR}$ (CD_3OD , 400 MHz) δ ppm 1.19-1.46 (m, 3H), 1.46-1.63 (m, 2H), 1.74 (d, $J=11.6\text{Hz}$, 1H), 1.88 (t, $J=14.0\text{Hz}$, 4H), 2.40 (tt, $J=11.6\text{Hz}$, $J=3.6\text{Hz}$, 1H), 8.34 (d, $J=2.0\text{Hz}$, 1H), 8.44 (t, $J=2.0\text{Hz}$, 1H), 8.64 (d, $J=2.0\text{Hz}$, 1H); ESIMS found $\text{C}_{12}\text{H}_{15}\text{BrN}_2\text{O}$ m/z 285.1 ($\text{Br}^{81}\text{M}+\text{H}$).



XXXIV

[0725] N-(5-bromopyridin-3-yl)-2-cyclohexylacetamide (XXXIV): Yellow solid (261 mg, 0.878 mmol, 84.4% yield). ESIMS found $C_{13}H_{17}BrN_2O$ m/z 297.1 ($Br^{81}M+H$).

[0726] Preparation of intermediate 5-bromo-N,N-dimethylpyridin-3-amine (XXXVI) is depicted below in Scheme 5.

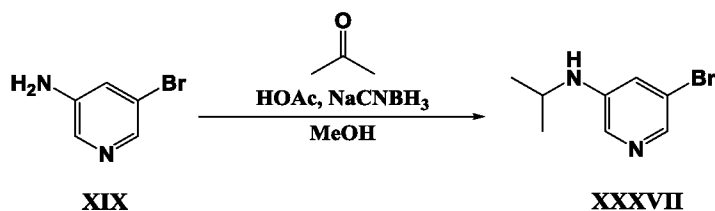


Scheme 5

Step 1

[0727] To a solution of 3,5-dibromopyridine (XXXV) (2.37 g, 10.0 mmol) in dry DMF (20.0 mL) was added K_2CO_3 (4.5 g, 33 mmol) and dimethylamino hydrochloride (1.79 g, 22 mmol). The mixture was heated overnight at 200°C in a sealed tube. The solution was cooled to room temperature and excess DMF was removed under vacuum. The residue was partitioned between EtOAc and water. The organic phase was separated. The aqueous phase was washed with EtOAc and the combined organic phases were dried over $MgSO_4$, and concentrated to afford 5-bromo-N,N-dimethylpyridin-3-amine (XXXVI) as an off-white solid (1.78g, 8.85 mmol, 88% yield). 1H NMR ($DMSO-d_6$, 500 MHz) δ ppm 2.94 (s, 6H), 7.25 (t, $J=2$ Hz, 1H), 7.91 (d, $J=2$ Hz, 1H), 8.07 (d, $J=2$ Hz, 1H); ESIMS found $C_7H_9BrN_2$ m/z 201.1 (M+H).

[0728] Preparation of intermediate 5-bromo-N-isopropylpyridin-3-amine (XXXVII) is depicted below in Scheme 6.

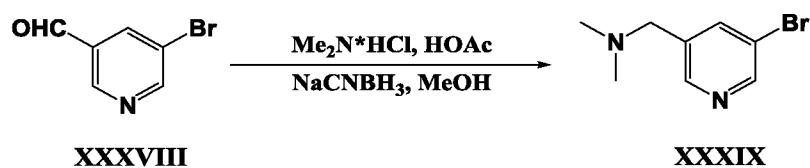


Scheme 6

Steps 1

[0729] To a solution of 5-bromopyridin-3-amine (**XIX**) (535 mg, 3.09 mmol) in MeOH (62 mL) was added acetone (296 μ L, 4.02 mL). The pH was adjusted to 4 using HOAc and stirred for 30 min. NaCNBH₃ (272 mg, 4.33 mmol) was added and stirred at room temperature overnight. The MeOH was removed under vacuum and the residue was partitioned between EtOAc and saturated aqueous NaHCO₃. The organic layer was dried over MgSO₄ and evaporated under vacuum. The crude product was purified on a silica gel column (100% hexane \rightarrow 90:10 hexane:EtOAc) to produce 5-bromo-N-isopropylpyridin-3-amine (**XXXVII**) as an oil which slowly solidified into an off-white solid (309 mg, 1.44 mmol, 47% yield). ¹H NMR (DMSO-d₆, 500 MHz) δ ppm 1.12 (d, $J=6.3$ Hz, 6H), 3.55-3.59 (m, 1H), 6.03 (d, $J=7.9$ Hz, 1H), 7.05-7.06 (m, 1H), 7.75 (d, $J=2$ Hz, 1H), 7.90 (d, $J=2$ Hz, 1H); ESIMS found C₈H₁₁BrN₂ m/z 215.1 (M+H).

[0730] Preparation of intermediate 1-(5-bromopyridin-3-yl)-N,N-dimethylmethanamine (**XXXIX**) is depicted below in Scheme 7.

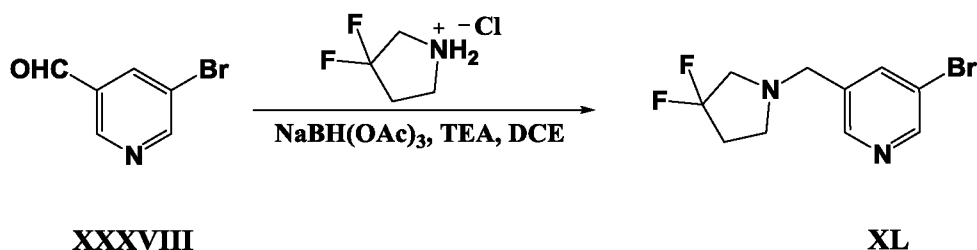


Scheme 7

Steps 1

[0731] Preparation of 1-(5-bromopyridin-3-yl)-N,N-dimethylmethanamine (**XXXIX**) was performed following the procedure listed in Scheme 6, Step 1. Brown oil (1.20 g, 5.59 mmol, 45% yield). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 2.15 (s, 6H), 3.43 (s, 2H), 7.94 (s, 1H), 8.47 (d, $J=1.1$ Hz, 1H), 8.59 (d, $J=2.2$ Hz, 1H); ESIMS found C₈H₁₁BrN₂ m/z 215 (M^{Br79}+H) and 217 (M^{Br81}+H).

[0732] Preparation of intermediate 3-bromo-5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridine (**XL**) is depicted below in Scheme 8.

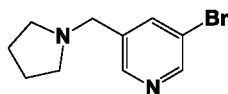


Scheme 8

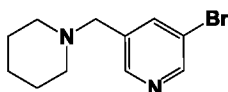
Steps 1

[0733] To a mixture of 5-bromopyridine-3-carbaldehyde (**XXXVIII**) (6.00 g, 32.26 mmol, 1.0 eq), 3,3-difluoropyrrolidine (5.56 g, 38.71 mmol, 1.20 eq) and TEA (5.39 mL, 38.71 mmol, 1.2 Eq) in DCE (200 mL) was stirred at room temperature for 30 min, then added sodium triacetoxyborohydride (10.25 g, 48.38 mmol, 1.50 eq) in one portion at room temperature under N₂. The mixture was stirred at room temperature for 6 hours. TLC showed the reaction was complete. The reaction was quenched with 1N NaOH (100 mL), extracted with DCE (100 mL X 2). The combined organic layers were washed with brine (100 mL), dried and concentrated. The residue was purified by silica gel chromatography (column height: 50 mm, diameter: 50 mm, 300-400 mesh silica gel, DCM/MeOH=30/1→20/1) to give 3-bromo-5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridine (**XL**): Yellow oil (8.00 g, 28.9 mmol, 89.5% yield). ¹H NMR (CDCl₃, 400 MHz) δ ppm 2.30 (spt, *J*=7.2 Hz, 2H), 2.75 (t, *J*=6.8 Hz, 2H), 2.91 (t, *J*=13.2 Hz, 2H), 7.85 (s, 1H), 8.45 (s, 1H), 8.59 (d, *J*=2 Hz, 1H); ESIMS found for C₁₀H₁₁BrF₂N₂ *m/z* 277.0 (M+H).

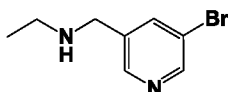
[0734] The following intermediates were prepared in accordance with the procedure described in the above Scheme 7 or Scheme 8.

**XLI**

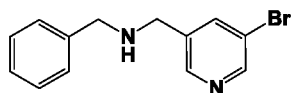
[0735] 3-Bromo-5-(pyrrolidin-1-ylmethyl)pyridine (**XLI**): Golden liquid (1.35 g, 97% yield). ¹H NMR (DMSO-*d*₆) 1.68-1.71 (m, 4H), 2.42-2.44 (m, 4H), 3.60 (s, 2H), 7.96 (s, 1H), 8.48 (d, *J*=2 Hz, 1H), 8.58 (d, *J*=3 Hz, 1H); ESIMS found for C₁₀H₁₃BrN₂ *m/z* 242.2 (M+H).

**XLII**

[0736] 3-Bromo-5-(piperidin-1-ylmethyl)pyridine (**XLII**): Brown liquid (13.1 g, 94% yield). ¹H NMR (DMSO-*d*₆) 1.36-1.39 (m, 2H), 1.46-1.51 (m, 4H), 2.31-2.32 (m, 4H), 3.46 (s, 2H), 7.94 (s, 1H), 8.47 (d, *J*=2 Hz, 1H), 8.58 (d, *J*=3 Hz, 1H); ESIMS found for C₁₁H₁₅BrN₂ *m/z* 257.0 (M+H).

**XLIII**

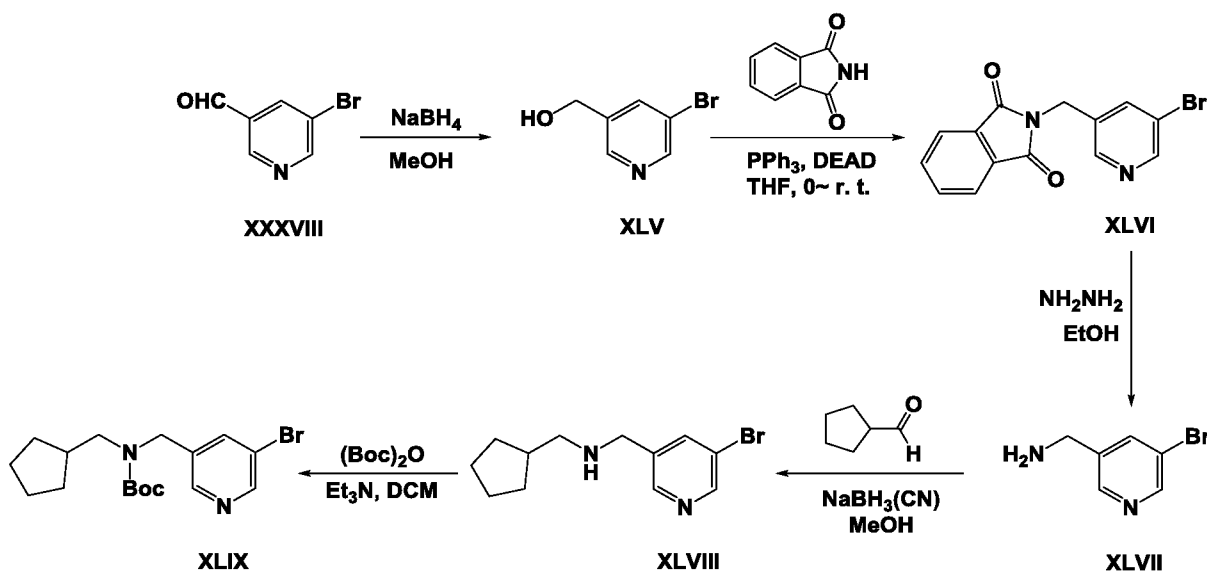
[0737] N-((5-Bromopyridin-3-yl)methyl)ethanamine (XLIII): Golden liquid (1.29 g, 6.00 mmol, 60% yield). $^1\text{H NMR}$ (CDCl_3 , 400 MHz) δ ppm 1.14 (t, $J=7.2\text{Hz}$, 3H), 2.67 (q, $J=7.2\text{Hz}$, 2H), 3.79 (s, 2H), 7.85 (t, $J=2\text{Hz}$, 1H), 8.46 (d, $J=1.6\text{Hz}$, 1H), 8.56 (d, $J=2.4\text{Hz}$, 1H); ESIMS found for $\text{C}_8\text{H}_{11}\text{BrN}_2$ m/z 215.1 (M+H).



XLIV

[0738] N-Benzyl-1-(5-bromopyridin-3-yl)methanamine (XLIV): Yellow oil (8.0 g, 28.9 mmol, 89.5% yield). $^1\text{H NMR}$ ($\text{DMSO}-d_6$, 400 MHz) δ ppm 3.71 (s, 2H), 3.74 (s, 2H), 7.18-7.28 (m, 1H), 7.28-7.40 (m, 4H), 8.04 (s, 1H), 8.52 (s, 1H), 8.58 (s, 1H); ESIMS found for $\text{C}_{13}\text{H}_{13}\text{BrN}_2$ m/z 277.1 (M+H).

[0739] Preparation of intermediate *tert*-butyl (5-bromopyridin-3-yl)methyl (cyclopentylmethyl)carbamate (XLIX) is depicted below in Scheme 9.



Scheme 9

Step 1

[0740] To a solution of 5-bromonicotinaldehyde (XXXVIII) (2.0 g, 10.8 mmol, 1 eq) in MeOH (20 mL) was added NaBH_4 (2.4 g, 64.9 mmol, 6 eq) and the reaction mixture was stirred at room temperature for 3 h. The mixture was concentrated in vacuo and the residue was diluted in water (15 mL), the aqueous phase was extracted with DCM (10 mL x 3). The combined organic layers were dried over MgSO_4 , filtered and concentrated in vacuo to afford (5-bromopyridin-3-yl)methanol (XLV) (1.8 g, 9.57 mmol, 90.0% yield) as a colorless oil. $^1\text{H NMR}$ (CDCl_3 , 500 MHz)

δ ppm 4.73 (s, 2H), 7.90 (s, 1H), 8.47 (s, 1H), 8.57 (s, 1H). ESIMS found for C_6H_6BrNO m/z 188.0 (M+H).

Step 2

[0741] To a stirred solution of (5-bromopyridin-3-yl)methanol (**XLV**) (1.60 g, 8.5 mmol, 1 eq), phthalimide (1.24 g, 8.5 mmol, 1 eq) and PPh_3 (3.33 g, 12.75 mmol, 1.5 eq) in anhydrous THF (15 mL) was added DEAD (2.21 g, 12.75 mmol, 1.5 eq) dropwise at 0°C under N_2 . Then the reaction mixture was stirred at room temperature for 6 h. The mixture was washed with saturated $NaHCO_3$ solution (15 mL), water (15 mL) and brine (15 mL) subsequently. The organic layers were dried over $MgSO_4$, concentrated under reduced pressure, the resultant residue was purified by flash chromatography on silica gel (PE:EtOAc = 4:1) to give 2-((5-bromopyridin-3-yl)methyl)isoindoline-1,3-dione (**XLVI**) (2.5 g, 7.88 mmol, 82.3% yield) as a white solid. ESIMS found for $C_{14}H_9BrN_2O_2$ m/z 317.1 (M+H).

Step 3

[0742] A solution of 2-((5-bromopyridin-3-yl)methyl)isoindoline-1,3-dione (**XLVI**) (1.9 g, 6.0 mmol, 1 eq) and hydrazine hydrate (2.0 g, 40 mmol, 6 eq) in EtOH (20 mL) was heated at 70°C for 3 h. The mixture was filtered through a Celite[®] pad and the filtrate was concentrated in vacuo, the crude product was dissolved in 1N HCl solution (15 mL) and concentrated to dryness, then it was washed with acetone (10 mL x 3), the precipitate was collected by filtration, dried in vacuo to give (5-bromopyridin-3-yl)methanamine (**XLVII**) (1.3 g, 6.95 mmol, 97.7% yield) as a white solid. 1H NMR (D_2O , 500 MHz) δ ppm 4.34 (s, 2H), 8.56 (s, 1H), 8.75 (d, $J=1.2$ Hz, 1H), 8.91 (d, $J=1.6$ Hz, 1H). ESIMS found for $C_6H_7BrN_2$ m/z 187.0 (M+H).

Step 4

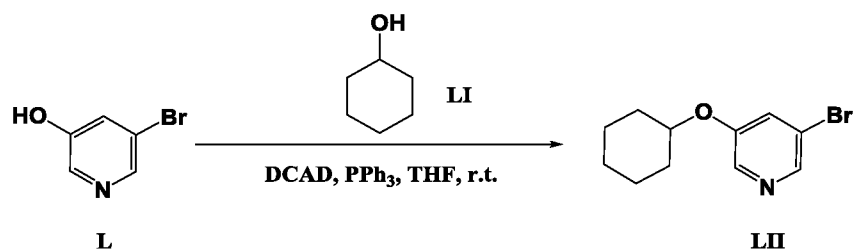
[0743] A solution of (5-bromopyridin-3-yl)methanamine (**XLVII**) (1.30 g, 5.8 mmol, 1.0 eq), cyclopentanecarbaldehyde (0.57 g, 5.8 mmol, 1.0 eq) and TEA (0.60 g, 5.8 mmol, 1.0 eq) in MeOH (15 mL) was stirred at room temperature for 2 h. Then $NaBH_3CN$ (1.98 g, 34.6 mmol, 6.0 eq) was added and the mixture was stirred at the same temperature for another 3 h. The solvent was removed under reduced pressure and the residue was diluted in water (20 mL) and extracted with DCM (10 mL x 3), combined organic layers were dried over $MgSO_4$ and concentrated in vacuo to give 1-(5-bromopyridin-3-yl)-N-(cyclopentylmethyl)methanamine (**XLVIII**) (1.23 g, 4.57 mmol, 79.3% yield) as a yellow oil. 1H NMR ($CDCl_3$, 400 MHz) δ ppm 1.07-1.23 (m, 2H), 1.47-

1.67 (m, 4H), 1.70-1.84 (m, 2H), 2.02 (spt, $J=7.6$ Hz, 1H), 2.53 (d, $J=7.2$ Hz, 2H), 3.80 (s, 2H), 7.86 (s, 1H), 8.47 (s, 1H), 8.56 (d, $J=2.0$ Hz, 1H); ESIMS found for $C_{12}H_{17}BrN_2$ m/z 269.1 (M+H).

Step 5

[0744] To a solution of 1-(5-bromopyridin-3-yl)-N-(cyclopentylmethyl) methanamine (**XLVIII**) (1.00 g, 3.7 mmol, 1 eq) and TEA (0.93 g, 9.2 mmol, 2.5 eq) in DCM (20 mL) was added portionwise $(Boc)_2O$ (0.85 g, 4.0 mmol, 1.1 eq) at $0^\circ C$, the reaction mixture was stirred at room temperature for 1 h. The mixture was washed with water (10 mL), brine (10 mL), the organic layer was separated, dried over $MgSO_4$ and concentrated in vacuo to give *tert*-butyl (5-bromopyridin-3-yl)methyl (cyclopentylmethyl)carbamate (**XLIX**) (1.25 g, 3.38 mmol, 91.9% yield) as a white solid. ESIMS found for $C_{17}H_{25}BrN_2O_2$ m/z 369.1 (M+H).

[0745] Preparation of intermediate 3-bromo-5-(cyclohexyloxy)pyridine (**LII**) is depicted below in Scheme 10.

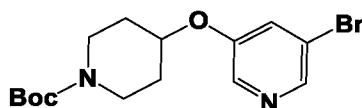


Scheme 10

Step 1

[0746] To a solution of 5-bromopyridin-3-ol (**L**) (523 mg, 3.01 mmol) in THF (30 mL) cooled to $0^\circ C$ were added triphenylphosphine (867 mg, 3.31 mmol) and cyclohexanol (**LI**) (331 mg, 3.31 mmol) followed by (*E*)-bis(4-chlorobenzyl) diazene-1,2-dicarboxylate (1.21 g, 3.31 mmol), added portionwise. The reaction mixture was then stirred at $25^\circ C$ overnight. The reaction was worked-up with a EtOAc- $NaHCO_3$ extraction and the solid filtered off. The solvent was removed and the residue was purified by Isco (20% EtOAc-Hexanes) to give 3-bromo-5-(cyclohexyloxy)pyridine (**LII**) (209 mg, 0.82 mmol, 27.2% yield) as a yellow oil. 1H NMR (DMSO- d_6 , 500 MHz) δ ppm 1.21 - 1.31 (m, 1 H) 1.34 - 1.48 (m, 4 H) 1.49 - 1.57 (m, 1 H) 1.70 (br dd, $J=9.74, 4.25$ Hz, 2 H) 1.88 - 1.96 (m, 2 H) 2.50 (dt, $J=3.70, 1.72$ Hz, 5 H) 4.46 - 4.54 (m, 1 H) 7.72 (t, $J=2.20$ Hz, 1 H) 8.24 (d, $J=1.92$ Hz, 1 H) 8.27 (d, $J=2.47$ Hz, 1 H).

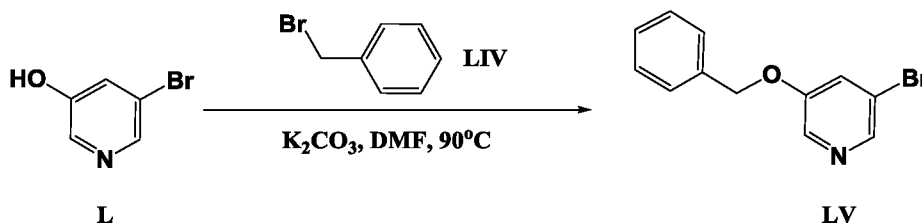
[0747] The following intermediate was prepared in accordance with the procedure described in the above Scheme 10.



LIII

[0748] *tert*-Butyl 4-((5-bromopyridin-3-yl)oxy)piperidine-1-carboxylate (LIII): Yellow oil (244 mg, 0.683 mmol, 23.2% yield). ESIMS found for $C_{15}H_{21}BrN_2O_3$ m/z 358.3 (M+H).

[0749] Preparation of intermediate 3-(benzyloxy)-5-bromopyridine (LV) is depicted below in Scheme 11.

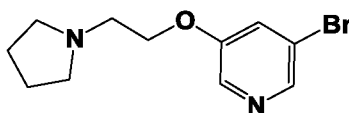


Scheme 11

Step 1

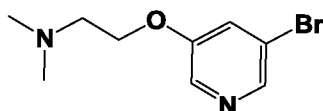
[0750] To a solution of 5-bromopyridin-3-ol (L) (174 mg, 1.0 mmol) in DMF (3 mL) was added potassium carbonate (415 mg, 3.0 mmol). The slurry was heated at 90°C for 1 hour and then cooled to 25°C. The (bromomethyl)benzene (LIV) (171 mg, 1.0 mmol) was added and the mixture was stirred at 25°C overnight. The reaction was worked-up using a saturated sodium bicarbonate and ethyl acetate extraction. The product was purified by ISCO column eluted with 40-100% EtOAc-Hexanes. The 3-(benzyloxy)-5-bromopyridine (LV) (105 mg, 0.398 mmol, 39.8 % yield) was obtained as yellow oil. MS: 266.1. ESIMS found for $C_{12}H_{10}BrNO$ m/z 266.1 (M+H).

[0751] The following intermediates were prepared in accordance with the procedure described in the above Scheme 11.



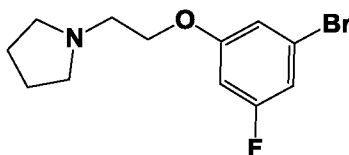
LVI

[0752] 3-Bromo-5-(2-(pyrrolidin-1-yl)ethoxy)pyridine (LVI): Yellow oil (97 mg, 0.358 mmol, 15.56% yield). ESIMS found for $C_{11}H_{15}BrN_2O$ m/z 272.2 (M+H).



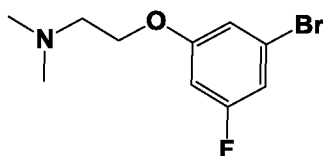
LVII

[0753] 2-((5-bromopyridin-3-yl)oxy)-N,N-dimethylethan-1-amine (LVII): Yellow oil (97 mg, 0.396 mmol, 28.9% yield). ESIMS found for $C_9H_{13}BrN_2O$ m/z 245.1 (M+H).



LVIII

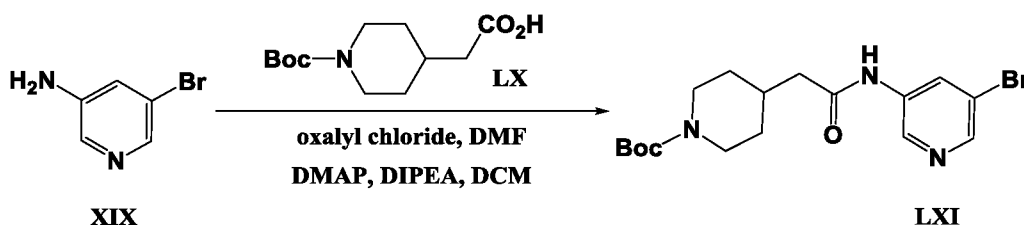
[0754] 1-(2-(3-bromo-5-fluorophenoxy)ethyl)pyrrolidine (LVIII): Yellow oil (370 mg, 1.284 mmol, 85.8% yield). ESIMS found for $C_{12}H_{15}BrFNO$ m/z 289.0 (M+H).



LIX

[0755] 2-(3-bromo-5-fluorophenoxy)-N,N-dimethylethan-1-amine (LIX): Yellow oil (364 mg, 1.389 mmol, 50.2% yield). ESIMS found for $C_{10}H_{13}BrFNO$ m/z 263.9 (M+H).

[0756] Preparation of intermediate *tert*-butyl 4-(2-((5-bromopyridin-3-yl)amino)-2-oxoethyl)piperidine-1-carboxylate (LXI) is depicted below in Scheme 12.



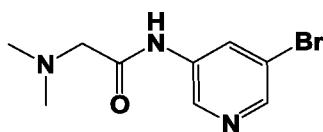
Scheme 12

Step 1

[0757] To a solution of 2-(1-(*tert*-butoxycarbonyl)piperidin-4-yl)acetic acid (LX) (3.4 g, 13.97 mmol) in DCM (10 mL) was added DMF (1 mL). The solution was cooled in ice-water to 0°C. Oxalyl chloride (1.835 mL, 20.96 mmol) was then added dropwise. The mixture was stirred for one hour at 25°C. The organic volatile was then removed under vacuum. The residue was dissolved in DCM (10 mL). DMAP (0.171 g, 1.397 mmol) and 5-bromopyridin-3-amine (XIX) (2.418 g, 13.97 mmol) were added to the solution and cooled to 0°C. DIEA (4.88 mL, 27.9 mmol) was then added dropwise and the mixture was stirred for 2 hours at 25°C. The reaction was worked-up with DCM and saturated $NaHCO_3$. The product was purified by ISCO eluted with 0-100% EtOAc-Hexanes. The *tert*-butyl 4-(2-((5-bromopyridin-3-yl)amino)-2-oxoethyl)piperidine-1-

carboxylate (**LXI**) (2.82 g, 7.08 mmol, 50.7 % yield) was obtained as yellow oil. ESIMS found for $C_{17}H_{24}BrN_3O_3$ m/z 343.1 (M-56).

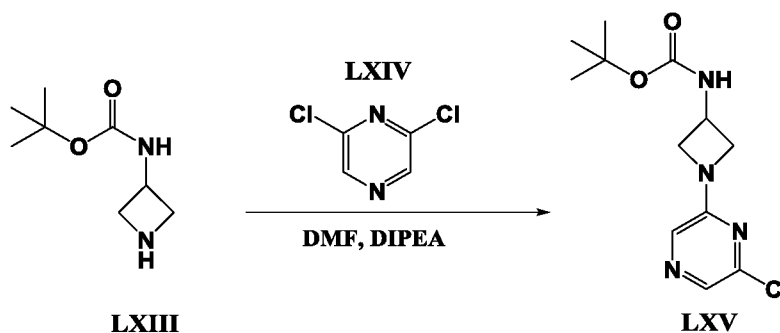
[0758] The following intermediate was prepared in accordance with the procedure described in the above Scheme 12.



LXII

[0759] N-(5-Bromopyridin-3-yl)-2-(dimethylamino)acetamide (**LXII**): Yellow oil (528 mg, 2.05 mmol, 19.0% yield). ESIMS found for $C_9H_{12}BrN_3O$ m/z 259.3 (M+H).

[0760] Preparation of *tert*-butyl (1-(6-chloropyrazin-2-yl)azetidin-3-yl)carbamate (**LXV**) is depicted below in Scheme 13.

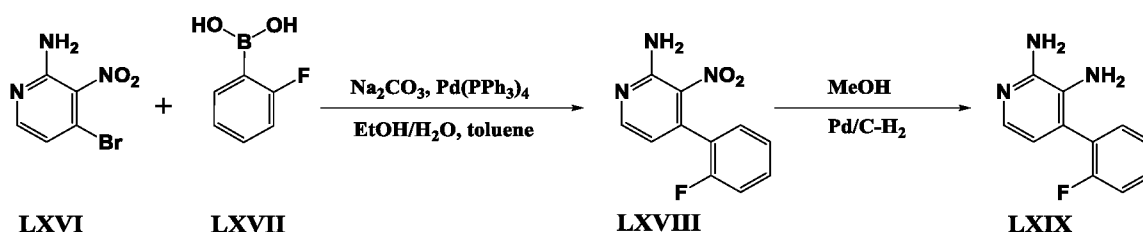


Scheme 13

Step 1

[0761] To a solution of *tert*-butyl azetidin-3-ylcarbamate hydrochloride (**LXIII**) (2 g, 9.58 mmol) in dry DMF (19.2 mL) was added DIPEA (8.37 ml, 47.9 mmol). To this mixture was added 2,6-dichloropyrazine (**LXIV**) (1.428 g, 9.58 mmol) and the reaction was stirred at 95°C for 3 hours. The reaction was quenched with water (20 mL) and extracted with EtOAc. The organic layer was dried over anhydrous Na_2SO_4 , filtered and concentrated. The residue was purified by silica gel column chromatography (40g) (100% hexanes→hexanes:EtOAc 1:1) to yield *tert*-butyl (1-(6-chloropyrazin-2-yl)azetidin-3-yl)carbamate (**LXV**) (2.2882 g, 8.04 mmol, 84 % yield) as a white solid. ESIMS found for $C_{12}H_{17}ClN_4O_2$ m/z 285.1 (M+H).

[0762] Preparation of intermediate 4-(2-fluorophenyl)-3-nitropyridin-2-amine (**LXIX**) is depicted below in Scheme 14.



Scheme 14

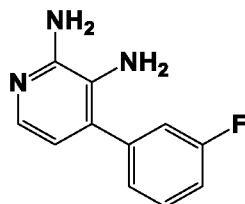
Step 1

[0763] A solution of 4-bromo-3-nitropyridin-2-amine (**LXVI**) (5.00 g, 22.9 mmol, 1.00 eq), (2-fluorophenyl)boronic acid (**LXVII**) (3.82 g, 27.5 mmol, 1.20 eq), Pd(PPh₃)₄ (1.32 g, 1.14 mmol, 0.05 eq), and Na₂CO₃ (4.85 g, 45.8 mmol, 2 eq) in a mixture of toluene (25 mL), H₂O (9 mL) and EtOH (6 mL) was stirred at 75°C for 15 h under nitrogen atmosphere. The reaction mixture was washed with brine (50 mL) and dried over anhydrous Na₂SO₄, filtered and concentrated in vacuo. The resultant residue was purified by chromatography on silica gel (PE:EtOAc = 3:1) to give 4-(2-fluorophenyl)-3-nitropyridin-2-amine (**LXVIII**) (4.0 g, 17.15 mmol, 74.9%) as a yellow solid. ¹H NMR (CDCl₃, 400 MHz) δ ppm 6.29 (brs, 2H), 6.68 (d, *J*=4.8Hz, 1H), 7.14 (t, *J*=5.2Hz, 1H), 7.23-7.50 (m, 3H), 8.32 (d, *J*=4.8Hz, 1H); ESIMS found C₁₁H₈FN₃O₂ m/z 234.2 (M+H).

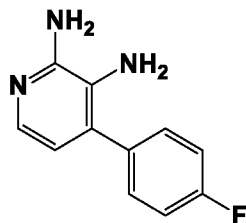
Step 2

[0764] A mixture of 4-(2-fluorophenyl)-3-nitropyridin-2-amine (**LXVIII**) (2.8 g, 12.0 mmol, 1 eq) and Pd/C (0.2 g) in MeOH (200 mL) was stirred under 50 psi of H₂ at room temperature overnight. The reaction was monitored by TLC. The mixture was filtered and the filtrate was concentrated in vacuo to produce 4-(2-fluorophenyl)pyridine-2,3-diamine (**LXIX**) (1.55g, 7.63 mmol, 63.6% yield) as a black solid. ¹H NMR (CDCl₃, 400 MHz) δ ppm 3.38 (brs, 2H), 4.40 (brs, 2H), 6.64 (d, *J*=4.8Hz, 1H), 7.11-7.53 (m, 4H), 7.72 (d, *J*=4.8Hz, 1H); ESIMS found C₁₁H₁₀FN₃ m/z 204.2 (M+H).

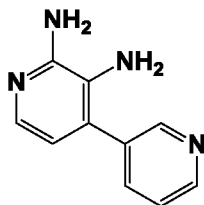
[0765] The following intermediates were prepared in accordance with the procedure described in the above Scheme 14.

**LXX**

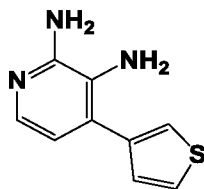
[0766] 4-(3-Fluorophenyl)pyridine-2,3-diamine (**LXX**): Grey solid, (1.55 g, 7.63 mmol, 86.0% yield). ¹H NMR (CDCl₃, 400 MHz) δ ppm 3.50 (brs, 2H), 4.36 (brs, 2H), 6.63 (d, *J*=3.6Hz, 1H), 7.3-7.37 (m, 3H), 7.47 (d, *J*=6Hz, 1H), 7.72 (d, *J*=3.6Hz, 1H); ESIMS found C₁₁H₁₀FN₃ *m/z* 204.2 (M+H).

**LXXI**

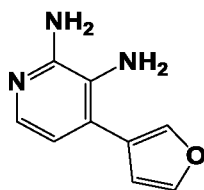
[0767] 4-(4-Fluorophenyl)pyridine-2,3-diamine (**LXXI**): Grey solid, (1.55 g, 7.63 mmol, 60.0% yield). ¹H NMR (CDCl₃, 400 MHz) δ ppm 3.46 (brs, 2H), 4.36 (brs, 2H), 6.62 (s, 1H), 7.19 (s, 2H), 7.43 (s, 2H), 7.70 (d, *J*=3.2Hz, 1H); ESIMS found C₁₁H₁₀FN₃ *m/z* 204.1 (M+H).

**LXXII**

[0768] 4-(4-Fluorophenyl)pyridine-2,3-diamine (**LXXII**): Grey solid, (1.55 g, 7.63 mmol, 60.0% yield). ¹H NMR (CDCl₃, 400 MHz) δ ppm 3.46 (brs, 2H), 4.36 (brs, 2H), 6.62 (s, 1H), 7.19 (s, 2H), 7.43 (s, 2H), 7.70 (d, *J*=3.2Hz, 1H); ESIMS found C₁₁H₁₀FN₃ *m/z* 204.1 (M+H).

**LXXIII**

[0769] 4-(Thiophen-3-yl)pyridine-2,3-diamine (**LXXIII**): Yellow solid, (1.9 g, 9.94 mmol, 84.9% yield). ¹H NMR (CDCl₃, 400 MHz) δ ppm 3.80 (brs, 2H), 4.34 (brs, 2H), 6.77 (s, 1H), 7.18 (s, 1H), 7.27 (s, 2H), 7.44 (s, 1H), 7.68 (s, 1H); ESIMS found C₉H₉N₃S *m/z* 192.2 (M+H).



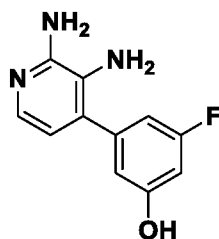
LXXIV

[0770] 4-(Furan-3-yl)pyridine-2,3-diamine (LXXIV): Black solid, (1.9 g, 10.84 mmol, 89.0% yield). ^1H NMR (CDCl_3 , 400 MHz) δ ppm 3.64 (brs, 2H), 4.32 (brs, 2H), 6.65 (s, 1H), 6.69 (d, $J=4.8\text{Hz}$, 1H), 7.58 (s, 1H), 7.67 (d, $J=4.8\text{Hz}$, 1H), 7.71 (s, 1H); ESIMS found $\text{C}_9\text{H}_9\text{N}_3\text{O}$ m/z 176.3 (M+H).



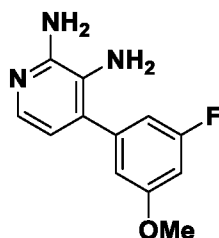
LXXV

[0771] 4-(Thiophen-2-yl)pyridine-2,3-diamine (LXXV): Yellow solid, (0.90 g, 4.71 mmol, 96.5% yield). ^1H NMR (CDCl_3 , 400 MHz) δ ppm 3.63 (brs, 2H), 4.35 (brs, 2H), 6.71 (s, 1H), 7.27 (s, 1H), 7.45 (s, 1H), 7.49 (s, 1H), 7.69 (s, 1H); ESIMS found $\text{C}_9\text{H}_9\text{N}_3\text{S}$ m/z 192.1 (M+H).



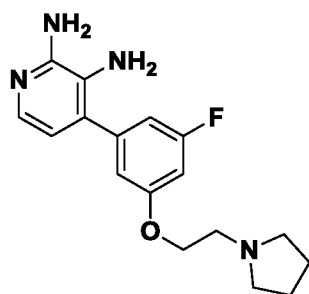
LXXVI

[0772] 3-(2,3-Diaminopyridin-4-yl)-5-fluorophenol (LXXVI): White solid (303 mg, 1.38 mmol, 84.4% yield). ESIMS found for $\text{C}_{11}\text{H}_{10}\text{FN}_3\text{O}$ m/z 220.1 (M+H).



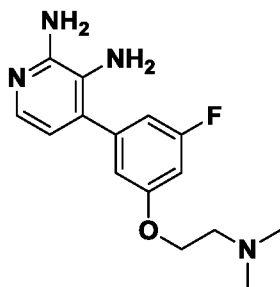
LXXVII

[0773] 4-(3-Fluoro-5-methoxyphenyl)pyridine-2,3-diamine (LXXVII): White solid (404 mg, 1.73 mmol, 75.1% yield). ESIMS found for $\text{C}_{12}\text{H}_{12}\text{FN}_3\text{O}$ m/z 234.1 (M+H).



LXXVIII

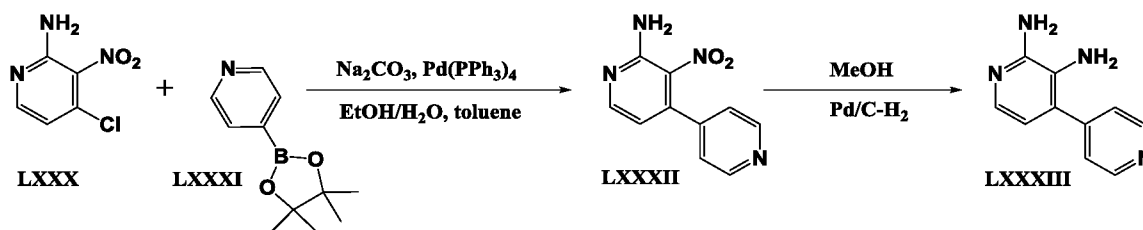
[0774] 4-(3-Fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)pyridine-2,3-diamine (LXXVIII): Black oil (368 mg, 1.163 mmol, 95.0% yield). ESIMS found for $C_{17}H_{21}FN_4O$ m/z 317.1 (M+H).



LXXIX

[0775] 4-(3-(2-(Dimethylamino)ethoxy)-5-fluorophenyl)pyridine-2,3-diamine (LXXIX): Black oil (352 mg, 1.212 mmol, 83.6% yield). ESIMS found for $C_{15}H_{19}FN_4O$ m/z 291.1 (M+H).

[0776] Preparation of intermediate 3-nitro-[4,4'-bipyridin]-2-amine (LXXXIII) is depicted below in Scheme 15.



Scheme 15

Step 1

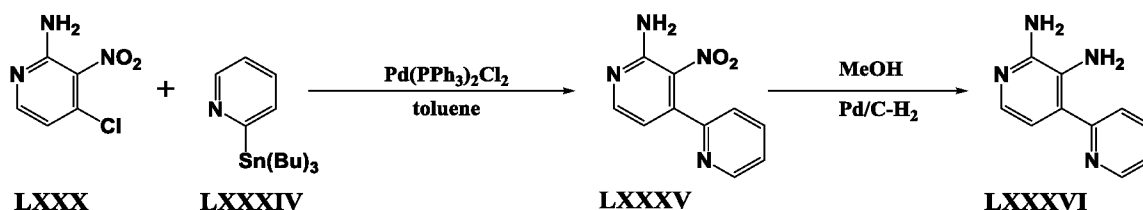
[0777] To a solution of 4-chloro-3-nitropyridin-2-amine (LXXX) (5.00 g, 28.9 mmol, 1.00 eq) and 4-(4,4,5,5-tetramethyl-1,3,2-dioxaborolan-2-yl)pyridine (LXXXI) (7.1 g, 34.6 mmol, 1.20 eq) in a mixture of toluene (30 mL), H_2O (18 mL) and EtOH (6 mL) was added $Pd(PPh_3)_4$ (1.0 g, 0.87 mmol, 0.03 eq) and Na_2CO_3 (6.1 g, 57.6 mmol, 2 eq). The mixture was stirred at $75^\circ C$ for

15 h under a nitrogen atmosphere. The reaction mixture was then poured into brine (100 mL) and extracted with EtOAc (30 mL x 3), the combined organic layers were dried over anhydrous Na₂SO₄, filtered and concentrated in vacuo. The resultant residue was purified by chromatography on silica gel (PE:EtOAc = 5:1→1:1) to afford 3-nitro-[4,4'-bipyridin]-2-amine (**LXXXII**) (1.80 g, 8.33 mmol, 28.8% yield) as a yellow solid. ESIMS found C₁₀H₈N₄O₂ m/z 217.1 (M+H).

Step 2

[0778] To a solution of 3-nitro-[4,4'-bipyridin]-2-amine (**LXXXII**) (1.80 g, 8.33 mol, 1 eq) in MeOH (50 mL) was added Pd/C (0.5 g) under a nitrogen atmosphere. The mixture was stirred under 50 psi of H₂ for 6 h at room temperature. Then the mixture was filtered through a Celite pad and the filtrate was concentrated in vacuo to afford [4,4'-bipyridine]-2,3-diamine (**LXXXIII**) (1.4 g, 7.52 mmol, 90.4% yield) as a black solid. ESIMS found C₁₀H₁₀N₄ m/z 186.0 (M+H).

[0779] Preparation of intermediate 3-nitro-[4,4'-bipyridin]-2-amine (**LXXXVI**) is depicted below in Scheme 16.



Scheme 16

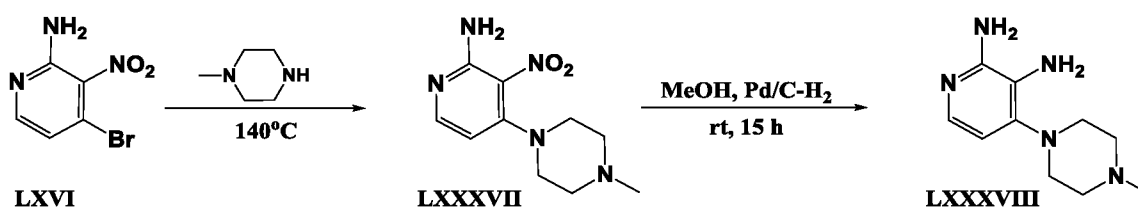
Step 1

[0780] A solution of 4-chloro-3-nitropyridin-2-amine (**LXXX**) (5.00 g, 28.9 mmol, 1.00 eq), 2-(tributylstannyl)pyridine (**LXXXIV**) (15.9 g, 43.4 mmol, 1.50 eq), and Pd(PPh₃)₂Cl₂ (1.05 g, 1.44 mmol, 0.05 eq) in a mixture of toluene (25 mL) and H₂O (9 mL) was stirred at 75°C for 16 h under a nitrogen atmosphere. The reaction mixture was then poured into brine (80 mL) and extracted with EtOAc (30 mL x 3). The combined organic layers were dried over anhydrous Na₂SO₄, filtered and concentrated in vacuo. The resultant residue was purified by chromatography on silica gel (PE:EtOAc = 5:1→1:1) to afford 3'-nitro-[2,4'-bipyridin]-2'-amine (**LXXXV**) (1.6 g, 7.40 mmol, 25.6% yield) as a yellow solid. ESIMS found C₁₀H₈N₄O₂ m/z 217.1 (M+H).

Step 2

[0781] To a solution of compound 3'-nitro-[2,4'-bipyridin]-2'-amine (**LXXXV**) (1.6 g, 7.4 mmol, 1.0 eq) in MeOH (50 mL) was added Pd/C (0.5 g) under a nitrogen atmosphere. The mixture was stirred under 50 psi of H₂ for 6 h at room temperature. The mixture was then filtered and concentrated in vacuo to afford [2,4'-bipyridine]-2',3'-diamine (**LXXXVI**) (1.1 g, 5.91 mmol, 79.8%) as a black solid. ESIMS found C₁₀H₁₀N₄ m/z 186.1 (M+H).

[0782] Preparation of intermediate 4-(4-methylpiperazin-1-yl)pyridine-2,3-diamine (**LXXXVIII**) is depicted below in Scheme 17.

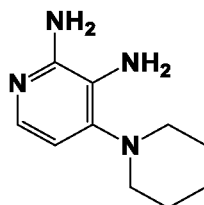
Step 1

[0783] A solution of 4-bromo-3-nitropyridin-2-amine (**LXVI**) (2.50 g, 11.47 mmol) in 1-methylpiperazine (4 mL, 34.5 mmol) was heated at 140°C overnight. The reaction was poured into an EtOAc/H₂O mixture; the organic layer was separated, dried over MgSO₄ and concentrated under vacuum. The crude product was purified on a silica gel column (100% CHCl₃ → 3:97 MeOH(7N NH₃):CHCl₃) to give 4-(4-methylpiperazin-1-yl)-3-nitropyridin-2-amine (**LXXXVII**) as a yellow solid (1.80 g, 7.59 mmol, 66.1% yield). ¹H NMR (CDCl₃, 400 MHz) δ ppm 2.36 (s, 3H), 2.54 (t, J=4.8Hz, 4H), 3.25 (t, J=5Hz, 4H), 6.18 (s, 2H), 6.22 (d, J=6Hz, 1H), 7.85 (d, J=6Hz, 1H); ESIMS found C₁₀H₁₅N₅O₂ m/z 238.0 (M+H).

Step 2

[0784] To a solution of 4-(4-methylpiperazin-1-yl)-3-nitropyridin-2-amine (**LXXXVII**) (1.80 g, 7.59 mmol) in MeOH (52 mL) was added 10% Pd/C (0.5 g). The solution was purged with hydrogen and stirred at room temperature under hydrogen for 4 h. The suspension was filtered through Celite[®] and the concentrated under vacuum to produce 4-(4-methylpiperazin-1-yl)pyridine-2,3-diamine (**LXXXVIII**) as black solid (1.4 g, 6.75 mmol, 89.0% yield). ¹H NMR (CDCl₃, 400 MHz): δ ppm 2.38 (s, 3H), 2.60 (brs, 4H), 2.99 (s, 4H), 3.49 (brs, 2H), 4.12 (brs, 2H), 6.52 (d, J=5.6Hz, 1H), 7.64 (d, J=5.6Hz, 1H); ESIMS found C₁₀H₁₇N₅ m/z 208.1 (M+H).

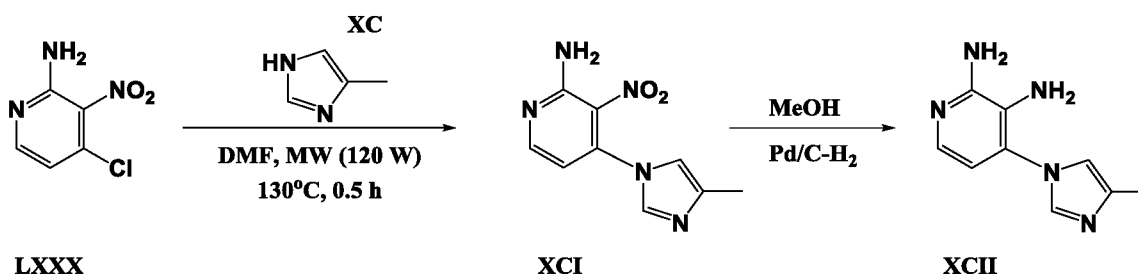
[0785] The following intermediate was prepared in accordance with the procedure described in the above Scheme 17.



LXXXIX

[0786] 4-(Piperidin-1-yl)pyridine-2,3-diamine (**LXXXIX**): Black solid, (2.40 g, 12.48 mmol, 92.5% yield). ¹H NMR (CDCl₃, 400 MHz): δ ppm 1.61 (brs, 2H), 1.73 (s, 4H), 2.88 (s, 4H), 3.48 (brs, 2H), 4.13 (brs, 2H), 6.50 (d, *J*=5.2Hz, 1H), 7.63 (d, *J*=4.8Hz, 1H); ESIMS found C₁₀H₁₆N₄ *m/z* 193.1 (M+H).

[0787] Preparation of intermediate 4-(2-fluorophenyl)-3-nitropyridin-2-amine (**XCII**) is depicted below in Scheme 18.



Scheme 18

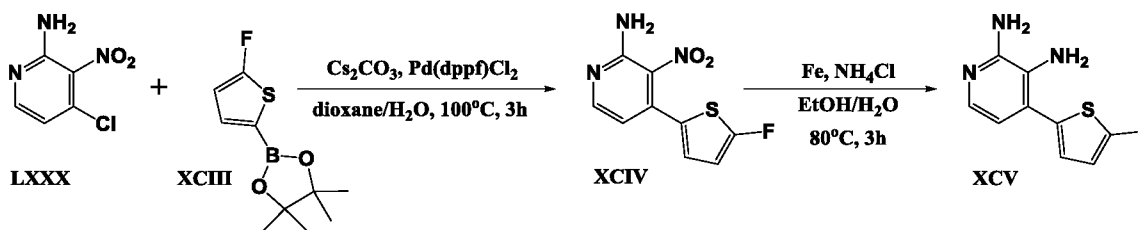
Step 1

[0788] 4-Chloro-3-nitro-pyridin-2-amine (**LXXX**) (3.00 g, 17.3 mmol, 1.0 eq) and 4-methyl-1H-imidazole (**XC**) (2.84 g, 34.6 mmol, 2.0 eq) were taken up into a microwave tube in DMF (20 mL). The sealed tube was heated at 130°C for 30 min under microwave. TLC showed the starting material was consumed, LC/MS showed the desired product was found. 10% NH₄Cl (60 mL) were added. The aqueous layer was extracted with DCM (2 x 100 mL). The combined organic layers were washed with brine (100 mL), dried over Na₂SO₄, concentrated in vacuum. The residue was purified by chromatography on silica gel (PE:THF = 1:1) to give the product 4-(4-methylimidazol-1-yl)-3-nitro-pyridin-2-amine (**XCI**) (1.20 g, 5.47 mmol, 31.6% yield) as a yellow solid. ¹H NMR (CDCl₃, 400 MHz) δ ppm 2.30 (d, *J*=0.75Hz, 3H), 5.96-6.26 (m, 2H), 6.68 (d, *J*=5.02Hz, 1H), 6.72-6.75 (m, 1H), 7.57 (d, *J*=1.0Hz, 1H), 8.32 (d, *J*=5.14Hz, 1H); ESIMS found C₉H₉N₅O₂ *m/z* 219.1 (M+H).

Step 2

[0789] To a solution of 4-(4-methylimidazol-1-yl)-3-nitro-pyridin-2-amine (**XCI**) (900.0 mg, 4.11 mmol, 1.0 eq) in MeOH (50 mL) was added Pd/C (100.0 mg, 4.11 mmol, 1.0 eq) at room temperature. The mixture was stirred at 20 °C under H₂ for 2 hr. TLC (DCM:MeOH = 20:1) showed that starting the material was consumed completely. The mixture was filtered and concentrated to afford 4-(4-methyl-1H-imidazol-1-yl)pyridine-2,3-diamine (**XCII**) (750.0 mg, 3.96 mmol, 96.4% yield) as a brown solid. ¹H NMR (CDCl₃, 400 MHz) δ ppm 2.17 (s, 3H), 4.58 (brs, 2H), 5.85 (brs, 2H), 6.36 (d, *J*=5.4Hz, 1H), 7.06 (s, 1H), 7.34 (s, 1H), 7.34 (s, 1H), 7.69 (d, *J*=0.88Hz, 1H); ESIMS found C₉H₁₁N₅ m/z 190.1 (M+H).

[0790] Preparation of intermediate 4-(5-fluorothiophen-2-yl)pyridine-2,3-diamine (**XCIV**) is depicted below in Scheme 19.

Step 1

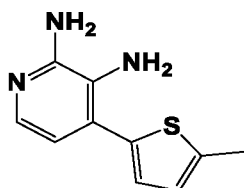
[0791] A solution of 2-chloro-3-nitro-pyridin-4-amine (**LXXX**) (1.5 g, 8.64 mmol), 2-(5-fluorothiophen-2-yl)-4,4,5,5-tetramethyl-1,3,2-dioxaborolane (**XCIII**) (2.96 g, 12.96 mmol), Pd(dppf)Cl₂ (632 mg, 0.86 mmol) and Cs₂CO₃ (5.63 g, 17.29 mmol) in dioxane (30 mL) and H₂O (5 mL) was de-gassed and then heated to 100°C under N₂ for 3 h. TLC (PE:EtOAc = 1:1) showed the starting material was consumed completely. The mixture was concentrated in vacuum to give a residue, which was purified by column chromatography to afford 4-(5-fluorothiophen-2-yl)-3-nitropyridin-2-amine (**XCIV**) (800 mg, 3.34 mmol, 38.7% yield). ESIMS found C₉H₆FN₃O₂S m/z 240.1 (M+H).

Step 2

[0792] A solution of 4-(5-fluorothiophen-2-yl)-3-nitropyridin-2-amine (**XCIV**) (700 mg, 2.93 mmol), Fe (817 mg, 14.63 mmol) and NH₄Cl (939 mg, 17.56 mmol) in EtOH (18 mL) and H₂O (6 mL) was heated to 80°C for 2 h. TLC (PE:EtOAc = 1:1) showed the starting material was consumed completely. The mixture was filtered, washed with HCl/EtOH, concentrated, basified to pH=7~8, extracted with EtOAc and H₂O, the organic layer was concentrated to give 4-

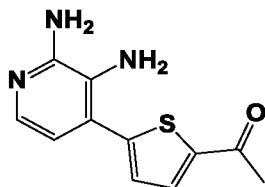
(5-fluorothiophen-2-yl)pyridine-2,3-diamine (**XCIV**) (550 mg, 2.63 mmol, 89.7% yield). ESIMS found $C_9H_8FN_3S$ m/z 210.0 (M+H).

[0793] The following intermediates were prepared in accordance with the procedure described in the above Scheme 19.



XCVI

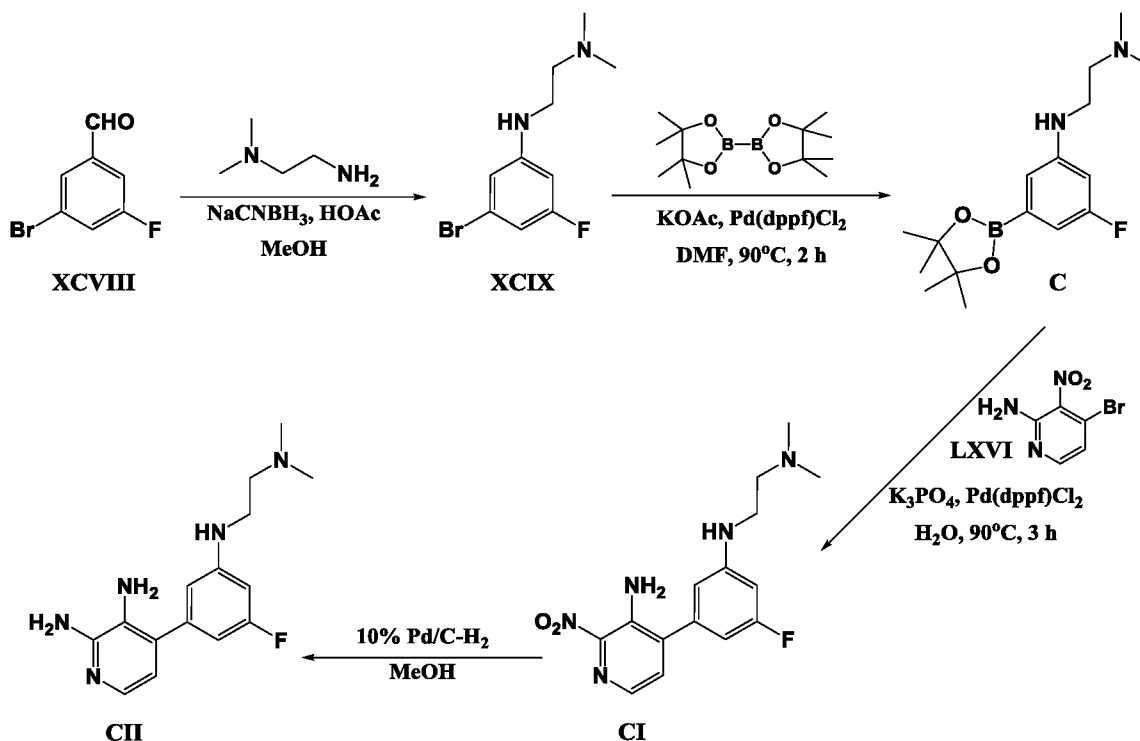
[0794] 4-(5-Methylthiophen-2-yl)pyridine-2,3-diamine (**XCVI**): Brown solid, (1.20 g, 5.85 mmol, 86.0% yield). 1H NMR (CD_3OD , 400 MHz) δ ppm 2.54 (s, 3H), 6.63 (d, $J=4.8$ Hz, 1H), 6.85 (s, 1H), 7.12 (s, 1H), 7.38 (d, $J=5.2$ Hz, 1H); ESIMS found $C_{10}H_{11}N_3S$ m/z 206.2 (M+H).



XCVII

[0795] 1-(5-(2,3-Diaminopyridin-4-yl)thiophen-2-yl)ethan-1-one (**XCVII**): Brown solid, (500 mg, 2.14 mmol, 56.4% yield). 1H NMR ($DMSO-d_6$, 400 MHz) δ ppm 2.55 (s, 3H), 4.89 (brs, 2H), 5.80 (brs, 2H), 6.52 (d, $J=5.2$ Hz, 1H), 7.33 (d, $J=5.2$ Hz, 1H), 7.46 (d, $J=4.0$ Hz, 1H), 7.96 (d, $J=4.0$ Hz, 1H); ESIMS found $C_{11}H_{11}N_3OS$ m/z 234 (M+H).

[0796] Preparation of intermediate 4-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)pyridine-2,3-diamine (CII) is depicted below in Scheme 20.



Scheme 20

Step 1

[0797] A solution of 3-bromo-5-fluorobenzaldehyde (XCVIII) (20.0 g, 98.2 mmol, 1.0 eq) in MeOH (1.8 L) was added N¹,N¹-dimethylethane-1,2-diamine (21.5 mL, 196.4 mmol, 2.0 eq). The pH was adjusted to 6 using HOAc and stirred for 1 h. NaCNBH₃ (8.6 g, 137.5 mmol, 1.4 eq) was added and stirred at room temperature overnight. The MeOH was removed under vacuum and the residue was partitioned between CHCl₃ and saturated aqueous NaHCO₃. The organic layer was dried over MgSO₄ and evaporated under vacuum. The crude product was purified on a silica gel column (100% CHCl₃ → 3:97 MeOH[7N NH₃]:CHCl₃) to produce N¹-(3-bromo-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine (XCIX) as a yellow oil (13.0 g, 49.9 mmol, 51% yield). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 1.28 (s, 6H), 2.39 (t, *J*=4Hz, 2H), 3.07 (q, *J*=6Hz, 2H), 6.10 (t, *J*=5Hz, 1H), 6.38 (td, *J*=12Hz, *J*=2Hz, 1H), 6.51 (td, *J*=8.6Hz, *J*=2Hz, 1H), 6.61 (t, *J*=2Hz, 1H); ESIMS found C₁₀H₁₄BrFN₂ *m/z* 261.0 (M+H).

Step 2

[0798] A solution of N¹-(3-bromo-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine (XCIX) (13.0 g, 49.9 mmol, 1.0 eq), bis(pinacolato)diboron (12.6 g, 59.9 mmol, 1.2 eq),

KOAc (12.1 g, 124.3 mmol, 2.5 eq) and dioxane (600 mL) was purged with argon. Pd(dppf)Cl₂ (2.0 g, 2.47 mmol, 0.05 eq) was added to the reaction and purged again with argon. The solution was heated at 90°C for 2 h. Once TLC showed the disappearance of (XCIX), the solution was cooled to room temperature and then concentrated under reduced pressure to produce crude N¹-(3-fluoro-5-(4,4,5,5-tetramethyl-1,3,2-dioxaborolan-2-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine (C) (7.4 g, 24.0 mmol, 48.2% yield).

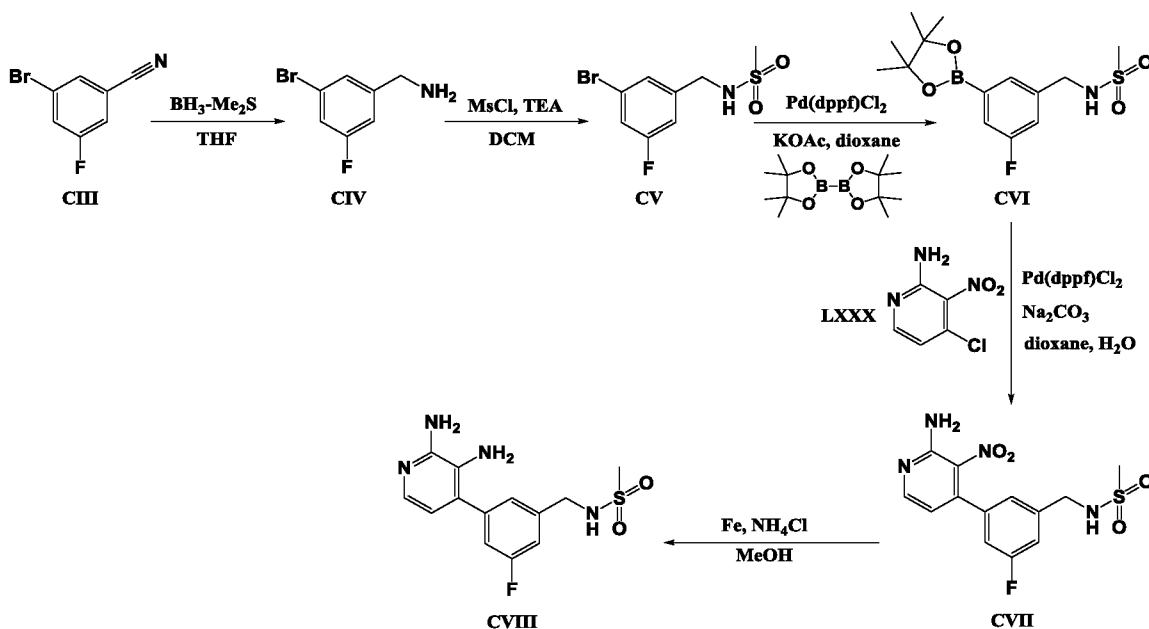
Step 3

[0799] To a solution of N¹-(3-fluoro-5-(4,4,5,5-tetramethyl-1,3,2-dioxaborolan-2-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine (C) (5.0 g, 16.22 mmol, 1.2 eq) in water (25 mL) was added K₂CO₃ (448 mg, 3.24 mmol, 2.0 eq), 4-bromo-3-nitropyridin-2-amine (LXVI) (3.5g, 16.22 mmol, 1.0 eq) and Pd(dppf)Cl₂ (1.0 g, 81.0 μmol, 0.05 eq). The solution was purged with argon and heated at 100°C for 48 h. The solution was cooled to room temperature and then concentrated under reduced pressure. The residue was purified on a silica gel column (100% CHCl₃ → 2:98 MeOH[7N NH₃]:CHCl₃) to give N¹-(3-(3-amino-2-nitropyridin-4-yl)-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine (CI) as a yellow amorphous solid (4.5 g, 14.1 mmol, 86.9% yield). ESIMS found for C₁₅H₁₈FN₅O₂ *m/z* 220.1(M+H).

Step 4

[0800] To a solution of N¹-(3-(3-amino-2-nitropyridin-4-yl)-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine (CI) (4.50 g, 10.43 mmol, 1.0 eq) in MeOH (15 mL) was added 10% Pd/C (540 mg, 15% by wt). The solution was purged with hydrogen and stirred for 48 h at room temperature under hydrogen (15 psi). The suspension was filtered through Celite[®] and concentrated under vacuum and purified by silica gel chromatography (MeOH:DCM = 10:1) to produce 4-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)pyridine-2,3-diamine (CII) as a tan solid (750.0 mg, 2.59 mmol, 24.9% yield). ¹H NMR (CD₃OD, 400 MHz) δ ppm 2.32 (s, 6H), 2.60 (t, *J*=6.8Hz, 2H), 3.26 (t, *J*=6.8Hz, 2H), 6.34-6.43 (m, 2H), 6.47 (d, *J*=5.6Hz, 1H), 7.58 (s, 1H), 7.75 (s, 1H); ESIMS found C₁₅H₂₀FN₅ *m/z* 290.1 (M+H).

[0801] Preparation of intermediate N-(3-(2,3-diaminopyridin-4-yl)-5-fluorobenzyl)methanesulfonamide (CVIII) is depicted below in Scheme 21.



Scheme 21

Step 1

[0802] A solution of 3-bromo-5-fluorobenzonitrile (CIII) (44.0 g, 220.0 mmol, 1.0 eq) was dissolved in THF (30 mL). $\text{BH}_3\text{-Me}_2\text{S}$ (33.43 g, 440.0 mmol, 2.0 eq) was added to the solution at 20°C. Then it was stirred at 80°C for 2 h, HCl (6 N, 100 mL) was added to the mixture slowly at 20°C. The mixture was stirred at 80°C for 1 h, then it was washed with EtOAc (300 ml). The water phase was basified with 50% aqueous NaOH and it was extracted with EtOAc (300 mL x 3). The combined organic layers were dried over anhydrous Na_2SO_4 and concentrated in vacuo to produce (3-bromo-5-fluoro-phenyl)methanamine (CIV) (24.0 g, 117.62 mmol, 53.5% yield). ¹H NMR (CDCl_3 , 300 MHz) 3.86 (s, 2H), 7.01 (d, $J=8\text{Hz}$, 1H), 7.12 (d, $J=8\text{Hz}$, 1H), 7.28 (s, 1H); ESIMS found $\text{C}_7\text{H}_7\text{BrFN}$ m/z 203.9 ($\text{Br}^{79}\text{M}+\text{H}$).

Step 2

[0803] A solution of (3-bromo-5-fluoro-phenyl)methanamine (CIV) (23.0 g, 112.7 mmol, 1.0 eq) was dissolved in DCM (15 mL), TEA (34.22 g, 338.2 mmol, 3.0 eq) was added to the mixture. Then MsCl (13.44 g, 117.3 mmol, 1.04 eq) was added slowly to the solution at 0°C. It was stirred at 0-30°C for 2 h. The reaction was washed with water and extracted with EtOAc. The combined organic layers were dried over anhydrous Na_2SO_4 and concentrated to give N-(3-bromo-5-fluorobenzyl) methanesulfonamide (CV) (34.0 g, 102.44 mmol, 90.9% yield, 85% purity) as an

oil. $^1\text{H NMR}$ (CDCl_3 , 300 MHz) 2.88 (s, 3H), 4.24 (d, $J=4.5\text{Hz}$, 2H), 6.99 (d, $J=9\text{Hz}$, 1H), 7.13 (dt, $J=8.1\text{Hz}$, $J=2\text{Hz}$, 1H), 7.25 (s, 1H); ESIMS found $\text{C}_8\text{H}_9\text{BrFNO}_2\text{S}$ m/z 282.0 ($\text{Br}^{79}\text{M}+\text{H}$).

Step 3

[0804] A solution of N-(3-bromo-5-fluorobenzyl)methanesulfonamide (**CV**) (34.0 g, 102.4 mmol, 1.0 eq) and 4,4,5,5-tetramethyl-2-(4,4,5,5-tetramethyl-1,3,2-dioxaborolan-2-yl)-1,3,2-dioxaborolane (52.02 g, 204.9 mmol, 2.0 eq), KOAc (20.11 g, 204.9 mmol, 2.0 eq) was dissolved in dioxane (20 mL). Then $\text{Pd}(\text{dppf})\text{Cl}_2$ (7.60 g, 10.2 mmol, 0.1 eq) was added to the mixture. It was stirred at 90°C for 2 h. Then the solvent was removed to get the residue which was purified by silica gel column (PE:EtOAc = 10:1 \rightarrow 100% EtOAc) to get N-(3-fluoro-5-(4,4,5,5-tetramethyl-1,3,2-dioxaborolan-2-yl)benzyl)methanesulfonamide (**CVI**) (30.0 g, crude). $^1\text{H NMR}$ (CDCl_3 , 400 MHz) 1.37 (s, 12H), 2.92 (s, 3H), 4.34 (d, $J=6.3\text{Hz}$, 2H), 7.19 (dt, $J=9.3\text{Hz}$, $J=2.1\text{Hz}$, 1H), 7.44 (dd, $J=8.7\text{Hz}$, $J=2.4\text{Hz}$, 1H), 7.54 (s, 1H); ESIMS found $\text{C}_{14}\text{H}_{21}\text{BFNO}_4\text{S}$ m/z 330.1 ($\text{M}+\text{H}$).

Step 4

[0805] A solution of N-(3-fluoro-5-(4,4,5,5-tetramethyl-1,3,2-dioxaborolan-2-yl)benzyl)methanesulfonamide (**CVI**) (6.83 g, 20.75 mmol, 1.2 eq), 4-chloro-3-nitropyridin-2-amine (**LXXX**) (3.0 g, 17.29 mmol, 1.0 eq), Na_2CO_3 (6.41 g, 60.52 mmol) and $\text{Pd}(\text{dppf})\text{Cl}_2$ (641.27 mg, 864.50 μmol) in dioxane (40 mL) and H_2O (8 mL) was de-gassed and then heated to 80°C overnight under N_2 . TLC (PE:EtOAc = 1:1) showed the starting material was consumed completely. The reaction mixture was poured into H_2O (300 mL). The mixture was extracted with EtOAc (3 x 250 mL). The organic phase was washed with saturated brine (300 mL), dried over anhydrous NaSO_4 , concentrated in vacuum to give a residue. The crude product was purified by silica gel chromatography (PE:EtOAc = 10:1) to give N-(3-(2-amino-3-nitropyridin-4-yl)-5-fluorobenzyl)methanesulfonamide (**CVII**) (2.2 g, 6.46 mmol, 37.4% yield) as brown solid. ESIMS found $\text{C}_{13}\text{H}_{13}\text{FN}_4\text{O}_4\text{S}$ m/z 341.1 ($\text{M}+\text{H}$).

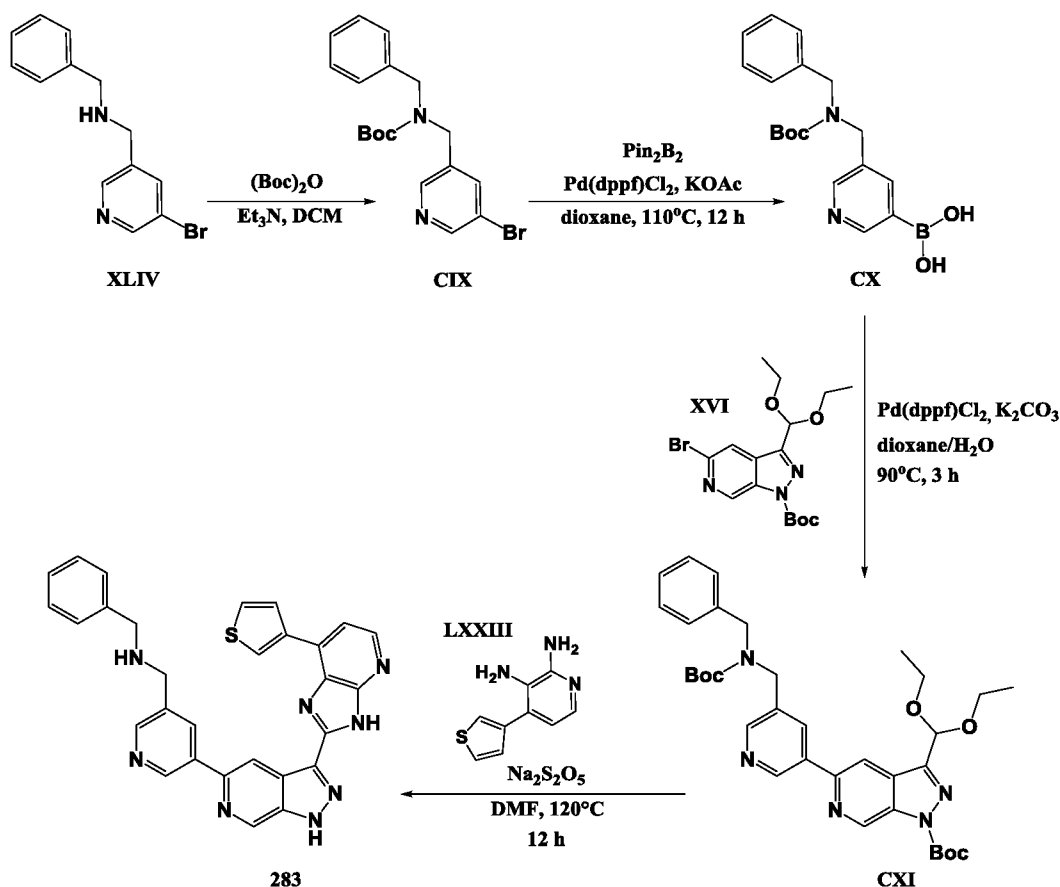
Step 5

[0806] A solution of N-[[3-(4-amino-3-nitro-2-pyridyl)-5-fluoro-phenyl]methyl]methanesulfonamide (**CVII**) (2.2 g, 6.46 mmol, 1.0 eq), Fe (1.44 g, 25.84 mmol, 4.0 eq) and NH_4Cl (2.8 g, 51.68 mmol, 8.0 eq) was dissolved in MeOH (30 mL). The mixture was stirred at 80°C for 16 h. The mixture was cooled to room temperature and concentrated in reduced pressure

at 60°C. The combined organic phase was washed with saturated brine (100 mL x 2), dried over anhydrous Na₂SO₄, filtered and concentrated in vacuo to get crude N-(3-(3,4-diaminopyridin-2-yl)-5-fluorobenzyl) methanesulfonamide (CVIII) (1.60 g, 5.16 mmol, 79.8% yield) as brown solid. ¹H NMR (CD₃OD, 400 MHz) 2.96 (s, 3H), 3.37 (s, 2H), 6.53 (d, *J*=4Hz, 1H), 7.13 (d, *J*=8.8Hz, 1H), 7.20 (d, *J*=9.2Hz, 1H), 7.31 (s, 1H), 7.44 (d, *J*=4Hz, 1H); ESIMS found C₁₃H₁₅FN₄O₂S *m/z* 311.1 (M+H).

Example 1.

[0807] Preparation of N-isopropyl-5-(3-(7-(pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-indazol-5-yl)pyridin-3-amine (90) is depicted below in Scheme 22.



Scheme 22

Step 1

[0808] To a solution of N-benzyl-1-(5-bromopyridin-3-yl)methanamine (XLIV) (8.32 g, 30.0 mmol, 1 eq) and TEA (10.45 mL, 75.0 mmol, 2.5 eq) in DCM (160 mL) was added portionwise (Boc)₂O (7.2 g, 33.0 mmol, 1.1 eq) at 0°C, the reaction mixture was stirred at room

temperature for 1 h. The mixture was washed with water (80 mL), brine (80 mL), the organic layer was separated, dried over MgSO₄ and concentrated in vacuo to give *tert*-butyl benzyl((5-bromopyridin-3-yl)methyl)carbamate (**CIX**) (10.38 g, 27.51 mmol, 91.7% yield) as a white solid. ESIMS found for C₁₈H₂₁BrN₂O₂ *m/z* 377.1 (M+H).

Step 2

[0809] A solution of *tert*-butyl benzyl((5-bromopyridin-3-yl)methyl)carbamate (**CIX**) (10.00 g, 26.5 mmol, 1.0 eq), Pd(dppf)Cl₂ (0.97 g, 1.33 mmol, 0.05 Eq), KOAc (5.20 g, 53.0 mmol, 2.0 eq) and Pin₂B₂ (6.7 g, 26.5 mmol, 1.0 eq) in dioxane (130 mL) was de-gassed and then heated to 110°C under N₂ for 12 hr. TLC (PE:EtOAc=2: 1) showed the starting material was consumed completely. The reaction mixture was filtered and concentrated to afford the crude (5-((benzyl(*tert*-butoxycarbonyl)amino)methyl)pyridin-3-yl)boronic acid (**CX**). The crude product was used in the next step without purification.

Step 3

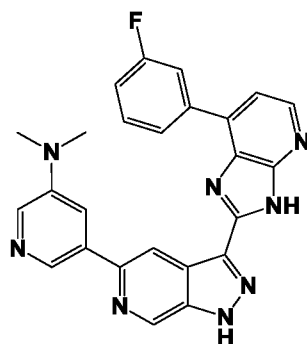
[0810] To a solution of *tert*-butyl 5-bromo-3-(diethoxymethyl)-1H-pyrazolo [3,4-*c*]pyridine-1-carboxylate (**XVI**) (3.50 g, 8.74 mmol, 1.0 eq) and (5-((benzyl(*tert*-butoxycarbonyl)amino)methyl)pyridin-3-yl)boronic acid (**CX**) (4.49 g, 13.12 mmol, 1.5 eq) in dioxane (30 mL) and H₂O (5 mL) was added Pd(dppf)Cl₂ (129.66 mg, 174.8 μmol, 0.02 eq), K₂CO₃ (2.42 g, 17.48 mmol, 2.0 eq) at room temperature. The mixture was stirred at 90°C for 3 hr. TLC (PE:EtOAc=0:1) showed that the starting material was consumed completely. The mixture was added water (50 mL) and extracted with EtOAc (80 mL x 2). The organic layers were washed with brine (60 mL), concentrated. The residue was purified by chromatography on silica gel (PE:EtOAc=1:1→1:5) to afford *tert*-butyl 5-(5-((benzyl(*tert*-butoxycarbonyl)amino)methyl)pyridin-3-yl)-3-(diethoxymethyl)-1H-pyrazolo[3,4-*c*]pyridine-1-carboxylate (**CXI**) (1.90 g, 3.95 mmol, 45.2% yield) as a brown oil. ¹H NMR (CDCl₃, 400 MHz) δ ppm 1.32 (t, *J*=6.8Hz, 6H), 1.61 (s, 9H), 1.79 (s, 9H), 3.63-3.75 (m, 2H), 3.80-3.92 (m, 2H), 4.41-4.50 (m, 2H), 4.50-4.61 (m, 2H), 5.87 (s, 1H), 7.21-7.33 (m, 2H), 7.33-7.42 (m, 3H), 8.25 (brs, 1H), 8.36 (s, 1H), 8.45-8.56 (m, 1H), 9.18 (s, 1H), 9.55 (s, 1H); ESIMS found for C₃₄H₄₃N₅O₆ *m/z* 618.1 (M+H).

Step 4

[0811] A mixture of *tert*-butyl 5-(5-((benzyl(*tert*-butoxycarbonyl)amino)methyl)pyridin-3-yl)-3-(diethoxymethyl)-1H-pyrazolo[3,4-*c*]pyridine-1-carboxylate (**CXI**) (100.0 mg, 0.16 mmol, 1 eq), 4-(thiophen-3-yl)pyridine-2,3-diamine (**LXXIII**) (31 mg, 0.16 mmol, 1.0

eq) and $\text{Na}_2\text{S}_2\text{O}_5$ (37 mg, 0.19 mmol, 1.2 eq) in DMF (2 mL) was stirred at 120°C for 24 h. LC/MS showed **CXI** was consumed. Water (5 mL) was added in dropwise and the mixture was filtered. The filtrate was washed by MeOH (0.5 mL) and purified by pre-HPLC (HCl) to give N-benzyl-1-(5-(3-(7-(thiophen-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl) methanamine (**283**) (12.1 mg, 0.02 mmol, 14.5%) as a white solid. ^1H NMR (DMSO-*d*₆, 400 MHz) δ ppm 4.32 (t, $J=5.90$ Hz, 2H), 4.48 (t, $J=4.96$ Hz, 2H), 7.38-7.51 (m, 3H), 7.60 (d, $J=1.51$ Hz, 1H), 7.61 (d, $J=1.88$ Hz, 1H), 7.72 (d, $J=5.27$ Hz, 1H), 7.81 (dd, $J=5.02, 3.01$ Hz, 1H), 8.23 (d, $J=4.77$ Hz, 1H), 8.42 (d, $J=5.27$ Hz, 1H), 8.95 (d, $J=1.51$ Hz, 1H), 9.04 (brs, 2H), 9.11 (d, $J=1.00$ Hz, 1H), 9.38 (d, $J=1.13$ Hz, 1H), 9.50 (d, $J=2.01$ Hz, 1H), 9.89 (brs, 2H), 14.74 (brs, 1H); ESIMS found for $\text{C}_{29}\text{H}_{22}\text{N}_8\text{S}$ m/z 515.1 (M+1).

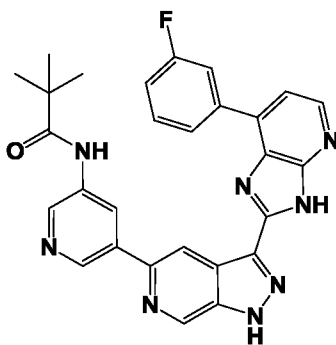
[0812] The following compound was prepared in accordance with the procedure described in the above Example 1.



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[0813] 5-(3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)-N,N-dimethylpyridin-3-amine **7**.

[0814] White solid (53.5 mg, 0.12 mmol). ^1H NMR (DMSO-*d*₆, 400 MHz) δ ppm 3.19 (s, 6H), 7.38 (tdd, $J=5.64\text{Hz}, J=2.28\text{Hz}, J=0.76\text{Hz}$, 1H), 7.66-7.75 (m, 2H), 8.28 (d, $J=2.64$ Hz, 1H), 8.30-8.36 (m, 2H), 8.40 (d, $J=10.16$ Hz, 1H), 8.48 (d, $J=5.15$ Hz, 1H), 8.72 (s, 1H), 9.04 (d, $J=1.25$ Hz, 1H), 9.38 (d, $J=1.25$ Hz, 1H), 14.76 (brs, 1H); ESIMS found for $\text{C}_{25}\text{H}_{19}\text{FN}_8$ m/z 451.2 (M+1).

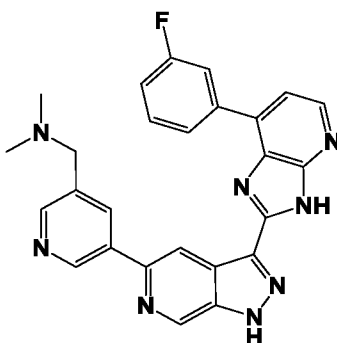


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8

[0815] N-(5-(3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide **8**.

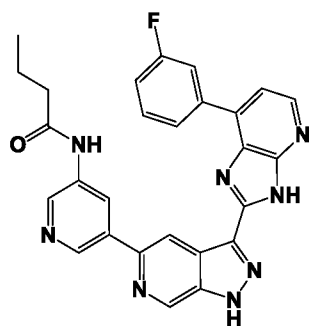
[0816] White solid (47.9 mg, 0.09 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.30 (s, 9H), 7.31-7.40 (m, 1H), 7.65-7.74 (m, 2H), 8.37 (brs, 1H), 8.44 (d, *J*=5.14 Hz, 1H), 8.59 (brs, 1H), 8.86 (s, 1H), 8.94 (d, *J*=1.51 Hz, 1H), 8.98 (s, 1H), 9.03 (s, 1H), 9.32 (s, 1H), 9.60 (s, 1H); ESIMS found for C₂₈H₂₃FN₈O *m/z* 507.1 (M+1).



13

[0817] 1-(5-(3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine **13**.

[0818] White solid (12.7 mg, 0.03 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 2.81 (s, 3H), 2.82 (s, 3H), 4.51 (d, *J*=4.89 Hz, 2H), 7.43 (td, *J*=8.91, 2.63 Hz, 1H), 7.66-7.77 (m, 2H), 8.24-8.32 (m, 1H), 8.48 (d, *J*=5.15 Hz, 1H), 8.54-8.66 (m, 1H), 8.83 (d, *J*=1.88 Hz, 1H), 8.87 (d, *J*=1.76 Hz, 1H), 9.07 (d, *J*=1.00 Hz, 1H), 9.37 (d, *J*=1.13 Hz, 1H), 9.45 (d, *J*=2.01 Hz, 1H), 10.55 (brs, 1H), 14.67 (brs, 1H); ESIMS found for C₂₆H₂₁FN₈ *m/z* 465.1 (M+1).

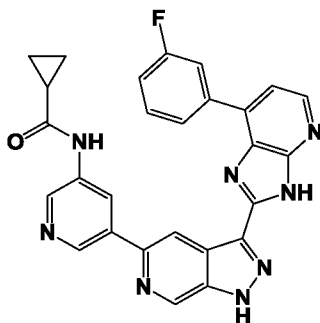


17

[0819] N-(5-(3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide **17**.

[0820] White solid (97.1 mg, 0.20 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.95 (t, *J*=7.34 Hz, 3H), 1.67 (sxt, *J*=7.28Hz, 2H), 2.47 (t, 2H), 7.37-7.47 (m, 1H), 7.76 (q, *J*=6.52Hz,

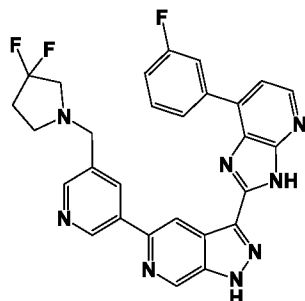
1H), 7.80 (d, $J=5.5$ Hz, 1H), 8.05-8.16 (m, 2H), 8.60 (d, $J=5.27$ Hz, 1H), 8.97 (s, 1H), 9.18 (s, 1H), 9.28 (s, 1H), 9.39 (s, 1H), 9.40 (s, 1H), 11.59 (s, 1H); ESIMS found for $C_{27}H_{21}FN_8O$ m/z 493.1 (M+1).



19

[0821] N-(5-(3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide **19**.

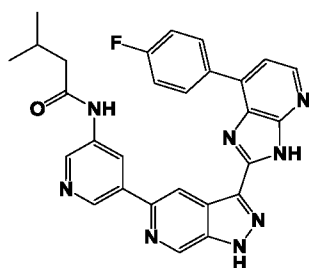
[0822] White solid (68.6 mg, 0.14 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 0.84-0.93 (m, 4H), 1.83-1.91 (m, 1H), 7.38 (td, $J=8.28$ Hz, $J=2.52$ Hz, 1H), 7.65-7.71 (m, 1H), 7.73 (d, $J=5.27$ Hz, 1H), 8.39 (d, $J=7.78$ Hz, 1H), 8.45 (d, $J=5.14$ Hz, 1H), 8.58 (d, $J=10.79$ Hz, 1H), 8.77 (d, $J=2.38$ Hz, 1H), 8.91 (s, 1H), 9.00 (d, $J=2.51$ Hz, 2H), 9.34 (s, 1H), 10.59 (s, 1H), 14.09 (s, 1H), 14.49 (s, 1H); ESIMS found for $C_{27}H_{20}FN_7O$ m/z 491.1 (M+1).



25

[0823] 5-(5-((3,3-Difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(3-fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridine **25**.

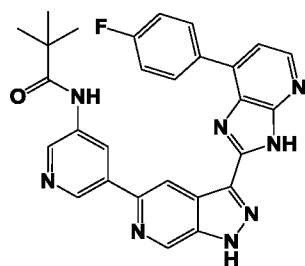
[0824] White solid (59.1 mg, 0.11 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 2.54-2.65 (m, 2H), 3.44-3.60 (m, 4H), 4.60 (brs, 2H), 7.41 (td, $J=8.28$, 2.26 Hz, 1H), 7.67-7.76 (m, 2H), 8.29 (d, $J=6.53$ Hz, 1H), 8.49 (d, $J=5.14$ Hz, 1H), 8.51-8.59 (m, 1H), 8.97 (s, 2H), 9.08 (s, 1H), 9.38 (d, $J=1.00$ Hz, 1H), 9.46 (s, 1H), 14.74 (brs, 1H); ESIMS found for $C_{28}H_{21}F_3N_8$ m/z 527.1 (M+1).



28

[0825] N-(5-(3-(7-(4-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide **28**.

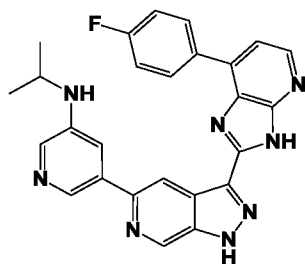
[0826] White solid (36.5 mg, 0.07 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.99 (d, *J*=6.65 Hz, 7H), 2.14-2.24 (m, 1H), 2.32 (d, *J*=7.28 Hz, 2H), 7.51 (t, *J*=8.91 Hz, 2H), 7.64 (d, *J*=5.27 Hz, 1H), 8.43 (d, *J*=5.15 Hz, 1H), 8.64 (dd, *J*=8.91, 5.52 Hz, 2H), 8.71 (d, *J*=2.26 Hz, 1H), 8.94 (s, 1H), 8.98 (t, *J*=2.07 Hz, 1H), 9.04 (d, *J*=1.88 Hz, 1H), 9.34 (d, *J*=1.13 Hz, 1H), 10.31 (s, 1H), 14.02 (s, 1H), 14.48 (s, 1H); ESIMS found for C₂₈H₂₃FN₈O *m/z* 507.1 (M+1).



34

[0827] N-(5-(3-(7-(4-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide **34**.

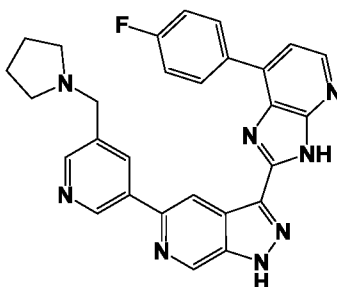
[0828] White solid (69.2 mg, 0.14 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.33 (s, 9H), 7.53 (t, *J*=8.85 Hz, 2H), 7.65 (d, *J*=5.27 Hz, 1H), 8.46 (d, *J*=5.14 Hz, 1H), 8.61 (brs, 2H), 9.01 (d, *J*=1.00 Hz, 1H), 9.22 (d, *J*=1.76 Hz, 1H), 9.25 (d, *J*=1.88 Hz, 1H), 9.35 (s, 1H), 9.39 (d, *J*=1.13 Hz, 1H), 10.24 (s, 1H), 14.71 (brs, 1H); ESIMS found for C₂₈H₂₃FN₈O *m/z* 507.2 (M+1).



38

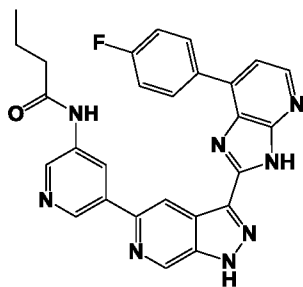
[0829] 5-(3-(7-(4-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)-N-isopropylpyridin-3-amine **38**.

[0830] White solid (63.2 mg, 0.14 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.22 (d, *J*=6.27 Hz, 6H), 3.71-3.83 (m, 1H), 6.44 (brs, 1H), 7.46 (t, *J*=8.85 Hz, 2H), 7.62 (d, *J*=4.02 Hz, 1H), 7.90 (brs, 1H), 8.06 (d, *J*=2.38 Hz, 1H), 8.43 (d, *J*=5.02 Hz, 1H), 8.56 (s, 1H), 8.61 (brs, 2H), 8.94 (s, 1H), 9.32 (s, 1H), 14.02 (brs, 1H), 14.49 (brs, 1H); ESIMS found for C₂₆H₂₁FN₈ *m/z* 465.1 (M+1).

**40**

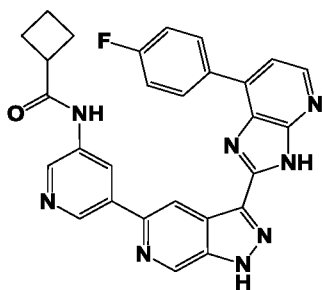
[0831] 3-(7-(4-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridine **40**.

[0832] White solid (21.2 mg, 0.04 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.87-1.97 (m, 2H), 2.01-2.12 (m, 2H), 3.11-3.23 (m, 2H), 3.41-3.52 (m, 3H), 4.63 (d, *J*=5.65 Hz, 2H), 7.53 (t, *J*=8.91 Hz, 2H), 7.64 (d, *J*=5.27 Hz, 1H), 8.47 (d, *J*=5.27 Hz, 1H), 8.54 (brs, 2H), 9.02-9.08 (m, 3H), 9.38 (d, *J*=1.13 Hz, 1H), 9.48 (s, 1H), 11.47 (brs, 1H), 14.83 (brs, 1H); ESIMS found for C₂₈H₂₃FN₈ *m/z* 491.1 (M+1).

**43**

[0833] N-(5-(3-(7-(4-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide **43**.

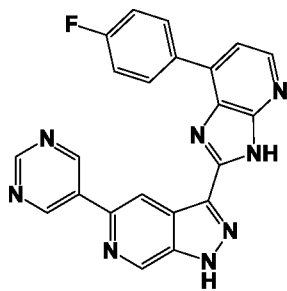
[0834] White solid (93.7 mg, 0.19 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.95 (t, *J*=7.34 Hz, 3H), 1.68 (sxt, *J*=7.38 Hz, 2H), 2.47 (t, 2H), 7.53 (t, *J*=8.78 Hz, 2H), 7.77 (d, *J*=5.77 Hz, 1H), 8.21-8.32 (m, 2H), 8.59 (d, *J*=5.77 Hz, 1H), 8.90 (s, 1H), 9.18 (s, 1H), 9.25 (s, 1H), 9.39 (d, *J*=8.28 Hz, 2H), 11.60 (s, 1H); ESIMS found for C₂₇H₂₁FN₈O *m/z* 493.1 (M+1).



46

[0835] N-(5-(3-(7-(4-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide **46**.

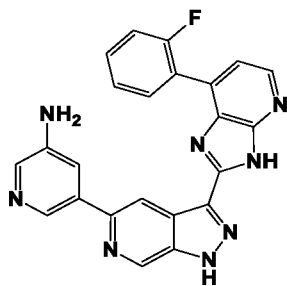
[0836] White solid (27.1 mg, 0.05 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.82-1.93 (m, 1H), 2.01 (sxt, *J*=9.28Hz, 1H), 2.16-2.27 (m, 2H), 2.33 (quin, *J*=9.52 Hz, 2H), 3.42 (quin, *J*=8.49 Hz, 2H), 7.55 (t, *J*=8.78 Hz, 2H), 7.65 (d, *J*=5.27 Hz, 1H), 8.47 (d, *J*=5.14 Hz, 1H), 8.55 (brs, 2H), 8.99 (s, 1H), 9.15 (s, 1H), 9.20 (s, 1H), 9.34 (brs, 1H), 9.38 (s, 1H), 11.02 (s, 1H), 14.82 (brs, 1H); ESIMS found for C₂₈H₂₁FN₈O *m/z* 505.2 (M+1).



52

[0837] 3-(7-(4-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1H-pyrazolo[3,4-*c*]pyridine **52**.

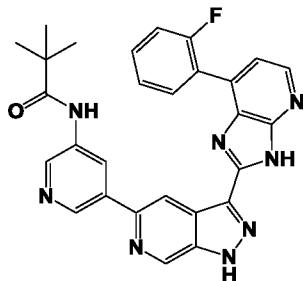
[0838] White solid (87.8 mg, 0.21 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 7.46 (t, *J*=8.91 Hz, 2H), 7.66 (d, *J*=5.27 Hz, 1H), 8.49 (d, *J*=5.27 Hz, 3H), 9.03 (d, *J*=0.88 Hz, 1H), 9.28 (s, 1H), 9.38 (d, *J*=1.13 Hz, 1H), 9.50 (s, 2H), 14.72 (brs, 1H); ESIMS found for C₂₂H₁₃FN₈ *m/z* 409.0 (M+1).



55

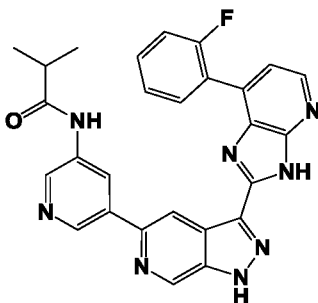
[0839] 5-(3-(7-(2-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine **55**.

[0840] White solid (9.8 mg, 0.02 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 7.46-7.57 (m, 3H), 7.59-7.68 (m, 1H), 8.05 (brs, 1H), 8.09 (d, *J*=2.38 Hz, 1H), 8.37 (t, *J*=1.76 Hz, 1H), 8.56 (d, *J*=5.27 Hz, 1H), 8.58 (s, 1H), 8.90 (d, *J*=1.13 Hz, 1H), 9.36 (d, *J*=1.13 Hz, 1H), 14.89 (brs, 1H); ESIMS found for C₂₃H₁₅FN₈ *m/z* 423.1 (M+1).

**60**

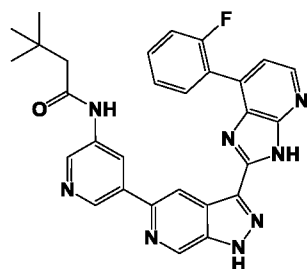
[0841] N-(5-(3-(7-(2-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide **60**.

[0842] White solid (49.8 mg, 0.10 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.31 (s, 9H), 7.41-7.54 (m, 3H), 7.55-7.63 (m, 1H), 8.46 (d, *J*=3.26 Hz, 1H), 8.82 (brs, 1H), 8.87 (s, 1H), 8.92 (brs, 1H), 8.96 (brs, 1H), 9.31 (s, 1H), 9.63 (s, 1H), 14.06 (brs, 1H), 14.39 (brs, 1H); ESIMS found for C₂₈H₂₃FN₈O *m/z* 507.1 (M+1).

**61**

[0843] N-(5-(3-(7-(2-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide **61**.

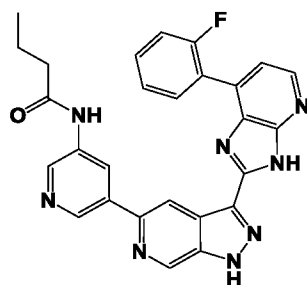
[0844] White solid (58.8 mg, 0.12 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.19 (d, *J*=6.90 Hz, 6H), 2.73 (spt, *J*=6.87 Hz, 1H), 7.43-7.54 (m, 3H), 7.57-7.65 (m, 1H), 8.12 (brs, 1H), 8.51 (d, *J*=4.89 Hz, 1H), 8.93 (d, *J*=0.88 Hz, 1H), 9.04 (s, 1H), 9.09 (s, 2H), 9.36 (d, *J*=1.25 Hz, 1H), 10.71 (s, 1H), 14.68 (brs, 1H); ESIMS found for C₂₇H₂₁FN₈O *m/z* 493.1 (M+1).



68

[0845] N-(5-(3-(7-(2-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide **68**.

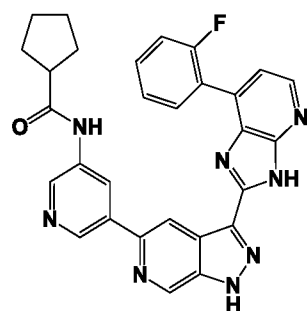
[0846] White solid (39.7 mg, 0.08 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.08 (s, 9H), 2.36 (s, 2H), 7.44-7.57 (m, 3H), 7.59-7.67 (m, 1H), 8.07 (brs, 1H), 8.55 (d, *J*=5.27 Hz, 1H), 8.95 (d, *J*=1.13 Hz, 1H), 9.14 (s, 1H), 9.21 (s, 2H), 9.39 (d, *J*=1.13 Hz, 1H), 11.08 (s, 1H), 14.89 (brs, 1H); ESIMS found for C₂₉H₂₅FN₈O *m/z* 521.3 (M+1).



69

[0847] N-(5-(3-(7-(2-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide **69**.

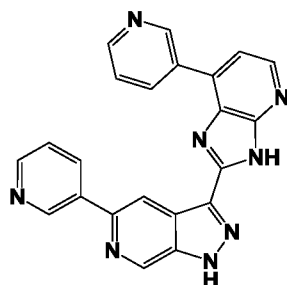
[0848] White solid (66.8 mg, 0.14 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.91 (t, *J*=7.40 Hz, 3H), 1.59-1.68 (sxt, *J*=7.40 Hz, 2H), 2.47 (t, *J*=6.88 Hz, 3H), 7.44-7.53 (m, 2H), 7.68 (q, *J*=6.48 Hz, 1H), 7.76-7.84 (m, 2H), 8.74 (d, *J*=6.02 Hz, 1H), 8.90 (s, 1H), 9.17 (s, 1H), 9.29 (s, 1H), 9.39 (s, 1H), 9.42 (s, 1H), 11.70 (s, 1H); ESIMS found for C₂₇H₂₁FN₈O *m/z* 493.1 (M+1).



73

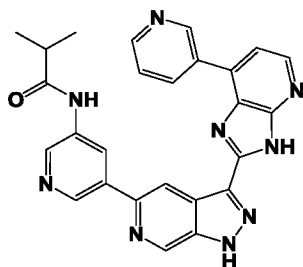
[0849] N-(5-(3-(7-(2-Fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide **73**.

[0850] White solid (21.2 mg, 0.04 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.53-1.66 (m, 2H), 1.66-1.86 (m, 4H), 1.88-2.02 (m, 2H), 2.89 (quin, *J*=7.81 Hz, 1H), 7.39-7.53 (m, 3H), 7.55-7.64 (m, 1H), 8.28 (s, 1H), 8.48 (d, *J*=4.77 Hz, 1H), 8.84 (brs, 1H), 8.88 (s, 1H), 8.93 (brs, 1H), 8.98 (brs, 1H), 9.33 (s, 1H), 10.40 (brs, 1H), 14.52 (brs, 1H); ESIMS found for C₂₉H₂₃FN₈O *m/z* 519.1 (M+1).

**82**

[0851] 5-(Pyridin-3-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridine **82**.

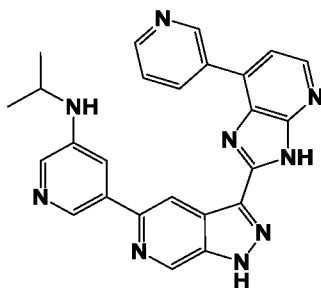
[0852] White solid (50.4 mg, 0.13 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 7.92 (d, *J*=5.14 Hz, 1H), 8.20-8.26 (m, 1H), 8.30 (brs, 1H), 8.59 (d, *J*=5.14 Hz, 1H), 8.97 (d, *J*=5.90 Hz, 1H), 9.05 (d, *J*=5.27 Hz, 1H), 9.17 (s, 1H), 9.29-9.35 (m, 1H), 9.43 (d, *J*=1.13 Hz, 1H), 9.50 (brs, 1H), 9.66 (d, *J*=2.01 Hz, 1H), 10.09 (s, 1H), 14.88 (brs, 1H); ESIMS found for C₂₂H₁₄N₈ *m/z* 391.0 (M+1).

**87**

[0853] N-(5-(3-(7-(Pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide **87**.

[0854] White solid (10.0 mg, 0.02 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.19 (d, *J*=6.80Hz, 6H), 2.71 (spt, *J*=6.80Hz, 1H), 7.46 (t, *J*=7.80Hz, 1H), 7.69 (d, *J*=8.40Hz, 2H), 7.92 (d, *J*=8.00Hz, 1H), 8.76 (d, *J*=1.2Hz, 1H), 8.86 (d, *J*=2.4Hz, 1H), 8.94 (d, *J*=4.0Hz, 2H), 9.04 (s,

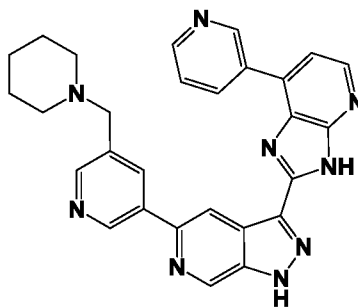
1H), 9.12 (brs, 1H), 9.34 (d, $J=1.2$ Hz, 1H), 9.61 (brs, 1H), 10.31 (s, 1H), 13.57 (brs, 1H), 14.45 (s, 1H); ESIMS found for $C_{26}H_{21}N_9O$ m/z 476.1 (M+1).



90

[0855] N-Isopropyl-5-(3-(7-(pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine **90**.

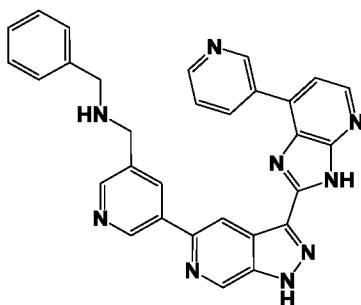
[0856] White solid (103.0 mg, 0.23 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 1.21 (d, $J=6.15$ Hz, 7H), 3.68-3.79 (m, 1H), 5.89 (d, $J=8.28$ Hz, 1H), 7.57 (s, 1H), 7.64 (dd, $J=7.97$, 5.46 Hz, 1H), 7.72 (d, $J=5.02$ Hz, 1H), 8.03 (d, $J=1.25$ Hz, 1H), 8.47 (d, $J=5.02$ Hz, 1H), 8.49 (s, 1H), 8.74 (d, $J=3.26$ Hz, 1H), 8.84 (d, $J=8.16$ Hz, 1H), 8.90 (s, 1H), 9.31 (d, $J=1.00$ Hz, 1H), 9.76 (s, 1H), 14.10 (brs, 1H), 14.44 (brs, 1H); ESIMS found for $C_{25}H_{21}N_9$ m/z 448.1 (M+1).



93

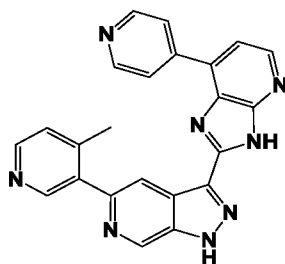
[0857] 5-(5-(Piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridine **93**.

[0858] White solid (17.8 mg, 0.04 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 1.31-1.45 (m, 1H), 1.62-1.74 (m, 1H), 1.74-1.86 (m, 4H), 2.90-3.05 (m, 2H), 3.38-3.46 (m, 2H), 4.53 (brs, 2H), 7.74-7.81 (m, 1H), 7.84-7.91 (m, 1H), 8.51 (d, $J=5.15$ Hz, 1H), 8.82 (s, 1H), 8.85-8.91 (m, 2H), 9.02 (brs, 1H), 9.06 (s, 1H), 9.36 (d, $J=1.13$ Hz, 1H), 9.44 (d, $J=2.13$ Hz, 1H), 9.90 (brs, 1H), 10.80 (brs, 1H), 14.73 (s, 1H); ESIMS found for $C_{28}H_{25}N_9$ m/z 488.1 (M+1).

**101**

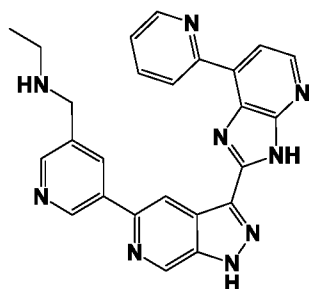
[0859] N-Benzyl-1-(5-(3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)methanamine **101**.

[0860] White solid (16.2 mg, 0.03 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 4.34 (brs, 2H), 4.62 (brs, 2H), 7.38-7.48 (m, 3H), 7.62 (d, *J*=1.38 Hz, 1H), 7.64 (d, *J*=1.88 Hz, 1H), 7.92 (d, *J*=5.14 Hz, 1H), 8.34 (dd, *J*=8.03, 5.65 Hz, 1H), 8.59 (d, *J*=5.15 Hz, 1H), 9.04 (s, 1H), 9.09 (d, *J*=5.02 Hz, 1H), 9.14 (d, *J*=1.13 Hz, 1H), 9.38 (brs, 1H), 9.41 (d, *J*=1.26 Hz, 1H), 9.49 (d, *J*=7.15 Hz, 1H), 9.61 (d, *J*=1.76 Hz, 1H), 10.10 (brs, 2H), 10.18 (brs, 1H), 14.93 (brs, 1H); ESIMS found for C₃₀H₂₃N₉ *m/z* 510.2 (M+1).

**109**

[0861] 5-(4-Methylpyridin-3-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine **109**.

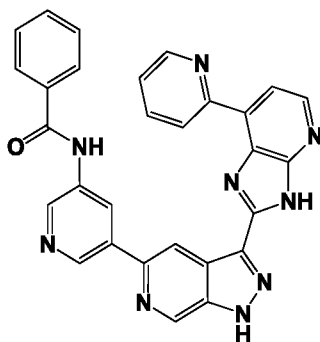
[0862] White solid (23.0 mg, 0.06 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 2.74 (s, 3H), 7.91 (d, *J*=5.14 Hz, 1H), 8.06 (d, *J*=6.02 Hz, 1H), 8.58 (d, *J*=5.15 Hz, 1H), 8.81 (s, 1H), 8.86 (d, *J*=6.02 Hz, 1H), 8.91 (brs, 2H), 9.00 (s, 1H), 9.01 (brs, 1H), 9.18 (s, 1H), 9.41 (d, *J*=1.13 Hz, 1H), 14.83 (brs, 1H); ESIMS found for C₂₃H₁₆N₈ *m/z* 405.1 (M+1).

**198**

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[0863] N-((5-(3-(7-(Pyridin-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine **136**.

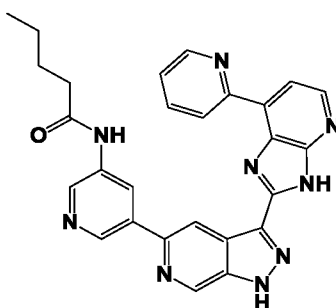
[0864] White solid (11.2 mg, 0.03 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.28 (t, *J*=7.28 Hz, 3H), 3.01-3.13 (m, 2H), 4.41 (t, *J*=5.58 Hz, 2H), 7.59-7.66 (m, 1H), 8.14 (d, *J*=5.15 Hz, 1H), 8.23 (td, *J*=7.81, 1.82 Hz, 1H), 8.57 (d, *J*=5.27 Hz, 1H), 8.88-8.95 (m, 2H), 8.99 (s, 1H), 9.11 (s, 1H), 9.20 (brs, 1H), 9.34-9.45 (m, 3H), 9.50 (d, *J*=2.01 Hz, 1H), 14.79 (brs, 1H); ESIMS found for C₂₅H₂₁N₉ *m/z* 448.2 (M+1).



141

[0865] N-(5-(3-(7-(Pyridin-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide **141**.

[0866] White solid (28.6 mg, 0.06 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 7.37 (brs, 1H), 7.55-7.72 (m, 4H), 8.07-8.19 (m, 4H), 8.49 (d, *J*=5.65 Hz, 1H), 8.79-8.87 (m, 1H), 9.02 (brs, 2H), 9.09 (brs, 1H), 9.15 (d, *J*=1.63 Hz, 1H), 9.35 (s, 1H), 10.76 (s, 1H); ESIMS found for C₂₉H₁₉N₉O *m/z* 510.1 (M+1).

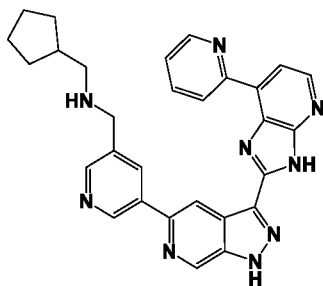


148

[0867] N-(5-(3-(7-(Pyridin-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide **148**.

[0868] White solid (73.6 mg, 0.15 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.94 (t, *J*=7.34 Hz, 3H), 1.39 (sxt, *J*=7.64 Hz, 2H), 1.66 (quin, *J*=7.43 Hz, 2H), 2.44 (t, *J*=7.47 Hz, 2H),

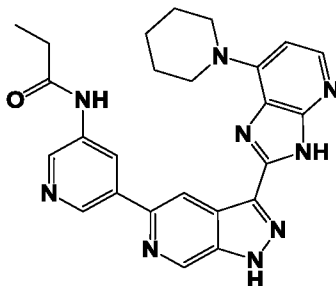
7.51 (dd, $J=6.84, 5.21$ Hz, 1H), 8.15-8.24 (m, 2H), 8.49 (d, $J=5.14$ Hz, 1H), 8.70 (brs, 1H), 8.83 (d, $J=4.64$ Hz, 1H), 8.98 (s, 1H), 9.01 (s, 1H), 9.06 (brs, 1H), 9.34 (s, 1H), 9.48 (d, $J=8.16$ Hz, 1H), 10.32 (s, 1H), 14.11 (brs, 1H), 14.51 (brs, 1H); ESIMS found for $C_{27}H_{23}N_9O$ m/z 490.1 (M+1).



154

[0869] 1-Cyclopentyl-N-((5-(3-(7-(pyridin-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine **154**.

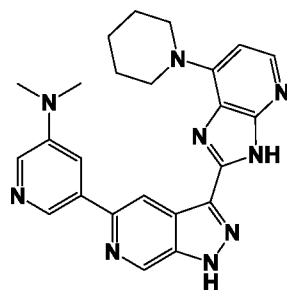
[0870] White solid (29.4 mg, 0.06 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 1.22-1.34 (m, 2H), 1.46-1.65 (m, 4H), 1.76-1.86 (m, 2H), 2.27 (spt, $J=7.64$ Hz, 1H), 2.96-3.04 (m, 2H), 4.42-4.48 (m, 2H), 7.63-7.68 (m, 1H), 8.15 (d, $J=5.14$ Hz, 1H), 8.31 (td, $J=7.78, 1.63$ Hz, 1H), 8.58 (d, $J=5.14$ Hz, 1H), 8.93 (d, $J=4.14$ Hz, 1H), 9.05 (d, $J=1.63$ Hz, 1H), 9.13 (d, $J=1.13$ Hz, 1H), 9.18 (brs, 2H), 9.40 (d, $J=1.26$ Hz, 1H), 9.50 (brs, 2H), 9.53 (d, $J=1.88$ Hz, 1H), 14.89 (brs, 1H); ESIMS found for $C_{29}H_{27}N_9$ m/z 502.2 (M+1).



157

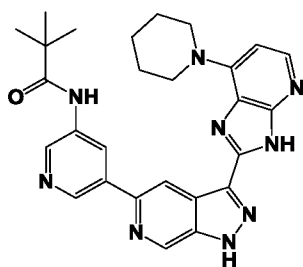
[0871] N-(5-(3-(7-(Piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide **157**.

[0872] White solid (15.9 mg, 0.03 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 1.09-1.19 (m, 3H), 1.65-1.81 (m, 6H), 2.35-2.46 (m, 2H), 3.95-4.16 (m, 4H), 6.48-6.60 (m, 1H), 7.90-8.00 (m, 1H), 8.68 (brs, 1H), 8.86 (brs, 1H), 8.94 (brs, 2H), 9.27 (brs, 1H), 10.27 (brs, 1H), 13.48 (brs, 1H), 14.23 (brs, 1H); ESIMS found for $C_{25}H_{25}N_9O$ m/z 468.1 (M+1).

**163**

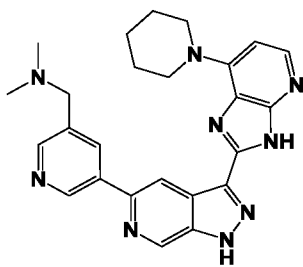
[0873] N,N-Dimethyl-5-(3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine **163**.

[0874] White solid (93.1 mg, 0.21 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.73 (brs, 6H), 3.05 (s, 6H), 4.05 (brs, 4H), 6.54 (d, *J*=5.90 Hz, 1H), 7.69-7.74 (m, 1H), 7.95 (d, *J*=5.90 Hz, 1H), 8.16 (d, *J*=2.76 Hz, 1H), 8.56 (s, 1H), 8.83 (d, *J*=1.13 Hz, 1H), 9.27 (d, *J*=1.26 Hz, 1H), 14.17 (brs, 1H); ESIMS found for C₂₄H₂₅N₉ *m/z* 440.1 (M+1).

**164**

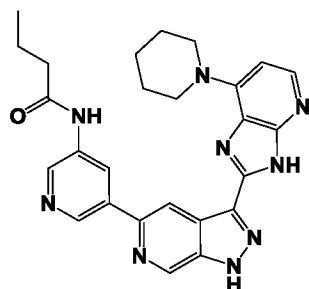
[0875] N-(5-(3-(7-(Piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide **164**.

[0876] White solid (45.6 mg, 0.09 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.29 (s, 9H), 1.73 (d, *J*=4.89 Hz, 6H), 4.06 (brs, 4H), 6.54 (d, *J*=5.52 Hz, 1H), 7.95 (d, *J*=5.27 Hz, 1H), 8.83 (brs, 1H), 8.86 (s, 1H), 8.88 (brs, 1H), 8.98 (s, 1H), 9.27 (s, 1H), 9.62 (s, 1H), 13.45 (brs, 1H), 14.19 (brs, 1H); ESIMS found for C₂₇H₂₉N₉O *m/z* 496.2 (M+1).

**169**

[0877] N,N-Dimethyl-1-(5-(3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine **169**.

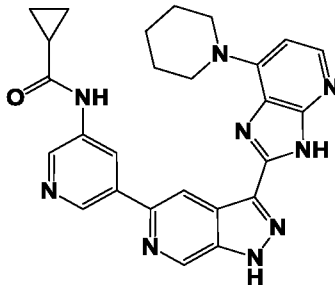
[0878] White solid (11.8 mg, 0.03 mmol). ^1H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.80 (brs, 6H), 2.79 (brs, 3H), 2.80 (brs, 3H), 4.47 (brs, 2H), 6.94-7.01 (m, 1H), 7.99 (brs, 1H), 8.76 (s, 1H), 8.78-8.83 (m, 2H), 9.30 (s, 1H), 9.37 (s, 1H), 10.69 (brs, 1H), 14.72 (brs, 1H); ESIMS found for C₂₅H₂₇N₉ *m/z* 454.2 (M+1).



173

[0879] N-(5-(3-(7-(Piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide **173**.

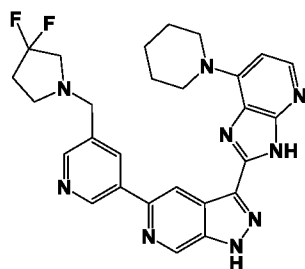
[0880] White solid (104.3 mg, 0.22 mmol). ^1H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.94 (t, *J*=7.34 Hz, 3H), 1.66 (sxt, *J*=7.40Hz, 3H), 1.73 (brs, 6H), 2.48 (t, 2H), 4.25 (brs, 4H), 6.87 (d, *J*=7.40 Hz, 1H), 7.91 (d, *J*=7.28 Hz, 1H), 8.59 (s, 1H), 9.03 (s, 1H), 9.13 (s, 1H), 9.25 (s, 1H), 9.42 (s, 1H), 11.74 (s, 1H); ESIMS found for C₂₆H₂₇N₉O *m/z* 482.2 (M+1).



175

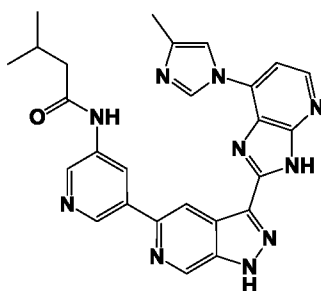
[0881] N-(5-(3-(7-(Piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide **175**.

[0882] White solid (54.5 mg, 0.11 mmol). ^1H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.83-0.93 (m, 4H), 1.66-1.81 (m, 6H), 1.86-1.91 (m, 1H), 4.10 (brs, 4H), 6.56 (d, *J*=6.00Hz, 1H), 7.96 (d, *J*=6.00Hz, 1H), 8.71 (s, 1H), 8.85 (s, 1H), 8.88 (s, 1H), 8.97 (s, 1H), 9.25 (s, 1H), 10.30 (s, 1H), 14.00 (brs, 1H); ESIMS found for C₂₆H₂₅N₉O *m/z* 480.3 (M+1).

**181**

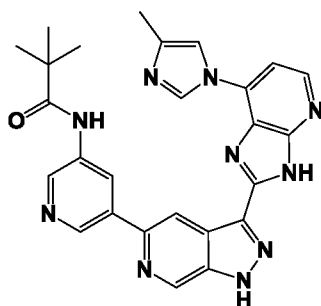
[0883] 5-(5-((3,3-Difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridine **181**.

[0884] White solid (19.5 mg, 0.04 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.79 (brs, 6H), 2.57-2.67 (m, 2H), 3.50-3.64 (m, 4H), 3.96-4.09 (m, 4H), 4.63 (brs, 2H), 6.97 (d, *J*=7.40 Hz, 1H), 8.00 (brs, 1H), 8.81 (s, 1H), 8.99 (s, 1H), 9.02 (s, 1H), 9.34-9.40 (m, 2H), 14.90 (brs, 1H); ESIMS found for C₂₇H₂₇F₂N₉ *m/z* 516.2 (M+1).

**184**

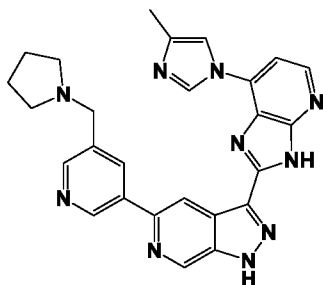
[0885] 3-Methyl-N-(5-(3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide **184**.

[0886] White solid (11.7 mg, 0.02 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.97 (d, *J*=6.65 Hz, 6H), 2.09-2.21 (m, 1H), 2.42 (d, *J*=7.15 Hz, 2H), 7.86 (d, *J*=5.27 Hz, 1H), 8.58 (d, *J*=5.14 Hz, 1H), 8.69 (s, 1H), 9.08 (s, 1H), 9.33-9.41 (m, 3H), 9.47 (s, 1H), 10.33 (s, 1H), 11.70 (s, 1H); ESIMS found for C₂₆H₂₄N₁₀O *m/z* 493.2 (M+1).

**190**

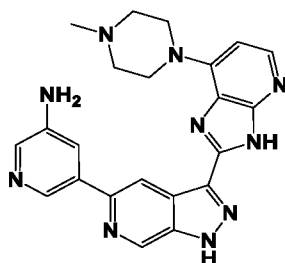
[0887] N-(5-(3-(7-(4-Methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide **190**.

[0888] White solid (22.4 mg, 0.05 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.31 (s, 9H), 7.86 (d, *J*=5.52 Hz, 1H), 8.56 (d, *J*=5.52 Hz, 1H), 8.70 (s, 1H), 9.08 (s, 1H), 9.36-9.40 (m, 2H), 9.45 (d, *J*=1.51 Hz, 1H), 9.73 (s, 1H), 10.31 (s, 1H), 10.83 (s, 1H); ESIMS found for C₂₆H₂₄N₁₀O *m/z* 493.2 (M+1).

**196**

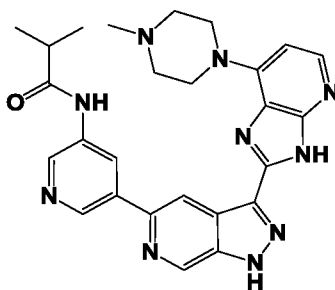
[0889] 3-(7-(4-Methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridine **196**.

[0890] White solid (6.8 mg, 0.01 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.77 (brs, 4H), 2.30 (s, 3H), 2.58-2.69 (m, 2H), 3.81-3.99 (m, 2H), 7.63 (d, *J*=5.40 Hz, 1H), 8.24 (s, 1H), 8.41 (d, *J*=5.40 Hz, 1H), 8.49 (brs, 1H), 8.61 (brs, 1H), 8.97 (s, 1H), 9.08 (s, 1H), 9.25 (s, 1H), 9.34 (d, *J*=1.00 Hz, 1H); ESIMS found for C₂₆H₂₄N₁₀ *m/z* 477.1 (M+1).

**211**

[0891] 5-(3-(7-(4-Methylpiperazin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine **211**.

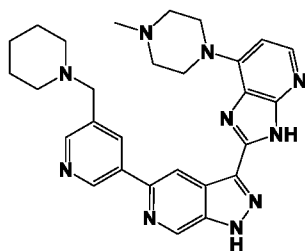
[0892] White solid (25.0 mg, 0.06 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 2.71 (brs, 3H), 3.23 (d, *J*=18.20 Hz, 4H), 4.02-4.47 (m, 4H), 6.62-6.70 (m, 1H), 7.67 (brs, 1H), 7.98 (brs, 1H), 8.05 (d, *J*=4.89 Hz, 1H), 8.45 (brs, 1H), 8.74 (s, 1H), 9.25 (s, 1H), 14.26 (brs, 1H); ESIMS found for C₂₂H₂₂N₁₀ *m/z* 427.1 (M+1).



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[0893] N-(5-(3-(7-(4-Methylpiperazin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide **217**.

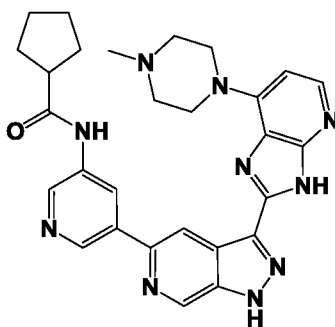
[0894] White solid (56.4 mg, 0.11 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.17 (d, *J*=6.90 Hz, 6H), 2.83-2.94 (m, 3H), 3.44-3.57 (m, 4H), 3.65 (brs, 2H), 5.11 (brs, 2H), 6.71 (d, *J*=5.90 Hz, 1H), 8.09 (d, *J*=5.27 Hz, 1H), 8.75-8.81 (m, 2H), 8.88 (s, 1H), 8.98 (d, *J*=2.01 Hz, 1H), 9.31 (s, 1H), 10.23 (s, 1H), 14.34 (brs, 1H); ESIMS found for C₂₆H₂₈N₁₀O *m/z* 497.2 (M+1).



223

[0895] 3-(7-(4-Methylpiperazin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridine **223**.

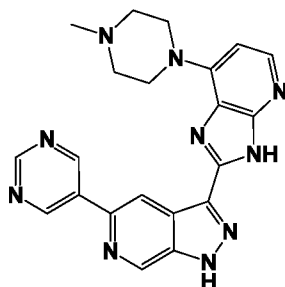
[0896] White solid (96.8 mg, 0.19 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.34-1.45 (m, 2H), 1.52 (d, *J*=4.89 Hz, 4H), 2.30 (s, 3H), 2.36-2.45 (m, 4H), 2.61 (d, *J*=3.76 Hz, 4H), 3.59 (s, 2H), 4.04 (d, *J*=1.76 Hz, 4H), 6.57 (d, *J*=5.77 Hz, 1H), 7.99 (d, *J*=6.15 Hz, 1H), 8.30 (s, 1H), 8.54 (s, 1H), 8.85 (s, 1H), 9.20 (s, 1H), 9.29 (s, 1H); ESIMS found for C₂₈H₃₂N₁₀ *m/z* 509.3 (M+1).



229

[0897] N-(5-(3-(7-(4-Methylpiperazin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide **229**.

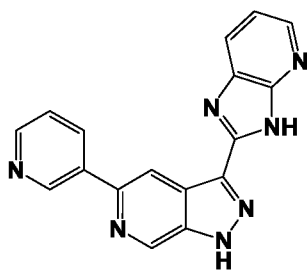
[0898] White solid (22.3 mg, 0.04 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.55-1.65 (m, 3H), 1.66-1.84 (m, 6H), 1.86-1.96 (m, 3H), 2.80-2.91 (m, 1H), 3.32 (s, 3H), 6.62 (d, *J*=3.39 Hz, 1H), 8.02 (d, *J*=5.52 Hz, 1H), 8.74 (s, 1H), 8.79 (s, 1H), 8.91 (brs, 1H), 8.95 (s, 1H), 9.29 (s, 1H), 10.27 (s, 1H), 13.57 (s, 1H), 14.28 (brs, 1H); ESIMS found for C₂₈H₃₀N₁₀O *m/z* 523.2 (M+1).



234

[0899] 3-(7-(4-Methylpiperazin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1H-pyrazolo[3,4-*c*]pyridine **234**.

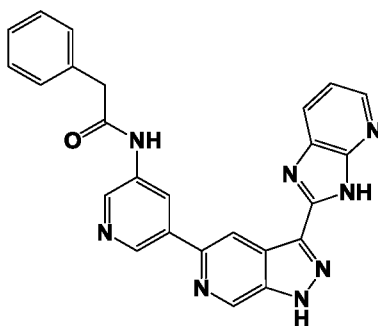
[0900] White solid (26.9 mg, 0.07 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 2.87 (brs, 3H), 3.66 (d, *J*=11.49 Hz, 4H), 3.73-3.86 (m, 2H), 5.28 (brs, 2H), 6.97 (brs, 1H), 8.17 (brs, 1H), 8.84 (s, 1H), 9.27 (s, 1H), 9.40 (s, 1H), 9.50 (s, 2H), 14.71 (brs, 2H); ESIMS found for C₂₁H₂₀N₁₀ *m/z* 413.1 (M+1).



238

[0901] 3-(3H-Imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridine **238**.

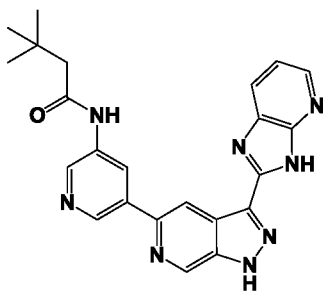
[0902] White solid (57.5 mg, 0.18 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 7.50 (t, *J*=6.53 Hz, 1H), 8.06-8.13 (m, 1H), 8.29 (d, *J*=7.78 Hz, 1H), 8.54 (d, *J*=5.02 Hz, 1H), 8.91 (d, *J*=4.77 Hz, 1H), 9.05 (s, 1H), 9.10-9.16 (m, 1H), 9.42 (d, *J*=1.00 Hz, 1H), 9.56 (s, 1H), 14.93 (brs, 1H); ESIMS found for C₁₇H₁₁N₇ *m/z* 314.0 (M+1).



244

[0903] N-(5-(3-(3H-Imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide **244**.

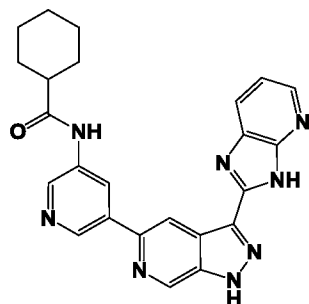
[0904] White solid (30.7 mg, 0.07 mmol). ^1H NMR (CD_3OD , 400 MHz) δ ppm 3.89 (s, 2H), 7.26-7.33 (m, 1H), 7.38 (t, $J=7.40$ Hz, 2H), 7.42-7.46 (m, 2H), 7.85 (dd, $J=8.03$, 6.02 Hz, 1H), 8.61-8.72 (m, 2H), 9.14 (d, $J=1.26$ Hz, 1H), 9.30 (d, $J=2.13$ Hz, 1H), 9.33 (d, $J=1.25$ Hz, 1H), 9.38 (d, $J=1.13$ Hz, 1H), 9.50 (t, $J=1.94$ Hz, 1H); ESIMS found for $\text{C}_{25}\text{H}_{18}\text{N}_8\text{O}$ m/z 447.1 (M+1).



250

[0905] N-(5-(3-(3H-Imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide **250**.

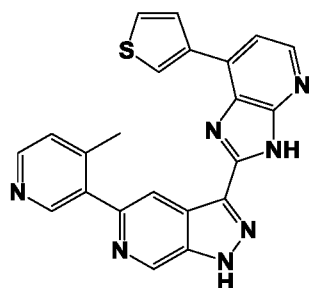
[0906] White solid (7.4 mg, 0.02 mmol). ^1H NMR (CD_3OD , 400 MHz) δ ppm 1.16 (s, 9H), 2.45 (s, 2H), 7.85 (dd, $J=7.97$, 5.96 Hz, 1H), 8.66-8.73 (m, 2H), 9.13 (d, $J=1.13$ Hz, 1H), 9.33 (s, 1H), 9.37 (d, $J=1.13$ Hz, 1H), 9.42 (s, 1H), 9.43-9.46 (m, 1H); ESIMS found for $\text{C}_{23}\text{H}_{22}\text{N}_8\text{O}$ m/z 427.1 (M+1).



256

[0907] N-(5-(3-(3H-Imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide **256**.

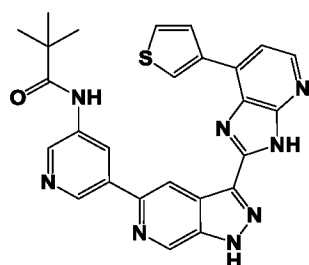
[0908] White solid (24.0 mg, 0.05 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.17-1.39 (m, 3H), 1.40-1.53 (m, 2H), 1.63-1.72 (m, 1H), 1.73-1.83 (m, 2H), 1.90 (d, *J*=12.30 Hz, 2H), 2.36-2.46 (m, 1H), 7.49-7.58 (m, 1H), 8.27-8.36 (m, 1H), 8.56 (d, *J*=4.89 Hz, 1H), 8.96 (s, 1H), 9.19 (d, *J*=1.13 Hz, 2H), 9.23 (s, 1H), 9.40 (d, *J*=1.00 Hz, 1H), 11.04 (brs, 1H), 15.03 (brs, 1H); ESIMS found for C₂₄H₂₂N₈O *m/z* 439.0 (M+1).



265

[0909] 5-(4-Methylpyridin-3-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridine **265**.

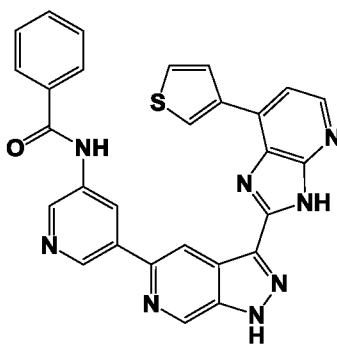
[0910] White solid (31.6 mg, 0.08 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 2.77 (s, 3H), 7.69 (d, *J*=5.27 Hz, 1H), 7.77 (dd, *J*=5.08, 3.07 Hz, 1H), 8.09 (d, *J*=6.15 Hz, 1H), 8.21 (d, *J*=3.89 Hz, 1H), 8.39 (d, *J*=5.27 Hz, 1H), 8.83 (d, *J*=1.00 Hz, 1H), 8.86 (d, *J*=5.90 Hz, 1H), 8.93 (d, *J*=1.63 Hz, 1H), 9.18 (s, 1H), 9.40 (d, *J*=1.26 Hz, 1H), 14.70 (brs, 1H); ESIMS found for C₂₂H₁₅N₇S *m/z* 410.0 (M+1).



268

[0911] N-(5-(3-(7-(Thiophen-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide **268**.

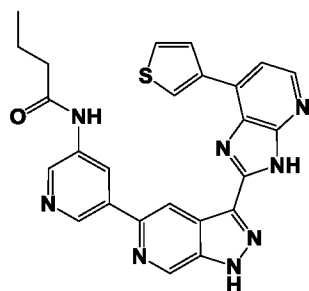
[0912] White solid (53.7 mg, 0.11 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.31 (s, 9H), 7.68 (d, *J*=5.14 Hz, 1H), 7.81 (dd, *J*=5.02, 3.01 Hz, 1H), 8.27 (d, *J*=4.89 Hz, 1H), 8.37 (d, *J*=5.14 Hz, 1H), 8.88-8.92 (m, 2H), 8.95 (brs, 1H), 8.97 (d, *J*=0.88 Hz, 1H), 9.06 (d, *J*=1.63 Hz, 1H), 9.32 (d, *J*=1.00 Hz, 1H), 9.63 (s, 1H); ESIMS found for C₂₆H₂₂N₈OS *m/z* 495.1 (M+1).



271

[0913] N-(5-(3-(7-(Thiophen-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide **271**.

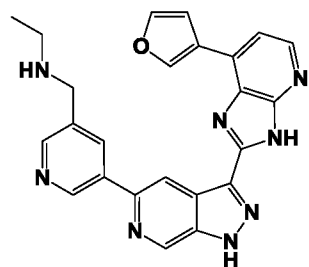
[0914] White solid (29.2 mg, 0.06 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 7.63 (t, *J*=7.16 Hz, 2H), 7.69 (d, *J*=7.15 Hz, 1H), 7.72 (d, *J*=5.27 Hz, 1H), 7.81 (dd, *J*=4.96, 2.95 Hz, 1H), 8.11-8.16 (m, 2H), 8.24 (d, *J*=5.02 Hz, 1H), 8.42 (d, *J*=5.27 Hz, 1H), 9.00 (d, *J*=2.01 Hz, 1H), 9.11 (s, 1H), 9.30 (d, *J*=3.39 Hz, 2H), 9.41 (d, *J*=1.13 Hz, 1H), 9.48 (brs, 1H), 11.16 (brs, 1H), 14.73 (brs, 1H); ESIMS found for C₂₈H₁₈N₈OS *m/z* 515.1 (M+1).



277

[0915] N-(5-(3-(7-(Thiophen-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide **277**.

[0916] White solid (88.9 mg, 0.18 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.95 (t, *J*=7.40 Hz, 3H), 1.68 (sxt, *J*=7.35 Hz, 2H), 2.44-2.50 (m, 2H), 7.87-7.94 (m, 2H), 8.12 (d, *J*=5.27 Hz, 1H), 8.53 (d, *J*=6.02 Hz, 1H), 8.88 (brs, 1H), 8.93 (s, 1H), 9.19 (s, 1H), 9.24 (d, *J*=1.63 Hz, 1H), 9.38 (s, 1H), 9.44 (s, 1H), 11.59 (s, 1H); ESIMS found for C₂₅H₂₀N₈OS *m/z* 481.1 (M+1).

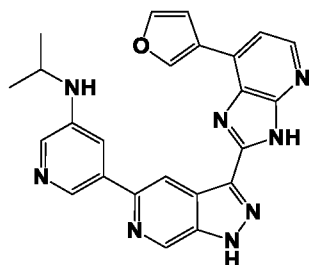


209

292

[0917] N-((5-(3-(7-(Furan-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo [3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine **292**.

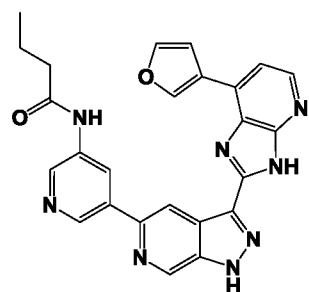
[0918] White solid (46.5 mg, 0.11 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.29 (t, *J*=7.22 Hz, 3H), 3.03-3.14 (m, 2H), 4.41 (t, *J*=5.58 Hz, 2H), 7.48 (d, *J*=1.38 Hz, 1H), 7.63 (d, *J*=5.14 Hz, 1H), 7.96 (t, *J*=1.69 Hz, 1H), 8.41 (d, *J*=5.27 Hz, 1H), 8.96 (d, *J*=1.76 Hz, 1H), 9.07 (s, 1H), 9.08-9.12 (m, 2H), 9.39 (d, *J*=1.25 Hz, 1H), 9.48 (brs, 2H), 9.51 (d, *J*=1.88 Hz, 1H), 14.78 (brs, 1H); ESIMS found for C₂₄H₂₀N₈O *m/z* 437.2 (M+1).



298

[0919] 5-(3-(7-(Furan-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)-N-isopropylpyridin-3-amine **298**.

[0920] White solid (111.0 mg, 0.25 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.23 (d, *J*=6.27 Hz, 6H), 3.68-3.80 (m, 1H), 5.93 (d, *J*=8.03 Hz, 1H), 7.46 (d, *J*=1.13 Hz, 1H), 7.56-7.62 (m, 2H), 7.94 (s, 1H), 8.02 (d, *J*=2.38 Hz, 1H), 8.36 (d, *J*=5.02 Hz, 1H), 8.52 (s, 1H), 8.89 (s, 1H), 9.01 (s, 1H), 9.31 (s, 1H), 13.93 (brs, 1H), 14.42 (brs, 1H); ESIMS found for C₂₄H₂₀N₈O *m/z* 437.1 (M+1).

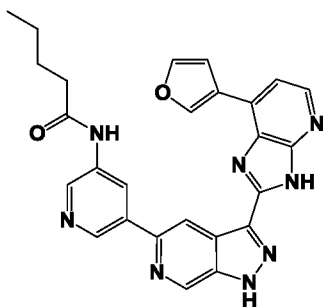


303

[0921] N-(5-(3-(7-(Furan-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo [3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide **303**.

[0922] White solid (84.0 mg, 0.18 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.94 (t, *J*=7.34 Hz, 3H), 1.67 (sxt, *J*=7.40 Hz, 2H), 2.48 (t, *J*=7.20 Hz, 2H), 7.50 (d, *J*=1.00 Hz, 1H), 7.91 (d, *J*=6.15 Hz, 1H), 8.01 (s, 1H), 8.56 (d, *J*=6.27 Hz, 1H), 8.93 (s, 1H), 9.10 (s, 1H), 9.22 (s, 1H),

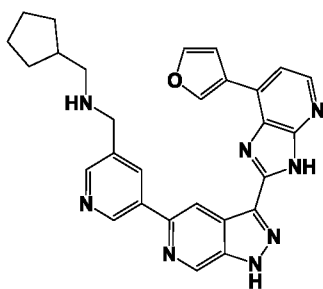
9.27 (d, $J=1.76$ Hz, 1H), 9.40 (s, 1H), 9.43 (s, 1H), 11.66 (s, 1H); ESIMS found for $C_{25}H_{20}N_8O_2$ m/z 465.1 (M+1).



304

[0923] N-(5-(3-(7-(Furan-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo [3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide **304**.

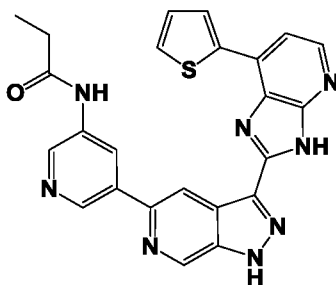
[0924] White solid (15.6 mg, 0.03 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 0.93 (t, $J=7.34$ Hz, 3H), 1.38 (sxt, $J=7.52$ Hz, 2H), 1.65 (quin, $J=7.28$ Hz, 2H), 2.47 (t, $J=7.56$ Hz, 2H), 7.50 (d, $J=1.25$ Hz, 1H), 7.62 (d, $J=5.14$ Hz, 1H), 7.94 (t, $J=1.51$ Hz, 1H), 8.39 (d, $J=5.14$ Hz, 1H), 9.03 (s, 2H), 9.10 (s, 1H), 9.23-9.28 (m, 2H), 9.38 (d, $J=1.00$ Hz, 1H), 10.95 (s, 1H), 14.73 (brs, 1H); ESIMS found for $C_{26}H_{22}N_8O_2$ m/z 479.1 (M+1).



310

[0925] 1-Cyclopentyl-N-((5-(3-(7-(furan-3-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine **310**.

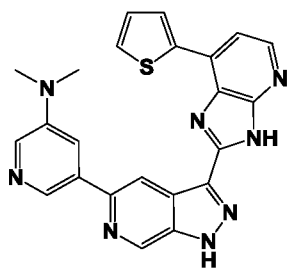
[0926] White solid (36.2 mg, 0.07 mmol). 1H NMR (CD $_3$ OD, 400 MHz) δ ppm 1.32-1.45 (m, 2H), 1.61-1.81 (m, 4H), 1.92-2.04 (m, 2H), 2.30-2.43 (m, 1H), 3.24 (d, $J=7.53$ Hz, 2H), 4.63 (brs, 2H), 7.33 (d, $J=1.13$ Hz, 1H), 7.88-7.95 (m, 2H), 8.59 (d, $J=6.02$ Hz, 1H), 8.88 (s, 1H), 9.01-9.08 (m, 1H), 9.25 (s, 1H), 9.38 (s, 1H), 9.47-9.60 (m, 1H), 9.61-9.68 (m, 1H); ESIMS found for $C_{28}H_{26}N_8O$ m/z 491.2 (M+1).



313

[0927] N-(5-(3-(7-(Thiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide **313**.

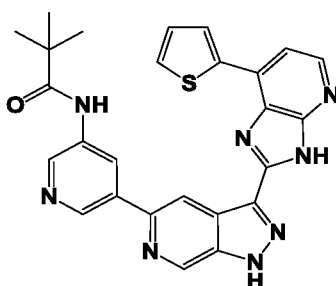
[0928] White solid (6.0 mg, 0.01 mmol). ^1H NMR (DMSO- d_6 , 400 MHz) δ ppm 1.17 (t, $J=7.53$ Hz, 4H), 2.46 (q, $J=7.88$ Hz, 6H), 7.34-7.39 (m, 1H), 7.69 (d, $J=5.14$ Hz, 1H), 7.89 (d, $J=5.14$ Hz, 1H), 8.35 (d, $J=5.02$ Hz, 2H), 8.70 (brs, 1H), 9.04 (brs, 1H), 9.07 (brs, 1H), 9.11 (s, 1H), 9.34 (s, 1H), 10.32 (s, 1H), 14.01 (brs, 1H), 14.48 (s, 1H); ESIMS found for $\text{C}_{24}\text{H}_{18}\text{N}_8\text{OS}$ m/z 467.1 (M+1).



319

[0929] N,N-Dimethyl-5-(3-(7-(thiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine **319**.

[0930] White solid (31.9 mg, 0.07 mmol). ^1H NMR (DMSO- d_6 , 400 MHz) δ ppm 3.22 (s, 6H), 7.34 (dd, $J=5.08, 3.70$ Hz, 1H), 7.69 (d, $J=5.27$ Hz, 1H), 7.91 (dd, $J=5.02, 1.00$ Hz, 1H), 8.30 (d, $J=2.51$ Hz, 1H), 8.33 (s, 1H), 8.36 (d, $J=5.27$ Hz, 1H), 8.39 (dd, $J=3.76, 1.00$ Hz, 1H), 8.71 (s, 1H), 9.21 (d, $J=1.13$ Hz, 1H), 9.37 (d, $J=1.25$ Hz, 1H), 14.08 (brs, 1H), 14.65 (s, 1H); ESIMS found for $\text{C}_{23}\text{H}_{18}\text{N}_8\text{S}$ m/z 439.1 (M+1).

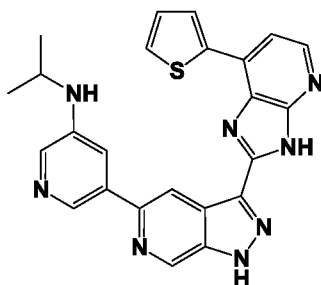


212

320

[0931] N-(5-(3-(7-(Thiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide **320**.

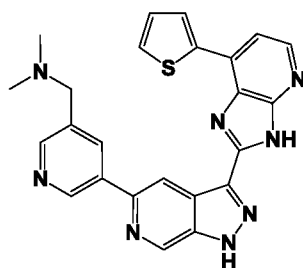
[0932] White solid (43.1 mg, 0.09 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.31 (s, 9H), 7.36 (t, *J*=4.33 Hz, 1H), 7.68 (d, *J*=5.02 Hz, 1H), 7.87 (d, *J*=5.14 Hz, 1H), 8.35 (d, *J*=5.27 Hz, 1H), 8.38 (d, *J*=2.26 Hz, 1H), 8.89 (brs, 1H), 8.94 (br. s., 1H), 9.08 (brs, 1H), 9.11 (s, 1H), 9.34 (s, 1H), 9.65 (s, 1H), 14.02 (brs, 1H), 14.48 (brs, 1H); ESIMS found for C₂₆H₂₂N₈OS *m/z* 495.1 (M+1).



324

[0933] N-Isopropyl-5-(3-(7-(thiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine **324**.

[0934] White solid (82.6 mg, 0.18 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.22 (d, *J*=5.02 Hz, 6H), 3.74 (brs, 2H), 6.30 (brs, 1H), 7.34 (d, *J*=1.76 Hz, 1H), 7.69 (d, *J*=3.39 Hz, 1H), 7.80 (brs, 1H), 7.82-7.89 (m, 1H), 8.06 (brs, 1H), 8.36 (brs, 2H), 8.55 (brs, 1H), 9.08 (brs, 1H), 9.33 (d, *J*=0.75 Hz, 1H), 14.03 (brs, 1H), 14.48 (brs, 1H); ESIMS found for C₂₄H₂₀N₈S *m/z* 453.1 (M+1).

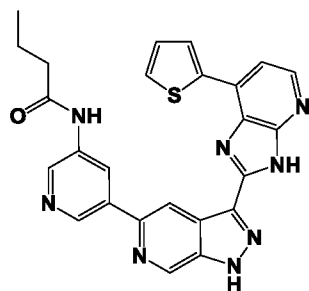


325

[0935] N,N-Dimethyl-1-(5-(3-(7-(thiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine **325**.

[0936] White solid (7.5 mg, 0.02 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ 2.82 (s, 3H), 2.83 (s, 3H), 4.59 (brs, 2H), 7.35-7.40 (m, 1H), 7.70 (d, *J*=5.27 Hz, 1H), 8.00 (d, *J*=4.77 Hz, 1H), 8.37 (d, *J*=5.27 Hz, 1H), 8.40 (d, *J*=3.26 Hz, 1H), 9.07 (s, 1H), 9.10 (s, 1H), 9.19 (s, 1H), 9.38

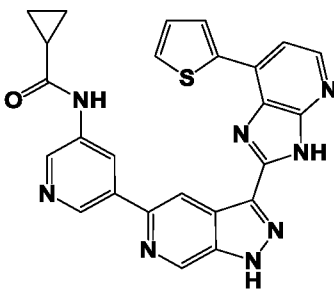
(d, $J=0.88$ Hz, 1H), 9.48 (s, 1H), 11.22 (brs, 1H), 14.82 (brs, 1H); ESIMS found for $C_{24}H_{20}N_8S$ m/z 453.1 (M+1).



329

[0937] N-(5-(3-(7-(Thiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide **7**.

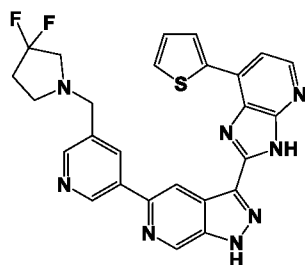
[0938] White solid (87.7 mg, 0.18 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 0.96 (t, $J=7.28$ Hz, 3H), 1.68 (sxt, $J=7.28$ Hz, 2H), 7.35 (t, $J=4.02$ Hz, 1H), 7.80 (d, $J=5.65$ Hz, 1H), 8.16 (d, $J=4.39$ Hz, 1H), 8.32 (brs, 1H), 8.34-8.39 (m, 1H), 8.92 (brs, 1H), 8.98 (brs, 1H), 9.15 (s, 1H), 9.26 (s, 1H), 9.42 (brs, 1H), 11.63 (s, 1H); ESIMS found for $C_{25}H_{20}N_8OS$ m/z 481.1 (M+1).



331

[0939] N-(5-(3-(7-(Thiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide **331**.

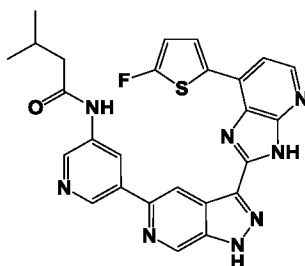
[0940] White solid (65.9 mg, 0.14 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 0.90 (d, $J=6.02$ Hz, 4H), 1.89 (quin, $J=6.80$ Hz, 1H), 7.35 (t, $J=5.52$ Hz, 1H), 7.69 (d, $J=5.27$ Hz, 1H), 7.85 (d, $J=5.02$ Hz, 1H), 8.33-8.40 (m, 2H), 8.72 (s, 1H), 8.99 (s, 1H), 9.07 (d, $J=1.00$ Hz, 1H), 9.11 (s, 1H), 9.34 (d, $J=0.88$ Hz, 1H), 10.64 (s, 1H), 14.01 (s, 1H), 14.48 (s, 1H); ESIMS found for $C_{25}H_{18}N_8OS$ m/z 479.1 (M+1).



337

[0941] 5-(5-((3,3-Difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridine **337**.

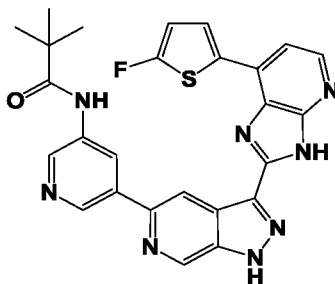
[0942] White solid (44.5 mg, 0.09 mmol). ^1H NMR (DMSO-*d*₆, 400 MHz) δ ppm 2.55-2.67 (m, 2H), 3.61 (brs, 2H), 3.83-3.96 (m, 2H), 4.70 (brs, 2H), 7.38 (dd, $J=4.96$, 3.83 Hz, 1H), 7.71 (d, $J=5.27$ Hz, 1H), 7.98-8.03 (m, 1H), 8.38 (d, $J=5.27$ Hz, 1H), 8.42 (d, $J=2.76$ Hz, 1H), 9.10 (d, $J=1.38$ Hz, 1H), 9.17 (s, 1H), 9.22 (d, $J=0.88$ Hz, 1H), 9.40 (d, $J=1.00$ Hz, 1H), 9.48 (d, $J=1.76$ Hz, 1H), 14.81 (brs, 1H); ESIMS found for C₂₆H₂₀F₂N₈S m/z 515.2 (M+1).



340

[0943] N-(5-(3-(7-(5-Fluorothiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide **340**.

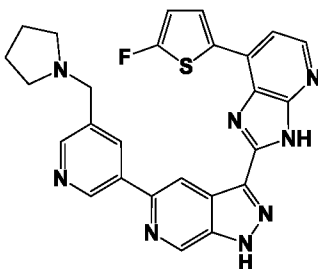
[0944] White solid (4.0 mg, 0.01 mmol). ^1H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.98 (d, $J=6.65$ Hz, 6H), 2.15 (non, $J=6.80$ Hz, 1H), 2.34 (d, $J=7.15$ Hz, 2H), 7.00 (dd, $J=4.27$, 1.88 Hz, 1H), 7.65 (d, $J=5.14$ Hz, 1H), 8.06 (t, $J=3.95$ Hz, 1H), 8.34 (d, $J=5.27$ Hz, 1H), 9.09 (s, 2H), 9.16 (s, 1H), 9.26 (s, 1H), 9.36 (s, 1H), 11.00 (s, 1H), 14.73 (brs, 1H); ESIMS found for C₂₆H₂₁FN₈OS m/z 513.2 (M+1).



346

[0945] N-(5-(3-(7-(5-Fluorothiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide **346**.

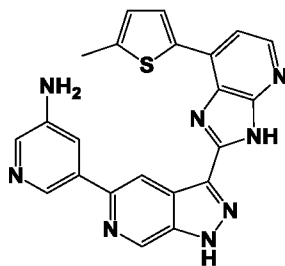
[0946] White solid (9.8 mg, 0.02 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.31 (s, 9H), 6.99 (d, *J*=2.64 Hz, 1H), 7.67 (d, *J*=4.77 Hz, 1H), 8.07 (t, *J*=3.83 Hz, 1H), 8.35 (d, *J*=5.14 Hz, 1H), 9.12 (s, 1H), 9.20 (s, 1H), 9.27 (s, 1H), 9.38 (s, 1H), 9.43 (s, 1H), 10.31 (s, 1H), 14.75 (brs, 1H); ESIMS found for C₂₆H₂₁FN₈OS *m/z* 513.2 (M+1).



352

[0947] 3-(7-(5-Fluorothiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridine **352**.

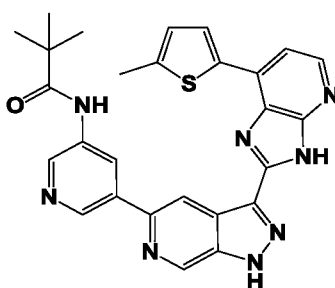
[0948] White solid (9.0 mg, 0.02 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.85-1.94 (m, 2H), 2.00-2.09 (m, 2H), 3.11-3.23 (m, 2H), 3.40-3.52 (m, 2H), 4.58 (d, *J*=4.89 Hz, 2H), 7.00 (d, *J*=4.02 Hz, 1H), 7.73 (d, *J*=5.40 Hz, 1H), 8.02 (t, *J*=4.16 Hz, 1H), 8.36 (d, *J*=5.40 Hz, 1H), 8.88 (s, 1H), 8.93 (s, 1H), 9.18 (s, 1H), 9.38 (s, 1H), 9.44 (s, 1H), 10.85 (brs, 1H), 14.66 (brs, 1H); ESIMS found for C₂₆H₂₁FN₈S *m/z* 497.1 (M+1).



367

[0949] 5-(3-(7-(5-Methylthiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine **367**.

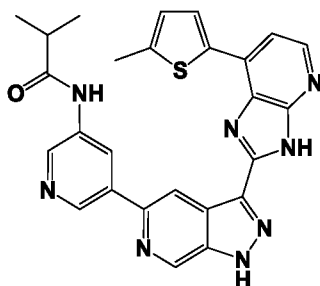
[0950] White solid (9.8 mg, 0.02 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 2.62 (s, 3H), 7.07 (d, *J*=2.51 Hz, 1H), 7.61 (d, *J*=5.40 Hz, 1H), 8.11 (d, *J*=2.01 Hz, 1H), 8.19 (d, *J*=3.51 Hz, 1H), 8.33 (d, *J*=5.27 Hz, 1H), 8.38 (s, 1H), 8.63 (s, 1H), 9.11 (s, 1H), 9.37 (s, 1H), 14.72 (brs, 1H); ESIMS found for C₂₂H₁₆N₈S *m/z* 425.0 (M+1).



372

[0951] N-(5-(3-(7-(5-Methylthiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide **372**.

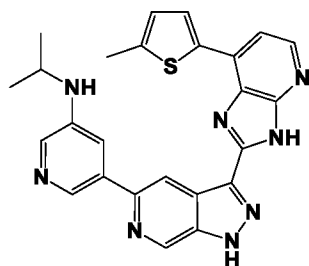
[0952] White solid (43.0 mg, 0.08 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.29 (s, 9H), 2.63 (s, 3H), 7.05 (d, *J*=2.76 Hz, 1H), 7.60 (d, *J*=5.14 Hz, 1H), 8.15 (d, *J*=3.51 Hz, 1H), 8.31 (d, *J*=5.27 Hz, 1H), 8.94 (s, 2H), 9.10 (d, *J*=1.51 Hz, 1H), 9.25 (s, 1H), 9.33 (s, 1H), 9.65 (s, 1H), 14.02 (brs, 1H), 14.41 (brs, 1H); ESIMS found for C₂₇H₂₄N₈OS *m/z* 509.1 (M+1).



373

[0953] N-(5-(3-(7-(5-Methylthiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide **373**.

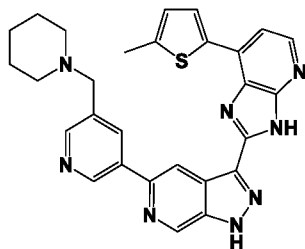
[0954] White solid (20.6 mg, 0.04 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.17 (d, *J*=6.90 Hz, 6H), 2.64 (s, 3H), 2.67-2.71 (m, 1H), 7.07 (d, *J*=2.63 Hz, 1H), 7.62 (d, *J*=5.27 Hz, 1H), 8.15 (d, *J*=3.39 Hz, 1H), 8.32 (d, *J*=5.27 Hz, 1H), 8.83 (d, *J*=2.38 Hz, 1H), 8.96 (t, *J*=2.13 Hz, 1H), 9.10 (d, *J*=1.88 Hz, 1H), 9.25 (d, *J*=1.13 Hz, 1H), 9.34 (d, *J*=1.26 Hz, 1H), 10.27 (s, 1H), 13.99 (s, 1H); ESIMS found for C₂₆H₂₂N₈OS *m/z* 495.1 (M+1).



376

[0955] N-Isopropyl-5-(3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine **376**.

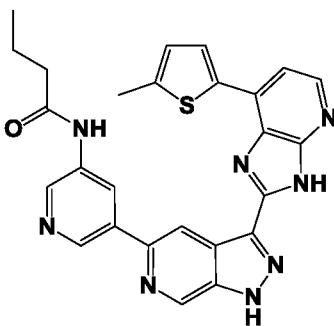
[0956] White solid (41.4 mg, 0.09 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.21 (d, *J*=6.27 Hz, 6H), 2.64 (s, 3H), 3.67-3.76 (m, 1H), 6.23 (brs, 1H), 7.00 (d, *J*=2.76 Hz, 1H), 7.61 (d, *J*=5.27 Hz, 1H), 7.85 (brs, 1H), 8.04 (d, *J*=2.38 Hz, 1H), 8.10 (d, *J*=3.51 Hz, 1H), 8.31 (d, *J*=5.27 Hz, 1H), 8.61 (d, *J*=1.26 Hz, 1H), 9.19 (s, 1H), 9.32 (d, *J*=0.88 Hz, 1H), 13.99 (brs, 1H), 14.43 (brs, 1H); ESIMS found for C₂₅H₂₂N₈S *m/z* 467.1 (M+1).



379

[0957] 3-(7-(5-Methylthiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridine **379**.

[0958] White solid (32.3 mg, 0.06 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.30-1.45 (m, 1H), 1.65-1.86 (m, 5H), 2.88-3.02 (m, 2H), 3.35-3.44 (m, 2H), 4.48 (d, *J*=5.02 Hz, 2H), 7.04 (dd, *J*=3.58, 1.07 Hz, 1H), 7.64 (d, *J*=5.27 Hz, 1H), 8.12 (d, *J*=3.51 Hz, 1H), 8.32 (d, *J*=5.27 Hz, 1H), 8.84 (s, 1H), 8.89 (d, *J*=2.01 Hz, 1H), 9.34 (s, 1H), 9.35 (d, *J*=1.26 Hz, 1H), 9.47 (d, *J*=2.01 Hz, 1H), 10.58 (brs, 1H), 14.02 (brs, 1H), 14.64 (s, 1H); ESIMS found for C₂₈H₂₆N₈S *m/z* 507.1 (M+1).

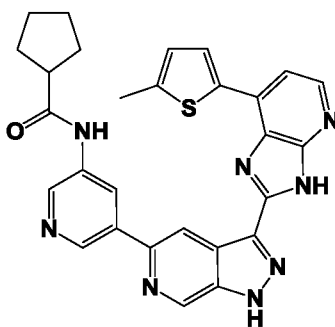


381

[0959] N-(5-(3-(7-(5-Methylthiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide **381**.

[0960] White solid (92.5 mg, 0.19 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.93 (t, *J*=7.40 Hz, 3H), 1.63 (sxt, *J*=7.24 Hz, 2H), 2.46 (t, *J*=7.40 Hz, 4H), 6.98 (d, *J*=2.89 Hz, 1H), 7.56 (d, *J*=5.90 Hz, 1H), 8.08 (d, *J*=3.26 Hz, 1H), 8.23 (d, *J*=6.02 Hz, 1H), 8.54 (brs, 1H), 8.87 (s, 1H),

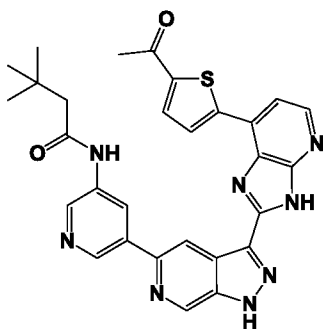
9.12 (s, 1H), 9.15 (s, 1H), 9.18 (brs, 1H), 11.55 (s, 1H); ESIMS found for $C_{26}H_{22}N_8OS$ m/z 495.1 (M+1).



385

[0961] N-(5-(3-(7-(5-Methylthiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide **385**.

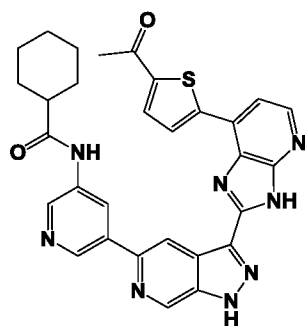
[0962] White solid (24.6 mg, 0.05 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 1.54-1.64 (m, 2H), 1.65-1.84 (m, 4H), 1.86-1.97 (m, 2H), 2.63 (s, 3H), 2.80-2.94 (m, 1H), 7.06 (d, $J=3.14$ Hz, 1H), 7.61 (d, $J=5.02$ Hz, 1H), 8.15 (d, $J=3.26$ Hz, 1H), 8.32 (d, $J=5.27$ Hz, 1H), 8.82 (d, $J=1.76$ Hz, 1H), 8.98 (brs, 1H), 9.09 (s, 1H), 9.24 (s, 1H), 9.34 (s, 1H), 10.33 (s, 1H), 13.99 (brs, 1H); ESIMS found for $C_{28}H_{24}N_8OS$ m/z 521.1 (M+1).



406

[0963] N-(5-(3-(7-(5-Acetylthiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide **406**.

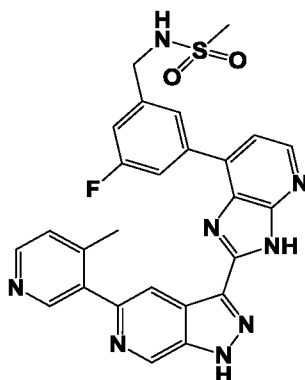
[0964] White solid (8.1 mg, 0.01 mmol). 1H NMR (DMSO- d_6 , 400 MHz) δ ppm 1.05 (s, 9H), 2.28 (s, 2H), 2.62 (s, 3H), 7.79 (d, $J=5.14$ Hz, 1H), 8.09 (d, $J=4.02$ Hz, 1H), 8.41 (d, $J=5.15$ Hz, 1H), 8.45 (d, $J=3.64$ Hz, 1H), 8.82 (d, $J=2.13$ Hz, 1H), 8.88 (s, 1H), 9.02 (s, 1H), 9.05 (d, $J=1.63$ Hz, 1H), 9.36 (s, 1H), 10.17 (s, 1H), 14.16 (brs, 1H), 14.55 (brs, 1H); ESIMS found for $C_{29}H_{26}N_8O_2S$ m/z 551.2 (M+1).



412

[0965] N-(5-(3-(7-(5-Acetylthiophen-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide **412**.

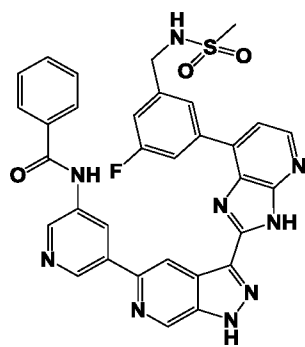
[0966] White solid (28.0 mg, 0.05 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.09-1.33 (m, 3H), 1.34-1.47 (m, 2H), 1.59-1.68 (m, 1H), 1.69-1.80 (m, 2H), 1.81-1.93 (m, 2H), 2.53 (s, 3H), 7.55 (d, *J*=1.88 Hz, 1H), 7.82-7.94 (m, 1H), 8.02-8.14 (m, 1H), 8.19 (brs, 1H), 8.44-8.57 (m, 1H), 8.68-8.80 (m, 1H), 8.96-9.08 (m, 2H), 9.21 (brs, 1H), 11.24 (brs, 1H); ESIMS found for C₃₀H₂₆N₈O₂S *m/z* 563.2 (M+1).



421

[0967] N-(3-Fluoro-5-(2-(5-(4-methylpyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide **421**.

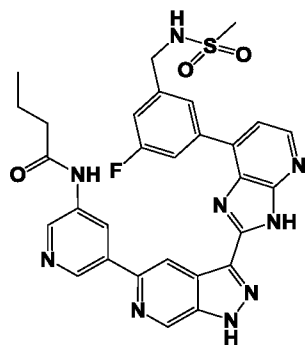
[0968] White solid (16.5 mg, 0.03 mmol). ¹H NMR (CD₃OD, 400 MHz) δ ppm 2.81 (s, 3H), 2.96 (s, 3H), 4.43 (s, 2H), 7.44 (d, *J*=10.29 Hz, 1H), 7.78-7.87 (m, 2H), 7.94 (s, 1H), 8.15 (d, *J*=6.02 Hz, 1H), 8.63 (d, *J*=5.77 Hz, 1H), 8.80 (d, *J*=6.78 Hz, 1H), 8.85 (d, *J*=1.13 Hz, 1H), 9.06 (s, 1H), 9.38 (d, *J*=1.25 Hz, 1H); ESIMS found for C₂₆H₂₁FN₈O₂S *m/z* 529.2 (M+1).



427

[0969] N-(5-(3-(7-(3-Fluoro-5-(methylsulfonamidomethyl)phenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide **427**.

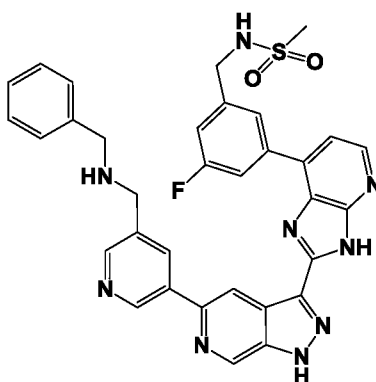
[0970] White solid (47.0 mg, 0.07 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 2.92 (s, 3H), 4.34 (d, *J*=6.27 Hz, 2H), 7.22 (d, *J*=9.16 Hz, 1H), 7.57-7.64 (m, 2H), 7.64-7.71 (m, 2H), 7.76 (t, *J*=6.40 Hz, 1H), 8.03-8.09 (m, 2H), 8.18 (brs, 1H), 8.48 (d, *J*=5.15 Hz, 1H), 8.53-8.69 (m, 1H), 9.02-9.05 (m, 1H), 9.05-9.08 (m, 1H), 9.10 (d, *J*=1.76 Hz, 1H), 9.12 (d, *J*=1.63 Hz, 1H), 9.37 (d, *J*=1.26 Hz, 1H), 10.68 (s, 1H), 14.13 (brs, 1H), 14.54 (s, 1H); ESIMS found for C₃₂H₂₄FN₉O₃S *m/z* 634.1 (M+1).



433

[0971] N-(5-(3-(7-(3-Fluoro-5-(methylsulfonamidomethyl)phenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide **433**.

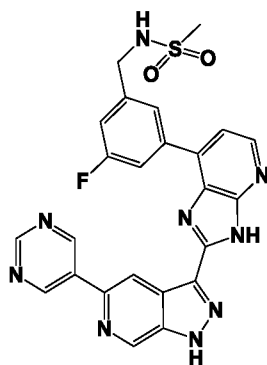
[0972] White solid (21.1 mg, 0.04 mmol). ¹H NMR (CD₃OD, 400 MHz) δ ppm 1.07 (t, *J*=7.40 Hz, 3H), 1.82 (sxt, *J*=7.40 Hz, 2H), 2.55 (t, *J*=7.40 Hz, 2H), 3.00 (s, 3H), 4.47 (s, 2H), 7.45-7.54 (m, 1H), 7.75-7.82 (m, 1H), 7.84-7.88 (m, 1H), 7.88-7.91 (m, 1H), 8.70 (d, *J*=6.15 Hz, 1H), 9.16 (d, *J*=1.25 Hz, 1H), 9.30 (d, *J*=1.13 Hz, 1H), 9.32 (d, *J*=1.88 Hz, 1H), 9.38 (d, *J*=1.38 Hz, 1H), 9.44 (t, *J*=2.01 Hz, 1H); ESIMS found for C₂₉H₂₆FN₉O₃S *m/z* 600.1 (M+1).



439

[0973] N-(3-(2-(5-(5-((Benzylamino)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide **439**.

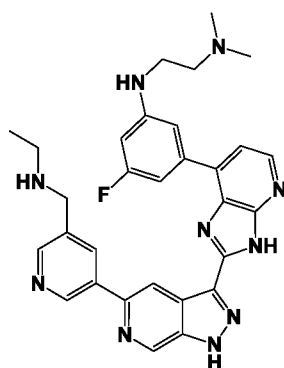
[0974] White solid (8.2 mg, 0.01 mmol). ^1H NMR (DMSO-*d*₆, 400 MHz) δ ppm 2.93 (s, 3H), 4.29 (brs, 2H), 4.37 (d, $J=4.52$ Hz, 2H), 4.50 (brs, 2H), 7.36-7.48 (m, 4H), 7.60 (s, 1H), 7.62 (brs, 1H), 7.71 (d, $J=5.27$ Hz, 1H), 7.84 (brs, 1H), 8.12 (brs, 1H), 8.41 (brs, 1H), 8.52 (d, $J=5.02$ Hz, 1H), 9.05 (s, 1H), 9.07 (brs, 1H), 9.16 (brs, 1H), 9.38 (s, 1H), 9.50 (s, 1H), 10.09 (brs, 2H), 14.95 (brs, 1H); ESIMS found for C₃₃H₂₈FN₉O₂S m/z 634.2 (M+1).



442

[0975] N-(3-Fluoro-5-(2-(5-(pyrimidin-5-yl)-1H-pyrazolo[3,4-*c*]pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide **442**.

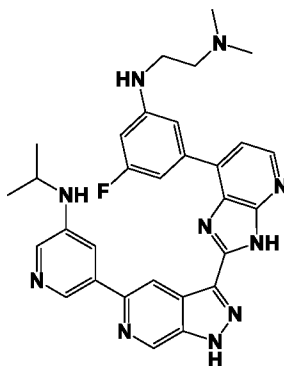
[0976] White solid (101.9 mg, 0.20 mmol). ^1H NMR (DMSO-*d*₆, 400 MHz) δ ppm 2.95 (s, 3H), 4.37 (d, $J=4.27$ Hz, 2H), 7.39 (d, $J=9.91$ Hz, 1H), 7.72 (d, $J=5.14$ Hz, 1H), 7.80 (brs, 1H), 8.13 (brs, 1H), 8.40-8.48 (m, 1H), 8.52 (d, $J=4.89$ Hz, 1H), 9.02 (s, 1H), 9.27 (s, 1H), 9.38 (s, 1H), 9.49 (s, 2H), 14.76 (brs, 1H); ESIMS found for C₂₄H₁₈FN₉O₂S m/z 516.1 (M+1).



448

[0977] N¹-(3-(2-(5-(5-((Ethylamino)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine **448**.

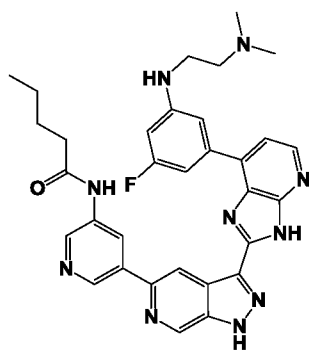
[0978] White solid (11.7 mg, 0.02 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm ¹H NMR (400 MHz, DMSO-*d*₆) δ 1.27 (t, *J*=7.15 Hz, 3H), 2.83 (s, 3H), 2.84 (s, 3H), 3.00-3.11 (m, 2H), 3.27-3.36 (m, 2H), 3.53-3.61 (m, 2H), 4.36 (d, *J*=4.52 Hz, 2H), 6.69 (d, *J*=11.04 Hz, 1H), 7.46 (brs, 1H), 7.62 (d, *J*=5.40 Hz, 2H), 8.46 (d, *J*=5.40 Hz, 1H), 8.87 (s, 1H), 8.89 (d, *J*=1.76 Hz, 1H), 9.07 (s, 1H), 9.38 (s, 3H), 9.43 (d, *J*=1.88 Hz, 1H), 10.44 (brs, 1H), 14.72 (brs, 1H); ESIMS found for C₃₀H₃₁FN₁₀ *m/z* 551.2 (M+1).



454

[0979] N¹-(3-Fluoro-5-(2-(5-(5-(isopropylamino)pyridin-3-yl)-1H-pyrazolo [3,4-*c*]pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine **454**.

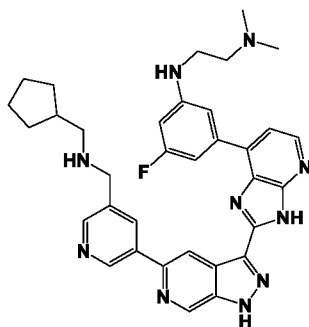
[0980] White solid (25.3 mg, 0.05 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 1.23 (d, *J*=6.27 Hz, 6H), 2.80 (s, 3H), 2.81 (s, 3H), 3.25 (q, *J*=4.39 Hz, 2H), 3.59 (t, *J*=6.34 Hz, 2H), 3.81 (spt, *J*=6.40 Hz, 3H), 6.66 (d, *J*=11.54 Hz, 1H), 7.34 (brs, 1H), 7.51 (brs, 1H), 7.62 (d, *J*=5.40 Hz, 1H), 8.13 (d, *J*=2.26 Hz, 1H), 8.28 (s, 1H), 8.48 (d, *J*=5.27 Hz, 1H), 8.59 (s, 1H), 8.93 (d, *J*=1.00 Hz, 1H), 9.37 (d, *J*=1.13 Hz, 1H), 10.77 (brs, 1H), 14.92 (brs, 1H); ESIMS found for C₃₀H₃₁FN₁₀ *m/z* 551.2 (M+1).



460

[0981] N-(5-(3-(7-(3-((2-(Dimethylamino)ethyl)amino)-5-fluorophenyl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide **460**.

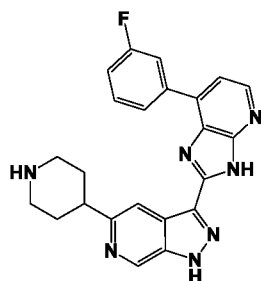
[0982] White solid (17.2 mg, 0.03 mmol). ¹H NMR (DMSO-*d*₆, 400 MHz) δ ppm 0.93 (t, *J*=7.34 Hz, 3H), 1.37 (sxt, *J*=7.43 Hz, 2H), 1.63 (quin, *J*=7.47 Hz, 2H), 2.44 (t, *J*=7.52 Hz, 2H), 2.82 (brs, 3H), 2.83 (brs, 3H), 3.25-3.34 (m, 3H), 6.65 (d, *J*=11.67 Hz, 1H), 7.42-7.50 (m, 1H), 7.50-7.59 (m, 1H), 7.61 (d, *J*=5.14 Hz, 1H), 8.46 (d, *J*=4.89 Hz, 1H), 9.02 (s, 1H), 9.06 (s, 2H), 9.13 (s, 1H), 9.38 (d, *J*=1.13 Hz, 1H), 10.22 (brs, 1H), 10.72 (brs, 1H), 14.69 (brs, 1H); ESIMS found for C₃₂H₃₃FN₁₀O *m/z* 593.3 (M+1).



466

[0983] N¹-(3-(2-(5-(5-(((Cyclopentylmethyl)amino)methyl)pyridin-3-yl)-3H-pyrazolo[3,4-*c*]pyridin-3-yl)-3H-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-N₂,N₂-dimethylethane-1,2-diamine **466**.

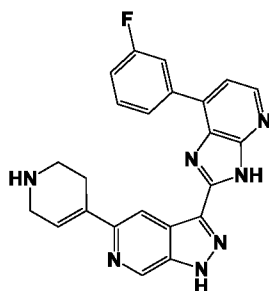
[0984] White solid (10.5 mg, 0.02 mmol). ¹H NMR (CD₃OD, 400 MHz) δ ppm 1.27-1.41 (m, 2H), 1.61-1.80 (m, 4H), 1.91-2.02 (m, 2H), 2.31 (spt, *J*=8.16 Hz, 1H), 2.99 (s, 6H), 3.19 (d, *J*=7.40 Hz, 2H), 3.46 (t, *J*=6.15 Hz, 2H), 3.69 (t, *J*=6.09 Hz, 2H), 4.50 (s, 2H), 6.75 (d, *J*=11.29 Hz, 1H), 7.16 (brs, 1H), 7.20-7.32 (m, 1H), 7.73 (d, *J*=5.90 Hz, 1H), 8.59 (d, *J*=5.52 Hz, 1H), 8.84 (s, 1H), 9.01 (s, 1H), 9.14 (s, 1H), 9.32 (d, *J*=1.13 Hz, 1H), 9.45 (d, *J*=1.76 Hz, 1H); ESIMS found for C₃₄H₃₇FN₁₀ *m/z* 605.3 (M+1).



937

[0985] 3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine **937**.

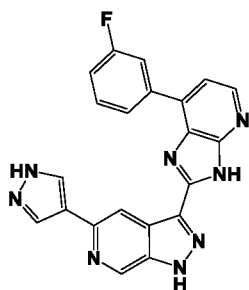
[0986] Off-white solid (35.0 mg, 0.09 mmol). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 1.93 - 2.06 (m, 2 H), 2.25 (br d, $J=12.90$ Hz, 2 H), 3.08 - 3.22 (m, 2 H), 3.45 (br d, $J=12.35$ Hz, 2 H), 7.39 (td, $J=8.44$, 2.33 Hz, 1 H), 7.64 - 7.74 (m, 2 H), 8.24 - 8.38 (m, 2 H), 8.43 (d, $J=4.94$ Hz, 1 H), 8.52 (br s, 1 H), 9.17 (d, $J=1.10$ Hz, 1 H); ESIMS found for $\text{C}_{23}\text{H}_{20}\text{FN}_7$ m/z 414.2 (M+1).



938

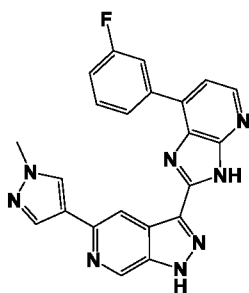
[0987] 3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine **938**.

[0988] Yellow solid (120.0 mg, 0.26 mmol). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 2.85 (br s, 2 H), 3.34 (br d, $J=5.76$ Hz, 2 H), 3.79 (br s, 2 H), 6.92 (br s, 1 H), 7.35 - 7.44 (m, 1 H), 7.62 - 7.70 (m, 1 H), 7.71 (br d, $J=4.12$ Hz, 1 H), 8.26 (br s, 1 H), 8.44 (d, $J=4.67$ Hz, 1 H), 8.58 (s, 1 H), 8.66 (br s, 1 H), 9.18 (s, 1 H); ESIMS found for $\text{C}_{23}\text{H}_{18}\text{FN}_7$ m/z 412.1 (M+1).

**939**

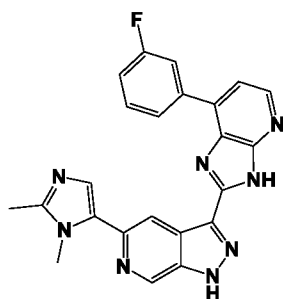
[0989] 3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine **939**.

[0990] Brown solid (16.4 mg, 0.04 mmol). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 7.37 (br s, 1 H), 7.59 (br s, 1 H), 7.65 (d, *J*=6.55 Hz, 1 H), 8.10 (br s, 2 H), 8.25 (br d, *J*=5.21 Hz, 1 H), 8.30 (br s, 1 H), 8.56 (br s, 1 H), 8.82 (br d, *J*=11.80 Hz, 1 H), 9.08 (br s, 1 H); ESIMS found for C₂₁H₁₃FN₈ *m/z* 397.1 (M+1).

**940**

[0991] 3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine **940**.

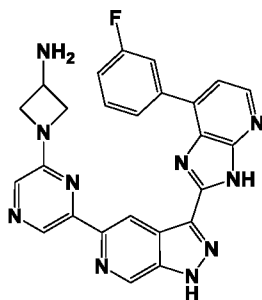
[0992] Yellow solid (9.5 mg, 0.02 mmol). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 3.94 (s, 3 H), 7.38 - 7.47 (m, 1 H), 7.65 - 7.75 (m, 2 H), 7.97 (s, 1 H), 8.21 (s, 1 H), 8.26 (br d, *J*=7.68 Hz, 1 H), 8.43 (d, *J*=5.21 Hz, 1 H), 8.62 (s, 1 H), 8.75 (br d, *J*=10.43 Hz, 1 H), 9.15 (s, 1 H), 13.98 (br s, 1 H), 14.26 (br s, 1 H); ESIMS found for C₂₂H₁₅FN₈ *m/z* 411.1 (M+1).



941

[0993] 5-(1,2-Dimethyl-1H-imidazol-5-yl)-3-(7-(3-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine **941**.

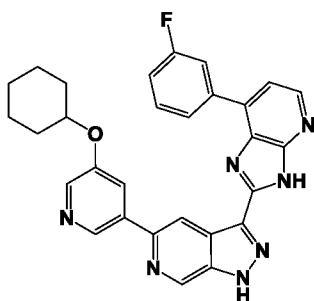
[0994] Brown solid (3.3 mg, 0.008 mmol). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 2.40 (s, 3 H), 3.85 (s, 3 H), 7.23 (s, 1 H), 7.36 - 7.43 (m, 1 H), 7.64 (br dd, $J=14.41$, 8.10 Hz, 1 H), 7.70 (br d, $J=5.21$ Hz, 1 H), 8.27 (br d, $J=8.78$ Hz, 1 H), 8.44 (br d, $J=5.21$ Hz, 1 H), 8.57 (br d, $J=11.25$ Hz, 1 H), 8.64 (s, 1 H), 9.26 (s, 1 H), 14.01 (br s, 1 H), 14.38 (br s, 1 H); ESIMS found for $\text{C}_{23}\text{H}_{17}\text{FN}_8$ m/z 425.1 (M+1).



942

[0995] 1-(6-(3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetidin-3-amine **942**.

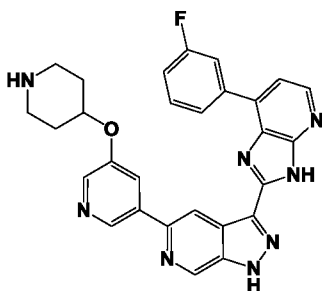
[0996] Brown solid 1 (6.8 mg, 0.014 mmol, 13.85% yield). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 3.67 - 3.77 (m, 2 H), 3.90 - 3.98 (m, 1 H), 4.27 (br t, $J=7.68$ Hz, 2 H), 7.31 - 7.39 (m, 1 H), 7.45 (br d, $J=5.49$ Hz, 1 H), 7.61 - 7.68 (m, 1 H), 7.77 (s, 1 H), 8.18 - 8.26 (m, 2 H), 8.32 (br d, $J=7.96$ Hz, 1 H), 8.84 (s, 1 H), 9.07 (br d, $J=3.02$ Hz, 2 H); ESIMS found for $\text{C}_{25}\text{H}_{19}\text{FN}_{10}$ m/z 479.2 (M+H).



943

[0997] 5-(5-(Cyclohexyloxy)pyridin-3-yl)-3-(7-(3-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine **943**.

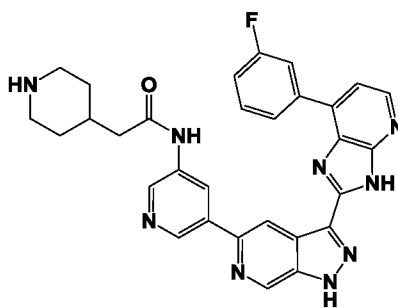
[0998] Brown solid (38 mg, 0.068 mmol, 11.46% yield). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 1.28 - 1.38 (m, 1 H), 1.38 - 1.49 (m, 2 H), 1.50 - 1.61 (m, 3 H), 1.69 - 1.80 (m, 2 H), 1.95 - 2.04 (m, 2 H), 4.54 - 4.63 (m, 1 H), 7.38 (td, *J*=8.58, 2.61 Hz, 1 H), 7.62 - 7.76 (m, 2 H), 7.98 - 8.04 (m, 1 H), 8.32 - 8.41 (m, 2 H), 8.45 (d, *J*=5.21 Hz, 1 H), 8.60 (br d, *J*=9.88 Hz, 1 H), 8.91 (d, *J*=1.65 Hz, 1 H), 9.00 (s, 1 H), 9.33 (d, *J*=1.37 Hz, 1 H), 14.08 (br s, 1 H), 14.42 (br s, 1 H); ESIMS found for C₂₉H₂₄FN₇O *m/z* 506.2 (M+1).



944

[0999] 3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine **944**.

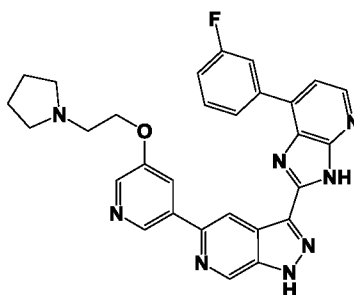
[01000] Amber colored solid (78 mg, 0.154 mmol, 77.0% yield). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 1.82 - 1.93 (m, 2 H), 2.14 (ddd, *J*=10.09, 6.93, 3.29 Hz, 2 H), 3.05 (ddd, *J*=12.62, 8.64, 3.43 Hz, 2 H), 3.22 - 3.30 (m, 3 H), 4.86 (tt, *J*=7.41, 3.57 Hz, 1 H), 7.40 (td, *J*=8.51, 2.20 Hz, 1 H), 7.66 (td, *J*=7.96, 6.59 Hz, 1 H), 7.70 (d, *J*=5.21 Hz, 1 H), 8.03 - 8.10 (m, 1 H), 8.32 (br d, *J*=7.68 Hz, 1 H), 8.43 (d, *J*=2.74 Hz, 1 H), 8.44 (d, *J*=5.21 Hz, 1 H), 8.59 (br d, *J*=10.70 Hz, 1 H), 8.97 (d, *J*=1.65 Hz, 1 H), 9.00 (d, *J*=1.10 Hz, 1 H), 9.32 (d, *J*=1.37 Hz, 1 H); ESIMS found for C₂₈H₂₃FN₈O *m/z* 507.2 (M+1).



945

[01001] N-(5-(3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo [3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide **945**.

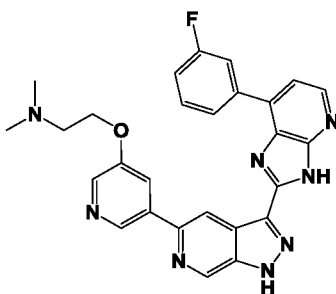
[01002] Brown solid (4.2 mg, 0.009 mmol). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 1.41 - 1.52 (m, 2 H), 1.88 (br d, $J=12.90$ Hz, 2 H), 2.05 - 2.16 (m, 1 H), 2.40 (d, $J=7.14$ Hz, 2 H), 2.84 - 2.97 (m, 2 H), 3.25 (br s, 2 H), 7.39 (br t, $J=7.41$ Hz, 1 H), 7.66 - 7.75 (m, 2 H), 8.37 (br d, $J=7.68$ Hz, 1 H), 8.42 - 8.52 (m, 2 H), 8.59 (br d, $J=9.88$ Hz, 1 H), 8.68 (br d, $J=7.96$ Hz, 1 H), 8.81 (s, 1 H), 8.91 (br s, 1 H), 9.00 (br d, $J=4.12$ Hz, 2 H), 9.33 (d, $J=1.10$ Hz, 1 H), 10.41 (s, 1 H), 14.07 (s, 1 H), 14.52 (s, 1 H); ESIMS found for $\text{C}_{30}\text{H}_{26}\text{FN}_9\text{O}$ m/z 548.2 (M+1).



946

[01003] 3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine **946**.

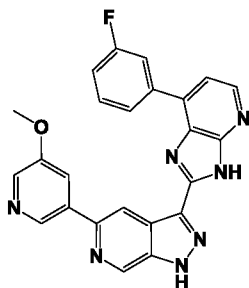
[01004] Yellow solid (7.6 mg, 0.015 mmol, 32.7% yield). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 1.92 (br d, $J=3.29$ Hz, 2 H), 2.07 (br s, 2 H), 3.18 (br d, $J=9.33$ Hz, 2 H), 3.70 (br s, 4 H), 4.55 (br s, 2 H), 7.37 - 7.49 (m, 1 H), 7.63 - 7.71 (m, 1 H), 7.73 (br d, $J=4.12$ Hz, 1 H), 8.16 (br s, 1 H), 8.34 (br d, $J=7.68$ Hz, 1 H), 8.42 - 8.51 (m, 2 H), 8.66 (br d, $J=10.43$ Hz, 1 H), 9.02 (d, $J=1.37$ Hz, 1 H), 9.05 (s, 1 H), 9.35 (s, 1 H), 14.10 (br s, 1 H), 14.54 (br s, 1 H); ESIMS found for $\text{C}_{29}\text{H}_{25}\text{FN}_8\text{O}$ m/z 521.2 (M+1).



947

[01005] 2-((5-(3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)-N,N-dimethylethan-1-amine **947**.

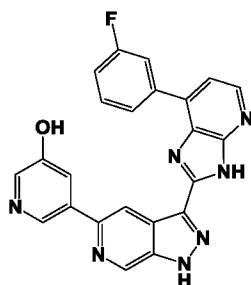
[01006] Brown solid (17.2 mg, 0.035 mmol, 33.0% yield). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 2.28 (s, 6 H), 2.74 (br s, 2 H), 4.27 (br t, *J*=5.35 Hz, 2 H), 7.38 (br t, *J*=8.23 Hz, 1 H), 7.62 - 7.74 (m, 2 H), 8.03 (br s, 1 H), 8.36 (br s, 2 H), 8.45 (br d, *J*=4.94 Hz, 1 H), 8.58 (br s, 1 H), 8.95 (s, 1 H), 9.00 (br s, 1 H), 9.33 (s, 1 H), 14.05 (br s, 2 H); ESIMS found for C₂₇H₂₃FN₈O *m/z* 495.2 (M+1).



948

[01007] 3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine **948**.

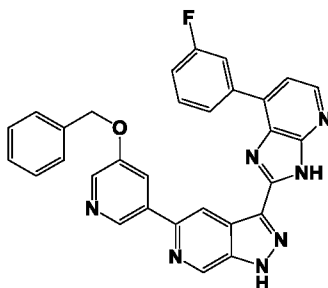
[01008] Orange solid (46.2 mg, 0.11 mmol). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 3.95 (s, 3 H), 7.31 - 7.38 (m, 1 H), 7.54 (d, *J*=5.49 Hz, 1 H), 7.59 - 7.67 (m, 1 H), 8.03 (br s, 1 H), 8.23 - 8.28 (m, 2 H), 8.31 (br d, *J*=7.96 Hz, 1 H), 8.71 (br d, *J*=5.21 Hz, 1 H), 8.84 (s, 1 H), 8.92 (s, 1 H), 9.15 (s, 1 H); ESIMS found for C₂₄H₁₆FN₇O *m/z* 438.1 (M+1).



949

[01009] 5-(3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol **949**.

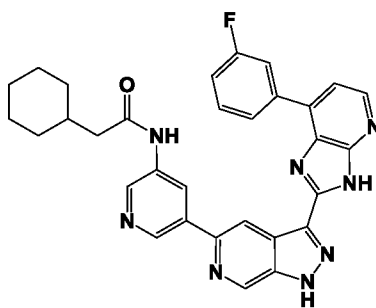
[01010] Brown solid (6.3 mg, 0.015 mmol, 11.11% yield). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 7.37 - 7.45 (m, 1 H), 7.64 - 7.71 (m, 1 H), 7.72 (br d, *J*=4.94 Hz, 1 H), 7.89 - 7.94 (m, 1 H), 8.20 (d, *J*=2.74 Hz, 1 H), 8.37 (br d, *J*=8.23 Hz, 1 H), 8.45 (d, *J*=5.21 Hz, 1 H), 8.60 (br d, *J*=10.70 Hz, 1 H), 8.81 (d, *J*=1.92 Hz, 1 H), 8.98 (s, 1 H), 9.32 (d, *J*=1.10 Hz, 1 H), 10.10 (s, 1 H), 14.07 (br s, 1 H), 14.46 (br s, 1 H); ESIMS found for C₂₃H₁₄FN₇O *m/z* 424.1 (M+1).



950

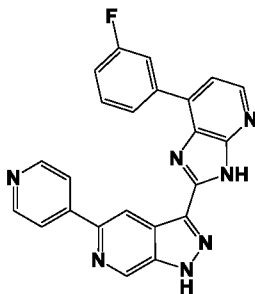
[01011] 5-(5-(Benzyloxy)pyridin-3-yl)-3-(7-(3-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine **950**.

[01012] Off-white solid (35 mg, 0.065 mmol, 32.0% yield). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 5.32 (s, 2 H), 7.19 - 7.28 (m, 1 H), 7.36 - 7.41 (m, 1 H), 7.45 (t, *J*=7.41 Hz, 2 H), 7.55 (d, *J*=7.41 Hz, 2 H), 7.61 - 7.68 (m, 1 H), 7.70 (br d, *J*=4.12 Hz, 1 H), 8.11 - 8.16 (m, 1 H), 8.35 (br s, 1 H), 8.45 (d, *J*=3.02 Hz, 2 H), 8.57 (br d, *J*=9.06 Hz, 1 H), 8.98 (d, *J*=1.37 Hz, 1 H), 9.03 (s, 1 H), 9.33 (d, *J*=1.10 Hz, 1 H), 14.08 (br s, 1 H), 14.42 (br s, 1 H); ESIMS found for C₃₀H₂₀FN₇O *m/z* 514.2 (M+1).

**951**

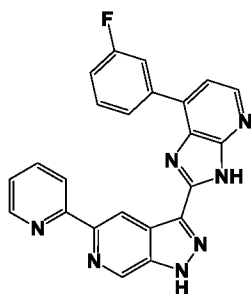
[01013] 2-Cyclohexyl-N-(5-(3-(7-(3-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide **951**.

[01014] Brown solid (50 mg, 0.09 mmol, 51.4% yield). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 0.95 - 1.07 (m, 2 H), 1.11 - 1.32 (m, 3 H), 1.63 (br d, $J=12.08$ Hz, 1 H), 1.68 (br d, $J=12.90$ Hz, 2 H), 1.74 (br d, $J=11.80$ Hz, 2 H), 1.82 (td, $J=7.14, 3.57$ Hz, 1 H), 2.28 (d, $J=7.14$ Hz, 2 H), 7.31 - 7.40 (m, 1 H), 7.65 - 7.76 (m, 2 H), 8.38 (br d, $J=7.96$ Hz, 1 H), 8.44 (d, $J=5.21$ Hz, 1 H), 8.57 (br d, $J=11.25$ Hz, 1 H), 8.78 (s, 1 H), 8.91 (br s, 1 H), 8.98 (br s, 2 H), 9.33 (s, 1 H), 10.23 (s, 1 H), 14.07 (br s, 1 H), 14.46 (br s, 1 H); ESIMS found for $\text{C}_{31}\text{H}_{27}\text{FN}_8\text{O}$ m/z 547.3 (M+1).

**952**

[01015] 3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine **952**.

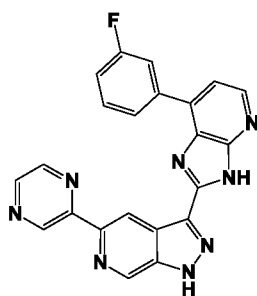
[01016] Yellow solid (5.0 mg, 0.01 mmol). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 7.37 - 7.45 (m, 1 H), 7.68 - 7.78 (m, 2 H), 8.31 (br d, $J=7.96$ Hz, 1 H), 8.49 (d, $J=5.21$ Hz, 1 H), 8.57 (br s, 1 H), 8.70 (d, $J=6.86$ Hz, 2 H), 9.03 (d, $J=6.31$ Hz, 2 H), 9.36 (s, 1 H), 9.44 (s, 1 H), 14.83 (br s, 1 H); ESIMS found for $\text{C}_{23}\text{H}_{14}\text{FN}_7$ m/z 408.1 (M+1).



953

[01017] 3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine **953**.

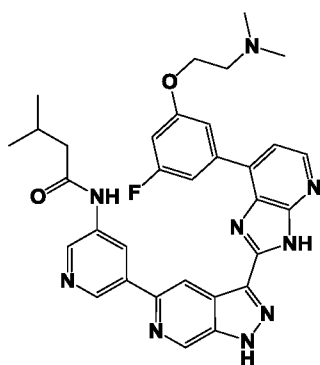
[01018] Brown solid (40 mg, 0.098 mmol, 50.8% yield). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 7.44 (br dd, $J=11.94$, 7.00 Hz, 2 H), 7.72 (br s, 2 H), 7.97 (br t, $J=7.55$ Hz, 1 H), 8.40 (br d, $J=7.41$ Hz, 1 H), 8.43 - 8.50 (m, 2 H), 8.59 (br d, $J=9.06$ Hz, 1 H), 8.75 (br s, 1 H), 9.28 (s, 1 H), 9.65 (br s, 1 H), 14.07 (br s, 1 H), 14.43 (br s, 1 H); ESIMS found for $\text{C}_{23}\text{H}_{14}\text{FN}_7$ m/z 408.1 (M+1).



954

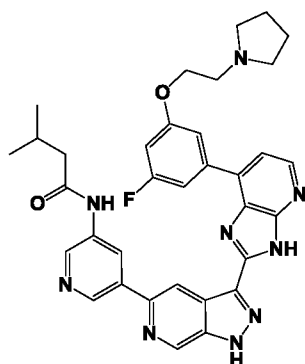
[01019] 3-(7-(3-Fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine **954**.

[01020] Brown solid (32 mg, 0.074 mmol, 37.4% yield). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 7.41 (td, $J=8.37$, 2.20 Hz, 1 H), 7.65 - 7.76 (m, 2 H), 8.38 (br s, 1 H), 8.45 (d, $J=4.94$ Hz, 1 H), 8.57 (br d, $J=7.96$ Hz, 1 H), 8.71 (d, $J=2.47$ Hz, 1 H), 8.79 (s, 1 H), 9.33 (d, $J=1.10$ Hz, 1 H), 9.61 (s, 1 H), 9.62 (d, $J=1.37$ Hz, 1 H), 14.10 (br s, 1 H), 14.49 (br s, 1 H); ESIMS found for $\text{C}_{22}\text{H}_{13}\text{FN}_8$ m/z 409.1 (M+1).

**1262**

[01021] N-(5-(3-(7-(3-(2-(Dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide **1262**.

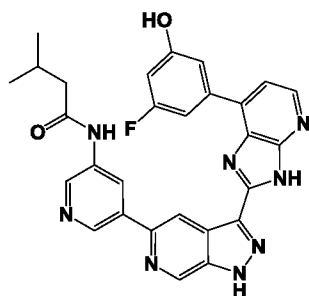
[01022] Brown solid (25.8 mg, 0.043 mmol, 36.1% yield). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 0.97 (d, $J=6.59$ Hz, 6 H), 2.07 - 2.19 (m, 7 H), 2.27 (d, $J=7.14$ Hz, 2 H), 2.54 - 2.60 (m, 2 H), 4.19 (t, $J=5.90$ Hz, 2 H), 7.00 (dt, $J=10.70, 2.20$ Hz, 1 H), 7.73 (br s, 1 H), 8.04 (br s, 2 H), 8.43 (d, $J=4.94$ Hz, 1 H), 8.80 - 8.84 (m, 1 H), 8.85 (br s, 1 H), 8.95 (s, 2 H), 9.34 (d, $J=1.10$ Hz, 1 H), 10.22 (s, 1 H), 14.09 (br s, 1 H), 14.35 (br s, 1 H); ESIMS found for $\text{C}_{32}\text{H}_{32}\text{FN}_9\text{O}_2$ m/z 594.3 (M+1).

**1306**

[01023] N-(5-(3-(7-(3-Fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide **1306**.

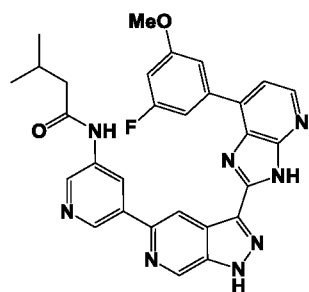
[01024] Yellow solid (2.8 mg, 4.52 μmol , 17.67% yield). ^1H NMR (DMSO- d_6 , 500 MHz) δ ppm 0.97 (d, $J=6.59$ Hz, 6 H), 1.61 (br s, 4 H), 2.08 - 2.17 (m, 1 H), 2.27 (d, $J=7.14$ Hz, 2 H), 2.45 (br s, 4 H), 2.73 (br s, 2 H), 4.21 (br t, $J=5.63$ Hz, 2 H), 7.00 (br d,

$J=10.70$ Hz, 1 H), 7.74 (br s, 1 H), 8.06 (br s, 1 H), 8.43 (br d, $J=5.21$ Hz, 1 H), 8.84 (br d, $J=10.98$ Hz, 2 H), 8.87 - 9.02 (m, 3 H), 9.34 (s, 1 H), 10.22 (s, 1 H), 14.07 (br s, 1 H), 14.46 (br s, 1 H); ESIMS found for $C_{34}H_{34}FN_9O_2$ m/z 620.3 (M+1).

**1350**

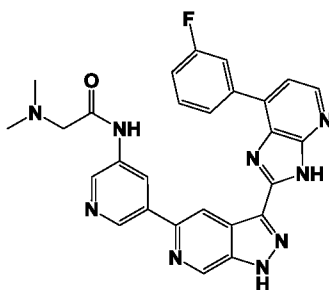
[01025] N-(5-(3-(7-(3-Fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide **1350**.

[01026] Yellow solid (4.1 mg, 7.85 μ mol, 12.53% yield). 1H NMR (DMSO- d_6 , 500 MHz) δ ppm 0.98 (d, $J=6.59$ Hz, 6 H), 2.07 - 2.19 (m, 1 H), 2.28 (d, $J=7.14$ Hz, 2 H), 6.74 (br d, $J=10.43$ Hz, 1 H), 7.59 (br s, 1 H), 7.77 - 8.00 (m, 2 H), 8.42 (d, $J=4.94$ Hz, 1 H), 8.84 (d, $J=2.20$ Hz, 1 H), 8.86 (s, 1 H), 9.00 (s, 1 H), 9.03 (d, $J=1.65$ Hz, 1 H), 9.33 (s, 1 H), 10.20 (s, 1 H), 14.28 (brs, 2 H); ESIMS found for $C_{28}H_{23}FN_8O_2$ m/z 523.2 (M+1).

**1394**

[01027] N-(5-(3-(7-(3-Fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide **1394**.

[01028] Yellow solid (7.7 mg, 0.014 mmol, 16.19% yield). 1H NMR (DMSO- d_6 , 500 MHz) δ ppm 0.97 (br s, 6 H), 2.13 (br s, 1 H), 2.27 (br s, 2 H), 3.94 (br s, 3 H), 6.97 (br d, $J=9.61$ Hz, 1 H), 7.65 (br s, 1 H), 7.98 (br s, 1 H), 8.24 (br s, 1 H), 8.33 (br s, 1 H), 8.83 (br s, 2 H), 8.90 (br s, 1 H), 8.99 (br s, 1 H), 9.24 (br s, 1 H), 10.23 (br s, 1 H); ESIMS found for $C_{29}H_{25}FN_8O_2$ m/z 537.2 (M+1).

**1437**

[01029] 2-(Dimethylamino)-N-(5-(3-(7-(3-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide **1437**.

[01030] Brown solid (20.0 mg, 0.04 mmol). ¹H NMR (DMSO-*d*₆, 500 MHz) δ ppm 2.82 (br s, 6 H), 4.03 (br s, 2 H), 7.37 - 7.46 (m, 1 H), 7.64 - 7.71 (m, 1 H), 7.73 (d, *J*=5.21 Hz, 1 H), 8.35 (br d, *J*=7.68 Hz, 1 H), 8.46 (d, *J*=5.21 Hz, 1 H), 8.65 (br d, *J*=10.43 Hz, 1 H), 8.85 (br d, *J*=2.20 Hz, 1 H), 8.88 (s, 1 H), 9.03 (s, 1 H), 9.07 (s, 1 H), 9.35 (d, *J*=1.37 Hz, 1 H), 10.81 (br s, 1 H), 14.08 (br s, 1 H), 14.51 (br s, 1 H); ESIMS found for C₂₇H₂₂FN₉O *m/z* 508.2 (M+1).

Example 2.

[01031] The screening assay for Wnt activity is described as follows. Reporter cell lines can be generated by stably transducing cancer cell lines (e.g., colon cancer) or primary cells (e.g., IEC-6 intestinal cells) with a lentiviral construct that includes a Wnt-responsive promoter driving expression of the firefly luciferase gene.

[01032] SW480 colon carcinoma cells were transduced with a lentiviral vector expressing luciferase with a human Sp5 promoter consisting of a sequence of eight TCF/LEF binding sites. SW480 cells stably expressing the Sp5-Luc reporter gene and a hygromycin resistance gene were selected by treatment with 150 μg/ml of hygromycin for 7 days. These stably transduced SW480 cells were expanded in cell culture and used for all further screening activities. For Sp5-Luc reporter gene assays, the cells were plated at 10,000 cells/well in 96-well plates with growth medium containing 10% fetal calf serum and incubated overnight at 37°C and 5% CO₂. Each compound was dissolved in DMSO as a 10 mM stock in standard j-vials and used to prepare compound source plates in dose-response format with 3-fold serial dilutions and a 10 mM top concentration. Compound transfer from serially diluted source plates to assay plates containing the cells was accomplished using a pintoole (Multimek 96, Beckman equipped with V&P Scientific FP1S50H pins) based liquid handling protocol. This protocol used a slotted pin to transfer 50 nl of

compound from a source plate well to an assay plate well containing 50 μ l of cells in growth medium. The 1000-fold dilution resulted in a final DMSO concentration of 0.1% on the cells in each well. Control wells received 50 nl of DMSO treatment for normalization and calculating IC₅₀ values. The treated cells were incubated at 37°C and 5% CO₂ for an additional forty-two hours. Following incubation, the growth medium was removed and 50 μ l of BrightGlo luminescence reagent (Promega) was added to each well of the 96-well assay plates. The plates were placed on an orbital shaker for 5 min and then luminescence was quantified on the Victor3 (Perkin Elmer) plate reader. Readings were normalized to DMSO only treated cells, and normalized activities were utilized for IC₅₀ calculations using the dose-response log (inhibitor) vs. response –variable slope (four parameters) nonlinear regression feature available in GraphPad Prism 5.0 or 6.0. Table 2 shows the measured activity for selected compounds of Formula I as described herein.

Table 2.

Compound	IC ₅₀ (μ M)	Compound	IC ₅₀ (μ M)	Compound	IC ₅₀ (μ M)
7	0.0854	196	>10	381	0.0095
8	0.1547	211	1.0440	385	0.0844
13	0.2897	217	>10	406	>10
17	0.0597	223	5.0810	412	>10
19	0.0465	229	>10	421	0.0233
25	0.0221	234	>10	427	>10
28	0.0597	238	0.0161	433	>10
34	0.0167	244	>10	439	>10
38	0.0103	250	>10	442	>10
40	0.3209	256	0.7957	448	>10
43	0.0212	265	0.0261	454	0.0947
46	0.0508	268	0.0097	460	0.1394
52	0.0653	271	0.0166	466	1.7390
55	0.0268	277	0.0249	937	1.856
60	0.0455	283	0.0747	938	>1.0
61	0.0728	292	0.4924	939	0.198
68	0.0244	298	0.0468	940	0.405
69	0.1275	303	0.0788	941	>10
73	0.0369	304	0.0607	942	1.491
82	>10	310	0.2431	943	0.950
87	>10	313	0.0717	944	>10
90	>10	319	0.1246	945	0.463
93	>10	320	0.0426	946	>10
101	>10	324	0.0585	947	1.600
109	0.8644	325	0.1367	948	0.850
136	0.7450	329	0.0175	949	0.111
141	>10	331	0.0831	950	4.430
148	>10	337	0.0248	951	>10
154	0.4639	340	0.0270	952	>1.0
157	>10	346	0.0621	953	0.276
163	1.0320	352	0.1558	954	0.661
164	0.0844	367	0.0226	1262	>10

169	0.2526	372	0.0248	1306	>10
173	0.1692	373	0.0164	1350	0.253
175	0.1554	376	0.0106	1394	>10
181	0.0167	379	0.1803	1437	0.646

Example 3.

[01033] The above synthesized compounds were screened using primary human mesenchymal stem cells (hMSCs) to determine their ability to induce chondrogenesis (process by which cartilage is developed).

[01034] *Human Mesenchymal Stem Cell Culture:* Primary human mesenchymal stem cells (hMSCs) were purchased from Lonza (Walkersville, MD) and expanded in Mesenchymal Stem Cell Growth Media (Lonza). Cells between passage 3 and 6 were used for the experiments.

[01035] *Compound Screening:* Each compound was dissolved in DMSO as a 10 mM stock and used to prepare compound source plates. Serial dilution (1:3, 6-point dose-response curves from 2700 nM to 10 nM) and compound transfer was performed using the ECHO 550 (Labcyte, Sunnyvale, CA) into 96-well clear bottom assay plates (Greiner Bio-One) with appropriate DMSO backfill for a final DMSO concentration of 0.03%. hMSCs were plated at 20,000 cells/well in 250 μ L/well Incomplete Chondrogenic Induction Medium (Lonza; DMEM, dexamethasone, ascorbate, insulin-transferrin-selenium [ITS supplement], gentamycin-amphotericin [GA-1000], sodium pyruvate, proline and L-glutamine). TGF- β 3 (10 ng/mL) was used as a positive control for differentiation while negative control wells were treated with 75 nL DMSO for normalization and calculating EC₅₀ values. Cells were incubated at 37°C and 5% CO₂ for 6 days. To image chondrogenic nodules, the cells were fixed using 4% formaldehyde (Electron Microscopy Sciences), and stained with 2 μ g/mL Rhodamine B (Sigma-Aldrich) and 20 μ M Nile Red (Sigma-Aldrich) [Johnson K., et.al, A Stem Cell-Based Approach to Cartilage Repair, *Science*, (2012), 336(6082), 717-721]. The nodules imaged (4 images per well at 4X magnification) by excitation at 531 nm and emission at 625 nm and quantified using the CellInsight CX5 (Thermo Scientific). Number of nodules in each well was normalized to the average of 3 DMSO treated wells on the same plate using Excel (Microsoft Inc.). The normalized averages (fold change over DMSO) of 3 replicate wells for each compound concentration were calculated. Due to solubility limitations of some of the compounds, curve fitting was incomplete leading to inaccurate EC₅₀ determinations.

[01036] Using TGF- β 3 as a positive control, the concentration of test compounds required to induce equivalent levels of chondrogenesis is reported. In addition, the maximum activity of each compound and the respective dose that each compound reached maximum

chondrogenesis activity is reported. Table 3 shows the activity of selected compounds as provided herein.

Table 3.

Compound	Conc (nM) of Max. activity	Max. Activity as % TGF- β 3 activity	Conc (nM) of 100% TGF- β 3 activity	Compound	Conc (nM) of Max. activity	Max. Activity as % TGF- β 3 activity	Conc (nM) of 100% TGF- β 3 activity
25	2700	N/A	111.9	329	300	300	109.8
38	300	300	74.9	381	900	N/A	59.6
68	300	N/A	9.3	421	30	30	70.0
181	900	N/A	25.6	945	10	N/A	54.7
238	100	30	113.0	949	100	100	142.9
268	900	N/A	25.8				

Example 4.

[01037] The above synthesized compounds were screened using primary human fibroblasts (derived from IPF patients) treated with TGF- β 1 to determine their ability to inhibit the fibrotic process.

[01038] *Human Fibroblast Cell Culture:* Primary human fibroblasts derived from IPF patients (LL29 cells) [Xiaoqiu Liu, et.al., “Fibrotic Lung Fibroblasts Show Blunted Inhibition by cAMP Due to Deficient cAMP Response Element-Binding Protein Phosphorylation”, *Journal of Pharmacology and Experimental Therapeutics* (2005), 315(2), 678-687; Watts, K. L., et.al., “RhoA signaling modulates cyclin D1 expression in human lung fibroblasts; implications for idiopathic pulmonary fibrosis”, *Respiratory Research* (2006), 7(1), 88] were obtained from American Type Culture Collection (ATCC) and expanded in F12 medium supplemented with 15% Fetal Bovine Serum and Penicillin/Streptomycin.

[01039] *Compound Screening:* Each compound was dissolved in DMSO as a 10 mM stock and used to prepare compound source plates. Serial dilution (1:2, 11-point dose-response curves from 10 μ M to 1.87 nM) and compound transfer was performed using the ECHO 550 (Labcyte, Sunnyvale, CA) into 384-well clear bottom assay plates (Greiner Bio-One) with appropriate DMSO backfill for a final DMSO concentration of 0.1%. LL29 cells were plated at 1,500 cells/well in 80 μ l/well F12 medium supplemented with 1% Fetal Bovine Serum. One hour after addition of the cells, TGF- β 1 (Peprotech; 20 ng/mL) was added to the plates to induce fibrosis (ref. 1 and 2 above). Wells treated with TGF- β 1 and containing DMSO were used as controls. Cells were incubated at 37°C and 5% CO₂ for 4 days. Following incubation for 4 days, SYTOX green nucleic acid stain (Life Technologies [Thermo Fisher Scientific]) was added to the wells at a final

concentration of 1 μ M and incubated at room temperature for 30 min. Cells were then fixed using 4% formaldehyde (Electron Microscopy Sciences), washed 3 times with PBS followed by blocking and permeabilization using 3% Bovine Serum Albumin (BSA; Sigma) and 0.3% Triton X-100 (Sigma) in PBS. Cells were then stained with antibody specific to α -smooth muscle actin (α SMA; Abcam) (ref. 1 and 2 above) in 3% Bovine Serum Albumin (BSA; Sigma) and 0.3% Triton X-100 (Sigma) in PBS, and incubated overnight at 4°C. Cells were then washed 3 times with PBS, followed by incubation with Alexa Flor-647 conjugated secondary antibody (Life Technologies [Thermo Fisher Scientific]) and DAPI at room temperature for 1 hour. Cells were then washed 3 times with PBS and plates were sealed for imaging. α SMA staining was imaged by excitation at 630nm and emission at 665 nm and quantified using the Compartmental Analysis program on the CellInsight CX5 (Thermo Scientific). Dead or apoptotic cells were excluded from analysis based on positive SYTOX green staining. % of total cells positive for α SMA were counted in each well and normalized to the average of 11 wells treated with TGF- β 1 on the same plate using Dotmatics' Studies Software. The normalized averages (fold change over untreated) of 3 replicate wells for each compound concentration were used to create dose-responses curves and EC₅₀ values were calculated using non-linear regression curve fit in the Dotmatics' Studies Software. The EC₅₀ values are reported.

[01040] Table 4 shows the activity of selected compounds as provided herein.

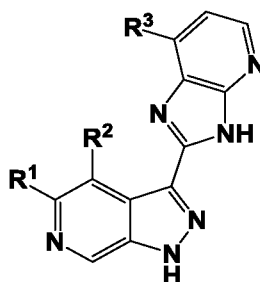
Table 4.

Compound	Inhibition of fibrosis EC ₅₀ (nM)	Compound	Inhibition of fibrosis EC ₅₀ (nM)	Compound	Inhibition of fibrosis EC ₅₀ (nM)
7	0.406	173	0.018	337	0.114
8	0.075	175	3.927	340	0.767
13	0.245	181	0.053	367	0.009
17	0.009	196	9.990	372	0.220
19	0.301	211	9.990	373	0.009
25	0.020	217	9.990	376	9.990
28	0.189	223	0.009	379	0.299
34	0.216	229	0.009	381	0.018
38	0.034	234	9.990	385	0.106
40	0.348	238	0.369	406	9.990
43	0.069	244	0.992	412	9.990
46	0.240	250	9.990	427	9.990
52	0.259	256	9.990	433	9.990
55	0.020	265	0.034	439	0.009
60	0.009	268	0.141	442	9.990
68	0.090	271	0.051	454	0.746
69	0.210	277	0.158	466	0.009
73	0.324	283	>10	939	0.014
82	0.009	292	2.300	940	0.353

87	9.990	298	0.472	941	1.005
90	9.990	303	0.020	942	1.470
101	0.009	304	3.654	945	0.040
109	9.990	310	0.641	947	9.990
136	0.009	313	0.039	948	0.632
141	9.990	319	0.075	949	0.031
148	0.009	320	0.025	952	9.990
154	0.596	324	9.990	953	0.213
157	9.990	325	0.254	954	9.990
164	0.009	329	0.034	1394	0.495
169	0.481	331	0.319		

WHAT IS CLAIMED IS:

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



I

wherein:

R^1 is selected from the group consisting of $-\text{heteroaryl}(R^4)_q$ and $-\text{heterocyclyl}(R^5)_h$;

R^2 is selected from the group consisting of H and halide;

R^3 is selected from the group consisting of H, $-\text{heteroaryl}(R^6)_q$, $-\text{heterocyclyl}(R^7)_h$, and $-\text{aryl}(R^8)_k$;

each R^4 is one substituent attached to the heteroaryl and is independently selected from the group consisting of halide, $-(C_{1-6} \text{ alkyl})$, $-(C_{1-4} \text{ alkylene})_p$, $\text{heterocyclyl}(R^9)_h$, $-(C_{1-4} \text{ alkylene})_p$, $\text{carbocyclyl}(R^{10})_j$, $-(C_{1-4} \text{ alkylene})_p$, $\text{aryl}(R^{11})_k$, $-\text{NHC}(=\text{O})R^{12}$, $-\text{NR}^{13}R^{14}$, $-(C_{1-6} \text{ alkylene})\text{NR}^{15}R^{16}$, and $-\text{OR}^{22}$;

each R^5 is one substituent attached to the heterocyclyl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-\text{CF}_3$, and $-\text{CN}$;

each R^6 is one substituent attached to the heteroaryl and is independently selected from the group consisting of $-(C_{1-6} \text{ alkyl})$, halide, $-\text{CF}_3$, $-\text{OCH}_3$, $-\text{CN}$, and $-\text{C}(=\text{O})R^{17}$;

each R^7 is one substituent attached to the heterocyclyl and is independently selected from the group consisting of $-(C_{1-6} \text{ alkyl})$, halide, $-\text{CF}_3$, $-\text{CN}$, and $-\text{OCH}_3$;

each R^8 is one substituent attached to the aryl and is independently selected from the group consisting of $-(C_{1-6} \text{ alkyl})$, halide, $-\text{CF}_3$, $-\text{CN}$, $-\text{OCH}_3$, $-(C_{1-6} \text{ alkylene})_p$, NHSO_2R^{17} , $-\text{NR}^{13}(C_{1-6} \text{ alkylene})\text{NR}^{13}R^{14}$, $-(C_{1-6} \text{ alkylene})_p$, $\text{NR}^{13}R^{14}$, and $-\text{OR}^{25}$;

each R^9 is one substituent attached to the heterocyclyl and is independently selected from the group consisting of amino, $-(C_{1-4} \text{ alkyl})$, halide, $-\text{CF}_3$, and $-\text{CN}$;

each R^{10} is one substituent attached to the carbocyclyl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-\text{CF}_3$, and $-\text{CN}$;

each R^{11} is one substituent attached to the aryl and is independently selected from the group consisting of $-(C_{1-4} \text{ alkyl})$, halide, $-\text{CF}_3$, and $-\text{CN}$;

each R¹² is independently selected from the group consisting of -(C₁₋₉ alkyl), -heteroaryl(R¹⁸)_q, -aryl(R¹⁹)_k, -CH₂aryl(R¹⁹)_k, -carbocyclyl(R²⁰)_j, -CH₂carbocyclyl(R²⁰)_j, -(C₁₋₄ alkylene)_pNR²³R²⁴, -heterocyclyl(R²¹)_h, and -CH₂heterocyclyl(R²¹)_h;

each R¹³ is independently selected from the group consisting of H and -(C₁₋₆ alkyl);

each R¹⁴ is independently selected from the group consisting of H, -(C₁₋₆ alkyl), -CH₂aryl(R¹⁹)_k, and -CH₂carbocyclyl(R²⁰)_j;

each R¹⁵ is independently selected from the group consisting of H and -(C₁₋₆ alkyl);

each R¹⁶ is independently selected from the group consisting of H, -(C₁₋₆ alkyl), -CH₂aryl(R¹⁹)_k, and -CH₂carbocyclyl(R²⁰)_j;

each R¹⁷ is a -(C₁₋₆ alkyl);

each R¹⁸ is one substituent attached to the heteroaryl and is independently selected from the group consisting of -(C₁₋₄ alkyl), halide, -CF₃, and -CN;

each R¹⁹ is one substituent attached to the aryl and is independently selected from the group consisting of -(C₁₋₄ alkyl), halide, -CF₃, and -CN;

each R²⁰ is one substituent attached to the carbocyclyl and is independently selected from the group consisting of -(C₁₋₄ alkyl), halide, -CF₃, and -CN;

each R²¹ is one substituent attached to the heterocyclyl and is independently selected from the group consisting of -(C₁₋₄ alkyl), halide, -CF₃, and -CN;

R²² is selected from the group consisting of H, -(C₁₋₆ alkyl), -(C₁₋₄ alkylene)_pheterocyclyl(R²¹)_h, -(C₁₋₄ alkylene)_pcarbocyclyl(R²⁰)_j, -(C₁₋₄ alkylene)_paryl(R¹⁹)_k, and -(C₁₋₆ alkylene)_pNR²³R²⁴;

each R²³ is independently selected from the group consisting of H and -(C₁₋₆ alkyl);

each R²⁴ is independently selected from the group consisting of H and -(C₁₋₆ alkyl);

R²⁵ is selected from the group consisting of H, -(C₁₋₆ alkyl), -(C₁₋₄ alkylene)_pheterocyclyl(R²¹)_h, and -(C₁₋₆ alkylene)_pNR²³R²⁴;

each p is independently 0 or 1;

each q is independently 0 to 4;

each h is independently 0 to 10;

each k is independently 0 to 5; and

each j is independently 0 to 12.

2. The compound of claim 1, wherein R² is H.

3. The compound of claim 1, wherein R² is F.

4. The compound of any one of the claims 1-3, wherein R¹ is -pyridinyl(R⁴)_q.

5. The compound of any one of the claims 1-4, wherein R¹ is -pyridin-3-yl(R⁴)_q.

6. The compound of any one of the claims 1-5, wherein R^1 is $-\text{pyridin-3-yl}(R^4)_q$ and q is 1.
7. The compound of any one of the claims 1-3, wherein R^1 is $-\text{pyrimidin-5-yl}(R^4)_q$ and q is 0.
8. The compound of any one of the claims 1-3, wherein R^1 is $-\text{pyrazolyl}(R^4)_q$.
9. The compound of any one of the claims 1-3 and 8, wherein R^1 is $-\text{pyrazol-4-yl}(R^4)_q$, q is 1, and R^4 is Me.
10. The compound of any one of the claims 1-3 and 8, wherein R^1 is $-\text{pyrazol-4-yl}(R^4)_q$ and q is 0.
11. The compound of any one of the claims 1-3, wherein R^1 is $-\text{imidazolyl}(R^4)_q$.
12. The compound of any one of the claims 1-3 and 11, wherein R^1 is $-\text{imidazol-5-yl}(R^4)_q$, q is 2, and both R^4 are Me.
13. The compound of any one of the claims 1-6, wherein R^4 is selected from the group consisting of $-(C_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{heterocyclyl}(R^9)_h$, $-\text{NHC}(=\text{O})R^{12}$, $-\text{NR}^{13}R^{14}$, $-\text{CH}_2\text{NR}^{15}R^{16}$, and $-\text{OR}^{22}$.
14. The compound of any one of the claims 1-6 and 13, wherein h is 0-2, and when h is 1 or 2 then each R^9 is independently selected from a halide.
15. The compound of any one of the claims 1-6 and 13, wherein R^{12} is selected from the group consisting of $-(C_{1-5} \text{ alkyl})$, $-\text{phenyl}(R^{19})_k$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, and $-\text{carbocyclyl}(R^{20})_j$.
16. The compound of any one of the claims 1-6 and 13, wherein R^{13} and R^{14} are independently selected from H and $-(C_{1-5} \text{ alkyl})$.
17. The compound of any one of the claims 1-6 and 13, wherein R^{15} and R^{16} are independently selected from H and $-(C_{1-5} \text{ alkyl})$.
18. The compound of any one of the claims 1-6 and 13, wherein R^{22} is selected from the group consisting of H, $-(C_{1-3} \text{ alkyl})$, $-\text{CH}_2\text{phenyl}(R^{19})_k$, $-\text{CH}_2\text{CH}_2\text{heterocyclyl}(R^{21})_h$, $-\text{CH}_2\text{carbocyclyl}(R^{20})_j$, and $-\text{CH}_2\text{CH}_2\text{NR}^{23}R^{24}$.
19. The compound of claim 15, wherein k and j are 0.
20. The compound of any one of the claims 1-19, wherein R^3 is $-\text{phenyl}(R^8)_k$.
21. The compound of claim 20, wherein k is 1 or 2 and each R^8 is independently selected from a halide.
22. The compound of claim 20, wherein k is 2 and one R^8 is halide and the other R^8 is $-\text{CH}_2\text{NH}\text{SO}_2R^{17}$.
23. The compound of claim 22, wherein R^{17} is $-(C_{1-3} \text{ alkyl})$.

24. The compound of claim 20, wherein k is 2 and one R⁸ is halide and one R⁸ is –NHCH₂CH₂NR¹³R¹⁴.

25. The compound of claim 24, wherein R¹³ and R¹⁴ are independently selected from H and –(C₁₋₃ alkyl).

26. The compound of any one of the claims 1-19, wherein R³ is –heteroaryl(R⁶)_q.

27. The compound any one of the claims 1-19 and 26, wherein R³ is selected from the group consisting of –pyridinyl(R⁶)_q, –imidazolyl(R⁶)_q, –furanlyl(R⁶)_q, and –thiophenyl(R⁶)_q.

28. The compound of any one of the claims 26-27, wherein q is 0 or 1, and when q is 1 then R⁶ is selected from the group consisting of halide, –(C₁₋₃ alkyl), and –C(=O)R¹⁷, wherein R¹⁷ is –(C₁₋₂ alkyl).

29. The compound of any one of the claims 1-19, wherein R³ is –heterocyclyl(R⁷)_h.

30. The compound any one of the claims 1-19 and 26, wherein R³ is selected from the group consisting of –piperidinyl(R⁷)_h, –morpholinyl(R⁷)_h, and –piperazinyl(R⁷)_h.

31. The compound of any one of the claims 29-30, wherein h is 0-2, and when h is 1 or 2 then each R⁷ is independently selected from a halide and –(C₁₋₃ alkyl).

32. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

N-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [1];

N-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [2];

5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [3];

3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [4];

3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [5];

N-((5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [6];

5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [7];

N-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [8];

- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [9];
- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [10];
- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [11];
- 5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [12];
- 1-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [13];
- 3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl) pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [14];
- 3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl) pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [15];
- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [16];
- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [17];
- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [18];
- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [19];
- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [20];
- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [21];
- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [22];
- N*-benzyl-1-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [23];
- 1-cyclopentyl-*N*-((5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [24];
- 5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [25];

- 3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [26];
- N*-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [27];
- N*-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [28];
- 5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [29];
- 3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [30];
- 3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [31];
- N*-((5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [32];
- 5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [33];
- N*-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [34];
- N*-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [35];
- N*-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [36];
- N*-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [37];
- 5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [38];
- 1-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [39];
- 3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [40];
- 3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [41];
- N*-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [42];

N-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [43];
N-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [44];
N-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [45];
N-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [46];
N-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [47];
N-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [48];
N-benzyl-1-(5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [49]; and
1-cyclopentyl-*N*-((5-(3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [50]; or a pharmaceutically acceptable salt thereof.

33. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [51];
3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [52];
N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [53];
N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [54];
5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [55];
3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [56];
3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [57];

N-((5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [58];

5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [59];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [60];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [61];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [62];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [63];

5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [64];

1-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [65];

3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl) pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [66];

3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl) pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [67];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [68];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [69];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [70];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [71];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [72];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [73];

N-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [74];

N-benzyl-1-(5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [75];

1-cyclopentyl-*N*-((5-(3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [76];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [77];

3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [78];

N-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [79];

3-methyl-*N*-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [80];

5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [81];

5-(pyridin-3-yl)-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [82];

5-(4-methylpyridin-3-yl)-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [83];

N-((5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [84];

N,N-dimethyl-5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [85];

N-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [86];

N-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [87];

2-phenyl-*N*-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)acetamide [88];

N-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [89];

N-isopropyl-5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [90];

N,N-dimethyl-1-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [91];

3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [92];

5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [93];

3,3-dimethyl-*N*-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [94];

N-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [95];

N-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [96];

N-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [97];

N-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [98];

N-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [99]; and

N-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [100]; or a pharmaceutically acceptable salt thereof.

34. The compound of claim 1, wherein the compound of Formula **I** is selected from the group consisting of:

N-benzyl-1-(5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [101];

1-cyclopentyl-*N*-((5-(3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [102];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [103];

3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [104];

N-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [105];

3-methyl-*N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [106];

5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [107];

- 5-(pyridin-3-yl)-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [108];
- 5-(4-methylpyridin-3-yl)-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [109];
- N*-((5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [110];
- N,N*-dimethyl-5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [111];
- N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [112];
- N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [113];
- 2-phenyl-*N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)acetamide [114];
- N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [115];
- N*-isopropyl-5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [116];
- N,N*-dimethyl-1-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [117];
- 3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [118];
- 5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [119];
- 3,3-dimethyl-*N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [120];
- N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [121];
- N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [122];
- N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [123];
- N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [124];

- N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [125];
- N*-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [126];
- N*-benzyl-1-(5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [127];
- 1-cyclopentyl-*N*-((5-(3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [128];
- 5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [129];
- 3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [130];
- N*-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [131];
- 3-methyl-*N*-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [132];
- 5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [133];
- 3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [134];
- 5-(4-methylpyridin-3-yl)-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [135];
- N*-((5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [136];
- N,N*-dimethyl-5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [137];
- N*-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [138];
- N*-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [139];
- 2-phenyl-*N*-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)acetamide [140];
- N*-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [141];

N-isopropyl-5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [142];

N,N-dimethyl-1-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [143];

3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [144];

5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [145];

3,3-dimethyl-*N*-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [146];

N-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [147];

N-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [148];

N-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [149]; and

N-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [150]; or a pharmaceutically acceptable salt thereof.

35. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

N-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [151];

N-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [152];

N-benzyl-1-(5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [153];

1-cyclopentyl-*N*-((5-(3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [154];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [155];

3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [156];

N-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [157];

- 3-methyl-*N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [158];
- 5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [159];
- 3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [160];
- 5-(4-methylpyridin-3-yl)-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [161];
- N*((5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [162];
- N,N*-dimethyl-5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [163];
- N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [164];
- N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [165];
- 2-phenyl-*N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)acetamide [166];
- N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [167];
- N*-isopropyl-5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [168];
- N,N*-dimethyl-1-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [169];
- 3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [170];
- 3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [171];
- 3,3-dimethyl-*N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [172];
- N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [173];
- N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [174];

- N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [175];
- N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [176];
- N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [177];
- N*-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [178];
- N*-benzyl-1-(5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [179];
- 1-cyclopentyl-*N*-((5-(3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [180];
- 5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [181];
- 3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [182];
- N*-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [183];
- 3-methyl-*N*-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [184];
- 5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [185];
- 3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [186];
- 3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [187];
- N*-((5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [188];
- N,N*-dimethyl-5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [189];
- N*-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [190];
- N*-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [191];

N-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [192];

N-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [193];

N-isopropyl-5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [194];

N,N-dimethyl-1-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [195];

3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [196];

3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [197];

3,3-dimethyl-*N*-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [198];

N-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [199]; and

N-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [200]; or a pharmaceutically acceptable salt thereof.

36. The compound of claim 1, wherein the compound of Formula **I** is selected from the group consisting of:

N-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [201];

N-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [202];

N-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [203];

N-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [204];

N-benzyl-1-(5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [205];

1-cyclopentyl-*N*-((5-(3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [206];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [207];

- 3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [208];
- N*-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [209];
- 3-methyl-*N*-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [210];
- 5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [211];
- 3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [212];
- 3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [213];
- N*-((5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [214];
- N,N*-dimethyl-5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [215];
- N*-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [216];
- N*-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [217];
- N*-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [218];
- N*-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [219];
- N*-isopropyl-5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [220];
- N,N*-dimethyl-1-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [221];
- 3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-yl)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [222];
- 3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-yl)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [223];
- 3,3-dimethyl-*N*-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [224];

N-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [225];

N-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [226];

N-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [227];

N-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [228];

N-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [229];

N-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [230];

N-benzyl-1-(5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [231];

1-cyclopentyl-*N*-((5-(3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [232];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [233];

3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [234];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [235];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [236];

5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [237];

3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [238];

3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [239];

N-((5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [240];

5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [241];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [242];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [243];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [244];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [245];

5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [246];

1-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [247];

3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [248];

3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [249]; and

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [250]; or a pharmaceutically acceptable salt thereof.

37. The compound of claim 1, wherein the compound of Formula **I** is selected from the group consisting of:

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [251];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [252];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [253];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [254];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [255];

N-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [256];

1-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N*-benzylmethanamine [257];

1-(5-(3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N*-(cyclopentylmethyl)methanamine [258];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [259];

3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [260];

N-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [261];

3-methyl-*N*-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [262];

5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [263];

5-(pyridin-3-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [264];

5-(4-methylpyridin-3-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [265];

N-((5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [266];

N,N-dimethyl-5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [267];

N-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [268];

N-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [269];

2-phenyl-*N*-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)acetamide [270];

N-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [271];

N-isopropyl-5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [272];

N,N-dimethyl-1-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [273];

5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [274];

5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [275];

3,3-dimethyl-*N*-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [276];

N-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [277];

N-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [278];

N-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [279];

N-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [280];

N-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [281];

N-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [282];

N-benzyl-1-(5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [283];

1-cyclopentyl-*N*-((5-(3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [284];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [285];

5-(pyrimidin-5-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [286];

N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [287];

N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [288];

5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [289];

3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [290];

3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [291];

N-((5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [292];

5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [293];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [294];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [295];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [296];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [297];
5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [298];
1-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [299]; and
3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [300]; or a pharmaceutically acceptable salt thereof.

38. The compound of claim 1, wherein the compound of Formula **I** is selected from the group consisting of:

3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [301];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [302];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [303];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [304];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [305];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [306];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [307];
N-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [308];

- N*-benzyl-1-(5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [309];
- 1-cyclopentyl-*N*-((5-(3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [310];
- 5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [311];
- 3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [312];
- N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [313];
- 3-methyl-*N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [314];
- 5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [315];
- 5-(pyridin-3-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [316];
- 5-(4-methylpyridin-3-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [317];
- N*-((5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [318];
- N,N*-dimethyl-5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [319];
- N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [320];
- N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [321];
- 2-phenyl-*N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)acetamide [322];
- N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [323];
- N*-isopropyl-5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [324];
- N,N*-dimethyl-1-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [325];

- 5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [326];
- 5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [327];
- 3,3-dimethyl-*N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [328];
- N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [329];
- N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [330];
- N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [331];
- N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [332];
- N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [333];
- N*-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [334];
- N*-benzyl-1-(5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [335];
- 1-cyclopentyl-*N*-((5-(3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [336];
- 5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [337];
- 5-(pyrimidin-5-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [338];
- N*-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [339];
- N*-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [340];
- 5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [341];
- 3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [342];

3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [343];
N-((5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [344];
5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [345];
N-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [346];
N-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [347];
N-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [348];
N-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [349]; and
5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [350]; or a pharmaceutically acceptable salt thereof.

39. The compound of claim 1, wherein the compound of Formula **I** is selected from the group consisting of:

1-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [351];
3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [352];
3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [353];
N-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [354];
N-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [355];
N-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [356];
N-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [357];
N-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [358];

- N*-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [359];
- N*-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [360];
- N*-benzyl-1-(5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [361];
- 1-cyclopentyl-*N*-((5-(3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [362];
- 5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [363];
- 3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [364];
- N*-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [365];
- 3-methyl-*N*-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [366];
- 5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [367];
- 3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [368];
- 5-(4-methylpyridin-3-yl)-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [369];
- N*-((5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [370];
- N,N*-dimethyl-5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [371];
- N*-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [372];
- N*-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [373];
- N*-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [374];
- N*-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [375];

N-isopropyl-5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [376];

N,N-dimethyl-1-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [377];

3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [378];

3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [379];

3,3-dimethyl-*N*-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butanamide [380];

N-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [381];

N-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [382];

N-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [383];

N-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [384];

N-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [385];

N-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [386];

N-benzyl-1-(5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [387];

1-cyclopentyl-*N*-((5-(3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [388];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [389];

3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [390];

N-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [391];

N-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [392];

1-(5-(2-(5-(5-aminopyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [393];

1-(5-(2-(5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [394];

1-(5-(2-(5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [395];

1-(5-(2-(5-(5-((ethylamino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [396];

1-(5-(2-(5-(5-(dimethylamino)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [397];

N-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [398];

N-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [399]; and

N-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [400]; or a pharmaceutically acceptable salt thereof.

40. The compound of claim 1, wherein the compound of Formula **I** is selected from the group consisting of:

N-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [401];

1-(5-(2-(5-(5-(isopropylamino)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [402];

1-(5-(2-(5-(5-((dimethylamino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [403];

1-(5-(2-(5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [404];

1-(5-(2-(5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [405];

N-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [406];

N-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [407];

N-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [408];

- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [409];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [410];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [411];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [412];
- 1-(5-(2-(5-(5-((benzylamino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [413];
- 1-(5-(2-(5-(5-(((cyclopentylmethyl)amino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [414];
- 1-(5-(2-(5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [415];
- 1-(5-(2-(5-(5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [416];
- N*-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [417];
- N*-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [418];
- N*-(3-(2-(5-(5-aminopyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [419];
- N*-(3-fluoro-5-(2-(5-(5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [420];
- N*-(3-fluoro-5-(2-(5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [421];
- N*-(3-(2-(5-(5-((ethylamino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [422];
- N*-(3-(2-(5-(5-(dimethylamino)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [423];
- N*-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [424];
- N*-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [425];

N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [426];

N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [427];

N-(3-fluoro-5-(2-(5-(5-(isopropylamino)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [428];

N-(3-(2-(5-(5-((dimethylamino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [429];

N-(3-fluoro-5-(2-(5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [430];

N-(3-fluoro-5-(2-(5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [431];

N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [432];

N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [433];

N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [434];

N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [435];

N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [436];

N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [437];

N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [438];

N-(3-(2-(5-(5-((benzylamino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [439];

N-(3-(2-(5-(5-(((cyclopentylmethyl)amino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [440];

N-(3-(2-(5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [441];

N-(3-fluoro-5-(2-(5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [442];

N-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [443];

N-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [444];

*N*¹-(3-(2-(5-(5-aminopyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [445];

*N*¹-(3-fluoro-5-(2-(5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [446];

*N*¹-(3-fluoro-5-(2-(5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [447];

*N*¹-(3-(2-(5-(5-((ethylamino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [448];

*N*¹-(3-(2-(5-(5-(dimethylamino)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [449]; and

N-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [450]; or a pharmaceutically acceptable salt thereof.

41. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

N-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [451];

N-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [452];

N-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [453];

*N*¹-(3-fluoro-5-(2-(5-(5-(isopropylamino)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [454];

*N*¹-(3-(2-(5-(5-((dimethylamino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [455];

*N*¹-(3-fluoro-5-(2-(5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [456];

*N*¹-(3-fluoro-5-(2-(5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [457];

- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [458];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [459];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [460];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [461];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [462];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [463];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [464];
- N*¹-(3-(2-(5-(5-((benzylamino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [465];
- N*¹-(3-(2-(5-(5-(((cyclopentylmethyl)amino)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [466];
- N*¹-(3-(2-(5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [467];
- N*¹-(3-fluoro-5-(2-(5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [468];
- N*-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [469];
- N*-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [470];
- 5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [471];
- 4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [472];
- 4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [473];

N-((5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [474];

5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [475];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [476];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [477];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [478];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [479];

5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [480];

1-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [481];

4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [482];

4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [483];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [484];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [485];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [486];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [487];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [488];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [489];

N-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [490];

N-benzyl-1-(5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [491];
1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [492];
5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [493];
4-fluoro-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [494];
N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [495];
N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [496];
5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [497];
4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [498];
4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [499]; and
N-((5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [500]; or a pharmaceutically acceptable salt thereof.

42. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [501];
N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [502];
N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [503];
N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [504];
N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [505];

5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [506];

1-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [507];

4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [508];

4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [509];

N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [510];

N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [511];

N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [512];

N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [513];

N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [514];

N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [515];

N-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [516];

N-benzyl-1-(5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [517];

1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [518];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [519];

4-fluoro-3-(7-(4-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [520];

N-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [521];

N-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [522];

- 5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [523];
- 4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [524];
- 4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [525];
- N*-((5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [526];
- 5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [527];
- N*-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [528];
- N*-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [529];
- N*-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [530];
- N*-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [531];
- 5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [532];
- 1-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [533];
- 4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [534];
- 4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [535];
- N*-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [536];
- N*-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [537];
- N*-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [538];
- N*-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [539];

N-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [540];

N-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [541];

N-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [542];

N-benzyl-1-(5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [543];

1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [544];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [545];

4-fluoro-3-(7-(2-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [546];

N-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [547];

N-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [548];

5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [549]; and

4-fluoro-5-(pyridin-3-yl)-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [550]; or a pharmaceutically acceptable salt thereof.

43. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

4-fluoro-5-(4-methylpyridin-3-yl)-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [551];

N-((5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [552];

5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [553];

N-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [554];

N-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [555];

- N*-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [556];
- N*-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [557];
- 5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [558];
- 1-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [559];
- 4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [560];
- 4-fluoro-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [561];
- N*-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [562];
- N*-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [563];
- N*-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [564];
- N*-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [565];
- N*-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [566];
- N*-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [567];
- N*-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [568];
- N*-benzyl-1-(5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [569];
- 1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [570];
- 5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [571];
- 4-fluoro-3-(7-(pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [572];

- N*-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [573];
- N*-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [574];
- 5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [575];
- 4-fluoro-5-(pyridin-3-yl)-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [576];
- 4-fluoro-5-(4-methylpyridin-3-yl)-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [577];
- N*-((5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [578];
- 5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [579];
- N*-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [580];
- N*-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [581];
- N*-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [582];
- N*-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [583];
- 5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [584];
- 1-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [585];
- 4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [586];
- 4-fluoro-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [587];
- N*-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [588];
- N*-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [589];

N-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [590];

N-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [591];

N-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [592];

N-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [593];

N-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [594];

N-benzyl-1-(5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [595];

1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [596];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [597];

4-fluoro-3-(7-(pyridin-4-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [598];

N-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [599]; and

N-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [600]; or a pharmaceutically acceptable salt thereof.

44. The compound of claim 1, wherein the compound of Formula **I** is selected from the group consisting of:

5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [601];

4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [602];

4-fluoro-5-(4-methylpyridin-3-yl)-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [603];

N-((5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [604];

5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [605];

- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [606];
- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [607];
- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [608];
- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [609];
- 5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [610];
- 1-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [611];
- 4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [612];
- 4-fluoro-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [613];
- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [614];
- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [615];
- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [616];
- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [617];
- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [618];
- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [619];
- N*-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [620];
- N*-benzyl-1-(5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [621];
- 1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(pyridin-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [622];

- 5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridine [623];
- 4-fluoro-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1H-pyrazolo[3,4-*c*]pyridine [624];
- N*-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [625];
- N*-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [626];
- 5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [627];
- 4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridine [628];
- 4-fluoro-5-(4-methylpyridin-3-yl)-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridine [629];
- N*-((5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [630];
- 5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [631];
- N*-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [632];
- N*-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [633];
- N*-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [634];
- N*-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [635];
- 5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [636];
- 1-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-1H-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [637];
- 4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridine [638];
- 4-fluoro-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-*c*]pyridine [639];

N-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [640];

N-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [641];

N-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [642];

N-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [643];

N-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [644];

N-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [645];

N-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [646];

N-benzyl-1-(5-(4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [647];

1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [648];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [649]; and

4-fluoro-3-(7-(piperidin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [650]; or a pharmaceutically acceptable salt thereof.

45. The compound of claim 1, wherein the compound of Formula **I** is selected from the group consisting of:

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [651];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [652];

5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [653];

4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [654];

4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [655];

N-((5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [656];

5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [657];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [658];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [659];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [660];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [661];

5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [662];

1-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [663];

4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [664];

4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [665];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [666];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [667];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [668];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [669];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [670];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [671];

N-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [672];

N-benzyl-1-(5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [673];

1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [674];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [675];

4-fluoro-3-(7-(4-methyl-1*H*-imidazol-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [676];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [677];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [678];

5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [679];

4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [680];

4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [681];

N-((5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [682];

5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [683];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [684];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [685];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [686];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [687];

5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [688];

1-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [689];

4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [690];

4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [691];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [692];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [693];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [694];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [695];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [696];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [697];

N-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [698];

N-benzyl-1-(5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [699]; and

1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [700]; or a pharmaceutically acceptable salt thereof.

46. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [701];

4-fluoro-3-(7-(4-methylpiperazin-1-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [702];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [703];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [704];

5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [705];

4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [706];

4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [707];

N-((5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [708];

5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [709];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [710];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [711];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [712];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [713];

5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [714];

1-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [715];

4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [716];

4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [717];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [718];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [719];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [720];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [721];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [722];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [723];

N-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [724];

N-benzyl-1-(5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [725];

1-cyclopentyl-*N*-((5-(4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [726];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [727];

4-fluoro-3-(3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [728];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [729];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [730];

5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [731];

4-fluoro-5-(pyridin-3-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [732];

4-fluoro-5-(4-methylpyridin-3-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [733];

N-((5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [734];

5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [735];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [736];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [737];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [738];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [739];

5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [740];

1-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [741];

4-fluoro-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [742];

4-fluoro-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [743];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [744];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [745];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [746];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [747];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [748];

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [749]; and

N-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [750]; or a pharmaceutically acceptable salt thereof.

47. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

N-benzyl-1-(5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [751];

1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [752];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [753];

- 4-fluoro-5-(pyrimidin-5-yl)-3-(7-(thiophen-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [754];
- N*-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [755];
- N*-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [756];
- 5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [757];
- 4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [758];
- 4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [759];
- N*-((5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [760];
- 5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [761];
- N*-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [762];
- N*-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [763];
- N*-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [764];
- N*-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [765];
- 5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [766];
- 1-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [767];
- 4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [768];
- 4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [769];
- N*-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [770];

N-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [771];

N-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [772];

N-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [773];

N-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [774];

N-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [775];

N-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [776];

N-benzyl-1-(5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [777];

1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [778];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [779];

4-fluoro-3-(7-(furan-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [780];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [781];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [782];

5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [783];

4-fluoro-5-(pyridin-3-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [784];

4-fluoro-5-(4-methylpyridin-3-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [785];

N-((5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [786];

5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [787];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [788];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [789];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [790];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [791];

5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [792];

1-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [793];

4-fluoro-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [794];

4-fluoro-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [795];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [796];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [797];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [798];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [799]; and

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [800]; or a pharmaceutically acceptable salt thereof.

48. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [801];

N-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [802];

N-benzyl-1-(5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [803];

1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [804];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [805];

4-fluoro-5-(pyrimidin-5-yl)-3-(7-(thiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [806];

N-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [807];

N-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [808];

5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine [809];

4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [810];

4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [811];

N-((5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [812];

5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [813];

N-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [814];

N-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [815];

N-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [816];

N-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [817];

5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [818];

1-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [819];

- 4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine **[820]**;
- 4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine **[821]**;
- N*-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide **[822]**;
- N*-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide **[823]**;
- N*-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide **[824]**;
- N*-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide **[825]**;
- N*-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide **[826]**;
- N*-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide **[827]**;
- N*-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide **[828]**;
- N*-benzyl-1-(5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine **[829]**;
- 1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine **[830]**;
- 5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine **[831]**;
- 4-fluoro-3-(7-(5-fluorothiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine **[832]**;
- N*-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide **[833]**;
- N*-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide **[834]**;
- 5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-amine **[835]**;
- 4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine **[836]**;

4-fluoro-5-(4-methylpyridin-3-yl)-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [837];
N-((5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [838];
5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N,N*-dimethylpyridin-3-amine [839];
N-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [840];
N-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [841];
N-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [842];
N-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [843];
5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)-*N*-isopropylpyridin-3-amine [844];
1-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-*N,N*-dimethylmethanamine [845];
4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [846];
4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [847];
N-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [848];
N-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [849]; and
N-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [850]; or a pharmaceutically acceptable salt thereof.

49. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

N-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [851];
N-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [852];

- N*-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [853];
- N*-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [854];
- N*-benzyl-1-(5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methanamine [855];
- 1-cyclopentyl-*N*-((5-(4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [856];
- 5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [857];
- 4-fluoro-3-(7-(5-methylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [858];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [859];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [860];
- 1-(5-(2-(5-(5-aminopyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [861];
- 1-(5-(2-(4-fluoro-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [862];
- 1-(5-(2-(4-fluoro-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [863];
- 1-(5-(2-(5-(5-((ethylamino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [864];
- 1-(5-(2-(5-(5-(dimethylamino)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [865];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [866];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [867];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [868];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [869];

- 1-(5-(2-(4-fluoro-5-(5-(isopropylamino)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [870];
- 1-(5-(2-(5-(5-((dimethylamino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [871];
- 1-(5-(2-(4-fluoro-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [872];
- 1-(5-(2-(4-fluoro-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [873];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [874];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [875];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [876];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [877];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [878];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [879];
- N*-(5-(3-(7-(5-acetylthiophen-2-yl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [880];
- 1-(5-(2-(5-(5-((benzylamino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [881];
- 1-(5-(2-(5-(5-(((cyclopentylmethyl)amino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [882];
- 1-(5-(2-(5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [883];
- 1-(5-(2-(4-fluoro-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)thiophen-2-yl)ethanone [884];
- N*-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [885];
- N*-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [886];

N-(3-(2-(5-(5-aminopyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [887];

N-(3-fluoro-5-(2-(4-fluoro-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [888];

N-(3-fluoro-5-(2-(4-fluoro-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [889];

N-(3-(2-(5-(5-((ethylamino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [890];

N-(3-(2-(5-(5-(dimethylamino)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [891];

N-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [892];

N-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [893];

N-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [894];

N-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [895];

N-(3-fluoro-5-(2-(4-fluoro-5-(5-(isopropylamino)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [896];

N-(3-(2-(5-(5-((dimethylamino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [897];

N-(3-fluoro-5-(2-(4-fluoro-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [898];

N-(3-fluoro-5-(2-(4-fluoro-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [899]; and

N-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [900]; or a pharmaceutically acceptable salt thereof.

50. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

N-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [901];

- N*-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [902];
- N*-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [903];
- N*-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [904];
- N*-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [905];
- N*-(5-(4-fluoro-3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [906];
- N*-(3-(2-(5-(5-((benzylamino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [907];
- N*-(3-(2-(5-(5-(((cyclopentylmethyl)amino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [908];
- N*-(3-(2-(5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [909];
- N*-(3-fluoro-5-(2-(4-fluoro-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)benzyl)methanesulfonamide [910];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)propionamide [911];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [912];
- N*¹-(3-(2-(5-(5-aminopyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [913];
- N*¹-(3-fluoro-5-(2-(4-fluoro-5-(pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [914];
- N*¹-(3-fluoro-5-(2-(4-fluoro-5-(4-methylpyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [915];
- N*¹-(3-(2-(5-(5-((ethylamino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [916];
- N*¹-(3-(2-(5-(5-(dimethylamino)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [917];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pivalamide [918];

- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)isobutyramide [919];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [920];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)benzamide [921];
- N*¹-(3-fluoro-5-(2-(4-fluoro-5-(5-(isopropylamino)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [922];
- N*¹-(3-(2-(5-(5-((dimethylamino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [923];
- N*¹-(3-fluoro-5-(2-(4-fluoro-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [924];
- N*¹-(3-fluoro-5-(2-(4-fluoro-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [925];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [926];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)butyramide [927];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)pentanamide [928];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [929];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [930];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [931];
- N*-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [932];
- N*¹-(3-(2-(5-(5-((benzylamino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [933];
- N*¹-(3-(2-(5-(5-(((cyclopentylmethyl)amino)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)-*N*²,*N*²-dimethylethane-1,2-diamine [934];

- N*¹-(3-(2-(5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-4-fluoro-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)-5-fluorophenyl)- *N*²,*N*²-dimethylethane-1,2-diamine [935];
- N*¹-(3-fluoro-5-(2-(4-fluoro-5-(pyrimidin-5-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-3-yl)-3*H*-imidazo[4,5-*b*]pyridin-7-yl)phenyl)- *N*²,*N*²-dimethylethane-1,2-diamine [936];
- 3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(piperidin-4-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [937];
- 3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [938];
- 3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(1*H*-pyrazol-4-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [939];
- 3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(1-methyl-1*H*-pyrazol-4-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [940];
- 5-(1,2-dimethyl-1*H*-imidazol-5-yl)-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [941];
- 1-(6-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [942];
- 5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [943];
- 3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [944];
- N*-(5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [945];
- 3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [946];
- 2-((5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-yl)oxy)-*N,N*-dimethylethan-1-amine [947];
- 3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-5-(5-methoxy-pyridin-3-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [948];
- 5-(3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridin-5-yl)pyridin-3-ol [949];
- 5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(3-fluorophenyl)-3*H*-imidazo[4,5-*b*]pyridin-2-yl)-1*H*-pyrazolo[3,4-*c*]pyridine [950];

- 2-cyclohexyl-N-(5-(3-(7-(3-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [951];
- 3-(7-(3-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [952];
- 3-(7-(3-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [953];
- 3-(7-(3-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [954];
- 3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [955];
- 3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [956];
- 3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [957];
- 3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [958];
- 5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [959];
- 1-(6-(3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetidin-3-amine [960];
- 5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [961];
- 3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [962];
- 3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [963];
- 3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [964];
- 2-((5-(3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)-N,N-dimethylethan-1-amine [965];
- 3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [966];
- 5-(3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [967];

5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [968];

2-cyclohexyl-N-(5-(3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [969];

3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [970];

3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [971];

3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [972];

3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [973];

3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [974];

3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [975];

3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [976];

5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [977];

1-(6-(3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetidin-3-amine [978];

5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [979];

3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [980];

N-(5-(3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [981];

3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [982];

2-((5-(3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)-N,N-dimethylethan-1-amine [983];

3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-methoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [984];

5-(3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [985];

5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [986];

2-cyclohexyl-N-(5-(3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [987];

3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [988];

3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [989];

3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [990];

5-(piperidin-4-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [991];

3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [992];

5-(1H-pyrazol-4-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [993];

5-(1-methyl-1H-pyrazol-4-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [994];

5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [995];

1-(6-(3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [996];

5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [997];

5-(5-(piperidin-4-yloxy)pyridin-3-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [998];

2-(piperidin-4-yl)-N-(5-(3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [999]; and

3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1000]; or a pharmaceutically acceptable salt thereof.

51. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

N,N-dimethyl-2-((5-(3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)ethan-1-amine [1001];

5-(5-methoxypyridin-3-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1002];

5-(3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1003];

5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1004];

2-cyclohexyl-N-(5-(3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1005];

3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1006];

5-(pyridin-2-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1007];

5-(pyrazin-2-yl)-3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1008];

5-(piperidin-4-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1009];

3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1010];

5-(1H-pyrazol-4-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1011];

5-(1-methyl-1H-pyrazol-4-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1012];

5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1013];

1-(6-(3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1014];

5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1015];

5-(5-(piperidin-4-yloxy)pyridin-3-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1016];

2-(piperidin-4-yl)-N-(5-(3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1017];

- 3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1018];
- N,N-dimethyl-2-((5-(3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)ethan-1-amine [1019];
- 5-(5-methoxypyridin-3-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1020];
- 5-(3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1021];
- 5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1022];
- 2-cyclohexyl-N-(5-(3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1023];
- 5-(pyridin-4-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1024];
- 5-(pyridin-2-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1025];
- 5-(pyrazin-2-yl)-3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1026];
- 5-(piperidin-4-yl)-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1027];
- 3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1028];
- 5-(1H-pyrazol-4-yl)-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1029];
- 5-(1-methyl-1H-pyrazol-4-yl)-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1030];
- 5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1031];
- 1-(6-(3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetidin-3-amine [1032];
- 5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1033];
- 5-(5-(piperidin-4-yloxy)pyridin-3-yl)-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1034];

- 2-(piperidin-4-yl)-N-(5-(3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1035];
- 3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1036];
- N,N-dimethyl-2-((5-(3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)ethan-1-amine [1037];
- 5-(5-methoxypyridin-3-yl)-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1038];
- 5-(3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1039];
- 5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1040];
- 2-cyclohexyl-N-(5-(3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1041];
- 3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1042];
- 5-(pyridin-2-yl)-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1043];
- 5-(pyrazin-2-yl)-3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1044];
- 3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1045];
- 3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1046];
- 3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1047];
- 5-(1-methyl-1H-pyrazol-4-yl)-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1048];
- 5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1049];
- 1-(6-(3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1050];
- 5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1051];

- 3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1052];
- N-(5-(3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1053];
- 3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1054];
- N,N-dimethyl-2-((5-(3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)ethan-1-amine [1055];
- 5-(5-methoxypyridin-3-yl)-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1056];
- 5-(3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1057];
- 5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1058];
- 2-cyclohexyl-N-(5-(3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1059];
- 3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1060];
- 3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1061];
- 3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1062];
- 3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1063];
- 3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1064];
- 3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1065];
- 3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1066];
- 5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1067];
- 1-(6-(3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1068];

5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1069];

3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1070];

N-(5-(3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1071];

3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1072];

N,N-dimethyl-2-((5-(3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)ethan-1-amine [1073];

5-(5-methoxypyridin-3-yl)-3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1074];

5-(3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1075];

5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1076];

2-cyclohexyl-N-(5-(3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1077];

3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1078];

3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1079];

3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1080];

3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1081];

3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1082];

3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1083];

5-(1-methyl-1H-pyrazol-4-yl)-3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1084];

5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1085];

1-(6-(3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1086];

5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1087];

3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1088];

N-(5-(3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1089];

3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1090];

N,N-dimethyl-2-((5-(3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)ethan-1-amine [1091];

5-(5-methoxypyridin-3-yl)-3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1092];

5-(3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1093];

5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1094];

2-cyclohexyl-N-(5-(3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1095];

3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1096];

3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1097];

3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1098];

3-(3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1099]; and

3-(3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1100]; or a pharmaceutically acceptable salt thereof.

52. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

3-(3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1101];

3-(3H-imidazo[4,5-b]pyridin-2-yl)-5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1102];

- 5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1103];
- 1-(6-(3-(3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1104];
- 5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1105];
- 3-(3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1106];
- N-(5-(3-(3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1107];
- 3-(3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1108];
- 2-((5-(3-(3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)-N,N-dimethylethan-1-amine [1109];
- 3-(3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1110];
- 5-(3-(3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1111];
- 5-(5-(benzyloxy)pyridin-3-yl)-3-(3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1112];
- N-(5-(3-(3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-cyclohexylacetamide [1113];
- 3-(3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1114];
- 3-(3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1115];
- 3-(3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1116];
- 5-(piperidin-4-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1117];
- 5-(1,2,3,6-tetrahydropyridin-4-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1118];
- 5-(1H-pyrazol-4-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1119];
- 5-(1-methyl-1H-pyrazol-4-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1120];
- 5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1121];

- 1-(6-(3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1122];
- 5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1123];
- 5-(5-(piperidin-4-yloxy)pyridin-3-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1124];
- 2-(piperidin-4-yl)-N-(5-(3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1125];
- 5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1126];
- N,N-dimethyl-2-((5-(3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)ethan-1-amine [1127];
- 5-(5-methoxypyridin-3-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1128];
- 5-(3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1129];
- 5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1130];
- 2-cyclohexyl-N-(5-(3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1131];
- 5-(pyridin-4-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1132];
- 5-(pyridin-2-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1133];
- 5-(pyrazin-2-yl)-3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1134];
- 3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1135];
- 3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1136];
- 3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1137];
- 3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1138];

- 5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1139];
- 1-(6-(3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1140];
- 5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1141];
- 3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1142];
- N-(5-(3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1143];
- 3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1144];
- 2-((5-(3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)-N,N-dimethylethan-1-amine [1145];
- 3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1146];
- 5-(3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1147];
- 5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1148];
- 2-cyclohexyl-N-(5-(3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1149];
- 3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1150];
- 3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1151];
- 3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1152];
- 5-(piperidin-4-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1153];
- 5-(1,2,3,6-tetrahydropyridin-4-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1154];
- 5-(1H-pyrazol-4-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1155];

- 5-(1-methyl-1H-pyrazol-4-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1156];
- 5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1157];
- 1-(6-(3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1158];
- 5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1159];
- 5-(5-(piperidin-4-yloxy)pyridin-3-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1160];
- 2-(piperidin-4-yl)-N-(5-(3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1161];
- 5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1162];
- N,N-dimethyl-2-((5-(3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)ethan-1-amine [1163];
- 5-(5-methoxypyridin-3-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1164];
- 5-(3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1165];
- 5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1166];
- 2-cyclohexyl-N-(5-(3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1167];
- 5-(pyridin-4-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1168];
- 5-(pyridin-2-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1169];
- 5-(pyrazin-2-yl)-3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1170];
- 3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1171];
- 3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1172];

- 3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1173];
- 3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1174];
- 5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1175];
- 1-(6-(3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1176];
- 5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1177];
- 3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1178];
- N-(5-(3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1179];
- 3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1180];
- 2-((5-(3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)-N,N-dimethylethan-1-amine [1181];
- 3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1182];
- 5-(3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1183];
- 5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1184];
- 2-cyclohexyl-N-(5-(3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1185];
- 3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1186];
- 3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1187];
- 3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1188];
- 3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1189];

3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1190];
3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1191];
5-(1-methyl-1H-pyrazol-4-yl)-3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1192];
5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1193];
1-(6-(3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1194];
5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1195];
3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1196];
N-(5-(3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1197];
3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1198];
N,N-dimethyl-2-((5-(3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)ethan-1-amine [1199]; and
5-(5-methoxypyridin-3-yl)-3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1200]; or a pharmaceutically acceptable salt thereof.

53. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

5-(3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1201];
5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1202];
2-cyclohexyl-N-(5-(3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1203];
3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1204];
3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1205];

- 3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1206];
- 1-(5-(2-(5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1207];
- 1-(5-(2-(5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1208];
- 1-(5-(2-(5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1209];
- 1-(5-(2-(5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1210];
- 1-(5-(2-(5-(1,2-dimethyl-1H-imidazol-5-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1211];
- 1-(5-(2-(5-(6-(3-aminoazetidin-1-yl)pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1212];
- 1-(5-(2-(5-(5-(cyclohexyloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1213];
- 1-(5-(2-(5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1214];
- N-(5-(3-(7-(5-acetylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1215];
- 1-(5-(2-(5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1216];
- 1-(5-(2-(5-(5-(2-(dimethylamino)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1217];
- 1-(5-(2-(5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1218];
- 1-(5-(2-(5-(5-hydroxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1219];
- 1-(5-(2-(5-(5-(benzyloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1220];
- N-(5-(3-(7-(5-acetylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-cyclohexylacetamide [1221];
- 1-(5-(2-(5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1222];

- 1-(5-(2-(5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1223];
- 1-(5-(2-(5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)thiophen-2-yl)ethan-1-one [1224];
- N-(3-fluoro-5-(2-(5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)benzyl)methanesulfonamide [1225];
- N-(3-fluoro-5-(2-(5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)benzyl)methanesulfonamide [1226];
- N-(3-(2-(5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [1227];
- N-(3-fluoro-5-(2-(5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)benzyl)methanesulfonamide [1228];
- N-(3-(2-(5-(1,2-dimethyl-1H-imidazol-5-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [1229];
- N-(3-(2-(5-(6-(3-aminoazetidin-1-yl)pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [1230];
- N-(3-(2-(5-(5-(cyclohexyloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [1231];
- N-(3-fluoro-5-(2-(5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)benzyl)methanesulfonamide [1232];
- N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1233];
- N-(3-fluoro-5-(2-(5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)benzyl)methanesulfonamide [1234];
- N-(3-(2-(5-(5-(2-(dimethylamino)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [1235];
- N-(3-fluoro-5-(2-(5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)benzyl)methanesulfonamide [1236];
- N-(3-fluoro-5-(2-(5-(5-hydroxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)benzyl)methanesulfonamide [1237];
- N-(3-(2-(5-(5-(benzyloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorobenzyl)methanesulfonamide [1238];
- 2-cyclohexyl-N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1239];

- N-(3-fluoro-5-(2-(5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)benzyl)methanesulfonamide [1240];
- N-(3-fluoro-5-(2-(5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)benzyl)methanesulfonamide [1241];
- N-(3-fluoro-5-(2-(5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)benzyl)methanesulfonamide [1242];
- N¹-(3-fluoro-5-(2-(5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine [1243];
- N¹-(3-fluoro-5-(2-(5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine [1244];
- N¹-(3-(2-(5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine [1245];
- N¹-(3-fluoro-5-(2-(5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine [1246];
- N¹-(3-(2-(5-(1,2-dimethyl-1H-imidazol-5-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine [1247];
- N¹-(3-(2-(5-(6-(3-aminoazetidin-1-yl)pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine [1248];
- N¹-(3-(2-(5-(5-(cyclohexyloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine [1249];
- N¹-(3-fluoro-5-(2-(5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine [1250];
- N-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1251];
- N¹-(3-fluoro-5-(2-(5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine [1252];
- N¹-(3-(2-(5-(5-(2-(dimethylamino)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine [1253];
- N¹-(3-fluoro-5-(2-(5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine [1254];
- 5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1255];
- N¹-(3-(2-(5-(5-(benzyloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenyl)-N²,N²-dimethylethane-1,2-diamine [1256];

2-cyclohexyl-N-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1257];

N¹-(3-fluoro-5-(2-(5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine [1258];

N¹-(3-fluoro-5-(2-(5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine [1259];

N¹-(3-fluoro-5-(2-(5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenyl)-N²,N²-dimethylethane-1,2-diamine [1260];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)propionamide [1261];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [1262];

5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-amine [1263];

2-(3-fluoro-5-(2-(5-(pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1264];

2-(3-fluoro-5-(2-(5-(4-methylpyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1265];

2-(3-(2-(5-(5-((ethylamino)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenoxy)-N,N-dimethylethan-1-amine [1266];

5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)-N,N-dimethylpyridin-3-amine [1267];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)pivalamide [1268];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)isobutyramide [1269];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [1270];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)benzamide [1271];

5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)-N-isopropylpyridin-3-amine [1272];

2-(3-(2-(5-(5-((dimethylamino)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenoxy)-N,N-dimethylethan-1-amine [1273];

2-(3-fluoro-5-(2-(5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1274];

2-(3-fluoro-5-(2-(5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1275];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [1276];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)butyramide [1277];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)pentanamide [1278];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [1279];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [1280];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [1281];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [1282];

2-(3-(2-(5-(5-((benzylamino)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenoxy)-N,N-dimethylethan-1-amine [1283];

2-(3-(2-(5-(5-(((cyclopentylmethyl)amino)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenoxy)-N,N-dimethylethan-1-amine [1284];

2-(3-(2-(5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenoxy)-N,N-dimethylethan-1-amine [1285];

2-(3-fluoro-5-(2-(5-(pyrimidin-5-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1286];

2-(3-fluoro-5-(2-(5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1287];

2-(3-fluoro-5-(2-(5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1288];

2-(3-fluoro-5-(2-(5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1289];

2-(3-(2-(5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenoxy)-N,N-dimethylethan-1-amine [1290];

2-(3-fluoro-5-(2-(5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1291];

2-(3-(2-(5-(1,2-dimethyl-1H-imidazol-5-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenoxy)-N,N-dimethylethan-1-amine [1292];

1-(6-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetidin-3-amine [1293];

2-(3-(2-(5-(5-(cyclohexyloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenoxy)-N,N-dimethylethan-1-amine [1294];

2-(3-fluoro-5-(2-(5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1295];

N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1296];

2-(3-fluoro-5-(2-(5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1297];

2-((5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)-N,N-dimethylethan-1-amine [1298];

2-(3-fluoro-5-(2-(5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1299]; and

5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1300]; or a pharmaceutically acceptable salt thereof.

54. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

2-(3-(2-(5-(5-(benzyloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenoxy)-N,N-dimethylethan-1-amine [1301];

2-cyclohexyl-N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1302];

2-(3-fluoro-5-(2-(5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1303];

2-(3-fluoro-5-(2-(5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenoxy)-N,N-dimethylethan-1-amine [1304];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)propionamide [1305];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [1306];

5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-amine [1307];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1308];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1309];

N-((5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [1310];

5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)-N,N-dimethylpyridin-3-amine [1311];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)pivalamide [1312];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)isobutyramide [1313];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [1314];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)benzamide [1315];

5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)-N-isopropylpyridin-3-amine [1316];

1-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-N,N-dimethylmethanamine [1317];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1318];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1319];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [1320];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)butyramide [1321];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)pentanamide [1322];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [1323];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [1324];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [1325];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [1326];

N-benzyl-1-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)methanamine [1327];

1-cyclopentyl-N-((5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [1328];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1329];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrimidin-5-yl)-1H-pyrazolo[3,4-c]pyridine [1330];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1331];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1332];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1333];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1334];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1335];

5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1336];

1-(6-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1337];

5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1338];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1339];

N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1340];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1341];

2-((5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)-N,N-dimethylethan-1-amine [1342];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1343];

5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1344];

5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1345];

2-cyclohexyl-N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1346];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1347];

3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1348];

N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)propionamide [1349];

N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [1350];

3-(2-(5-(5-aminopyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1351];

3-fluoro-5-(2-(5-(pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1352];

3-fluoro-5-(2-(5-(4-methylpyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1353];

3-(2-(5-(5-((ethylamino)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1354];

3-(2-(5-(5-(dimethylamino)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1355];

N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)pivalamide [1356];

- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)isobutyramide [1357];
- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [1358];
- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)benzamide [1359];
- 3-fluoro-5-(2-(5-(5-(isopropylamino)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1360];
- 3-(2-(5-(5-((dimethylamino)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1361];
- 3-fluoro-5-(2-(5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1362];
- 3-fluoro-5-(2-(5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1363];
- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-indazol-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [1364];
- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)butyramide [1365];
- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)pentanamide [1366];
- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [1367];
- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [1368];
- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [1369];
- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [1370];
- 3-(2-(5-(5-((benzylamino)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1371];
- 3-(2-(5-(5-(((cyclopentylmethyl)amino)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1372];
- 3-(2-(5-(5-(((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1373];

- 3-fluoro-5-(2-(5-(pyrimidin-5-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1374];
- 3-fluoro-5-(2-(5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1375];
- 3-fluoro-5-(2-(5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1376];
- 3-fluoro-5-(2-(5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1377];
- 3-(2-(5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1378];
- 3-fluoro-5-(2-(5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1379];
- 3-(2-(5-(1,2-dimethyl-1H-imidazol-5-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1380];
- 3-(2-(5-(6-(3-aminoazetid-1-yl)pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1381];
- 3-(2-(5-(5-(cyclohexyloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1382];
- 3-fluoro-5-(2-(5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1383];
- N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1384];
- 3-fluoro-5-(2-(5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1385];
- 3-(2-(5-(5-(2-(dimethylamino)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1386];
- 3-fluoro-5-(2-(5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1387];
- 5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1388];
- 3-(2-(5-(5-(benzyloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)-5-fluorophenol [1389];
- 2-cyclohexyl-N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1390];

3-fluoro-5-(2-(5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1391];

3-fluoro-5-(2-(5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)phenol [1392];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)propionamide [1393];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3-methylbutanamide [1394];

5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-amine [1395];

3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1396];

3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(4-methylpyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1397];

N-((5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)methyl)ethanamine [1398];

5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)-N,N-dimethylpyridin-3-amine [1399]; and

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)pivalamide [1400]; or a pharmaceutically acceptable salt thereof.

55. The compound of claim 1, wherein the compound of Formula I is selected from the group consisting of:

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)isobutyramide [1401];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-phenylacetamide [1402];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)benzamide [1403];

5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)-N-isopropylpyridin-3-amine [1404];

1-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-N,N-dimethylmethanamine [1405];

3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(pyrrolidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1406];

3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-1-ylmethyl)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1407];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-3,3-dimethylbutanamide [1408];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)butyramide [1409];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)pentanamide [1410];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclopropanecarboxamide [1411];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclobutanecarboxamide [1412];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclopentanecarboxamide [1413];

N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)cyclohexanecarboxamide [1414];

N-benzyl-1-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)methanamine [1415];

1-cyclopentyl-N-((5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)methyl)methanamine [1416];

5-(5-((3,3-difluoropyrrolidin-1-yl)methyl)pyridin-3-yl)-3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1417];

3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrimidin-5-yl)-1H-pyrazolo[3,4-c]pyridine [1418];

3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1419];

3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(piperidin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1420];

3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1,2,3,6-tetrahydropyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1421];

3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1422];

3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1423];

- 5-(1,2-dimethyl-1H-imidazol-5-yl)-3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1424];
- 1-(6-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyrazin-2-yl)azetid-3-amine [1425];
- 5-(5-(cyclohexyloxy)pyridin-3-yl)-3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1426];
- 3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(piperidin-4-yloxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1427];
- N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(piperidin-4-yl)acetamide [1428];
- 3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-(2-(pyrrolidin-1-yl)ethoxy)pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1429];
- 2-((5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)oxy)-N,N-dimethylethan-1-amine [1430];
- 3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-methoxypyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1431];
- 5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-ol [1432];
- 5-(5-(benzyloxy)pyridin-3-yl)-3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1433];
- 2-cyclohexyl-N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1434];
- 3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-4-yl)-1H-pyrazolo[3,4-c]pyridine [1435];
- 3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyrazin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1436];
- 2-(dimethylamino)-N-(5-(3-(7-(3-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1437];
- 2-(dimethylamino)-N-(5-(3-(7-(4-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1438];
- 2-(dimethylamino)-N-(5-(3-(7-(2-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1439];
- 2-(dimethylamino)-N-(5-(3-(7-(pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1440];

- 2-(dimethylamino)-N-(5-(3-(7-(pyridin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1441];
- 2-(dimethylamino)-N-(5-(3-(7-(pyridin-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1442];
- 2-(dimethylamino)-N-(5-(3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1443];
- 2-(dimethylamino)-N-(5-(3-(7-(4-methyl-1H-imidazol-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1444];
- 2-(dimethylamino)-N-(5-(3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1445];
- N-(5-(3-(3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(dimethylamino)acetamide [1446];
- 2-(dimethylamino)-N-(5-(3-(7-(thiophen-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1447];
- 2-(dimethylamino)-N-(5-(3-(7-(furan-3-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1448];
- 2-(dimethylamino)-N-(5-(3-(7-(thiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1449];
- 2-(dimethylamino)-N-(5-(3-(7-(5-fluorothiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1450];
- 2-(dimethylamino)-N-(5-(3-(7-(5-methylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1451];
- N-(5-(3-(7-(5-acetylthiophen-2-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)-2-(dimethylamino)acetamide [1452];
- 2-(dimethylamino)-N-(5-(3-(7-(3-fluoro-5-(methylsulfonamidomethyl)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1453];
- 2-(dimethylamino)-N-(5-(3-(7-(3-((2-(dimethylamino)ethyl)amino)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1454];
- 2-(dimethylamino)-N-(5-(3-(7-(3-(2-(dimethylamino)ethoxy)-5-fluorophenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1455];
- 2-(dimethylamino)-N-(5-(3-(7-(3-fluoro-5-(2-(pyrrolidin-1-yl)ethoxy)phenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1456];
- 2-(dimethylamino)-N-(5-(3-(7-(3-fluoro-5-hydroxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1457];

- 2-(dimethylamino)-N-(5-(3-(7-(3-fluoro-5-methoxyphenyl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridin-5-yl)pyridin-3-yl)acetamide [1458];
- 4-(2-(5-(pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)morpholine [1459];
- 3-(7-(4,4-difluoropiperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1460];
- 3-(7-(1-methylpiperidin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(pyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1461];
- 4-(2-(5-(5-fluoropyridin-3-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)morpholine [1462];
- 5-(5-fluoropyridin-3-yl)-3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1463];
- 3-(7-(4,4-difluoropiperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(5-fluoropyridin-3-yl)-1H-pyrazolo[3,4-c]pyridine [1464];
- 5-(5-fluoropyridin-3-yl)-3-(7-(1-methylpiperidin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1465];
- 5-(5-fluoropyridin-3-yl)-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1466];
- 4-(2-(5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)morpholine [1467];
- 3-(7-(4,4-difluoropiperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1-methyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1468];
- 5-(1-methyl-1H-pyrazol-4-yl)-3-(7-(1-methylpiperidin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1469];
- 4-(2-(5-(1-cyclopropyl-1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)morpholine [1470];
- 5-(1-cyclopropyl-1H-pyrazol-4-yl)-3-(7-(4-methylpiperazin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1471];
- 5-(1-cyclopropyl-1H-pyrazol-4-yl)-3-(7-(4,4-difluoropiperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1472];
- 5-(1-cyclopropyl-1H-pyrazol-4-yl)-3-(7-(1-methylpiperidin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1473];
- 5-(1-cyclopropyl-1H-pyrazol-4-yl)-3-(7-(piperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-1H-pyrazolo[3,4-c]pyridine [1474];

4-(2-(5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridin-3-yl)-3H-imidazo[4,5-b]pyridin-7-yl)morpholine [1475];
3-(7-(4,4-difluoropiperidin-1-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1476]; and
3-(7-(1-methylpiperidin-4-yl)-3H-imidazo[4,5-b]pyridin-2-yl)-5-(1H-pyrazol-4-yl)-1H-pyrazolo[3,4-c]pyridine [1477]; or a pharmaceutically acceptable salt thereof.

56. A pharmaceutical composition comprising a compound according to any one of the claims 1-55, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable excipient.

57. A method of treating or ameliorating in a patient a disorder or disease selected from the group consisting of: cancer, pulmonary fibrosis, idiopathic pulmonary fibrosis (IPF), degenerative disc disease, bone/osteoporotic fracture, bone or cartilage disease, and osteoarthritis, the method comprising administering to the patient a therapeutically effective amount of a compound according to any one of the claims 1-55, or a pharmaceutically acceptable salt thereof.

58. The method of claim 57, wherein the disorder or disease is cancer.

59. The method of claim 57, wherein the disorder or disease is pulmonary fibrosis.

60. The method of claim 57, wherein the disorder or disease is idiopathic pulmonary fibrosis (IPF).

61. The method of claim 57, wherein the disorder or disease is degenerative disc disease.

62. The method of claim 57, wherein the disorder or disease is a bone/osteoporotic fracture.

63. The method of claim 57, wherein the disorder or disease is a bone or cartilage disease.

64. The method of claim 57, wherein the disorder or disease is osteoarthritis.

65. The method of claim 57, wherein the patient is a human.

66. The method of claim 58, wherein the cancer is selected from the group consisting of: colon cancer, colorectal cancer, leukemia, breast cancer, skin cancer, prostate cancer, stomach (gastric) cancer, lung cancer, pancreatic cancer, and liver (hepatic) cancer.

67. The method of claim 57, wherein the compound inhibits one or more proteins in the Wnt pathway.

68. The method of claim 57, wherein the compound inhibits signaling induced by one or more Wnt proteins.

69. The method of claim 67 or 68, wherein the Wnt proteins are selected from the group consisting of: WNT1, WNT2, WNT2B, WNT3, WNT3A, WNT4, WNT5A, WNT5B, WNT6, WNT7A, WNT7B, WNT8A, WNT8B, WNT9A, WNT9B, WNT10A, WNT10B, WNT11, and WNT16.

70. The method of claim 57, wherein the compound inhibits a kinase activity.

INTERNATIONAL SEARCH REPORT

International application No.

PCT/US2015/048680

A. CLASSIFICATION OF SUBJECT MATTER

IPC(8) - A61K 31/4162 (2015.01)

CPC - A61K 31/4162 (2015.10)

According to International Patent Classification (IPC) or to both national classification and IPC

B. FIELDS SEARCHED

Minimum documentation searched (classification system followed by classification symbols)

IPC(8) - A61K 31/4162, 31/4188; C07D 403/04, 403/14 (2015.01)

CPC - A61K 31/4162, 31/4188; C07D 403/04, 403/14 (2015.10)

Documentation searched other than minimum documentation to the extent that such documents are included in the fields searched

USPC - 514/243, 393; 546/118; 548/305.1; IPC(8) - A61K 31/4162, 31/4188; C07D 403/04, 403/14; CPC - A61K 31/4162, 31/4188; C07D 403/04, 403/14 (keyword delimited)

Electronic data base consulted during the international search (name of data base and, where practicable, search terms used)

Orbit, Google Patents, Google, PubChem, SureChem, STN.

Search terms used: imidazo, pyridine, pyrazolo, azaindazole, Wnt proteins.

C. DOCUMENTS CONSIDERED TO BE RELEVANT

Category*	Citation of document, with indication, where appropriate, of the relevant passages	Relevant to claim No.
A	US 2009/0048249 A1 (CHIU et al) 19 February 2009 (19.02.2009) entire document	1-4, 7, 8, 11, 32-55
A	US 8,288,425 B2 (EDWARDS et al) 16 October 2012 (16.10.2012) entire document	1-4, 7, 8, 11, 32-55
A	US 2006/0079564 A1 (JANSEN et al) 13 April 2006 (13.04.2006) entire document	1-4, 7, 8, 11, 32-55
A	WO 2003/035065 A1 (AVENTIS PHARMACEUTICALS INC) 01 May 2003 (01.05.2003) entire document	1-4, 7, 8, 11, 32-55

 Further documents are listed in the continuation of Box C. See patent family annex.

* Special categories of cited documents:

"A" document defining the general state of the art which is not considered to be of particular relevance

"E" earlier application or patent but published on or after the international filing date

"L" document which may throw doubts on priority claim(s) or which is cited to establish the publication date of another citation or other special reason (as specified)

"O" document referring to an oral disclosure, use, exhibition or other means

"P" document published prior to the international filing date but later than the priority date claimed

"T" later document published after the international filing date or priority date and not in conflict with the application but cited to understand the principle or theory underlying the invention

"X" document of particular relevance; the claimed invention cannot be considered novel or cannot be considered to involve an inventive step when the document is taken alone

"Y" document of particular relevance; the claimed invention cannot be considered to involve an inventive step when the document is combined with one or more other such documents, such combination being obvious to a person skilled in the art

"&" document member of the same patent family

Date of the actual completion of the international search

21 October 2015

Date of mailing of the international search report

11 JAN 2016

Name and mailing address of the ISA/

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INTERNATIONAL SEARCH REPORT

International application No.

PCT/US2015/048680

Box No. II Observations where certain claims were found unsearchable (Continuation of item 2 of first sheet)

This international search report has not been established in respect of certain claims under Article 17(2)(a) for the following reasons:

- 1. Claims Nos.:
because they relate to subject matter not required to be searched by this Authority, namely:

- 2. Claims Nos.:
because they relate to parts of the international application that do not comply with the prescribed requirements to such an extent that no meaningful international search can be carried out, specifically:

- 3. Claims Nos.: 5, 6, 9, 10, 12-31, 56-70
because they are dependent claims and are not drafted in accordance with the second and third sentences of Rule 6.4(a).

Box No. III Observations where unity of invention is lacking (Continuation of item 3 of first sheet)

This International Searching Authority found multiple inventions in this international application, as follows:

- 1. As all required additional search fees were timely paid by the applicant, this international search report covers all searchable claims.
- 2. As all searchable claims could be searched without effort justifying additional fees, this Authority did not invite payment of additional fees.
- 3. As only some of the required additional search fees were timely paid by the applicant, this international search report covers only those claims for which fees were paid, specifically claims Nos.:

- 4. No required additional search fees were timely paid by the applicant. Consequently, this international search report is restricted to the invention first mentioned in the claims; it is covered by claims Nos.:

Remark on Protest

- The additional search fees were accompanied by the applicant's protest and, where applicable, the payment of a protest fee.
- The additional search fees were accompanied by the applicant's protest but the applicable protest fee was not paid within the time limit specified in the invitation.
- No protest accompanied the payment of additional search fees.