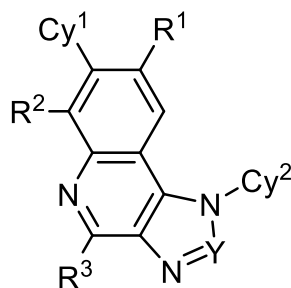


WHAT IS CLAIMED IS:

1. Use of:
 - (i) an FGFR1 inhibitor; and
 - (ii) a KRAS inhibitorin the manufacture of a medicament for the treatment of cancer in a patient.
2. The use of claim 1, wherein the FGFR1 inhibitor is selected from pemigatinib, futibatinib, erdafitinib and infigratinib, or a pharmaceutically acceptable salt thereof.
3. The use of claim 1 or 2, wherein the FGFR1 inhibitor is pemigatinib, or a pharmaceutically acceptable salt thereof.
4. The use of claim 1 or 2, wherein the FGFR1 inhibitor is futibatinib, or a pharmaceutically acceptable salt thereof.
5. The use of claim 1 or 2, wherein the FGFR1 inhibitor is erdafitinib, or a pharmaceutically acceptable salt thereof.
6. The use of claim 1 or 2, wherein the FGFR1 inhibitor is infigratinib, or a pharmaceutically acceptable salt thereof.
7. The use of claim 1, wherein the FGFR1 inhibitor further inhibits FGFR2, FGFR3, or a combination thereof.
8. The use of any one of claims 1-7, wherein the KRAS inhibitor is a compound of Formula (I):



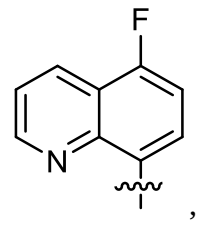
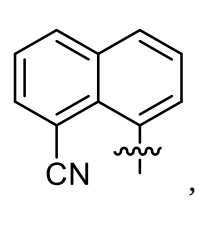
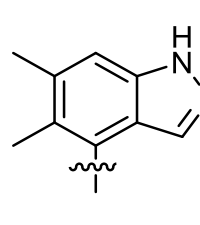
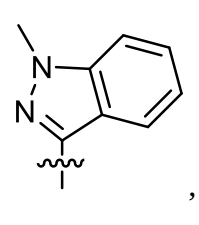
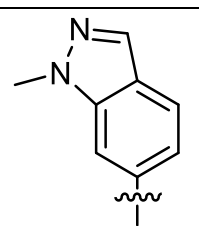
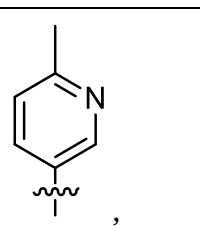
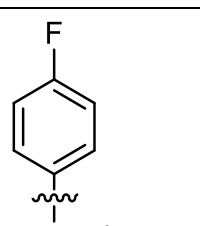
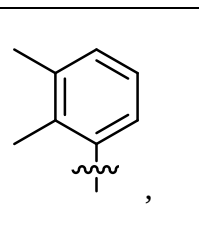
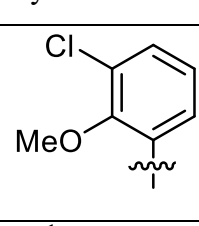
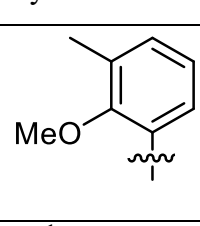
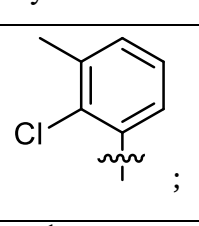
I

or a pharmaceutically acceptable salt thereof, wherein:

Y is N or CH;

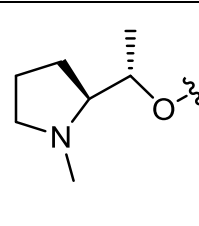
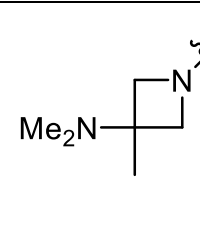
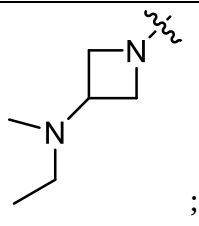
R¹ is selected from Cl, CH₃, CH₂F, CHF₂, and CF₃;

Cy¹ is selected from

			
Cy ¹ -a	Cy ¹ -b	Cy ¹ -c	Cy ¹ -d
			
Cy ¹ -e	Cy ¹ -f	Cy ¹ -g	Cy ¹ -h
		and	
Cy ¹ -i	Cy ¹ -j		Cy ¹ -k

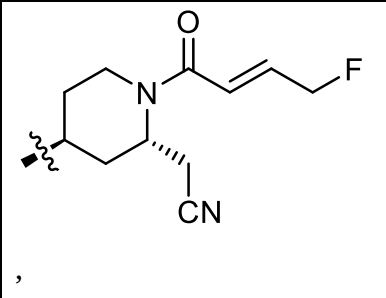
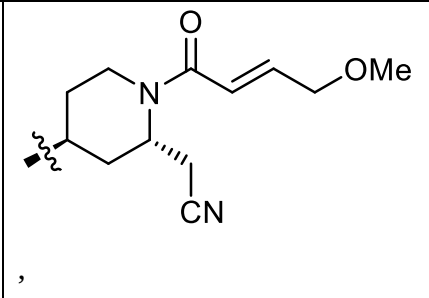
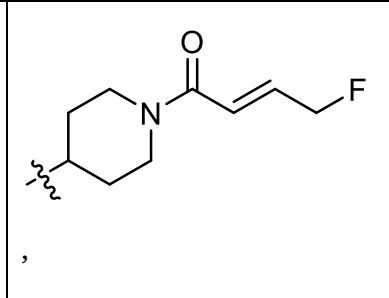
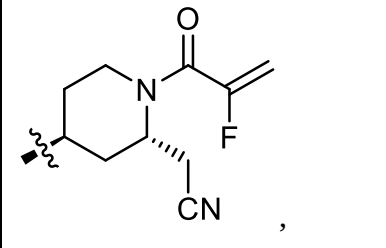
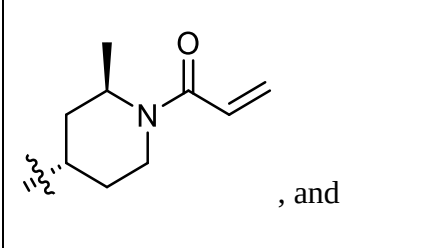
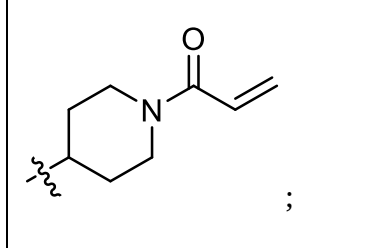
R² is selected from F and Cl;

R³ is selected from

		and	
R ³ -a	R ³ -b		R ³ -c

and,

Cy² is selected from

		
Cy ² -a	Cy ² -b	Cy ² -c
		
Cy ² -d	Cy ² -e	Cy ² -f

provided that the compound of Formula (I) is other than

2-((2S,4S)-4-(7-(2-chloro-3-methylphenyl)-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-8-methyl-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile and

2-((2S,4S)-4-(4-(3-(dimethylamino)-3-methylazetidin-1-yl)-7-(2,3-dimethylphenyl)-6-fluoro-8-methyl-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile.

9. The use of claim 8, wherein the compound of Formula (I) or the pharmaceutically acceptable salt thereof is selected from:

2-((2S,4S)-4-(4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-7-(5-fluoroquinolin-8-yl)-8-(trifluoromethyl)-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-7-(2-methoxy-3-methylphenyl)-8-(trifluoromethyl)-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(7-(3-chloro-2-methoxyphenyl)-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-8-(trifluoromethyl)-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6-fluoro-7-(5-fluoroquinolin-8-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-8-(trifluoromethyl)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

1-(4-(6-fluoro-7-(5-fluoroquinolin-8-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-8-(trifluoromethyl)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)piperidin-1-yl)prop-2-en-1-one;

2-((2S,4S)-4-(4-(3-(dimethylamino)-3-methylazetidin-1-yl)-7-(2,3-dimethylphenyl)-6-fluoro-8-methyl-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6-fluoro-8-methyl-7-(1-methyl-1*H*-indazol-6-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6-fluoro-8-methyl-7-(6-methylpyridin-3-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6-fluoro-8-methyl-7-(1-methyl-1*H*-indazol-3-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6-fluoro-7-(4-fluorophenyl)-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;

8-(1-(1-acryloylpiperidin-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-7-yl)-1-naphthonitrile;

2-((2S,4S)-4-(7-(2-chloro-3-methylphenyl)-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-8-methyl-1*H*-imidazo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(7-(2-chloro-3-methylphenyl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-imidazo[4,5-*c*]quinolin-1-yl)-1-(2-fluoroacryloyl)piperidin-2-yl)acetonitrile;

8-(1-((2*R*,4*S*)-1-acryloyl-2-methylpiperidin-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-7-yl)-1-naphthonitrile;

2-((2S,4S)-4-(7-(5,6-dimethyl-1*H*-indazol-4-yl)-4-(3-(ethyl(methyl)amino)azetidin-1-yl)-6-fluoro-8-methyl-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-(2-fluoroacryloyl)piperidin-2-yl)acetonitrile;

8-(6-fluoro-1-(1-((*E*)-4-fluorobut-2-enoyl)piperidin-4-yl)-8-methyl-4-((*S*)-1-((*S*)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-7-yl)-1-naphthonitrile;

8-(1-((2*S*,4*S*)-2-(cyanomethyl)-1-(2-fluoroacryloyl)piperidin-4-yl)-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-8-methyl-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-7-yl)-1-naphthonitrile;

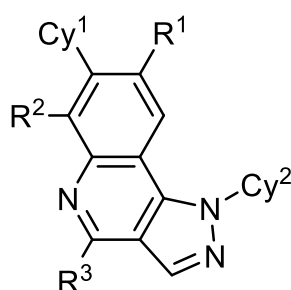
2-((2*S*,4*S*)-4-(6,8-dichloro-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-7-(5-fluoroquinolin-8-yl)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2*S*,4*S*)-4-(6,8-dichloro-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-7-(5-fluoroquinolin-8-yl)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile; and

2-((2*S*,4*S*)-4-(4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-7-(5-fluoroquinolin-8-yl)-8-(trifluoromethyl)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;

or a pharmaceutically acceptable salt thereof.

10. The use of any one of claims 1-7, wherein the KRAS inhibitor is a compound of Formula (II):



II

or a pharmaceutically acceptable salt thereof, wherein:

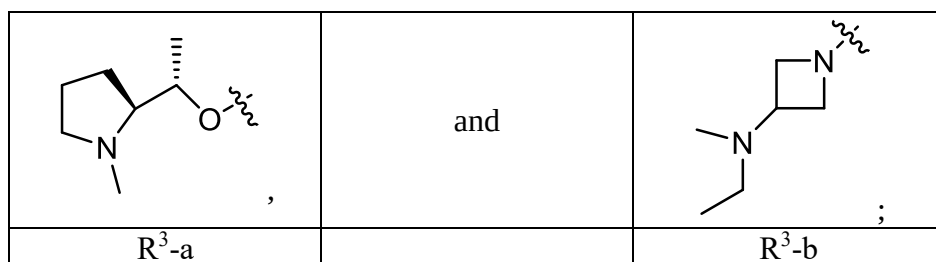
R¹ is selected from Cl, CH₃, CH₂F, CHF₂, and CF₃;

Cy¹ is selected from

		and	
Cy¹-a	Cy¹-b		Cy¹-c

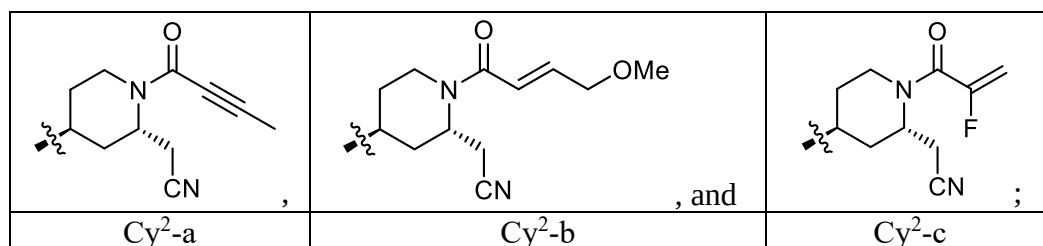
R² is selected from F and Cl;

R³ is selected from



and,

Cy² is selected from



provided that the compound of Formula (II) is other than,

8-(1-((2S,4S)-2-(cyanomethyl)-1-(2-fluoroacryloyl)piperidin-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile,
8-(1-((2S,4S)-2-(cyanomethyl)-1-((E)-4-methoxybut-2-enoyl)piperidin-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile,
2-((2S,4S)-4-(7-(5,6-dimethyl-1H-indazol-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-1-yl)-1-(2-fluoroacryloyl)piperidin-2-yl)acetonitrile,
2-((2S,4S)-1-(but-2-ynoyl)-4-(7-(5,6-dimethyl-1H-indazol-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-1-yl)piperidin-2-yl)acetonitrile, and
2-((2S,4S)-4-(7-(5,6-dimethyl-1H-indazol-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-1-yl)-1-((E)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile.

11. The use of claim 10, wherein the compound of Formula (II) or the pharmaceutically acceptable salt thereof is selected from:

1-(1-((2S,4S)-1-(but-2-ynoyl)-2-(cyanomethyl)piperidin-4-yl)-8-chloro-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)isoquinoline-8-carbonitrile;

1-(8-chloro-1-((2S,4S)-2-(cyanomethyl)-1-((E)-4-methoxybut-2-enoyl)piperidin-4-yl)-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)isoquinoline-8-carbonitrile;

1-(8-chloro-1-((2S,4S)-2-(cyanomethyl)-1-(2-fluoroacryloyl)piperidin-4-yl)-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)isoquinoline-8-carbonitrile;

2-((2S,4S)-4-(8-chloro-7-(5,6-dimethyl-1H-indazol-4-yl)-4-(3-(ethyl(methyl)amino)azetidin-1-yl)-6-fluoro-1H-pyrazolo[4,3-c]quinolin-1-yl)-1-(2-fluoroacryloyl)piperidin-2-yl)acetonitrile;

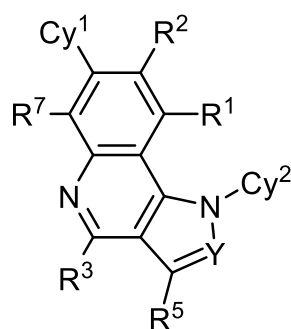
8-(8-chloro-1-((2S,4S)-2-(cyanomethyl)-1-(2-fluoroacryloyl)piperidin-4-yl)-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile;

8-(8-chloro-1-((2S,4S)-2-(cyanomethyl)-1-((E)-4-methoxybut-2-enoyl)piperidin-4-yl)-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile; and

8-(1-((2S,4S)-1-(but-2-ynoyl)-2-(cyanomethyl)piperidin-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile;

or a pharmaceutically acceptable salt thereof.

12. The use of any one of claims 1-7, wherein the KRAS inhibitor a compound of Formula (III):



III

or a pharmaceutically acceptable salt thereof, wherein:

Y is N or CR⁶;

R¹ is selected from H, C₁₋₃ alkyl, C₁₋₃ haloalkyl, cyclopropyl, halo, D, CN, and OR^{a1}; wherein said C₁₋₃ alkyl and cyclopropyl are each optionally substituted with 1 or 2 substituents independently selected from R^g;

R² is selected from H, C₁₋₃ alkyl, C₁₋₃ haloalkyl, 4-6 membered heterocycloalkyl, phenyl, 5-6 membered heteroaryl, 4-6 membered heterocycloalkyl-C₁₋₃ alkylene, phenyl-C₁₋₃

alkylene, 5-6 membered heteroaryl-C₁₋₃ alkylene, halo, D, CN, and OR^{a2}; wherein said C₁₋₃ alkyl, 4-6 membered heterocycloalkyl, phenyl, 5-6 membered heteroaryl, 4-6 membered heterocycloalkyl-C₁₋₃ alkylene, phenyl-C₁₋₃ alkylene, 5-6 membered heteroaryl-C₁₋₃ alkylene are each optionally substituted with 1 or 2 substituents independently selected from R^g;

Cy¹ is selected from C₃₋₁₀ cycloalkyl, 4-10 membered heterocycloalkyl, C₆₋₁₀ aryl and 6-10 membered heteroaryl; wherein the 4-10 membered heterocycloalkyl and 6-10 membered heteroaryl each has at least one ring-forming carbon atom and 1, 2, 3, or 4 ring-forming heteroatoms independently selected from N, O, and S; wherein a ring-forming carbon atom of 6-10 membered heteroaryl and 4-10 membered heterocycloalkyl is optionally substituted by oxo to form a carbonyl group; and wherein the C₃₋₁₀ cycloalkyl, 4-10 membered heterocycloalkyl, C₆₋₁₀ aryl and 6-10 membered heteroaryl are each optionally substituted with 1, 2, 3, or 4 substituents independently selected from R¹⁰;

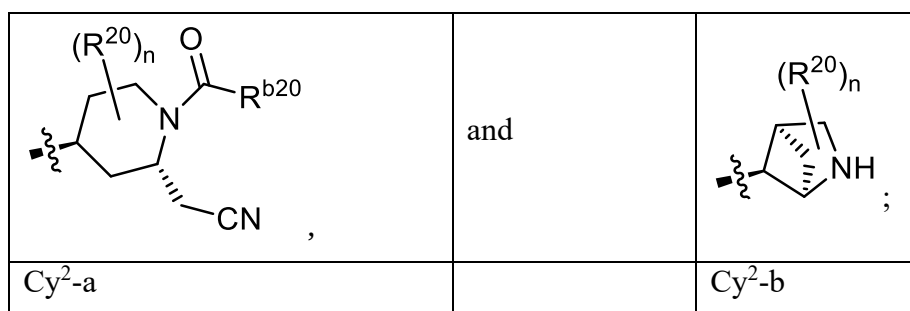
R³ is selected from H, C₁₋₃ alkyl, C₁₋₃ haloalkyl, C₃₋₆ cycloalkyl, 4-6 membered heterocycloalkyl, phenyl, 5-6 membered heteroaryl, C₃₋₆ cycloalkyl-C₁₋₃ alkylene, 4-6 membered heterocycloalkyl-C₁₋₃ alkylene, phenyl-C₁₋₃ alkylene, 5-6 membered heteroaryl-C₁₋₃ alkylene, halo, D, CN, OR^{f3}, C(O)NR^{c3}R^{d3}, NR^{c3}R^{j3}, and NR^{c3}C(O)R^{b3}; wherein said C₁₋₃ alkyl, C₃₋₆ cycloalkyl, 4-6 membered heterocycloalkyl, phenyl, 5-6 membered heteroaryl, C₃₋₆ cycloalkyl-C₁₋₃ alkylene, 4-6 membered heterocycloalkyl-C₁₋₃ alkylene, phenyl-C₁₋₃ alkylene, and 5-6 membered heteroaryl-C₁₋₃ alkylene are each optionally substituted with 1, 2, or 3 substituents independently selected from R³⁰;

R⁵ is selected from H, C₁₋₃ alkyl, C₁₋₃ haloalkyl, cyclopropyl, halo, D, CN, and OR^{a5}; wherein said C₁₋₃ alkyl and cyclopropyl are each optionally substituted with 1 or 2 substituents independently selected from R^g;

R⁶ is selected from H, C₁₋₃ alkyl, C₁₋₃ haloalkyl, C₃₋₆ cycloalkyl, 4-9 membered heterocycloalkyl, phenyl, 5-6 membered heteroaryl, C₃₋₆ cycloalkyl-C₁₋₃ alkylene, 4-6 membered heterocycloalkyl-C₁₋₃ alkylene, phenyl-C₁₋₃ alkylene, 5-6 membered heteroaryl-C₁₋₃ alkylene, halo, D, CN, OR^{a6}, and C(O)NR^{c6}R^{d6}; wherein said C₁₋₃ alkyl, C₃₋₆ cycloalkyl, 4-9 membered heterocycloalkyl, phenyl, 5-6 membered heteroaryl, C₃₋₆ cycloalkyl-C₁₋₃ alkylene, 4-6 membered heterocycloalkyl-C₁₋₃ alkylene, phenyl-C₁₋₃ alkylene, and 5-6 membered heteroaryl-C₁₋₃ alkylene are each optionally substituted with 1 or 2 substituents independently selected from R⁶⁰;

R⁷ is selected from H, C₁₋₃ alkyl, C₁₋₃ haloalkyl, cyclopropyl, halo, D, CN, and OR^{a7}; wherein said C₁₋₃ alkyl and cyclopropyl are each optionally substituted with 1 or 2 substituents independently selected from R^g;

Cy² is selected from



wherein n is 0, 1, or 2;

each R¹⁰ is independently selected from C₁₋₃ alkyl, C₁₋₃ haloalkyl, halo, D, CN, OR^{a10}, C(O)R^{b10}, C(O)NR^{c10}R^{d10}, C(O)OR^{a10}, NR^{c10}R^{d10}, and S(O)₂R^{b10};

each R²⁰ is independently selected from C₁₋₃ alkyl, C₁₋₃ haloalkyl, halo, D, CN, and OR^{a20};

each R³⁰ is independently selected from C₁₋₃ alkyl, C₁₋₃ haloalkyl, C₃₋₆ cycloalkyl, 4-6 membered heterocycloalkyl, phenyl, 5-6 membered heteroaryl, halo, D, CN, OR^{a30}, C(O)R^{b30}, C(O)NR^{c30}R^{d30}, C(O)OR^{a30}, NR^{c30}R^{d30}, and S(O)₂R^{b30}; wherein said C₁₋₃ alkyl, C₃₋₆ cycloalkyl, 4-6 membered heterocycloalkyl, phenyl, and 5-6 membered heteroaryl are each optionally substituted with 1 or 2 substituents independently selected from R³¹;

each R³¹ is independently selected from C₁₋₃ alkyl, C₁₋₃ haloalkyl, halo, D, CN, OR^{a31}, C(O)R^{b31}, C(O)NR^{c31}R^{d31}, C(O)OR^{a31}, NR^{c31}R^{d31}, and S(O)₂R^{b31};

each R³³ is independently selected from C₁₋₃ alkyl, C₁₋₃ haloalkyl, C₃₋₆ cycloalkyl, 4-membered heterocycloalkyl, 6-membered heterocycloalkyl, phenyl, 5-6 membered heteroaryl, halo, D, CN, OR^{a30}, C(O)NR^{c30}R^{d30}, and NR^{c30}R^{d30}; wherein said C₁₋₃ alkyl, C₃₋₆ cycloalkyl, 4-membered heterocycloalkyl, 6-membered heterocycloalkyl, phenyl, and 5-6 membered heteroaryl are each optionally substituted with 1 or 2 substituents independently selected from R³¹;

each R⁶⁰ is independently selected from C₁₋₃ alkyl, C₁₋₃ haloalkyl, 4-6 membered heterocycloalkyl, 5-6 membered heteroaryl, halo, D, CN, OR^{a60}, C(O)R^{b60}, C(O)NR^{c60}R^{d60}, NR^{c60}C(O)R^{b60}, C(O)OR^{a60}, NR^{c60}C(O)OR^{a60}, NR^{c60}R^{d60}, NR^{c60}S(O)₂R^{b60}, and S(O)₂R^{b60}; wherein said C₁₋₃ alkyl, 4-6 membered heterocycloalkyl, and 5-6 membered heteroaryl are each optionally substituted with 1 or 2 substituents independently selected from R⁶¹;

each R⁶¹ is independently selected from C₁₋₃ alkyl, C₁₋₃ haloalkyl, halo, D, CN, OR^{a61}, and NR^{c61}R^{d61};

R^{a1} is selected from H, C₁₋₃ alkyl, and C₁₋₃ haloalkyl;

each R^{a2} is independently selected from H, C₁₋₃ alkyl, and C₁₋₃ haloalkyl;

each R^{b3} , R^{c3} and R^{d3} is independently selected from H, C_{1-3} alkyl, C_{1-3} haloalkyl, C_{3-6} cycloalkyl, 4-6 membered heterocycloalkyl, phenyl and 5-6 membered heteroaryl; wherein said C_{1-3} alkyl, C_{3-6} cycloalkyl, 4-6 membered heterocycloalkyl, phenyl and 5-6 membered heteroaryl are each optionally substituted with 1, 2, or 3 substituents independently selected from R^{30} ;

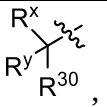
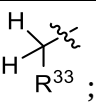
or R^{c3} and R^{d3} attached to the same N atom, together with the N atom to which they are attached, form a 4-, 5-, or 6-membered heterocycloalkyl group optionally substituted with 1, 2, or 3 substituents independently selected from R^{30} ;

R^{j3} is selected from C_{1-3} alkyl, C_{1-3} haloalkyl, C_{3-6} cycloalkyl, 4-6 membered heterocycloalkyl, phenyl and 5-6 membered heteroaryl; wherein said C_{1-3} alkyl, C_{3-6} cycloalkyl, 4-6 membered heterocycloalkyl, phenyl and 5-6 membered heteroaryl are each optionally substituted with 1, 2, or 3 substituents independently selected from R^{30} ;

or R^{c3} and R^{j3} attached to the same N atom, together with the N atom to which they are attached, form a 4-, 5-, or 6-membered heterocycloalkyl group optionally substituted with 1, 2, or 3 substituents independently selected from R^{30} ;

R^{f3} is selected from C_{1-3} haloalkyl, C_{3-6} cycloalkyl, 4-6 membered heterocycloalkyl, phenyl and 5-6 membered heteroaryl; wherein said C_{1-3} haloalkyl, C_{3-6} cycloalkyl, 4-6 membered heterocycloalkyl, phenyl, and 5-6 membered heteroaryl are each optionally substituted with 1, 2, or 3 substituents independently selected from R^{30} ; or

R^{f3} is selected from

	and	
R^{f3} -a		R^{f3} -b

wherein R^x is H or C_{1-2} alkyl and R^y is C_{1-2} alkyl;

or R^x and R^y , together with the C atom to which they are attached, form a 3-, or 4-membered cycloalkyl group;

R^{a5} is selected from H, C_{1-3} alkyl, and C_{1-3} haloalkyl;

each R^{a6} , R^{c6} and R^{d6} is independently selected from H, C_{1-3} alkyl, C_{1-3} haloalkyl, C_{3-6} cycloalkyl, 4-6 membered heterocycloalkyl, phenyl and 5-6 membered heteroaryl; wherein said C_{1-3} alkyl, C_{3-6} cycloalkyl, 4-6 membered heterocycloalkyl, phenyl and 5-6 membered heteroaryl are each optionally substituted with 1 or 2 substituents independently selected from R^{60} ;

R^{a7} is selected from H, C_{1-3} alkyl, and C_{1-3} haloalkyl;

each R^{a10}, R^{b10}, R^{c10} and R^{d10} is independently selected from H, C₁₋₃ alkyl, and C₁₋₃ haloalkyl;

each R^{a20} is independently selected from H, C₁₋₃ alkyl, and C₁₋₃ haloalkyl;

R^{b20} is selected from NH₂, C₁₋₃ alkyl, and C₁₋₃ haloalkyl;

each R^{a30}, R^{b30}, R^{c30} and R^{d30} is independently selected from H, C₁₋₃ alkyl, and C₁₋₃ haloalkyl;

each R^{a31}, R^{b31}, R^{c31} and R^{d31} is independently selected from H, C₁₋₃ alkyl, and C₁₋₃ haloalkyl;

each R^{a60}, R^{b60}, R^{c60} and R^{d60} is independently selected from H, C₁₋₃ alkyl, C₁₋₃ haloalkyl, C₃₋₆ cycloalkyl, 4-6 membered heterocycloalkyl, and 5-6 membered heteroaryl; wherein said C₁₋₃ alkyl, C₃₋₆ cycloalkyl, 4-6 membered heterocycloalkyl, and 5-6 membered heteroaryl are each optionally substituted with 1 or 2 substituents independently selected from R⁶¹;

or any R^{c60} and R^{d60} attached to the same N atom, together with the N atom to which they are attached, form a 4-, 5-, or 6-membered heterocycloalkyl group optionally substituted with 1 or 2 substituents independently selected from R⁶¹; and

each R^{a61}, R^{c61}, and R^{d61}, is independently selected from H, C₁₋₃ alkyl, and C₁₋₃ haloalkyl; and

each R^g is independently selected from D, OH, CN, halo, C₁₋₃ alkyl, C₁₋₃ haloalkyl, C₁₋₃ alkoxy, C₁₋₃ haloalkoxy, amino, C₁₋₃ alkylamino, and di(C₁₋₃ alkyl)amino;

provided that the compound of Formula (III) is other than,

3-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-4-ethoxy-6-fluoro-7-(3-hydroxynaphthalen-1-yl)-1H-pyrrolo[3,2-c]quinolin-2-yl)-N,N-dimethylpropanamide.

13. The use of claim 12, wherein the KRAS inhibitor is selected from:

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(7-chloro-3-hydroxynaphthalen-1-yl)-6-fluoro-2-methyl-4-(1H-1,2,4-triazol-1-yl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(5,7-difluoro-1H-indol-3-yl)-6-fluoro-2-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(6-fluoro-5-methyl-1*H*-indol-3-yl)-2-methyl-4-((*S*)-1-((*S*)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

3-(2-(3-(Azetidin-1-yl)-3-oxopropyl)-1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichlorophenyl)-6-fluoro-4-((*S*)-1-((*S*)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

3-((1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(3-hydroxynaphthalen-1-yl)-8-methyl-4-((*S*)-1-((*S*)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)methyl)oxazolidin-2-one;

8-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-2,8-dimethyl-4-((*S*)-1-((*S*)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-pyrrolo[3,2-*c*]quinolin-7-yl)-1-naphthonitrile;

1-((2*S*,4*S*)-1-Acetyl-2-(cyanomethyl)piperidin-4-yl)-7-(8-cyanonaphthalen-1-yl)-6-fluoro-4-((*S*)-1-((*S*)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-pyrazolo[4,3-*c*]quinoline-8-carbonitrile;

8-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-8-methyl-4-((*S*)-1-((*S*)-1-methylpyrrolidin-2-yl)ethoxy)-2-((3-oxomorpholino)methyl)-1*H*-pyrrolo[3,2-*c*]quinolin-7-yl)-1-naphthonitrile;

3-(7-(Benzo[*b*]thiophen-3-yl)-1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-4-((*S*)-1-((*S*)-1-methylpyrrolidin-2-yl)ethoxy)-2-((2-oxopyrrolidin-1-yl)methyl)-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

3-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-4-(((*S*)-1-(dimethylamino)propan-2-yl)oxy)-6-fluoro-7-(7-fluoronaphthalen-1-yl)-2-((2-oxopyrrolidin-1-yl)methyl)-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

8-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-6-fluoro-2-methyl-4-((*S*)-1-((*S*)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-pyrrolo[3,2-*c*]quinolin-7-yl)-1-naphthonitrile;

3-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichloro-5-hydroxyphenyl)-6-fluoro-2-methyl-4-((*S*)-1-((*S*)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

3-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-6-fluoro-4-((3-fluoro-1-methylazetidin-3-yl)methoxy)-7-(3-hydroxynaphthalen-1-yl)-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)-*N,N*-dimethylpropanamide;

3-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(3-hydroxynaphthalen-1-yl)-2-methyl-4-(5-methylpyrazin-2-yl)-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(7-fluoronaphthalen-1-yl)-4-methyl-2-((4-methyl-2-oxopiperazin-1-yl)methyl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichloro-5-hydroxyphenyl)-4-ethoxy-6-fluoro-2-((4-isopropyl-2-oxopiperazin-1-yl)methyl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-7-(7-fluoronaphthalen-1-yl)-2-((3-oxomorpholino)methyl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-4-ethoxy-6-fluoro-7-(3-hydroxynaphthalen-1-yl)-2-(1-(3-oxomorpholino)ethyl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((endo)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(3-hydroxynaphthalen-1-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-2-(pyridin-3-yl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(2-(3-(azetidin-1-yl)-3-oxopropyl)-1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-7-(7,8-difluoronaphthalen-1-yl)-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(2-(3-(azetidin-1-yl)-3-oxopropyl)-1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-7-(6,7-difluoronaphthalen-1-yl)-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(7-fluoro-3-hydroxynaphthalen-1-yl)-2-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

1-(1-((2S,4S)-1-Acetyl-2-(cyanomethyl)piperidin-4-yl)-8-chloro-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)isoquinoline-8-carbonitrile;

8-(1-((2S,4S)-1-acetyl-2-(cyanomethyl)piperidin-4-yl)-8-chloro-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrrolo[3,2-c]quinolin-7-yl)-1-naphthonitrile;

8-(1-((2S,4S)-1-acetyl-2-(cyanomethyl)piperidin-4-yl)-8-chloro-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(7-fluoro-3-hydroxynaphthalen-1-yl)-2-methyl-4-(1H-1,2,4-triazol-1-yl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(7-fluoronaphthalen-1-yl)-2-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

(2R)-2-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(1H-1,2,4-triazol-1-yl)-1H-pyrrolo[3,2-c]quinolin-2-yl)-N,N-dimethylpyrrolidine-1-carboxamide;

methyl (2R)-2-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-7-(2-chloro-3-methylphenyl)-8-(2-cyanoethyl)-6-fluoro-4-(1H-1,2,4-triazol-1-yl)-1H-pyrrolo[3,2-c]quinolin-2-yl)pyrrolidine-1-carboxylate;

Methyl (1S,3R,5S)-3-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-4-(6-(dimethylcarbamoyl)pyridin-3-yl)-6-fluoro-1H-pyrrolo[3,2-c]quinolin-2-yl)-2-azabicyclo[3.1.0]hexane-2-carboxylate;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-2-(5-oxo-1,2,3,5-tetrahydroindolizin-3-yl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

Methyl (2R)-2-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-4-(6-(dimethylcarbamoyl)pyridin-3-yl)-6-fluoro-1H-pyrrolo[3,2-c]quinolin-2-yl)pyrrolidine-1-carboxylate;

Methyl (2R)-2-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(6-(methylcarbamoyl)pyridin-3-yl)-1H-pyrrolo[3,2-c]quinolin-2-yl)pyrrolidine-1-carboxylate;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(2-chloro-3-fluorophenyl)-2-((R)-1-(cyclopropanecarbonyl)pyrrolidin-2-yl)-6-fluoro-4-methyl-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

8-(2-((R)-1-Acetylpyrrolidin-2-yl)-1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-8-methyl-4-(2-methylpyridin-4-yl)-1H-pyrrolo[3,2-c]quinolin-7-yl)-1,2,3,4-tetrahydronaphthalene-1-carbonitrile;

5-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(3-chloro-2-methylphenyl)-8-(2-cyanoethyl)-6-fluoro-2-((R)-1-(2-oxopyrazin-1(2H)-yl)ethyl)-1H-pyrrolo[3,2-c]quinolin-4-yl)-N-methylpicolinamide;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(7-fluoronaphthalen-1-yl)-4-(6-(2-hydroxypropan-2-yl)pyridin-3-yl)-2-((R)-1-(2-oxopyrazin-1(2H)-yl)ethyl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(3-chloro-2-methylphenyl)-6-fluoro-4-(5-methylpyrazin-2-yl)-2-((R)-1-(2-oxopyrazin-1(2H)-yl)ethyl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

Methyl (2R)-2-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(5-fluoro-6-(methylcarbamoyl)pyridin-3-yl)-1H-pyrrolo[3,2-c]quinolin-2-yl)pyrrolidine-1-carboxylate;

Methyl (2R)-2-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(6-(2-hydroxypropan-2-yl)pyridin-3-yl)-1H-pyrrolo[3,2-c]quinolin-2-yl)pyrrolidine-1-carboxylate;

Ethyl (2R)-2-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(6-(2-hydroxypropan-2-yl)pyridin-3-yl)-1H-pyrrolo[3,2-c]quinolin-2-yl)pyrrolidine-1-carboxylate;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichlorophenyl)-2-((R)-1-(3,3-difluoroazetidine-1-carbonyl)pyrrolidin-2-yl)-6-fluoro-4-(methyl-d3)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichlorophenyl)-2-((R)-1-(3,3-difluoroazetidine-1-carbonyl)pyrrolidin-2-yl)-6-fluoro-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(3-chloro-2-methylphenyl)-6-fluoro-4-(5-methylpyrazin-2-yl)-2-((R)-1-(3-oxomorpholino)ethyl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

5-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-6-fluoro-7-(7-fluoronaphthalen-1-yl)-2-((R)-1-(3-oxomorpholino)ethyl)-1H-pyrrolo[3,2-c]quinolin-4-yl)-N-methylpicolinamide;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(7-fluoronaphthalen-1-yl)-4-(6-(2-hydroxypropan-2-yl)pyridin-3-yl)-2-((R)-1-(3-oxomorpholino)ethyl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

3-(1-((1R,4R,5S)-2-Azabicyclo[2.1.1]hexan-5-yl)-6-fluoro-7-(7-fluoronaphthalen-1-yl)-4-(5-methylpyrazin-2-yl)-2-((R)-1-(3-oxomorpholino)ethyl)-1H-pyrrolo[3,2-c]quinolin-8-yl)propanenitrile;

Methyl (1R,3R,5R)-3-(1-((1R,4R,5S)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-4-(6-(dimethylcarbamoyl)pyridin-3-yl)-6-fluoro-1H-pyrrolo[3,2-c]quinolin-2-yl)-2-azabicyclo[3.1.0]hexane-2-carboxylate;

3-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-2-((1*R*,3*R*,5*R*)-2-(cyclopropanecarbonyl)-2-azabicyclo[3.1.0]hexan-3-yl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

3-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(6-(2-hydroxypropan-2-yl)pyridin-3-yl)-2-((*R*)-1-(2-oxopyrazin-1(2*H*)-yl)ethyl)-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

Methyl (2*R*,4*S*)-2-(1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(6-(2-hydroxypropan-2-yl)pyridin-3-yl)-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)-4-fluoropyrrolidine-1-carboxylate;

Methyl (2*R*,5*R*)-2-(1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(6-(2-hydroxypropan-2-yl)pyridin-3-yl)-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)-5-methylpyrrolidine-1-carboxylate;

Methyl (2*R*)-2-(1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-3-chloro-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(6-(2-hydroxypropan-2-yl)pyridin-3-yl)-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)pyrrolidine-1-carboxylate;

4-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-2-((*R*)-1-(2-oxopyrazin-1(2*H*)-yl)ethyl)-1*H*-pyrrolo[3,2-*c*]quinolin-4-yl)-2-fluoro-*N*-methylbenzamide;

Methyl ((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)carbamate;
N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)-2,2-difluoroacetamide;

N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)-2,2-difluoroacetamide;

(2*S*)-*N*-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)tetrahydrofuran-2-carboxamide;

N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)cyclopropanesulfonamide;

N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)thiazole-4-carboxamide;

N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)-*N*-methylcyclopropanecarboxamide;

N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(1-hydroxyethyl)-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)-1-methylcyclopropane-1-carboxamide;

3-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(1-hydroxyethyl)-2-((1*R*,3*R*,5*R*)-2-(1-methylcyclopropane-1-carbonyl)-2-azabicyclo[3.1.0]hexan-3-yl)-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

3-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichlorophenyl)-6-fluoro-2-((1*R*,3*R*,5*R*)-2-(1-fluorocyclopropane-1-carbonyl)-2-azabicyclo[3.1.0]hexan-3-yl)-4-(1-hydroxyethyl)-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

3-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichlorophenyl)-6-fluoro-2-((1*R*,3*R*,5*R*)-2-(1-fluorocyclopropane-1-carbonyl)-2-azabicyclo[3.1.0]hexan-3-yl)-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(1-hydroxyethyl)-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)-1-fluorocyclopropane-1-carboxamide;

N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-(1-hydroxyethyl)-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)-1-fluorocyclobutane-1-carboxamide;

3-(1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-7-(3-chloro-2-methylphenyl)-2-(1-(2,6-dimethyl-3-oxo-2,3-dihydropyridazin-4-yl)ethyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)pyrimidine-4-carboxamide;

N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)pyridazine-3-carboxamide;

N-((1*R*)-1-(1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)ethyl)-3,3-difluoroazetidine-1-carboxamide;

3-(1-((1*R*,4*R*,5*S*)-2-azabicyclo[2.1.1]hexan-5-yl)-7-(2,3-dichlorophenyl)-6-fluoro-4-methyl-2-((*R*)-1-((1-methyl-1*H*-pyrazol-4-yl)amino)ethyl)-1*H*-pyrrolo[3,2-*c*]quinolin-8-yl)propanenitrile;

5-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-6-fluoro-2-((*R*)-1-(1-fluorocyclopropane-1-carbonyl)pyrrolidin-2-yl)-1*H*-pyrrolo[3,2-*c*]quinolin-4-yl)-*N,N*-dimethylpicolinamide; and

methyl (2*R*)-2-(1-((1*R*,4*R*,5*S*)-2-Azabicyclo[2.1.1]hexan-5-yl)-8-(2-cyanoethyl)-7-(2,3-dichlorophenyl)-4-(4-((dimethylamino)methyl)-2,3-difluorophenyl)-6-fluoro-1*H*-pyrrolo[3,2-*c*]quinolin-2-yl)pyrrolidine-1-carboxylate,

or a pharmaceutically acceptable salt thereof.

14. The use of any one of claims 1-7, wherein the KRAS inhibitor is selected from sotorasib, adagrasib, Compound 2, Compound 3 and Compound 4, or a pharmaceutically acceptable salt thereof.

15. The use of any one of claims 1-7, wherein the KRAS inhibitor is selected from sotorasib, Compound 2, Compound 3 and Compound 4, or a pharmaceutically acceptable salt thereof.

16. The use of any one of claims 1-7, wherein the KRAS inhibitor is selected from Compound 5, Compound 6, and Compound 7.

17. The use of any one of claims 1-7, wherein the KRAS inhibitor is sotorasib, or a pharmaceutically acceptable salt thereof.

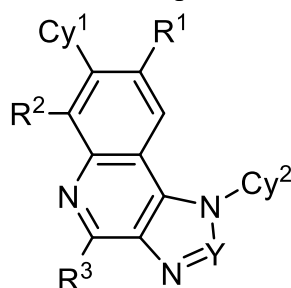
18. The use of any one of claims 1-7, wherein the KRAS inhibitor is adagrasib, or a pharmaceutically acceptable salt thereof.

19. The use of any one of claims 1-7, wherein the KRAS inhibitor is Compound 2, or a pharmaceutically acceptable salt thereof.

20. The use of any one of claims 1-7, wherein the KRAS inhibitor is Compound 3, or a pharmaceutically acceptable salt thereof.
21. The use of any one of claims 1-7, wherein the KRAS inhibitor is Compound 4, or a pharmaceutically acceptable salt thereof.
22. The use of any one of claims 1-11, 14, 15, and 17-21, wherein the KRAS inhibitor is a KRAS G12C inhibitor.
23. The use of any one of claims 1-7, 12, 13, and 16, wherein the KRAS inhibitor is a KRAS G12D inhibitor.

24. Use of:

- (i) pemigatinib, or a pharmaceutically acceptable salt thereof; and
- (ii) a KRAS inhibitor, which is a compound of Formula (I):



I

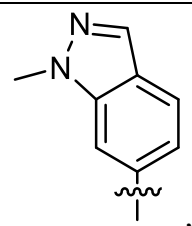
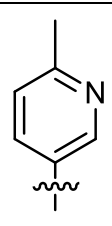
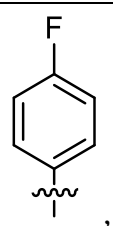
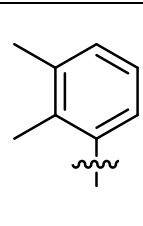
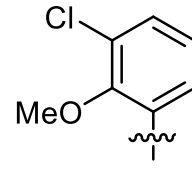
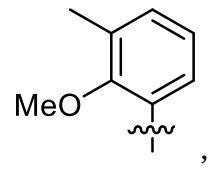
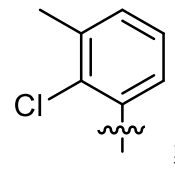
or a pharmaceutically acceptable salt thereof, wherein:

Y is N or CH;

R¹ is selected from Cl, CH₃, CH₂F, CHF₂, and CF₃;

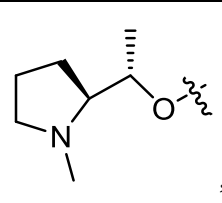
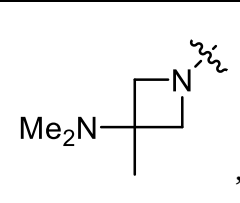
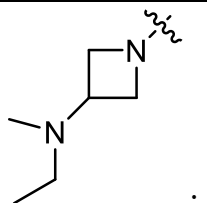
Cy¹ is selected from

Cy ¹ -a	Cy ¹ -b	Cy ¹ -c	Cy ¹ -d

			
Cy ¹ -e	Cy ¹ -f	Cy ¹ -g	Cy ¹ -h
		and	
Cy ¹ -i	Cy ¹ -j		Cy ¹ -k

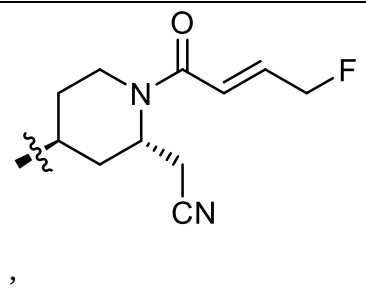
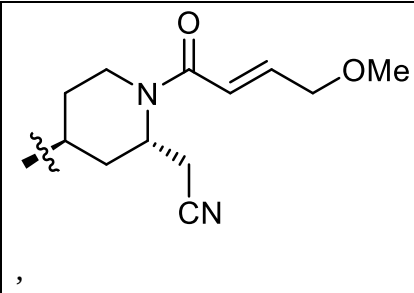
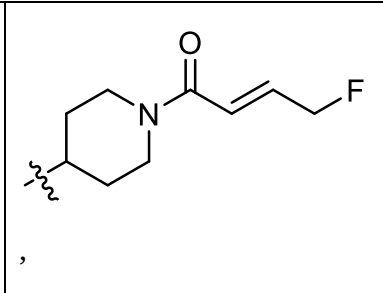
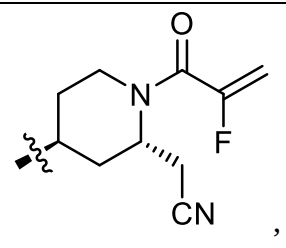
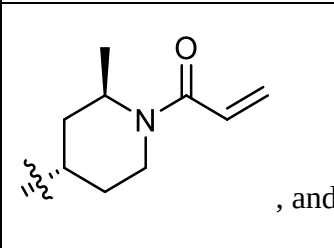
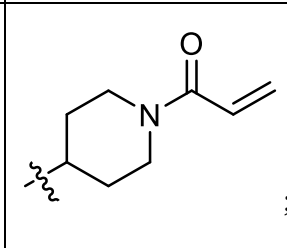
R² is selected from F and Cl;

R³ is selected from

		and	
R ³ -a	R ³ -b		R ³ -c

and,

Cy² is selected from

		
Cy ² -a	Cy ² -b	Cy ² -c
		
Cy ² -d	Cy ² -e	Cy ² -f

provided that the compound of Formula (I) is other than

2-((2S,4S)-4-(7-(2-chloro-3-methylphenyl)-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-8-methyl-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile and

2-((2S,4S)-4-(4-(3-(dimethylamino)-3-methylazetidin-1-yl)-7-(2,3-dimethylphenyl)-6-fluoro-8-methyl-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

in the manufacture of a medicament for the treatment of cancer in a patient.

25. The use of claim 24, wherein the compound of Formula (I) or the pharmaceutically acceptable salt thereof is selected from:

2-((2S,4S)-4-(4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-7-(5-fluoroquinolin-8-yl)-8-(trifluoromethyl)-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-7-(2-methoxy-3-methylphenyl)-8-(trifluoromethyl)-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(7-(3-chloro-2-methoxyphenyl)-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-8-(trifluoromethyl)-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6-fluoro-7-(5-fluoroquinolin-8-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-8-(trifluoromethyl)-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

1-(4-(6-fluoro-7-(5-fluoroquinolin-8-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-8-(trifluoromethyl)-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)piperidin-1-yl)prop-2-en-1-one;

2-((2S,4S)-4-(4-(3-(dimethylamino)-3-methylazetidin-1-yl)-7-(2,3-dimethylphenyl)-6-fluoro-8-methyl-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6-fluoro-8-methyl-7-(1-methyl-1H-indazol-6-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-[1,2,3]triazolo[4,5-c]quinolin-1-yl)-1-((E)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6-fluoro-8-methyl-7-(6-methylpyridin-3-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6-fluoro-8-methyl-7-(1-methyl-1*H*-indazol-3-yl)-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6-fluoro-7-(4-fluorophenyl)-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;

8-(1-(1-acryloylpiperidin-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-7-yl)-1-naphthonitrile;

2-((2S,4S)-4-(7-(2-chloro-3-methylphenyl)-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-8-methyl-1*H*-imidazo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(7-(2-chloro-3-methylphenyl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-imidazo[4,5-*c*]quinolin-1-yl)-1-(2-fluoroacryloyl)piperidin-2-yl)acetonitrile;

8-(1-((2*R*,4*S*)-1-acryloyl-2-methylpiperidin-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-7-yl)-1-naphthonitrile;

2-((2S,4S)-4-(7-(5,6-dimethyl-1*H*-indazol-4-yl)-4-(3-(ethyl(methyl)amino)azetidin-1-yl)-6-fluoro-8-methyl-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-(2-fluoroacryloyl)piperidin-2-yl)acetonitrile;

8-(6-fluoro-1-(1-((*E*)-4-fluorobut-2-enoyl)piperidin-4-yl)-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-7-yl)-1-naphthonitrile;

8-(1-((2S,4S)-2-(cyanomethyl)-1-(2-fluoroacryloyl)piperidin-4-yl)-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-8-methyl-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-7-yl)-1-naphthonitrile;

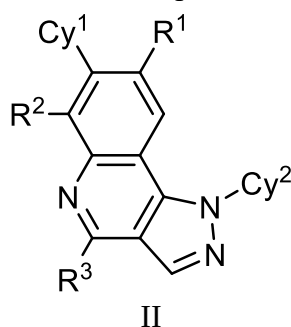
2-((2S,4S)-4-(6,8-dichloro-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-7-(5-fluoroquinolin-8-yl)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;

2-((2S,4S)-4-(6,8-dichloro-4-(3-(dimethylamino)-3-methylazetidin-1-yl)-7-(5-fluoroquinolin-8-yl)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-fluorobut-2-enoyl)piperidin-2-yl)acetonitrile; and

2-((2*S*,4*S*)-4-(4-(3-(dimethylamino)-3-methylazetidin-1-yl)-6-fluoro-7-(5-fluoroquinolin-8-yl)-8-(trifluoromethyl)-1*H*-[1,2,3]triazolo[4,5-*c*]quinolin-1-yl)-1-((*E*)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;
or a pharmaceutically acceptable salt thereof.

26. Use of:

- (i) pemigatinib, or a pharmaceutically acceptable salt thereof; and
- (ii) a KRAS inhibitor, which is a compound of Formula (II):



or a pharmaceutically acceptable salt thereof, wherein:

R^1 is selected from Cl, CH₃, CH₂F, CHF₂, and CF₃;

Cy^1 is selected from

		and	
Cy^1 -a	Cy^1 -b		Cy^1 -c

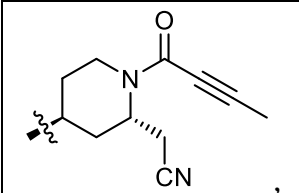
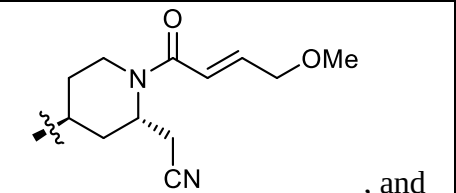
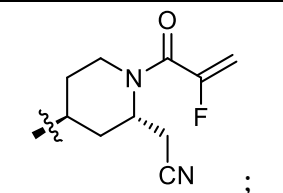
R^2 is selected from F and Cl;

R^3 is selected from

	and	
R^3 -a		R^3 -b

and,

Cy^2 is selected from

		
Cy ² -a	Cy ² -b	Cy ² -c

provided that the compound of Formula (II) is other than,

8-(1-((2S,4S)-2-(cyanomethyl)-1-(2-fluoroacryloyl)piperidin-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile,
8-(1-((2S,4S)-2-(cyanomethyl)-1-((E)-4-methoxybut-2-enoyl)piperidin-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile,
2-((2S,4S)-4-(7-(5,6-dimethyl-1H-indazol-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-1-yl)-1-(2-fluoroacryloyl)piperidin-2-yl)acetonitrile,
2-((2S,4S)-1-(but-2-ynoyl)-4-(7-(5,6-dimethyl-1H-indazol-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-1-yl)piperidin-2-yl)acetonitrile, and
2-((2S,4S)-4-(7-(5,6-dimethyl-1H-indazol-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-1-yl)-1-((E)-4-methoxybut-2-enoyl)piperidin-2-yl)acetonitrile;

in the manufacture of a medicament for the treatment of cancer in a patient.

27. The use of claim 26, wherein the compound of Formula (II) or the pharmaceutically acceptable salt thereof is selected from:

1-(1-((2S,4S)-1-(but-2-ynoyl)-2-(cyanomethyl)piperidin-4-yl)-8-chloro-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)isoquinoline-8-carbonitrile;

1-(8-chloro-1-((2S,4S)-2-(cyanomethyl)-1-((E)-4-methoxybut-2-enoyl)piperidin-4-yl)-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)isoquinoline-8-carbonitrile;

1-(8-chloro-1-((2S,4S)-2-(cyanomethyl)-1-(2-fluoroacryloyl)piperidin-4-yl)-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)isoquinoline-8-carbonitrile;

2-((2S,4S)-4-(8-chloro-7-(5,6-dimethyl-1H-indazol-4-yl)-4-(3-(ethyl(methyl)amino)azetidin-1-yl)-6-fluoro-1H-pyrazolo[4,3-c]quinolin-1-yl)-1-(2-fluoroacryloyl)piperidin-2-yl)acetonitrile;

8-(8-chloro-1-((2S,4S)-2-(cyanomethyl)-1-(2-fluoroacryloyl)piperidin-4-yl)-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile;

8-(8-chloro-1-((2S,4S)-2-(cyanomethyl)-1-((E)-4-methoxybut-2-enoyl)piperidin-4-yl)-6-fluoro-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile; and

8-(1-((2S,4S)-1-(but-2-ynoyl)-2-(cyanomethyl)piperidin-4-yl)-6-fluoro-8-methyl-4-((S)-1-((S)-1-methylpyrrolidin-2-yl)ethoxy)-1H-pyrazolo[4,3-c]quinolin-7-yl)-1-naphthonitrile;

or a pharmaceutically acceptable salt thereof.

28. Use of:

(i) an FGFR1 inhibitor selected from pemigatinib, futibatinib, erdafitinib and infigratinib, or a pharmaceutically acceptable salt thereof; and

(ii) a KRAS inhibitor selected from sotorasib, adagrasib, Compound 2, Compound 3 and Compound 4, or a pharmaceutically acceptable salt thereof;
in the manufacture of a medicament for the treatment of cancer in a patient.

29. The use of claim 28, wherein the FGFR1 inhibitor is pemigatinib, or a pharmaceutically acceptable salt thereof.

30. The use of claim 28, wherein the FGFR1 inhibitor is futibatinib, or a pharmaceutically acceptable salt thereof.

31. The use of claim 28, wherein the FGFR1 inhibitor is erdafitinib, or a pharmaceutically acceptable salt thereof.

32. The use of claim 28, wherein the FGFR1 inhibitor is infigratinib, or a pharmaceutically acceptable salt thereof.

33. The use of any one of claims 28-32, wherein the KRAS inhibitor is sotorasib, or a pharmaceutically acceptable salt thereof.

34. The use of any one of claims 28-32, wherein the KRAS inhibitor is adagrasib, or a pharmaceutically acceptable salt thereof.
35. The use of any one of claims 28-32, wherein the KRAS inhibitor is Compound 2, or a pharmaceutically acceptable salt thereof.
36. The use of any one of claims 28-32, wherein the KRAS inhibitor is Compound 3, or a pharmaceutically acceptable salt thereof.
37. The use of any one of claims 28-32, wherein the KRAS inhibitor is Compound 4, or a pharmaceutically acceptable salt thereof.
38. Use of:
- (i) pemigatinib, or a pharmaceutically acceptable salt thereof; and
 - (ii) a KRAS inhibitor selected from sotorasib, adagrasib, Compound 2, Compound 3 and Compound 4, or a pharmaceutically acceptable salt thereof;
- in the manufacture of a medicament for the treatment of cancer in a patient.
39. The use of claim 38, wherein the KRAS inhibitor is sotorasib, or a pharmaceutically acceptable salt thereof.
40. The use of claim 38, wherein the KRAS inhibitor is adagrasib, or a pharmaceutically acceptable salt thereof.
41. The use of claim 38, wherein the KRAS inhibitor is Compound 2, or a pharmaceutically acceptable salt thereof.
42. The use of claim 38, wherein the KRAS inhibitor is Compound 3, or a pharmaceutically acceptable salt thereof.
43. The use of claim 38, wherein the KRAS inhibitor is Compound 4, or a pharmaceutically acceptable salt thereof.

44. Use of pemigatinib, or a pharmaceutically acceptable salt thereof, in the manufacture of a medicament for the treatment of cancer in a patient;
wherein the medicament is to be administered with sotorasib, or a pharmaceutically acceptable salt thereof.
45. The use of claim 44, wherein the medicament comprising pemigatinib, or a pharmaceutically acceptable salt thereof, and sotorasib or a pharmaceutically acceptable salt thereof, are to be administered simultaneously.
46. The use of claim 44, wherein the medicament comprising pemigatinib, or a pharmaceutically acceptable salt thereof, and sotorasib or a pharmaceutically acceptable salt thereof, are to be administered sequentially.
47. The use of any one of claims 44-46, wherein the medicament comprising pemigatinib, or a pharmaceutically acceptable salt thereof, is to be administered orally.
48. The use of claim 47, wherein the medicament comprising pemigatinib, or a pharmaceutically acceptable salt thereof, is to be administered in the form of a tablet.
49. The use of any one of claims 44-48, wherein the medicament comprising pemigatinib, or a pharmaceutically acceptable salt thereof, is to be administered in a daily dose of about 1 mg to about 10 mg.
50. The use of any one of claims 44-48, wherein the medicament comprising pemigatinib, or a pharmaceutically acceptable salt thereof, is to be administered in a daily dose of about 1 mg to about 5 mg.
51. The use of any one of claims 44-48, wherein the medicament comprising pemigatinib, or a pharmaceutically acceptable salt thereof, is to be administered in a daily dose of about 2 mg.
52. The use of any one of claims 44-51, wherein sotorasib, or a pharmaceutically acceptable salt thereof, is to be administered orally.

53. The use of claim 52, wherein sotorasib, or a pharmaceutically acceptable salt thereof, is to be administered in the form of a tablet.
54. The use of any one of claims 44-53, wherein sotorasib, or a pharmaceutically acceptable salt thereof, is to be administered in a daily dose of about 50 mg to about 300 mg.
55. The use of any one of claims 44-53, wherein sotorasib, or a pharmaceutically acceptable salt thereof, is to be administered in a daily dose of about 100 mg to about 200 mg.
56. The use of any one of claims 44-53, wherein sotorasib, or a pharmaceutically acceptable salt thereof, is to be administered in a daily dose of about 133 mg.
57. The use of any one of claims 44-53, wherein sotorasib, or a pharmaceutically acceptable salt thereof, is to be administered in a daily dose of about 120 mg.
58. The use of claim 44, wherein the medicament comprising pemigatinib, or a pharmaceutically acceptable salt thereof, and sotorasib, or a pharmaceutically acceptable salt thereof, are to be orally administered simultaneously in a daily dose of about 2 mg and 120 mg, respectively.
59. The use of claim 44, wherein the medicament comprising pemigatinib, or a pharmaceutically acceptable salt thereof, and sotorasib, or a pharmaceutically acceptable salt thereof, are to be orally administered sequentially in a daily dose of about 2 mg and 120 mg, respectively.
60. The use of claim 58 or 59, wherein the medicament comprising pemigatinib, or a pharmaceutically acceptable salt thereof, and sotorasib, or a pharmaceutically acceptable salt thereof, are each to be administered in the form of a tablet.
61. The use of any one of claims 1-60, wherein the medicament is to be further administered with one or more additional therapeutic agents.

62. The use of claim 61, wherein the medicament is to be administered with one additional therapeutic agent.
63. The use of claim 62, wherein the additional therapeutic agent is an inhibitor of PD-1 or PD-L1.
64. The use of any one of claims 1-63, wherein the cancer comprises one or more KRAS mutations.
65. The use of claim 64, wherein the one or more KRAS mutations comprise mutations selected from G12C, G12D, C12V and combinations thereof.
66. The use of claim 64 or 65, wherein the cancer further comprises high FGFR1 expression.
67. The use of any one of claims 1-66, wherein the cancer is selected from carcinomas, pancreatic cancer, colorectal cancer, lung cancer, non-small cell lung cancer, ovarian cancer, bladder cancer, gastric cancer, esophageal cancer, breast cancer, head and neck cancer, cervical cancer, skin cancer, thyroid cancer, hematopoietic malignancies, multiple myeloma, acute myelogenous leukemia, myeloproliferative neoplasms, neoplasms, glioblastoma and sarcomas.
68. The use of any one of claims 1-67, wherein the cancer is lung cancer.
69. The use of claim 68, wherein the lung cancer is non-small cell lung cancer.
70. The use of any one of claims 1-67, wherein the cancer is colorectal cancer.
71. The use of any one of claims 1-67, wherein the cancer is pancreatic cancer.
72. The use of any one of claims 1-67, wherein the cancer is ovarian cancer.