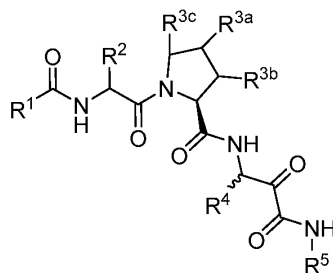


WHAT IS CLAIMED IS:

1. A compound of Formula (I):



(I)

or a pharmaceutically acceptable salt, solvate, solvate of the salt or prodrug thereof wherein:

R¹ is selected from the group consisting of:

- (a) -aryl and
- (b) -heteroaryl;

wherein the aryl and heteroaryl of choices (a) and (b) are each optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (i) -halogen,
- (ii) -CN,
- (iii) -C₁₋₆alkyl,
- (iv) -C₂₋₆alkenyl,
- (v) -C₂₋₆alkynyl,
- (vi) -C(O)R⁸,
- (vii) -CO₂R⁸,
- (viii) -CONR⁵R⁶,
- (ix) -OH,
- (x) -O-C₁₋₆alkyl,
- (xi) -SH,
- (xii) -S(O)_p-C₁₋₆alkyl,
- (xiii) -S(O)₂NR⁵R⁶,
- (xiv) -NO₂,
- (xv) -NR⁵R⁶,

- (xvi) $-\text{NHC}(\text{O})\text{R}^8$,
- (xvii) $-\text{NHC}(\text{O})\text{OR}^8$,
- (xviii) $-\text{NHC}(\text{O})\text{NR}^5\text{R}^6$, and
- (xix) $-\text{NHSO}_2\text{C}_{1-6}\text{alkyl}$,

wherein each of the alkyl group of choices (iii), (x), (xii) and (xix) is optionally substituted with 1 to 5 substituents independently selected from $-\text{halogen}$, $-\text{haloC}_{1-4}\text{alkyl}$, $-\text{COR}^8$, $-\text{CO}_2\text{R}^8$, $-\text{CONR}^5\text{R}^6$, $-\text{NR}^5\text{R}^6$, $-\text{OH}$, $-\text{O}-\text{C}_{1-4}\text{alkyl}$, $-\text{SH}$ and $-\text{S}-\text{C}_{1-4}\text{alkyl}$;

R^2 is selected from the group consisting of:

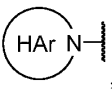
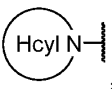
- (a) $-\text{C}_{3-8}\text{alkyl}$,
- (b) $-\text{C}_{0-6}\text{alkyl}-\text{R}^7$, and
- (c) $-(\text{CH}_2)_{1-6}-\text{N}(\text{R}^{13})(\text{R}^{13})$,

wherein each of the alkyl group of choices (a) and (b) is optionally substituted with 1 to 5 substituents independently selected from:

- (i) $-\text{halogen}$,
- (ii) $-\text{C}_{1-4}\text{alkyl}$,
- (iii) $-\text{haloC}_{1-4}\text{alkyl}$,
- (iv) $-\text{OH}$,
- (v) $-\text{O}-\text{C}_{1-4}\text{alkyl}$,
- (vi) $-\text{SH}$, and
- (vii) $-\text{S}-\text{C}_{1-4}\text{alkyl}$;

R^{3b} and R^{3c} together represent $-(\text{CH}_2)_{2-3}-$, and R^{3a} is H; or

R^{3b} and R^{3c} are each H, and R^{3a} is selected from the group consisting of:

- (a) $-\text{H}$,
- (b) $-\text{aryl}$,
- (c) , and
- (d) ,

wherein HAr is heteroaryl and Hcyl is heterocycle, wherein each of the aryl of choice (b), HAr and Hcyl is optionally substituted with 1 to 3 groups independently selected from the group consisting of:

- (i) -halogen,
- (ii) -OH,
- (iii) $-\text{CR}^{10}\text{R}^{11}\text{R}^{12}$,
- (iv) $-(\text{CH}_2)_{0-3}-\text{NHSO}_2-\text{C}_{1-4}\text{alkyl}$, and
- (v) $-(\text{CH}_2)_{0-3}-\text{SO}_2-\text{C}_{1-4}\text{alkyl}$;

R^4 is selected from the group consisting of:

- (a) $-\text{C}_{1-6}\text{alkyl}$,
- (b) $-\text{haloC}_{1-6}\text{alkyl}$,
- (c) $-\text{C}_{2-6}\text{alkenyl}$,
- (d) $-\text{C}_{2-6}\text{alkynyl}$,
- (e) $-\text{C}_{1-6}\text{alkyl}-\text{C}_{3-6}\text{cycloalkyl}$,
- (f) $-\text{C}_{1-6}\text{alkyl-aryl}$, wherein aryl is optionally substituted with nitro or $-\text{N}(\text{R}^{13})(\text{R}^{13})$,
- (g) $-\text{C}_{1-6}\text{alkyl}-\text{R}^9$, and
- (h) $-\text{haloC}_{1-6}\text{alkyl}-\text{R}^9$;

each R^5 and each R^6 are independently selected from the group consisting of:

- (a) -H,
- (b) $-\text{C}_{1-6}\text{alkyl}$,
- (c) $-\text{C}_{0-6}\text{alkyl}-\text{C}_{3-12}\text{cycloalkyl}$,
- (d) $-\text{C}_{0-6}\text{alkyl-heterocyclyl}$,
- (e) $-\text{C}_{0-6}\text{alkyl-heteroaryl}$, and
- (f) $-\text{C}_{0-6}\text{alkyl-aryl}$,

wherein the alkyl group of choices (b) – (f) is optionally substituted with 1 to 3 groups independently selected from:

- (i) -halogen,
- (ii) $-\text{C}(\text{O})\text{C}_{1-4}\text{alkyl}$,
- (iii) $-\text{C}(\text{O})\text{OC}_{1-4}\text{alkyl}$,
- (iv) -OH,

- (v) $-\text{OC}_{1-4}\text{alkyl}$,
- (vi) $-\text{SH}$,
- (vii) $-\text{SC}_{1-4}\text{alkyl}$,
- (viii) $-\text{NH}_2$,
- (ix) $-\text{NH}(\text{C}_{1-4}\text{alkyl})$, and
- (x) $-\text{N}(\text{C}_{1-4}\text{alkyl})(\text{C}_{1-4}\text{alkyl})$; or

R⁵, R⁶ and the nitrogen atom to which they are attached together form a 3- to 7-membered monocyclic or 6- to 11-membered bicyclic heterocycle optionally having an additional heteroatomic moiety selected from $-\text{O}-$, $-\text{S}(\text{O})_p-$, and $-\text{NR}^{13}-$, and wherein said heterocycle is optionally substituted with 1 to 2 groups independently selected from halogen, $-\text{haloC}_{1-4}\text{alkyl}$, $-\text{OH}$, $-\text{O}-\text{C}_{1-4}\text{alkyl}$, $-\text{SH}$ and $-\text{S}-\text{C}_{1-4}\text{alkyl}$;

R⁷ is selected from the group consisting of:

- (a) $-\text{C}_{3-10}\text{cycloalkyl}$, and
- (b) $-\text{C}_{4-10}\text{heterocyclyl}$,

wherein each of choices (a) and (b) is optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (i) $-\text{C}_{1-4}\text{alkyl}$,
- (ii) $-\text{halogen}$,
- (iii) $-\text{OH}$,
- (iv) $-\text{O}-\text{C}_{1-4}\text{alkyl}$,
- (v) $-\text{SH}$, and
- (vi) $-\text{S}-\text{C}_{1-4}\text{alkyl}$;

R⁸ is selected from the group consisting of:

- (a) $-\text{C}_{1-6}\text{alkyl}$,
- (b) $-\text{C}_{0-6}\text{alkyl}-\text{C}_{3-12}\text{cycloalkyl}$,
- (c) $-\text{C}_{0-6}\text{alkyl}-\text{heterocyclyl}$,
- (d) $-\text{C}_{0-6}\text{alkyl}-\text{heteroaryl}$, and
- (e) $-\text{C}_{0-6}\text{alkyl}-\text{aryl}$,

wherein each of the alkyl group of choices (a) – (e) is optionally substituted with 1 to 3 groups independently selected from:

- (i) $-\text{halogen}$,

- (ii) $-\text{OH}$,
- (iii) $-\text{OC}_{1-4}\text{alkyl}$,
- (iv) $-\text{SH}$, and
- (v) $-\text{SC}_{1-4}\text{alkyl}$;

R^9 is selected from the group consisting of:

- (a) $-\text{NH}_2$,
- (b) $-\text{NH}-\text{C}_{1-4}\text{alkyl}$,
- (c) $-\text{N}(\text{C}_{1-4}\text{alkyl})_2$,
- (d) $-\text{NH}-\text{C}(=\text{O})-\text{NH}_2$,
- (e) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{1-4}\text{alkyl}$,
- (f) $-\text{NH}-\text{C}(=\text{O})-\text{N}(\text{C}_{1-4}\text{alkyl})_2$,
- (g) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{3-5}\text{alkenyl}$,
- (h) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{3-5}\text{alkynyl}$,
- (i) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{3-6}\text{cycloalkyl}$,
- (j) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{aryl}$,
- (k) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{heterocycle}$,
- (l) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{heteroaryl}$,
- (m) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{SO}_2-\text{C}_{1-4}\text{alkyl}$,
- (n) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{SO}_2-\text{C}_{3-6}\text{cycloalkyl}$,
- (o) $-\text{NH}-\text{C}(=\text{O})-\text{O}-\text{C}_{1-4}\text{alkyl}$,
- (p) $-\text{NH}-\text{C}(=\text{O})-\text{O}-\text{C}_{1-4}\text{alkylaryl}$,
- (q) $-\text{NH}-\text{C}(=\text{O})-\text{C}_{1-4}\text{alkyl}$,
- (r) $-\text{NH}-\text{C}(=\text{O})-\text{C}_{3-6}\text{cycloalkyl}$,
- (s) $-\text{NH}-\text{C}(=\text{O})-\text{aryl}$,
- (t) $-\text{NH}-\text{C}(=\text{O})-\text{heterocycle}$,
- (u) $-\text{NH}-\text{C}(=\text{O})-\text{heteroaryl}$, and
- (v) $-\text{NH}-\text{SO}_2-\text{C}_{1-4}\text{alkyl}$,

wherein each of choices (b) to (v) is optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (i) $-\text{C}_{1-4}\text{alkyl}$,
- (ii) $-\text{halogen}$,

- (iii) $-\text{OH}$,
- (iv) $-\text{O}-\text{C}_{1-4}\text{alkyl}$,
- (v) $-\text{SH}$,
- (vi) $-\text{S}-\text{C}_{1-4}\text{alkyl}$;
- (vii) $-\text{NO}_2$, and
- (viii) $-\text{CN}$;

R^{10} , R^{11} , and R^{12} are independently selected from the group consisting of: H, halogen, $-\text{OH}$ and $-\text{C}_{1-6}\text{alkyl}$; or

R^{10} , R^{11} and the carbon atom to which they are attached together form a $\text{C}_{3-12}\text{cycloalkyl}$ or a heterocyclyl group;

R^{13} is selected from the group consisting of:

- (a) $-\text{H}$,
- (b) $-\text{C}_{1-4}\text{alkyl}$,
- (c) $-\text{C}(\text{O})-\text{C}_{1-4}\text{alkyl}$,
- (d) $-\text{C}(\text{O})\text{NH}_2$,
- (e) $-\text{C}(\text{O})-\text{NH}(\text{C}_{1-4}\text{alkyl})$,
- (f) $-\text{C}(\text{O})-\text{NH}(\text{C}_{3-6}\text{cycloalkyl})$,
- (g) $-\text{C}(\text{O})-\text{N}(\text{C}_{1-4}\text{alkyl})_2$,
- (h) $-\text{C}(\text{O})\text{O}-\text{C}_{1-4}\text{alkyl}$, and
- (i) $-\text{C}(\text{O})\text{O}-\text{C}_{1-4}\text{alkylaryl}$;

p is 0, 1 or 2.

2. A compound of Claim 1 wherein

R^1 is selected from the group consisting of:

- (a) $-\text{aryl}$ and
- (b) $-\text{heteroaryl}$;

wherein the aryl and heteroaryl of choices (a) and (b) are each optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (i) $-\text{halogen}$,
- (ii) $-\text{CN}$,
- (iii) $-\text{C}(\text{O})\text{R}^8$,

- (iv) $-\text{CONR}^5\text{R}^6$,
- (v) $-\text{OH}$,
- (vi) $-\text{O}-\text{C}_{1-6}\text{alkyl}$,
- (vii) $-\text{S}(\text{O})_p-\text{C}_{1-6}\text{alkyl}$,
- (viii) $-\text{S}(\text{O})_2\text{NR}^5\text{R}^6$,
- (ix) $-\text{NHC}(\text{O})\text{R}^8$,
- (x) $-\text{NHC}(\text{O})\text{OR}^8$,
- (xi) $-\text{NHSO}_2\text{C}_{1-6}\text{alkyl}$,

wherein each of the alkyl group of choices (iii), (x), (xii) and (xix) is optionally substituted with 1 to 5 substituents independently selected from $-\text{halogen}$, $-\text{haloC}_{1-4}\text{alkyl}$, $-\text{COR}^8$, $-\text{CO}_2\text{R}^8$, $-\text{CONR}^5\text{R}^6$, $-\text{NR}^5\text{R}^6$, $-\text{OH}$, $-\text{O}-\text{C}_{1-4}\text{alkyl}$, $-\text{SH}$ and $-\text{S}-\text{C}_{1-4}\text{alkyl}$.

3. A compound of Claim 1 or Claim 2 wherein

R^2 is selected from the group consisting of:

- (a) $-\text{C}_{1-6}\text{alkyl}-\text{R}^7$, and
- (b) $-(\text{CH}_2)_{1-6}-\text{N}(\text{R}^{13})(\text{R}^{13})$,

wherein the alkyl group of choice (a) is optionally substituted with 1 to 5 substituents independently selected from:

- (i) $-\text{halogen}$,
- (ii) $-\text{C}_{1-4}\text{alkyl}$,
- (iii) $-\text{haloC}_{1-4}\text{alkyl}$,
- (iv) $-\text{OH}$,
- (v) $-\text{O}-\text{C}_{1-4}\text{alkyl}$,
- (vi) $-\text{SH}$, and
- (vii) $-\text{S}-\text{C}_{1-4}\text{alkyl}$.

4. A compound of any one of Claims 1 - 3 wherein

R^2 is selected from the group consisting of:

- (a) $-(\text{CH}_2)_{1-6}-\text{R}^7$, and
- (b) $-(\text{CH}_2)_{1-6}-\text{N}(\text{R}^{13})(\text{R}^{13})$.

5. A compound of any one of Claims 1 - 4 wherein R^{3b} and R^{3c} are each H, and R^{3a} is selected from the group consisting of:



wherein HAr is heteroaryl and Hcyl is heterocycle, wherein each of HAr and Hcyl is optionally substituted with 1 to 3 groups independently selected from the group consisting of:

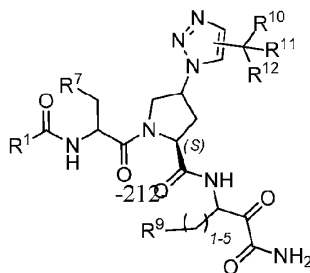
- (i) -halogen,
- (ii) -OH,
- (iii) $-CR^{10}R^{11}R^{12}$,
- (iv) $-(CH_2)_{0-3}-NHSO_2-C_{1-4}alkyl$, and
- (v) $-(CH_2)_{0-3}-SO_2-C_{1-4}alkyl$.

6. A compound of any one of Claims 1 - 5 wherein R^4 is selected from the group consisting of:

- (a) $-C_{1-6}alkyl$,
- (b) $-C_{2-6}alkenyl$,
- (c) $-C_{1-6}alkyl-C_{3-6}cycloalkyl$,
- (d) $-C_{1-6}alkyl-aryl$, wherein aryl is optionally substituted with nitro or $-N(R^{13})(R^{13})$, and
- (e) $-C_{1-6}alkyl-R^9$.

7. A compound of any one of Claims 1 - 6 wherein R^4 is $-C_{1-6}alkyl-R^9$.

8. A compound of Claim 1 having the Formula (Ia):



(Ia)

or a pharmaceutically acceptable salt, solvate, solvate of the salt or prodrug thereof; wherein R^1 is selected from the group consisting of:

- (a) -aryl and
- (b) -heteroaryl;

wherein the aryl and heteroaryl of choices (a) and (b) are each optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (i) -halogen,
- (ii) -CN,
- (iii) -C₁₋₆alkyl,
- (iv) -C₂₋₆alkenyl,
- (v) -C₂₋₆alkynyl,
- (vi) -C(O)R⁸,
- (vii) -CO₂R⁸,
- (viii) -CONR⁵R⁶,
- (ix) -OH,
- (x) -O-C₁₋₆alkyl,
- (xi) -SH,
- (xii) -S(O)_p-C₁₋₆alkyl,
- (xiii) -S(O)₂NR⁵R⁶,
- (xiv) -NO₂,
- (xv) -NR⁵R⁶,
- (xvi) -NHC(O)R⁸,
- (xvii) -NHC(O)OR⁸,
- (xviii) -NHC(O)NR⁵R⁶, and
- (xix) -NHSO₂C₁₋₆alkyl,

wherein each of the alkyl group of choices (iii), (x), (xii) and (xix) is optionally substituted with 1 to 5 substituents independently selected from -halogen, -haloC₁₋₄alkyl, -COR⁸, -CO₂R⁸, -CONR⁵R⁶, -NR⁵R⁶, -OH, -O-C₁₋₄alkyl, -SH and -S-C₁₋₄alkyl;

each R⁵ and each R⁶ are independently selected from the group consisting of:

- (a) -H,
- (b) -C₁₋₆alkyl,
- (c) -C₀₋₆alkyl-C₃₋₁₂cycloalkyl,
- (d) -C₀₋₆alkyl-heterocyclyl,
- (e) -C₀₋₆alkyl-heteroaryl, and
- (f) -C₀₋₆alkyl-aryl,

wherein each of the alkyl groups of choices (b) – (f) is optionally substituted with 1 to 3 groups independently selected from:

- (i) -halogen,
- (ii) -C(O)C₁₋₄alkyl,
- (iii) -C(O)OC₁₋₄alkyl,
- (iv) -OH,
- (v) -OC₁₋₄alkyl,
- (vi) -SH,
- (vii) -SC₁₋₄alkyl,
- (viii) -NH₂,
- (ix) -NH(C₁₋₄alkyl), and
- (x) -N(C₁₋₄alkyl)(C₁₋₄alkyl); or

R⁵, R⁶ and the nitrogen atom to which they are attached together form a 3- to 7-membered monocyclic or 6- to 11-membered bicyclic heterocycle optionally having an additional heteroatomic moiety selected from -O-, -S(O)_p-, and -NR¹³-, and wherein said heterocycle is optionally substituted with 1 to 2 groups independently selected from halogen, -haloC₁₋₄alkyl, -OH, -O-C₁₋₄alkyl, -SH and -S-C₁₋₄alkyl;

R⁷ is selected from the group consisting of:

- (a) -C₃₋₁₀cycloalkyl, and
- (b) -heterocyclyl,

wherein each of choices (a) and (b) is optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (i) -C₁₋₄alkyl,
- (ii) -halogen,
- (iii) -OH,

- (iv) $-\text{O}-\text{C}_{1-4}\text{alkyl}$,
- (v) $-\text{SH}$, and
- (vi) $-\text{S}-\text{C}_{1-4}\text{alkyl}$;

R^8 is selected from the group consisting of:

- (a) $-\text{C}_{1-6}\text{alkyl}$,
- (b) $-\text{C}_{0-6}\text{alkyl}-\text{C}_{3-12}\text{cycloalkyl}$,
- (c) $-\text{C}_{0-6}\text{alkyl}-\text{heterocyclyl}$,
- (d) $-\text{C}_{0-6}\text{alkyl}-\text{heteroaryl}$, and
- (e) $-\text{C}_{0-6}\text{alkyl}-\text{aryl}$,

wherein the alkyl group of choices (a) – (e) is optionally substituted with 1 to 3 groups independently selected from:

- (i) $-\text{halogen}$,
- (ii) $-\text{OH}$,
- (iii) $-\text{OC}_{1-4}\text{alkyl}$,
- (iv) $-\text{SH}$, and
- (v) $-\text{SC}_{1-4}\text{alkyl}$;

R^9 is selected from the group consisting of:

- (a) $-\text{NH}_2$,
- (b) $-\text{NH}-\text{C}_{1-4}\text{alkyl}$,
- (c) $-\text{N}(\text{C}_{1-4}\text{alkyl})_2$,
- (d) $-\text{NH}-\text{C}(=\text{O})-\text{NH}_2$,
- (e) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{1-4}\text{alkyl}$,
- (f) $-\text{NH}-\text{C}(=\text{O})-\text{N}(\text{C}_{1-4}\text{alkyl})_2$,
- (g) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{3-5}\text{alkenyl}$,
- (h) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{3-5}\text{alkynyl}$,
- (i) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{3-6}\text{cycloalkyl}$,
- (j) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{aryl}$,
- (k) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{heterocycle}$,
- (l) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{heteroaryl}$,
- (m) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{SO}_2-\text{C}_{1-4}\text{alkyl}$,
- (n) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{SO}_2-\text{C}_{3-6}\text{cycloalkyl}$,

- (o) $-\text{NH}-\text{C}(=\text{O})-\text{O}-\text{C}_{1-4}\text{alkyl}$,
- (p) $-\text{NH}-\text{C}(=\text{O})-\text{O}-\text{C}_{1-4}\text{alkylaryl}$,
- (q) $-\text{NH}-\text{C}(=\text{O})-\text{C}_{1-4}\text{alkyl}$,
- (r) $-\text{NH}-\text{C}(=\text{O})-\text{C}_{3-6}\text{cycloalkyl}$,
- (s) $-\text{NH}-\text{C}(=\text{O})-\text{aryl}$,
- (t) $-\text{NH}-\text{C}(=\text{O})-\text{heterocycle}$,
- (u) $-\text{NH}-\text{C}(=\text{O})-\text{heteroaryl}$, and
- (v) $-\text{NH}-\text{SO}_2-\text{C}_{1-4}\text{alkyl}$,

wherein each of choices (b) to (v) is optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (i) $-\text{C}_{1-4}\text{alkyl}$,
- (ii) $-\text{halogen}$,
- (iii) $-\text{OH}$,
- (iv) $-\text{O}-\text{C}_{1-4}\text{alkyl}$,
- (v) $-\text{SH}$,
- (vi) $-\text{S}-\text{C}_{1-4}\text{alkyl}$;
- (vii) $-\text{NO}_2$, and
- (viii) $-\text{CN}$;

R¹⁰, R¹¹, and R¹² are independently selected from the group consisting of: H, halogen, $-\text{OH}$ and $-\text{C}_{1-6}\text{alkyl}$; or

R¹⁰, R¹¹ and the carbon atom to which they are attached together form a $\text{C}_{3-12}\text{cycloalkyl}$ or a heterocyclyl group;

R¹³ is selected from the group consisting of:

- (a) $-\text{H}$,
- (b) $-\text{C}_{1-4}\text{alkyl}$,
- (c) $-\text{C}(\text{O})-\text{C}_{1-4}\text{alkyl}$,
- (d) $-\text{C}(\text{O})\text{NH}_2$,
- (e) $-\text{C}(\text{O})-\text{NH}(\text{C}_{1-4}\text{alkyl})$,
- (f) $-\text{C}(\text{O})-\text{NH}(\text{C}_{3-6}\text{cycloalkyl})$,
- (g) $-\text{C}(\text{O})-\text{N}(\text{C}_{1-4}\text{alkyl})_2$,
- (h) $-\text{C}(\text{O})\text{O}-\text{C}_{1-4}\text{alkyl}$, and

- (i) $-\text{C}(\text{O})\text{O}-\text{C}_{1-4}\text{alkylaryl}$; and

p is 0, 1 or 2.

9. A compound of Claim 8 wherein

R¹ is selected from the group consisting of:

- (a) $-\text{aryl}$, and
(b) $-\text{heteroaryl}$;

wherein the aryl and heteroaryl of choices (a) and (b) are each optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (i) $-\text{halogen}$,
(ii) $-\text{CN}$,
(iii) $-\text{C}(\text{O})\text{R}^8$,
(iv) $-\text{CONR}^5\text{R}^6$,
(v) $-\text{OH}$,
(vi) $-\text{O}-\text{C}_{1-6}\text{alkyl}$,
(vii) $-\text{S}(\text{O})_p-\text{C}_{1-6}\text{alkyl}$,
(viii) $-\text{S}(\text{O})_2\text{NR}^5\text{R}^6$,
(ix) $-\text{NHC}(\text{O})\text{R}^8$,
(x) $-\text{NHC}(\text{O})\text{OR}^8$, and
(xi) $-\text{NHSO}_2\text{C}_{1-6}\text{alkyl}$,

wherein each of the alkyl group of choices (vi), (vii), and (xi) is optionally substituted with 1 to 5 substituents independently selected from $-\text{halogen}$, $-\text{haloC}_{1-4}\text{alkyl}$, $-\text{COR}^8$, $-\text{CO}_2\text{R}^8$, $-\text{CONR}^5\text{R}^6$, $-\text{NR}^5\text{R}^6$, $-\text{OH}$, $-\text{O}-\text{C}_{1-4}\text{alkyl}$, $-\text{SH}$ and $-\text{S}-\text{C}_{1-4}\text{alkyl}$;

R⁷ is $-\text{C}_{3-10}\text{cycloalkyl}$;

R⁹ is selected from the group consisting of:

- (m) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{1-4}\text{alkyl}$,
(n) $-\text{NH}-\text{C}(=\text{O})-\text{N}(\text{C}_{1-4}\text{alkyl})_2$,
(o) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{3-5}\text{alkenyl}$,
(p) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{3-5}\text{alkynyl}$,
(q) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{C}_{3-6}\text{cycloalkyl}$,
(r) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{aryl}$,

- (s) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{heterocycle}$,
- (t) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{heteroaryl}$,
- (u) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{SO}_2-\text{C}_{1-4}\text{alkyl}$,
- (v) $-\text{NH}-\text{C}(=\text{O})-\text{NH}-\text{SO}_2-\text{C}_{3-6}\text{cycloalkyl}$,
- (w) $-\text{NH}-\text{C}(=\text{O})-\text{O}-\text{C}_{1-4}\text{alkyl}$, and
- (x) $-\text{NH}-\text{C}(=\text{O})-\text{O}-\text{C}_{1-4}\text{alkylaryl}$,

wherein each of choices (a) to (l) is optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

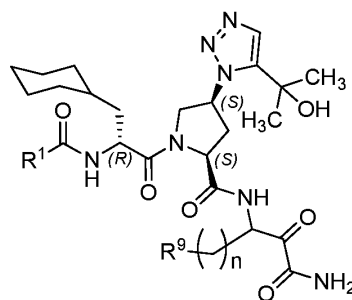
- (vii) $-\text{C}_{1-4}\text{alkyl}$,
- (viii) $-\text{halogen}$,
- (ix) $-\text{OH}$,
- (x) $-\text{O}-\text{C}_{1-4}\text{alkyl}$,
- (xi) $-\text{SH}$, and
- (xii) $-\text{S}-\text{C}_{1-4}\text{alkyl}$;

R^{10} and R^{11} are each $-\text{C}_{1-4}\text{alkyl}$, or

R^{10} , R^{11} and the carbon atom to which they are attached together form a $\text{C}_{3-6}\text{cycloalkyl}$ or a 4- to 6-membered heterocycle, and

R^{12} is $-\text{OH}$.

10. A compound of Claim 1 having the formula (Ib):



(Ib)

or a pharmaceutically acceptable salt, solvate, solvate of the salt or prodrug thereof, wherein

R^1 is selected from the group consisting of:

- (a) $-\text{aryl}$, and

(b) -heteroaryl;

wherein the aryl and heteroaryl of choices (a) and (b) are each optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (xiii) -halogen,
- (xiv) -CN,
- (xv) -C(O)R⁸,
- (xvi) -CONR⁵R⁶,
- (xvii) -OH,
- (xviii) -O-C₁₋₆alkyl,
- (xix) -S(O)_p-C₁₋₆alkyl,
- (xx) -S(O)₂NR⁵R⁶,
- (xxi) -NHC(O)R⁸,
- (xxii) -NHC(O)OR⁸,
- (xxiii) -NHSO₂C₁₋₆alkyl, and
- (xxiv) C₁₋₄alkyl;

wherein each of the alkyl group of choices (vi), (vii), and (xi) is optionally substituted with 1 to 5 substituents independently selected from -halogen, -haloC₁₋₄alkyl, -COR⁸, -CO₂R⁸, -CONR⁵R⁶, -NR⁵R⁶, -OH, -O-C₁₋₄alkyl, -SH and -S-C₁₋₄alkyl;

R⁹ is selected from the group consisting of:

- (y) -NH-C(=O)-NH-C₁₋₄alkyl,
- (z) -NH-C(=O)-N(C₁₋₄alkyl)₂,
- (aa) -NH-C(=O)-NH-C₃₋₅alkenyl,
- (bb) -NH-C(=O)-NH-C₃₋₅alkynyl,
- (cc) -NH-C(=O)-NH-C₃₋₆cycloalkyl,
- (dd) -NH-C(=O)-NH-aryl,
- (ee) -NH-C(=O)-NH-heterocycle,
- (ff) -NH-C(=O)-NH-heteroaryl,
- (gg) -NH-C(=O)-NH-SO₂-C₁₋₄alkyl,
- (hh) -NH-C(=O)-NH-SO₂-C₃₋₆cycloalkyl,
- (ii) -NH-C(=O)-O-C₁₋₄alkyl, and
- (jj) -NH-C(=O)-O-C₁₋₄alkylaryl,

wherein each of choices (a) to (l) is optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (xiii) $-C_{1-4}alkyl$,
- (xiv) $-halogen$,
- (xv) $-OH$,
- (xvi) $-O-C_{1-4}alkyl$,
- (xvii) $-SH$, and
- (xviii) $-S-C_{1-4}alkyl$; and

n is 1 to 5.

11. A compound of Claim 10 wherein n is 4.

12. A compound of Claim 10 or Claim 11 wherein

R^1 is phenyl optionally substituted with 1 to 3 substituents independently selected from the group consisting of:

- (ix) $-halogen$,
- (x) $-CN$,
- (xi) $-C(O)R^8$,
- (xii) $-CONR^5R^6$,
- (xiii) $-OH$,
- (xiv) $-O-C_{1-6}alkyl$,
- (xv) $-S(O)_p-C_{1-6}alkyl$, and
- (xvi) $-S(O)_2NR^5R^6$;

wherein each of the alkyl group of choices (vi) and (vii) is optionally substituted with 1 to 5 substituents independently selected from $-halogen$, $-haloC_{1-4}alkyl$, $-COR^8$, $-CO_2R^8$, $-CONH_2$, $-OH$, and $-O-C_{1-4}alkyl$.

13. A compound of Claim 10 or Claim 11 wherein R^1 is naphthyl.

14. A compound of Claim 10 or Claim 11 wherein R^1 is selected from the group consisting of:

(c) 5- or 6-membered monocyclic heteroaryl ring having a heteroatom selected from N, O and S, and optionally 1, 2 or 3 additional N atoms; and

(d) 8-, 9-, or 10-membered fused bicyclic heteroaryl ring having a heteroatom selected from N, O and S, and optionally 1, 2 or 3 additional N atoms;

wherein each of choices (a) and (b) is optionally substituted with a group selected from: OH and C₁₋₄alkyl.

15. A compound of Claim 1 selected from the group consisting of:

(2*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-4-methyl-1,2-dioxopentan-3-yl)pyrrolidine-2-carboxamide;

methyl (3-((*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)pyrrolidine-2-carboxamido)-4-methyl-2-oxopentanoyl)glycinate;

(2*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-((2-methoxyethyl)amino)-4-methyl-1,2-dioxopentan-3-yl)pyrrolidine-2-carboxamide;

(2*S*,4*R*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-4-methyl-1,2-dioxopentan-3-yl)-4-phenylpyrrolidine-2-carboxamide;

(2*S*,4*R*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-4-methyl-1,2-dioxopentan-3-yl)-4-(piperidin-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*R*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-cyclohexylureido)-2-hydroxy-1-oxoheptan-3-yl)-4-(piperidin-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*R*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(4-amino-1-cyclobutyl-3,4-dioxobutan-2-yl)-4-(piperidin-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*R*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-4-methyl-1,2-dioxohexan-3-yl)-4-(piperidin-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*R*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(4-amino-3,4-dioxo-1-phenylbutan-2-yl)-4-(piperidin-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(4-amino-1-cyclobutyl-3,4-dioxobutan-2-yl)-4-(piperidin-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-cyclopropylureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-isobutylureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-cyclohexylureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-1,2-dioxo-7-(3-(pyridin-4-yl)ureido)heptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-ethylureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-(1-methyl-1*H*-pyrazol-3-yl)ureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-(isoxazol-3-yl)ureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-1,2-dioxo-7-(3-(tetrahydro-2*H*-pyran-4-yl)ureido)heptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-1,2-dioxo-7-(3-(pyrimidin-5-yl)ureido)heptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-(1-oxidotetrahydro-2*H*-thiopyran-4-yl)ureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-1,2-dioxo-7-(3-(tetrahydro-2*H*-thiopyran-4-yl)ureido)heptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-1,2-dioxo-7-(3-(pyridin-3-yl)ureido)heptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-1,2-dioxo-7-(3-(pyridin-2-yl)ureido)heptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-1,2-dioxo-7-(3-(prop-2-yn-1-yl)ureido)heptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-(5-(tert-butyl)isoxazol-3-yl)ureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(7-(3-allylureido)-1-amino-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-(cyanomethyl)ureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-(cyclobutylureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-1,2-dioxo-7-((trifluoromethyl)sulfonamido)heptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-(cyclopropylsulfonyl)ureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

benzyl (5-((2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;

benzyl (5-((2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-4-(4-(2-hydroxypropan-2-

yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;
benzyl (5-((2*S*,4*R*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;
benzyl (5-((2*S*,4*R*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-4-(4-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;
(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-1,2-dioxohex-5-en-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;
N-((2*R*)-1-((2*S*,4*S*)-2-((7-(3-propargylureido)-1-amino-1,2-dioxoheptan-3-yl)carbamoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidin-1-yl)-3-cyclohexyl-1-oxopropan-2-yl)isonicotinamide;
(2*S*,4*S*)-*N*-(7-(3-propargylureido)-1-amino-1,2-dioxoheptan-3-yl)-1-((*R*)-3-cyclohexyl-2-(4-(methylsulfonyl)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;
(2*S*,4*S*)-*N*-(1-amino-1,2-dioxo-7-(3-(prop-2-yn-1-yl)ureido)heptan-3-yl)-1-((*R*)-3-cyclohexyl-2-(4-(2,2,2-trifluoro-1,1-dihydroxyethyl)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;
benzyl (5-((2*S*,4*S*)-1-((*R*)-2-(1*H*-benzo[*d*]imidazole-2-carboxamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(4-(methylsulfonyl)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (5-((2*S*,4*S*)-1-((*R*)-2-(1*H*-benzo[*d*]imidazole-4-carboxamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;
benzyl (5-((2*S*,4*S*)-1-((*R*)-2-(1*H*-benzo[*d*]imidazole-6-carboxamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1,6-naphthyridine-5-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-

carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1,6-naphthyridine-8-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1,8-naphthyridine-4-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(quinoxaline-5-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(quinazoline-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(quinoxaline-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1,5-naphthyridine-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1,6-naphthyridine-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1,8-naphthyridine-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1,7-naphthyridine-3-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;
benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1,5-naphthyridine-3-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-

carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(quinoxaline-6-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(4-oxo-3,4-dihydroquinazoline-7-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1,5-naphthyridine-4-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1-methyl-1*H*-imidazole-5-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1-methyl-1*H*-imidazole-4-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1-methyl-1*H*-imidazole-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1*H*-pyrrolo[3,2-*c*]pyridine-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1*H*-indole-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1*H*-indole-5-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1*H*-indole-6-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-

dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1*H*-indazole-4-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1*H*-indazole-5-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1*H*-indazole-6-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1*H*-indazole-7-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (5-((2*S*,4*S*)-1-((*R*)-2-(1-naphthamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(isoxazole-4-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(oxazole-4-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(oxazole-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-2-(3-carbamoylbenzamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-2-(4-carbamoylbenzamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-2-(2-cyanobenzamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-2-(3-cyanobenzamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-2-(4-cyanobenzamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(2-fluorobenzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(4-fluorobenzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(2-methoxybenzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(3-methoxybenzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(4-methoxybenzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(2-((methylperoxy)thio)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(3-((methylperoxy)thio)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(4-((difluoromethyl)sulfonyl)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(4-((2-methoxyethyl)sulfonyl)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-2-(4-((2-amino-2-oxoethyl)sulfonyl)benzamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(4-(isopropylsulfonyl)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(nicotinamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(picolinamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(isonicotinamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(pyrimidine-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(isoquinoline-1-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(quinoline-4-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(quinoline-5-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(isoquinoline-5-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(isoquinoline-8-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(quinoline-8-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(quinoline-2-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1-oxo-1,2-dihydroisoquinoline-3-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(isoquinoline-3-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(quinoline-3-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(2-oxo-1,2-dihydroquinoline-6-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(isoquinoline-6-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(isoquinoline-7-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(imidazo[1,2-*a*]pyridine-6-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(quinoline-7-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(pyrido[2,3-*b*]pyrazine-8-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(pyrido[2,3-*b*]pyrazine-6-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(pyrido[2,3-*b*]pyrazine-7-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1*H*-1,2,3-triazole-4-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (7-amino-5-((2*S*,4*S*)-1-((*R*)-3-cyclohexyl-2-(1-methyl-1*H*-1,2,3-triazole-4-carboxamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-6,7-dioxoheptyl)carbamate;

(2*S*,4*S*)-*N*-(1-amino-7-(3-cyclopropylureido)-1,2-dioxoheptan-3-yl)-1-((*R*)-3-cyclohexyl-2-(4-(methylsulfonyl)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-*N*-(1-amino-7-(3-cyclopropylureido)-1,2-dioxoheptan-3-yl)-1-((*R*)-2-(4-cyanobenzamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

N-((2*R*)-1-((2*S*,4*S*)-2-((1-amino-7-(3-cyclopropylureido)-1,2-dioxoheptan-3-yl)carbamoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidin-1-yl)-3-cyclohexyl-1-oxopropan-2-yl)imidazo[1,2-*a*]pyridine-6-carboxamide;

N-((2*R*)-1-((2*S*,4*S*)-2-((1-amino-7-(3-cyclopropylureido)-1,2-dioxoheptan-3-yl)carbamoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidin-1-yl)-3-cyclohexyl-1-oxopropan-2-yl)quinoline-3-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(4-((2-amino-2-oxoethyl)sulfonyl)benzamido)-3-cyclohexylpropanoyl)-*N*-(1-amino-7-(3-cyclopropylureido)-1,2-dioxoheptan-3-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-*N*-(1-amino-7-(3-cyclopropylureido)-1,2-dioxoheptan-3-yl)-1-((*R*)-3-cyclohexyl-2-(4-((difluoromethyl)sulfonyl)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

N-((2*R*)-1-((2*S*,4*S*)-2-((1-amino-7-(3-cyclopropylureido)-1,2-dioxoheptan-3-yl)carbamoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidin-1-yl)-3-cyclohexyl-1-oxopropan-2-yl)-1*H*-indazole-7-carboxamide;

(2*S*,4*S*)-*N*-(1-amino-7-(3-cyclopropylureido)-1,2-dioxoheptan-3-yl)-1-((*R*)-3-cyclohexyl-2-(4-((2-methoxyethyl)sulfonyl)benzamido)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(4-amino-1-(4-nitrophenyl)-3,4-dioxobutan-2-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(1-(4-acetamidophenyl)-4-amino-3,4-dioxobutan-2-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

(2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-*N*-(4-amino-1-(4-(3-cyclohexylureido)phenyl)-3,4-dioxobutan-2-yl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamide;

benzyl (2-((2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-cyclohexylpropanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-4-amino-3,4-dioxobutyl)carbamate;

benzyl (7-amino-5-((3*S*)-2-((*R*)-3-cyclohexyl-2-(4-(methylsulfonyl)benzamido)propanoyl)-2-azabicyclo[2.2.1]heptane-3-carboxamido)-6,7-dioxoheptyl)carbamate;

benzyl (5-((2*S*,4*S*)-1-(2-(2-naphthamido)-4-((1*S*,4*R*)-bicyclo[2.2.1]heptan-2-yl)butanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;

benzyl (5-((2*S*,4*S*)-1-((*R*)-2-(2-naphthamido)-3-(((benzyloxy)carbonyl)amino)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;

benzyl (5-((2*S*,4*S*)-1-(2-(2-naphthamido)-3-(tetrahydro-2*H*-pyran-4-yl)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;

benzyl (5-((2*S*,4*S*)-1-(2-(2-naphthamido)-3-((3*S*,5*S*,7*S*)-adamantan-1-yl)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;

benzyl (5-((2*S*,4*S*)-1-(2-(2-naphthamido)-3-((1*S*,2*R*,5*S*)-6,6-dimethylbicyclo[3.1.1]heptan-2-yl)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate; and

benzyl (5-((2*S*,4*S*)-1-(2-(2-naphthamido)-3-(3-methyloxetan-3-yl)propanoyl)-4-(5-(2-hydroxypropan-2-yl)-1*H*-1,2,3-triazol-1-yl)pyrrolidine-2-carboxamido)-7-amino-6,7-dioxoheptyl)carbamate;

or a pharmaceutically acceptable salt, solvate, salt of the solvate or prodrug thereof.

16. A pharmaceutical composition comprising a compound of any of Claim 1-15 and a pharmaceutically acceptable carrier.

17. A method of preventing, or treating a disease of the eye selected from dry-AMD, wet-AMD, geographic atrophy, diabetic retinopathy, retinopathy of prematurity, polypoidal choroidal vasculopathy, and degeneration of retinal or photoreceptor cells, comprising: administering to a subject in need thereof a therapeutically effective amount of a compound according to Claim 1-15 or a pharmaceutically acceptable salt, solvate, solvate of the salt or prodrug thereof, or the pharmaceutical composition of Claim 16.

18. The method of preventing a disease of the eye, according to Claim 17 wherein the method of prevention is selected from delaying the onset of disease and reducing the risk of developing a disease of the eye, wherein the disease of the eye is selected from dry-AMD, wet-AMD, geographic atrophy, diabetic retinopathy, retinopathy of prematurity, polypoidal choroidal vasculopathy, and degeneration of retinal or photoreceptor cells.

19. The method of treating a disease of the eye according to Claim 17 wherein the method is selected from controlling, alleviating, and slowing the progression of, wherein the disease is selected from dry-AMD, wet-AMD, geographic atrophy, diabetic retinopathy, retinopathy of prematurity, polypoidal choroidal vasculopathy, and degeneration of retinal or photoreceptor cells.

20. The method according to any of Claims 17 to 19 wherein the disease is geographic atrophy.

21. A method of inhibiting HTRA1 protease activity in an eye comprising administering to a subject in need thereof a therapeutically effective amount of compound according to Claims 1-15 or a pharmaceutically acceptable salt, solvate, solvate of the salt or prodrug thereof, or a pharmaceutical composition of Claim 16.