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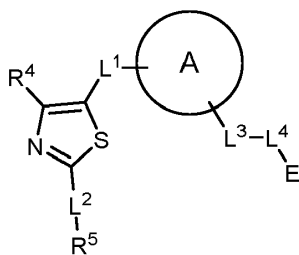
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(54)	Title	COMPOUNDS AND METHODS FOR MODULATING INTERLEUKIN-2-INDUCIBLE T-CELL KINASE
(56)	References Cited:	US-A1- 2005 176 789 US-A1- 2010 179 121 US-A1- 2003 069 290 US-A1- 2010 113 520 US-B2- 6 706 717 WO-A2-2017/044858 CN-A- 105 884 711 WO-A1-2014/117292

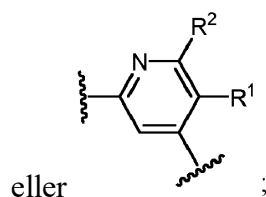
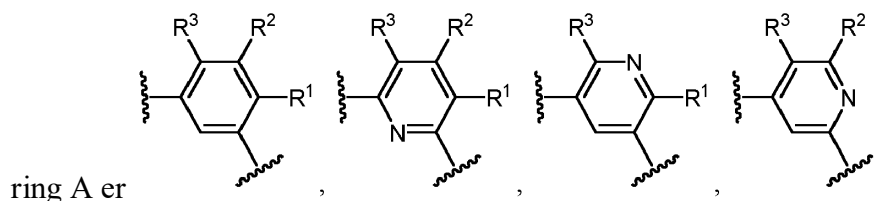
Enclosed is a translation of the patent claims in Norwegian. Please note that as per the Norwegian Patents Acts, section 66i the patent will receive protection in Norway only as far as there is agreement between the translation and the language of the application/patent granted at the EPO. In matters concerning the validity of the patent, language of the application/patent granted at the EPO will be used as the basis for the decision. The patent documents published by the EPO are available through Espacenet (<http://worldwide.espacenet.com>) or via the search engine on our website here: <https://search.patentstyret.no/>

Patentkrav

1. Forbindelse eller et farmasøytisk akseptabelt salt derav, som har formelen:



5 hvori,



- 10 R^1 er uavhengig $-OR^{1D}$, hydrogen, halogen, $-CX^1_3$, $-CHX^1_2$, $-CH_2X^1$, $-OCX^1_3$, $-OCH_2X^1$, $-OCHX^1_2$, $-CN$, $-SO_{n1}R^{1D}$, $-SO_{v1}NR^{1A}R^{1B}$, $-NHC(O)NR^{1A}R^{1B}$, $-N(O)_{m1}$, $-NR^{1A}R^{1B}$, $-C(O)R^{1C}$, $-C(O)-OR^{1C}$, $-C(O)NR^{1A}R^{1B}$, $-NR^{1A}SO_2R^{1D}$, $-NR^{1A}C(O)R^{1C}$, $-NR^{1A}C(O)OR^{1C}$,

- 15 $-NR^{1A}OR^{1C}$, substituert eller usubstituert alkyl, substituert eller usubstituert heteroalkyl, substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl, substituert eller usubstituert aryl eller substituert eller usubstituert heteroaryl;

- R^2 er uavhengig hydrogen, halogen, $-CX^2_3$, $-CHX^2_2$, $-CH_2X^2$, $-OCX^2_3$, $-OCH_2X^2$, $-OCHX^2_2$, $-CN$, $-SO_{n2}R^{2D}$, $-SO_{v2}NR^{2A}R^{2B}$, $-NHC(O)NR^{2A}R^{2B}$, $-N(O)_{m2}$, $-NR^{2A}R^{2B}$, $-C(O)R^{2C}$, $-C(O)-OR^{2C}$, $-C(O)NR^{2A}R^{2B}$, $-OR^{2D}$, $-NR^{2A}SO_2R^{2D}$, $-NR^{2A}C(O)R^{2C}$, $-NR^{2A}C(O)OR^{2C}$, $-NR^{2A}OR^{2C}$, substituert eller usubstituert alkyl, substituert eller usubstituert heteroalkyl, substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl, substituert eller usubstituert aryl eller substituert eller usubstituert heteroaryl;
- 20

R^3 er uavhengig usubstituert eller substituert alkyl, hydrogen, halogen, $-CX^3_3$, $-CHX^3_2$, $-CH_2X^3$, $-OCX^3_3$, $-OCH_2X^3$, $-OCHX^3_2$, $-CN$, $-SO_{n3}R^{3D}$, $-SO_{v3}NR^{3A}R^{3B}$, $-NHC(O)NR^{3A}R^{3B}$, $-N(O)_{m3}$, $-NR^{3A}R^{3B}$, $-C(O)R^{3C}$, $-C(O)-OR^{3C}$, $-C(O)NR^{3A}R^{3B}$, $-OR^{3D}$, $-NR^{3A}SO_2R^{3D}$, $-NR^{3A}C(O)R^{3C}$, $-NR^{3A}C(O)OR^{3C}$, $-NR^{3A}OR^{3C}$, substituert eller usubstituert heteroalkyl,

5 substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl, substituert eller usubstituert aryl eller substituert eller usubstituert heteroaryl;

R^4 er uavhengig hydrogen, halogen, $-CX^4_3$, $-CHX^4_2$, $-CH_2X^4$, $-OCX^4_3$, $-OCH_2X^4$, $-OCHX^4_2$, $-CN$, $-SO_{n4}R^{4D}$, $-SO_{v4}NR^{4A}R^{4B}$, $-NHC(O)NR^{4A}R^{4B}$, $-N(O)_{m4}$, $-NR^{4A}R^{4B}$, $-C(O)R^{4C}$,

10 $-C(O)-OR^{4C}$, $-C(O)NR^{4A}R^{4B}$, $-OR^{4D}$, $-NR^{4A}SO_2R^{4D}$, $-NR^{4A}C(O)R^{4C}$, $-NR^{4A}C(O)OR^{4C}$, $-NR^{4A}OR^{4C}$, substituert eller usubstituert alkyl, substituert eller usubstituert heteroalkyl, substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl, substituert eller usubstituert aryl eller substituert eller usubstituert heteroaryl;

R^5 er uavhengig usubstituert eller substituert sykloalkyl, substituert eller usubstituert alkyl, substituert eller usubstituert heteroalkyl, substituert eller usubstituert heterosykloalkyl, substituert eller usubstituert aryl eller substituert eller usubstituert heteroaryl;

L^1 er $-S-$, $-O-$ eller substituert eller usubstituert C_1 - C_2 alkylen eller substituert eller usubstituert 2-leddet heteroalkylen;

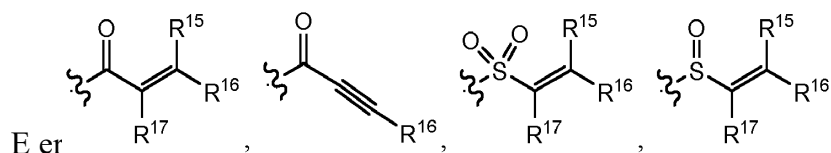
L^2 er $-NHC(O)-$, en binding eller $-NH-$;

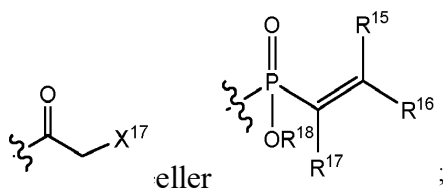
20 L^3 er $-C(O)-$, en

binding, $-S(O)_2-$, $-N(R^6)-$, $-O-$, $-S-$, $-C(O)N(R^6)-$, $-N(R^6)C(O)-$, $-N(R^6)C(O)NH-$, $-NHC(O)N(R^6)-$, $-C(O)O-$, $-OC(O)-$, substituert eller usubstituert alkylen, substituert eller usubstituert heteroalkylen, substituert eller usubstituert sykloalkylen, substituert eller usubstituert heterosykloalkylen, substituert eller usubstituert arylen eller substituert eller usubstituert heteroarylen;

R^6 er uavhengig hydrogen, $-CX^6_3$, $-CHX^6_2$, $-CH_2X^6$, $-CN$, $-C(O)R^{6C}$, $-C(O)OR^{6C}$, $-C(O)NR^{6A}R^{6B}$, substituert eller usubstituert alkyl, substituert eller usubstituert heteroalkyl, substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl, substituert eller usubstituert aryl eller substituert eller usubstituert heteroaryl;

30 L^4 er substituert eller usubstituert heterosykloalkylen;





R^{15} er uavhengig hydrogen, halogen, $-CX^{15}_3$, $-CHX^{15}_2$, $-CH_2X^{15}$, $-CN$, $-SO_{n15}R^{15D}$,
 $-SO_{v15}NR^{15A}R^{15B}$, $-NHN R^{15A}R^{15B}$, $-ONR^{15A}R^{15B}$, $-NHC=(O)NHN R^{15A}R^{15B}$,
 $-NHC(O)NR^{15A}R^{15B}$, $-N(O)_{m15}$, $-NR^{15A}R^{15B}$, $-C(O)R^{15C}$, $-C(O)-OR^{15C}$, $-$
 $C(O)NR^{15A}R^{15B}$,
 $-OR^{15D}$, $-NR^{15A}SO_2R^{15D}$, $-NR^{15A}C(O)R^{15C}$, $-NR^{15A}C(O)OR^{15C}$, $-NR^{15A}OR^{15C}$, $-OCX^{15}_3$,
 $-OCHX^{15}_2$, substituert eller usubstituert alkyl, substituert eller usubstituert heteroalkyl,
substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl,

substituert eller usubstituert aryl, substituert eller usubstituert heteroaryl;

R^{16} er uavhengig hydrogen, halogen, $-CX^{16}_3$, $-CHX^{16}_2$, $-CH_2X^{16}$, $-CN$, $-SO_{n16}R^{16D}$,
 $-SO_{v16}NR^{16A}R^{16B}$, $-NHN R^{16A}R^{16B}$, $-ONR^{16A}R^{16B}$, $-NHC=(O)NHN R^{16A}R^{16B}$,
 $-NHC(O)NR^{16A}R^{16B}$, $-N(O)_{m16}$, $-NR^{16A}R^{16B}$, $-C(O)R^{16C}$, $-C(O)-OR^{16C}$, $-$
 $C(O)NR^{16A}R^{16B}$,
 $-OR^{16D}$, $-NR^{16A}SO_2R^{16D}$, $-NR^{16A}C(O)R^{16C}$, $-NR^{16A}C(O)OR^{16C}$, $-NR^{16A}OR^{16C}$, $-OCX^{16}_3$,
 $-OCHX^{16}_2$, substituert eller usubstituert alkyl, substituert eller usubstituert heteroalkyl,
substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl,
substituert eller usubstituert aryl, substituert eller usubstituert heteroaryl;

R^{17} er uavhengig hydrogen, halogen, $-CX^{17}_3$, $-CHX^{17}_2$, $-CH_2X^{17}$, $-CN$, $-SO_{n17}R^{17D}$,
 $-SO_{v17}NR^{17A}R^{17B}$, $-NHN R^{17A}R^{17B}$, $-ONR^{17A}R^{17B}$, $-NHC=(O)NHN R^{17A}R^{17B}$,
 $-NHC(O)NR^{17A}R^{17B}$, $-N(O)_{m17}$, $-NR^{17A}R^{17B}$, $-C(O)R^{17C}$, $-C(O)-OR^{17C}$, $-$
 $C(O)NR^{17A}R^{17B}$,
 $-OR^{17D}$, $-NR^{17A}SO_2R^{17D}$, $-NR^{17A}C(O)R^{17C}$, $-NR^{17A}C(O)OR^{17C}$, $-NR^{17A}OR^{17C}$, $-OCX^{17}_3$,
 $-OCHX^{17}_2$, substituert eller usubstituert alkyl, substituert eller usubstituert heteroalkyl,

substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl,
substituert eller usubstituert aryl, substituert eller usubstituert heteroaryl;

R^{18} er uavhengig hydrogen, $-CX^{18}_3$, $-CHX^{18}_2$, $-CH_2X^{18}$, $-C(O)R^{18C}$, $-C(O)OR^{18C}$,
 $-C(O)NR^{18A}R^{18B}$, substituert eller usubstituert alkyl, substituert eller usubstituert
heteroalkyl, substituert eller usubstituert sykloalkyl, substituert eller usubstituert

heterosykloalkyl, substituert eller usubstituert aryl, substituert eller usubstituert heteroaryl;

hver R^{1A} , R^{1B} , R^{1C} , R^{1D} , R^{2A} , R^{2B} , R^{2C} , R^{2D} , R^{3A} , R^{3B} , R^{3C} , R^{3D} , R^{4A} , R^{4B} , R^{4C} , R^{4D} , R^{6A} , R^{6B} og R^{6C} er uavhengig usubstituert eller substituert alkyl, hydrogen, $-CX_3$, $-CN$, $-COOH$, $-CONH_2$, $-CHX_2$, $-CH_2X$, substituert eller usubstituert heteroalkyl, substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl, substituert eller usubstituert aryl eller substituert eller usubstituert heteroaryl;

R^{1A} - og R^{1B} -substituenten bundet til det samme nitrogenatomet kan eventuelt være forbundet for å danne et substituert eller usubstituert heterosykloalkyl eller substituert eller usubstituert heteroaryl; R^{2A} - og R^{2B} -substituenten bundet til det samme nitrogenatomet kan eventuelt være forbundet for å danne et substituert eller usubstituert heterosykloalkyl eller substituert eller usubstituert heteroaryl; R^{3A} - og R^{3B} -substituenten bundet til det samme nitrogenatomet kan eventuelt være forbundet for å danne et substituert eller usubstituert heterosykloalkyl eller substituert eller usubstituert heteroaryl; R^{4A} - og R^{4B} -substituenten bundet til det samme nitrogenatomet kan eventuelt være forbundet for å danne et substituert eller usubstituert heterosykloalkyl eller substituert eller usubstituert heteroaryl; R^{6A} - og R^{6B} -substituenten bundet til det samme nitrogenatomet kan eventuelt være forbundet for å danne et substituert eller usubstituert heterosykloalkyl eller substituert eller usubstituert heteroaryl;

R^{15A} , R^{15B} , R^{15C} , R^{15D} , R^{16A} , R^{16B} , R^{16C} , R^{16D} , R^{17A} , R^{17B} , R^{17C} , R^{17D} , R^{18A} , R^{18B} , R^{18C} , R^{18D} , er uavhengig hydrogen, $-CX_3$, $-CN$, $-COOH$, $-CONH_2$, $-CHX_2$, $-CH_2X$, substituert eller usubstituert alkyl, substituert eller usubstituert heteroalkyl, substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl, substituert eller usubstituert aryl eller substituert eller usubstituert heteroaryl; R^{15A} - og R^{15B} -substituenten bundet til det samme nitrogenatomet kan eventuelt være forbundet for å danne et substituert eller usubstituert heterosykloalkyl eller substituert eller usubstituert heteroaryl; R^{16A} - og R^{16B} -substituenten bundet til det samme nitrogenatomet kan eventuelt være forbundet for å danne et substituert eller usubstituert heterosykloalkyl eller substituert eller usubstituert heteroaryl; R^{17A} - og R^{17B} -substituenten bundet til det samme nitrogenatomet kan eventuelt være forbundet for å danne et substituert eller usubstituert heterosykloalkyl eller substituert eller usubstituert heteroaryl; R^{18A} - og R^{18B} -substituenten bundet til det samme nitrogenatomet kan eventuelt være forbundet for å danne et substituert eller usubstituert heterosykloalkyl eller substituert eller usubstituert heteroaryl;

hver X , X^1 , X^2 , X^3 , X^4 , X^6 , X^{15} , X^{16} , X^{17} og X^{18} er uavhengig $-F$, $-Cl$, $-Br$ eller $-I$;

n_1 , n_2 , n_3 , n_4 , n_{15} , n_{16} og n_{17} , er uavhengig et heltall fra 0 til 4; og

m_1 , m_2 , m_3 , m_4 , m_{15} , m_{16} , m_{17} , v_1 , v_2 , v_3 , v_4 , v_{15} , v_{16} og v_{17} er uavhengig 1 eller

2. Forbindelsen ifølge krav 1, hvori:

- (i) L^1 er -S-, -O- eller substituert eller usubstituert metylen; eller
- (ii) L^1 er -SCH₂-; eller
- (iii) L^1 er -CH(CH₃)-.

3. Forbindelsen ifølge krav 1 eller 2, hvori:

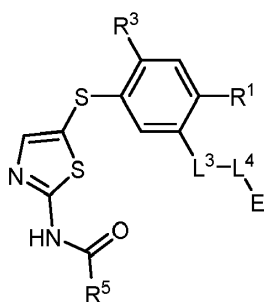
- (i) L^2 er -NHC(O)-; eller
- (ii) L^2 er -NH-; eller
- (iii) L^2 er en binding.

4. Forbindelsen ifølge et hvilket som helst av kravene 1 til 3, hvori R^4 er hydrogen.

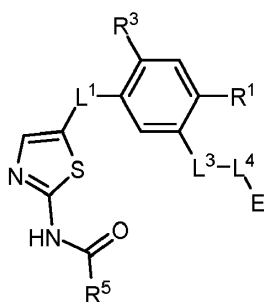
5. Forbindelsen ifølge et hvilket som helst av kravene 1 til 4, hvori:

- (i) R^2 er hydrogen; eller
- (ii) R^2 er usubstituert C₁-C₂alkyl; eller
- (iii) R^2 er usubstituert metyl.

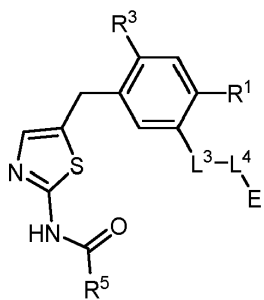
6. Forbindelsen ifølge krav 1, som har formelen:



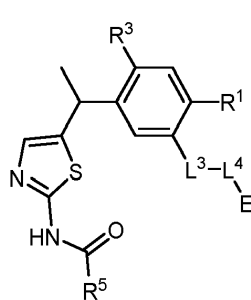
(IIA),



(II),

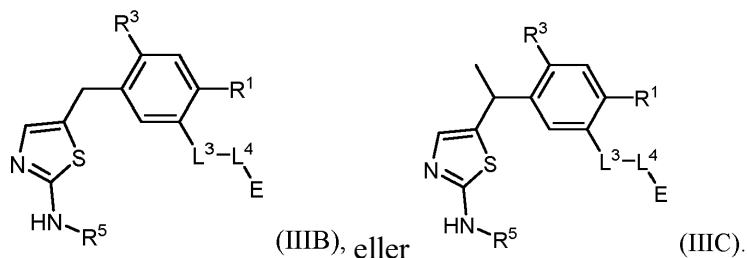
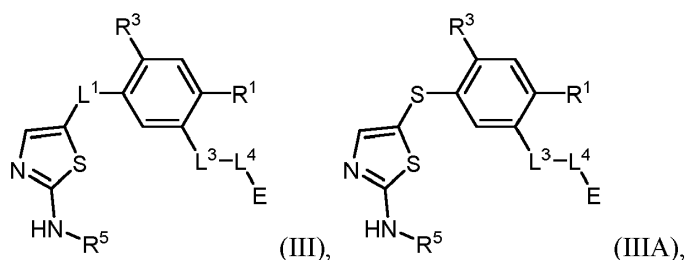


(IIB) eller



(IIC).

7. Forbindelsen ifølge krav 1, som har formelen:



- 5 **8.** Forbindelsen ifølge et hvilket som helst av kravene 1 til 7, hvori:
- (i) R^1 er uavhengig hydrogen, halogen, $-CX^1_3$, $-CHX^1_2$, $-CH_2X^1$, $-OCX^1_3$, $-OCH_2X^1$, $-OCHX^1_2$, substituert eller usubstituert alkyl eller substituert eller usubstituert heteroalkyl; eller
- (ii) R^1 er uavhengig hydrogen, halogen, $-CX^1_3$, $-CHX^1_2$, $-CH_2X^1$, $-OCX^1_3$, $-OCH_2X^1$, $-OCHX^1_2$, substituert eller usubstituert (C₁-C₄)alkyl eller substituert eller usubstituert 2- til 4-leddet heteroalkyl; eller
- 10 (iii) R^1 er hydrogen, halogen, $-CH_3$, $-CH_2CH_3$, $-CX^1_3$, $-CHX^1_2$, $-CH_2X^1$, $-OCH_3$, $-OCX^1_3$, $-OCH_2X^1$, $-OCHX^1_2$, $-SCH_3$, $-SCX^1_3$, $-SCH_2X^1$ eller $-SCHX^1_2$.
- 15
- 9.** Forbindelsen ifølge et hvilket som helst av kravene 1 til 8, hvori:
- (i) R^3 er uavhengig hydrogen, halogen, $-CX^3_3$, $-CHX^3_2$, $-CH_2X^3$, $-OCX^3_3$, $-OCH_2X^3$, $-OCHX^3_2$, $-CN$, substituert eller usubstituert alkyl eller substituert eller usubstituert heteroalkyl; eller
- 20 (ii) R^3 er uavhengig hydrogen, halogen, $-CX^3_3$, $-CHX^3_2$, $-CH_2X^3$, $-OCX^3_3$, $-OCH_2X^3$, $-OCHX^3_2$, $-CN$, substituert eller usubstituert (C₁-C₄)alkyl eller substituert eller usubstituert 2- til 4-leddet heteroalkyl; eller
- (iii) R^3 er hydrogen, halogen, $-CH_3$, $-CH_2CH_3$, $-CX^3_3$, $-CHX^3_2$, $-CH_2X^3$, $-OCH_3$, $-OCX^3_3$, $-OCH_2X^3$, $-OCHX^3_2$, $-CN$, $-SCH_3$, $-SCX^3_3$, $-SCH_2X^3$ eller $-SCHX^3_2$.
- 25

10. Forbindelsen ifølge et hvilket som helst av kravene 1 til 9, hvori:

(i) R^5 er uavhengig substituert eller usubstituert (C_1 - C_8)alkyl, substituert eller usubstituert 2- til 8-leddet heteroalkyl, substituert eller usubstituert (C_3 - C_6)sykloalkyl, substituert eller usubstituert 3- til 6-leddet heterosykloalkyl, substituert eller usubstituert fenyl eller substituert eller usubstituert 5- til 6-leddet heteroaryl; eller

(ii) R^5 er uavhengig substituert eller usubstituert (C_1 - C_4)alkyl; eller

(iii) R^5 er uavhengig usubstituert (C_1 - C_4)alkyl; eller

(iv) R^5 er uavhengig usubstituert metyl, usubstituert etyl, usubstituert isopropyl eller usubstituert tert-butyl; eller

(v) R^5 er uavhengig usubstituert metyl eller usubstituert etyl; eller

(vi) R^5 er uavhengig substituert eller usubstituert 2- til 8-leddet heteroalkyl; eller

(vii) R^5 er uavhengig substituert eller usubstituert 2- til 4-leddet heteroalkyl; eller

(viii) R^5 er uavhengig usubstituert 2- til 4-leddet heteroalkyl; eller

(ix) R^5 er uavhengig $-CH_2N(CH_3)_2$; eller

(x) R^5 er uavhengig substituert eller usubstituert (C_3 - C_6)sykloalkyl; eller

(xi) R^5 er uavhengig usubstituert (C_3 - C_6)sykloalkyl; eller

(xii) R^5 er uavhengig usubstituert syklopropyl, usubstituert syklobutyl eller usubstituert syklopentyl; eller

(xiii) R^5 er uavhengig usubstituert syklopropyl eller usubstituert syklobutyl; eller

(xiv) R^5 er uavhengig usubstituert syklopropyl; eller

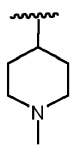
(xv) R^5 er uavhengig substituert eller usubstituert 3- til 6-leddet heterosykloalkyl; eller

(xvi) R^5 er uavhengig substituert eller usubstituert 5- til 6-leddet heterosykloalkyl; eller

(xvii) R^5 er uavhengig substituert eller usubstituert 6-leddet heterosykloalkyl; eller

(xviii) R^5 er uavhengig substituert eller usubstituert piperidinyll; eller

(xix) R^5 er uavhengig



; eller

(xx) R^5 er uavhengig substituert eller usubstituert fenyl; eller

(xxi) R^5 er uavhengig usubstituert fenyl; eller

(xxii) R^5 er uavhengig 2-substituert fenyl; eller

(xxiii) R^5 er uavhengig 3-substituert fenyl; eller

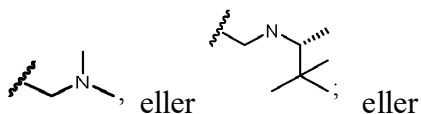
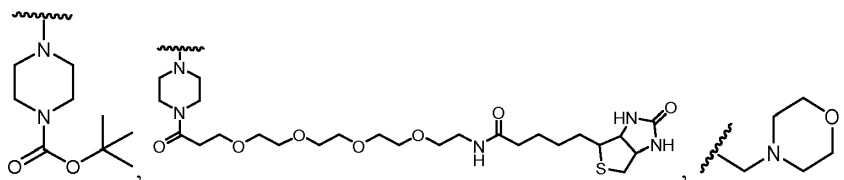
(xxiv) R^5 er uavhengig 4-substituert fenyl; eller

(xxv) R^5 er uavhengig fenyl substituert med halogen, substituert eller usubstituert alkyl, substituert eller usubstituert heteroalkyl, substituert eller usubstituert sykloalkyl, substituert eller usubstituert heterosykloalkyl, substituert eller usubstituert aryl eller substituert eller usubstituert heteroaryl; eller

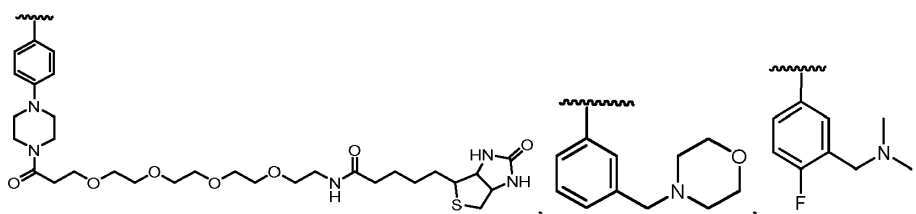
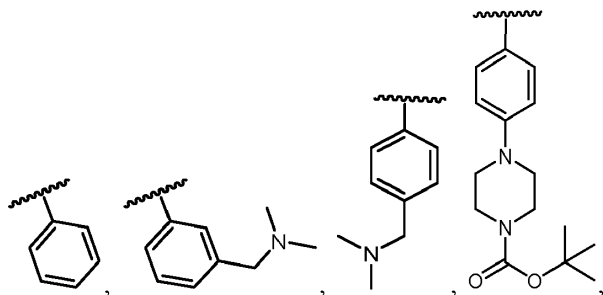
(xxvi) R^5 er uavhengig fenyl substituert med halogen, substituert eller usubstituert

(C_1 - C_8)alkyl, substituert eller usubstituert 2- til 8-leddet heteroalkyl, substituert eller usubstituert (C_3 - C_6)sykloalkyl, substituert eller usubstituert 3- til 6-leddet heterosykloalkyl, substituert eller usubstituert fenyl eller substituert eller usubstituert 5- til 6-leddet heteroaryl; eller

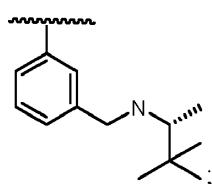
(xxvii) R^5 er uavhengig fenyl substituert med -F,



(xxviii) R^5 er uavhengig

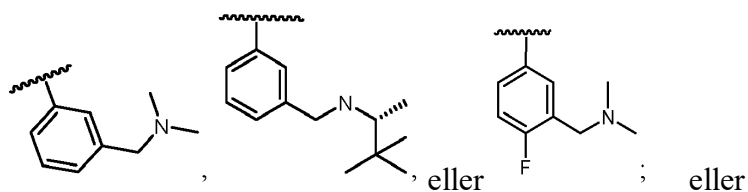


eller



eller

(xxix) R^5 er uavhengig



(xxx) R^5 er uavhengig substituert eller usubstituert heteroaryl; eller

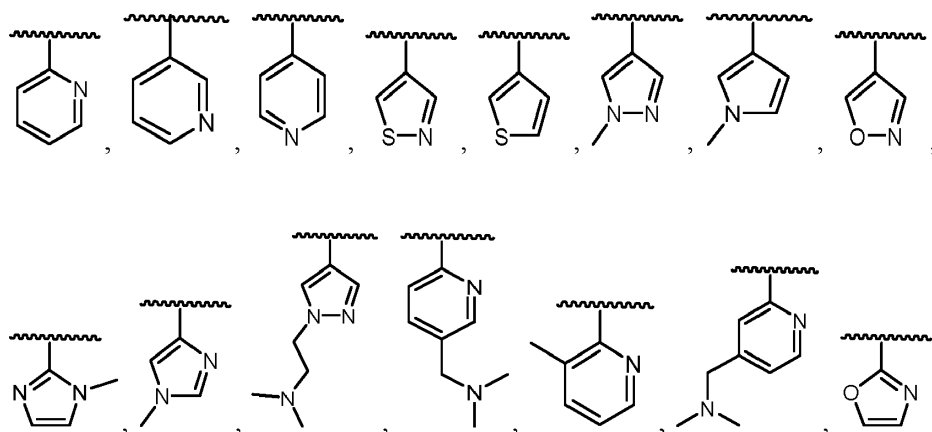
5 (xxxi) R^5 er uavhengig substituert eller usubstituert 5- til 6-leddet heteroaryl; eller

(xxxii) R^5 er uavhengig substituert eller usubstituert pyridyl, substituert eller

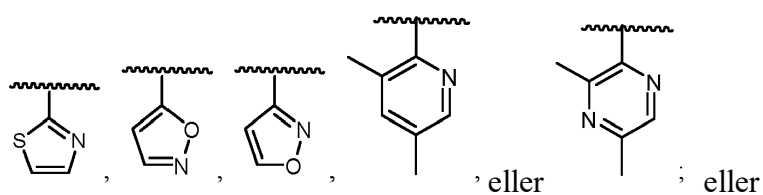
usubstituert tienyl, substituert eller usubstituert furanyl, substituert eller usubstituert pyrrolyl, substituert eller usubstituert imidazolyl, substituert eller usubstituert pyrazolyl, substituert eller usubstituert tiazolyl, substituert eller usubstituert isotiazolyl, substituert eller usubstituert

10 oksazolyl eller substituert eller usubstituert isoksazolyl; eller

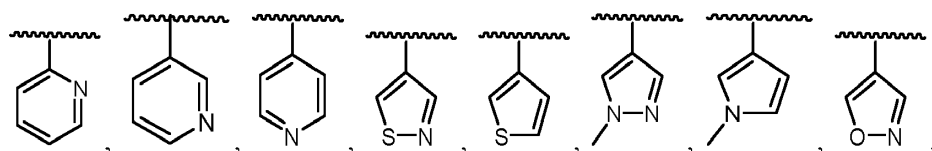
(xxxiii) R^5 er uavhengig

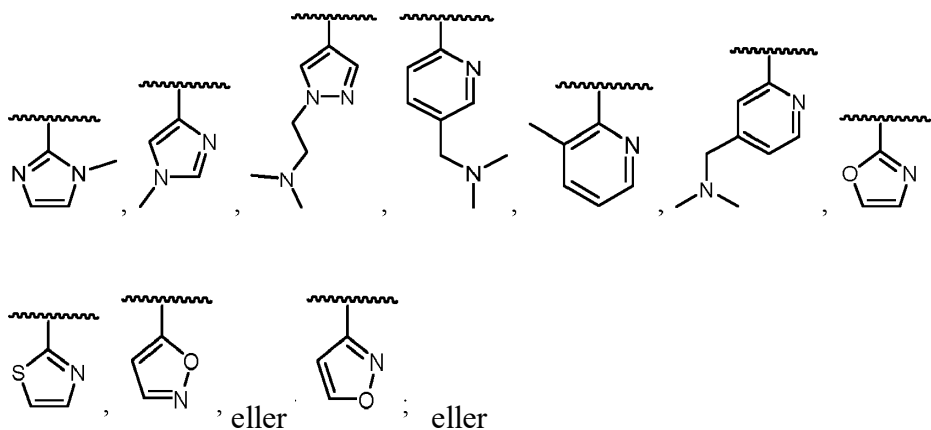


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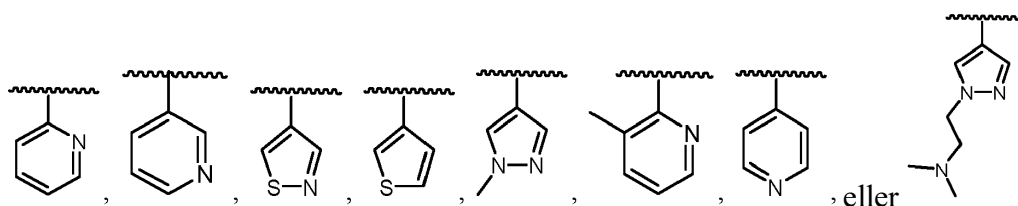


(xxxiv) R^5 er uavhengig





(xxxv) R^5 er uafhængig



11. Forbindelsen ifølge et hvilket som helst af kravene 1 til 10, hvori:

(i) L^3 er $-C(O)-$, en binding, $-N(R^6)-$ eller $-C(O)N(R^6)-$; og R^6 er uafhængig hydrogen, $-CX^6_3$, $-CHX^6_2$, $-CH_2X^6$ eller usubstitueret (C_1 - C_4)alkyl; eller

(ii) L^3 er $-C(O)-$, en binding, $-N(R^6)-$ eller $-C(O)N(R^6)-$; og R^6 er uafhængig hydrogen eller usubstitueret metyl; eller

(iii) L^3 er $-C(O)-$, en binding, $-C(O)N(CH_3)-$, $-N(CH_3)-$ eller $-NH-$; eller

(iv) L^3 er $-C(O)-$; eller

(v) L^3 er en binding; eller

(vi) L^3 er $-C(O)N(CH_3)-$.

12. Forbindelsen ifølge et hvilket som helst af kravene 1 til 11, hvori:

(i) L^4 er substitueret eller usubstitueret monosyklisk heterosykloalkylen, substitueret eller usubstitueret kondensert ringheterosykloalkylen, substitueret eller usubstitueret spirosyklisk heterosykloalkylen eller substitueret eller usubstitueret brodannet ringheterosykloalkylen; eller

(ii) L^4 er substitueret monosyklisk heterosykloalkylen, substitueret kondensert ringheterosykloalkylen, substitueret spirosyklisk heterosykloalkylen eller substitueret brodannet ringheterosykloalkylen; eller

(iii) L^4 er usubstitueret monosyklisk heterosykloalkylen, usubstitueret kondensert ringheterosykloalkylen, usubstitueret spirosyklisk heterosykloalkylen eller usubstitueret brodannet ringheterosykloalkylen; eller

(iv) L^4 er et substituert eller usubstituert 4- til 10-leddet monosyklisk heterosykloalkylen, substituert eller usubstituert 5- til 10-leddet kondensert ringheterosykloalkylen, substituert eller usubstituert 6- til 10-leddet spirosyklisk heterosykloalkylen eller substituert eller usubstituert 5- til 10-leddet brodannet ringheterosykloalkylen; eller

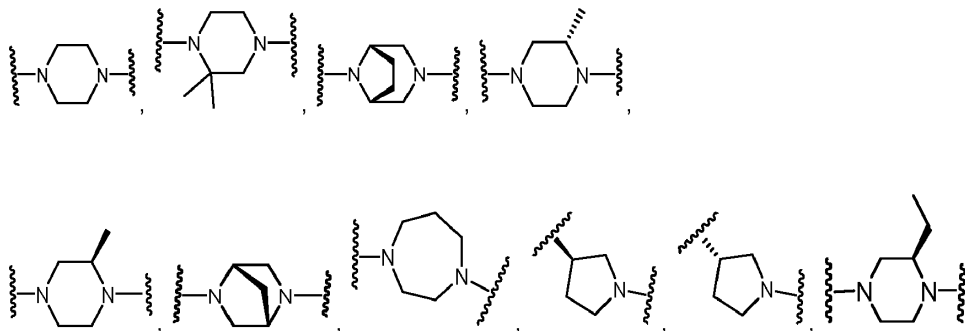
(v) L^4 er et substituert eller usubstituert 5- til 10-leddet monosyklisk heterosykloalkylen, substituert eller usubstituert 6- til 10-leddet kondensert ringheterosykloalkylen, substituert eller usubstituert 7- til 10-leddet spirosyklisk heterosykloalkylen eller substituert eller usubstituert 5- til 10-leddet brodannet ringheterosykloalkylen; eller

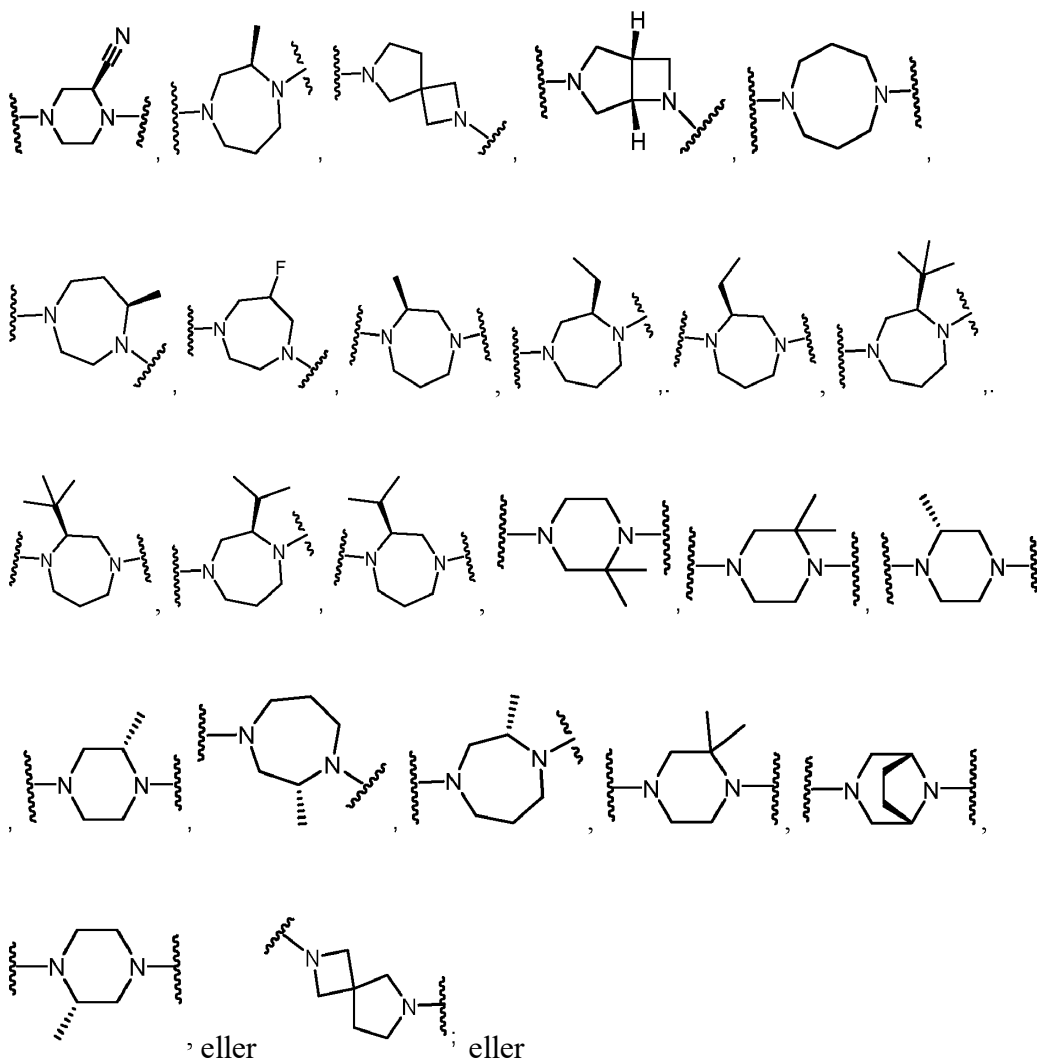
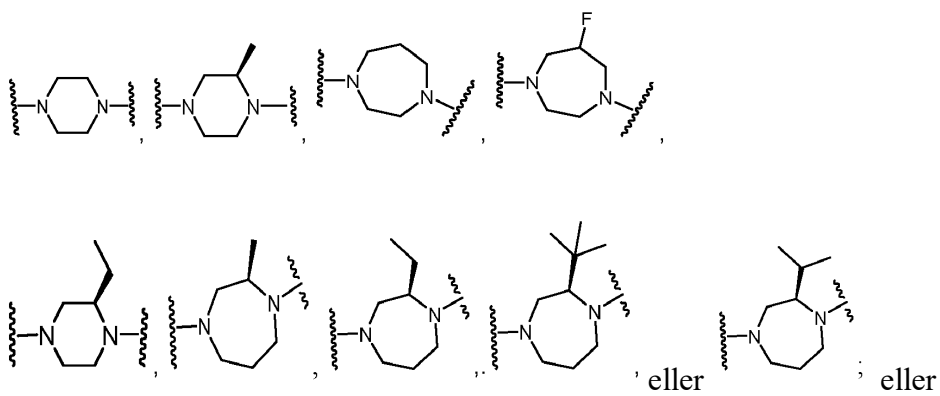
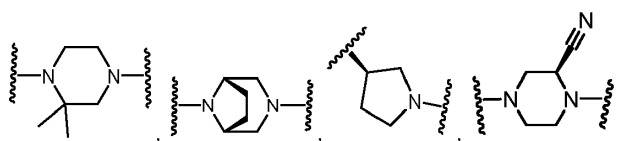
(vi) L^4 er et substituert eller usubstituert 5- til 8-leddet monosyklisk heterosykloalkylen, substituert eller usubstituert 7- til 8-leddet kondensert ringheterosykloalkylen, substituert eller usubstituert 7- til 8-leddet spirosyklisk heterosykloalkylen eller substituert eller usubstituert 7- til 8-leddet brodannet ringheterosykloalkylen; eller

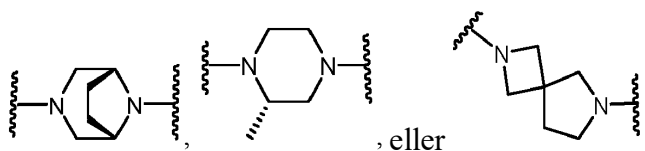
(vii) L^4 er et usubstituert 7- til 8-leddet brodannet ringheterosykloalkylen; eller
 (viii) L^4 er et usubstituert 7- til 8-leddet kondensert ringheterosykloalkylen; eller
 (ix) L^4 er et usubstituert 7- til 8-leddet spirosyklisk heterosykloalkylen; eller
 (x) L^4 er et metyl-substituert, etyl-substituert, cyano-substituert, halogen-substituert eller usubstituert 5- til 8-leddet monosyklisk heterosykloalkylen; eller

(xi) L^4 er et usubstituert 5- til 8-leddet monosyklisk heterosykloalkylen; eller
 (xii) L^4 er et metyl-substituert 5- til 8-leddet monosyklisk heterosykloalkylen, et etyl-substituert 5- til 8-leddet monosyklisk heterosykloalkylen, et isopropyl-substituert 5- til 8-leddet monosyklisk heterosykloalkylen eller et tert-butyl-substituert 5- til 8-leddet monosyklisk heterosykloalkylen; eller

(xiii) L^4 er



(xiv) L⁴ er(xv) L⁴ er



13. Forbindelsen ifølge et hvilket som helst av kravene 1 til 12, hvori:

Alternativ A:

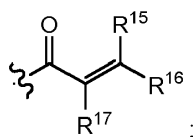
5 R^{15} er uavhengig hydrogen, halogen, CX^{15}_3 , $-CHX^{15}_2$, $-CH_2X^{15}$, $-CN$, $-SR^{15D}$,
 $-NR^{15A}R^{15B}$, $-C(O)R^{15C}$, $-C(O)-OR^{15C}$, $-C(O)NR^{15A}R^{15B}$, $-OR^{15D}$, $-NR^{15A}C(O)R^{15C}$,
 $-NR^{15A}C(O)OR^{15C}$, $-OCX^{15}_3$, $-OCHX^{15}_2$, substituert eller usubstituert alkyl, substituert
eller usubstituert heteroalkyl, substituert eller usubstituert sykloalkyl, substituert eller
usubstituert heterosykloalkyl, substituert eller usubstituert aryl, substituert eller usubstituert
10 heteroaryl;

R^{16} er uavhengig hydrogen, halogen, CX^{16}_3 , $-CHX^{16}_2$, $-CH_2X^{16}$, $-CN$, $-SR^{16D}$,
 $-NR^{16A}R^{16B}$, $-C(O)R^{16C}$, $-C(O)-OR^{16C}$, $-C(O)NR^{16A}R^{16B}$, $-OR^{16D}$, $-NR^{16A}C(O)R^{16C}$,
 $-NR^{16A}C(O)OR^{16C}$, $-OCX^{16}_3$, $-OCHX^{16}_2$, substituert eller usubstituert alkyl, substituert
eller usubstituert heteroalkyl, substituert eller usubstituert sykloalkyl, substituert eller
15 usubstituert heterosykloalkyl, substituert eller usubstituert aryl, substituert eller usubstituert
heteroaryl; og

R^{17} er uavhengig hydrogen, halogen, $-CX^{17}_3$, $-CHX^{17}_2$, $-CH_2X^{17}$, $-CN$, $-SR^{17D}$,
 $-NR^{17A}R^{17B}$, $-C(O)R^{17C}$, $-C(O)-OR^{17C}$, $-C(O)NR^{17A}R^{17B}$, $-OR^{17D}$, $-NR^{17A}C(O)R^{17C}$,
 $-NR^{17A}C(O)OR^{17C}$, $-OCX^{17}_3$, $-OCHX^{17}_2$, substituert eller usubstituert alkyl eller
20 substituert eller usubstituert heteroalkyl; eller

Alternativ B: R^{15} , R^{16} , R^{17} og R^{18} er hydrogen; eller

Alternativ C: E er:



hensiktsmessig i alternativ C

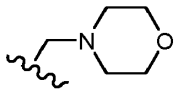
25 R^{15} er hydrogen;

R^{16} er hydrogen; og

R^{17} er hydrogen; eller

hensiktsmessig i alternativ C

R^{15} er hydrogen;

R^{16} er hydrogen, $-CH_3$, $-CH_2NR^{16A}R^{16B}$ eller  ;

R^{17} er hydrogen; og

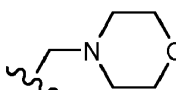
R^{16A} og R^{16B} er uavhengig hydrogen eller usubstituert alkyl, for eksempel er R^{16A} og

R^{16B} uavhengig usubstituert metyl; eller

5 hensiktsmessig i alternativ C

R^{15} er hydrogen;

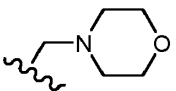
R^{16} er hydrogen;

R^{17} er hydrogen, $-CH_3$, $-CH_2NR^{17A}R^{17B}$ eller  ; og

R^{17A} og R^{17B} er uavhengig hydrogen eller usubstituert alkyl, for eksempel hvori R^{17A} og

10 R^{17B} er uavhengig usubstituert metyl; eller

hensiktsmessig i alternativ C

R^{15} er hydrogen, $-CH_3$, $-CH_2NR^{15A}R^{15B}$ eller  ;

R^{16} er hydrogen;

R^{17} er hydrogen; og

15 R^{15A} og R^{15B} er uavhengig hydrogen eller usubstituert alkyl, for eksempel hvori R^{15A} og

R^{15B} er uavhengig usubstituert metyl; eller

Alternativ D: E er

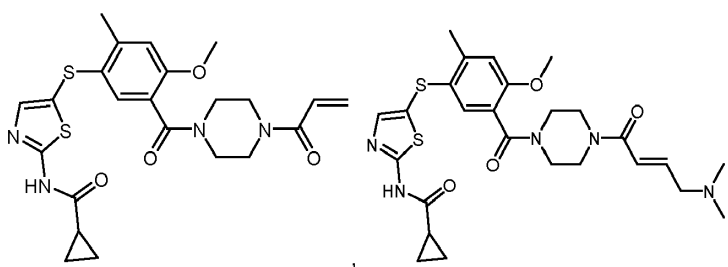
$-C(O)CH=CH_2$, $-C(O)CH=CHCH_2N(CH_3)_2$, $-C(O)C(=CH_2)CH_2N(CH_3)_2$,

$-C(O)C\equiv CCH_3$ eller $-C(O)C(=CH_2)CH_3$.

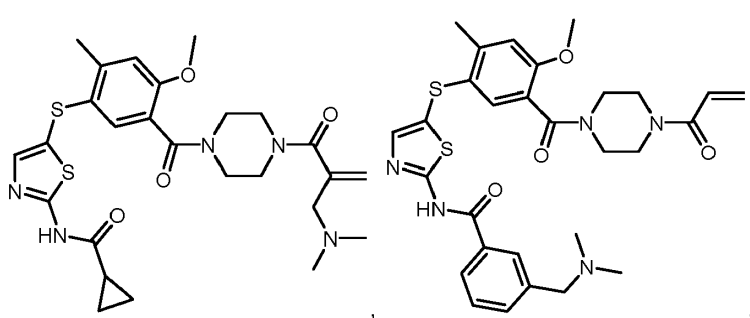
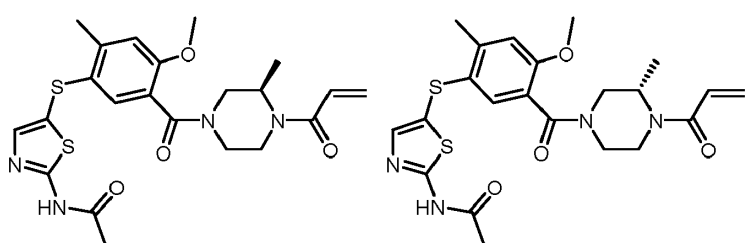
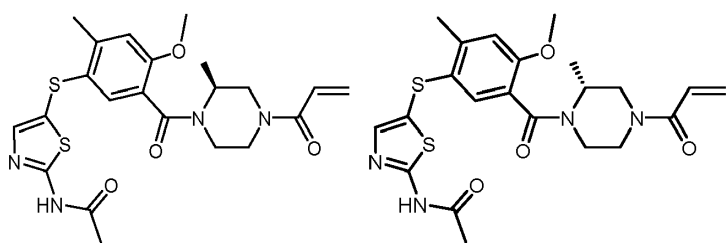
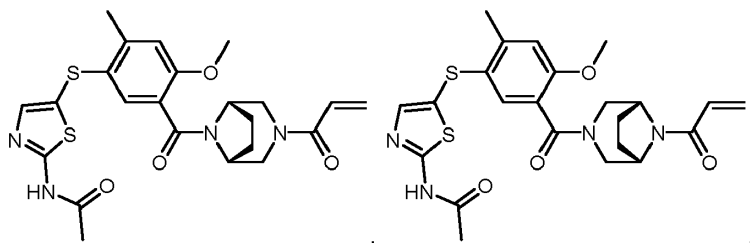
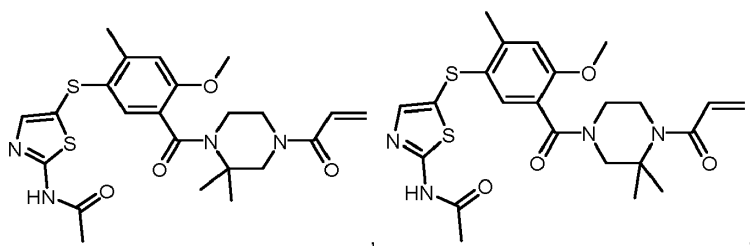
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14. Forbindelse ifølge krav 1, hvori forbindelsen er en forbindelse valgt fra én av forbindelsene oppført i alternativ A, B, C, D, E, F, G, H, I, J, K, L, M, N, O, P, Q, R, S eller T:

Alternativ A:

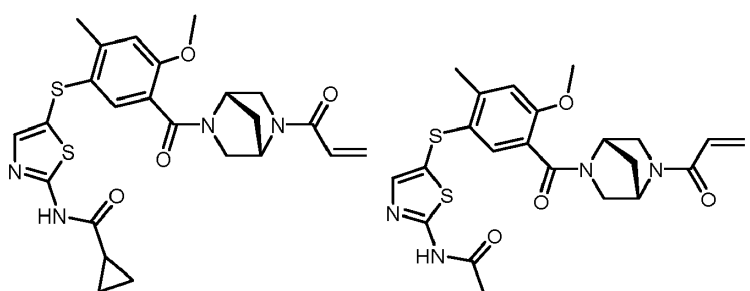
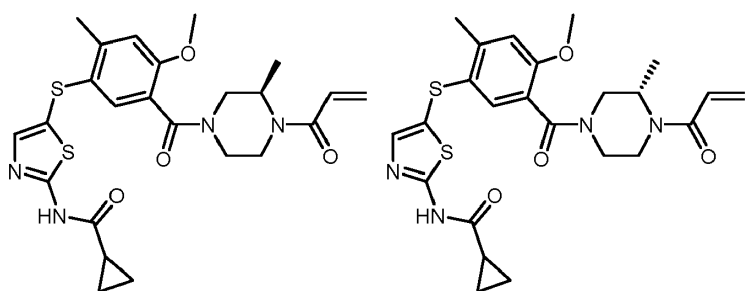
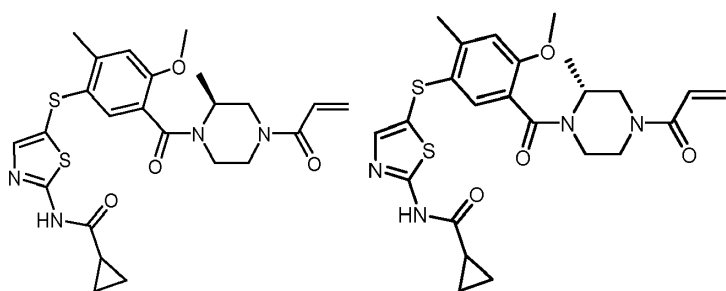
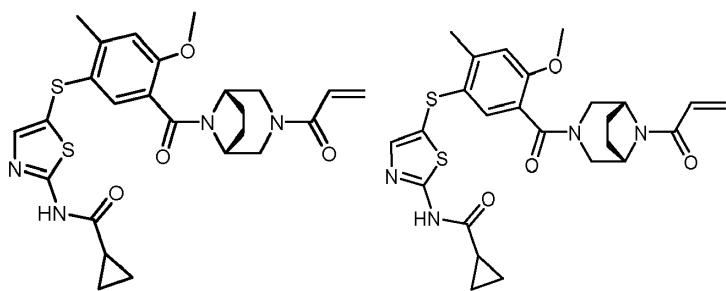
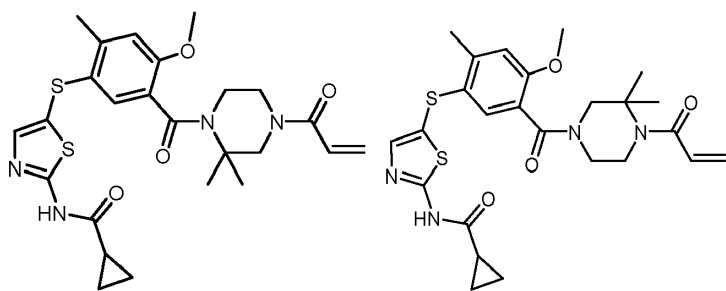


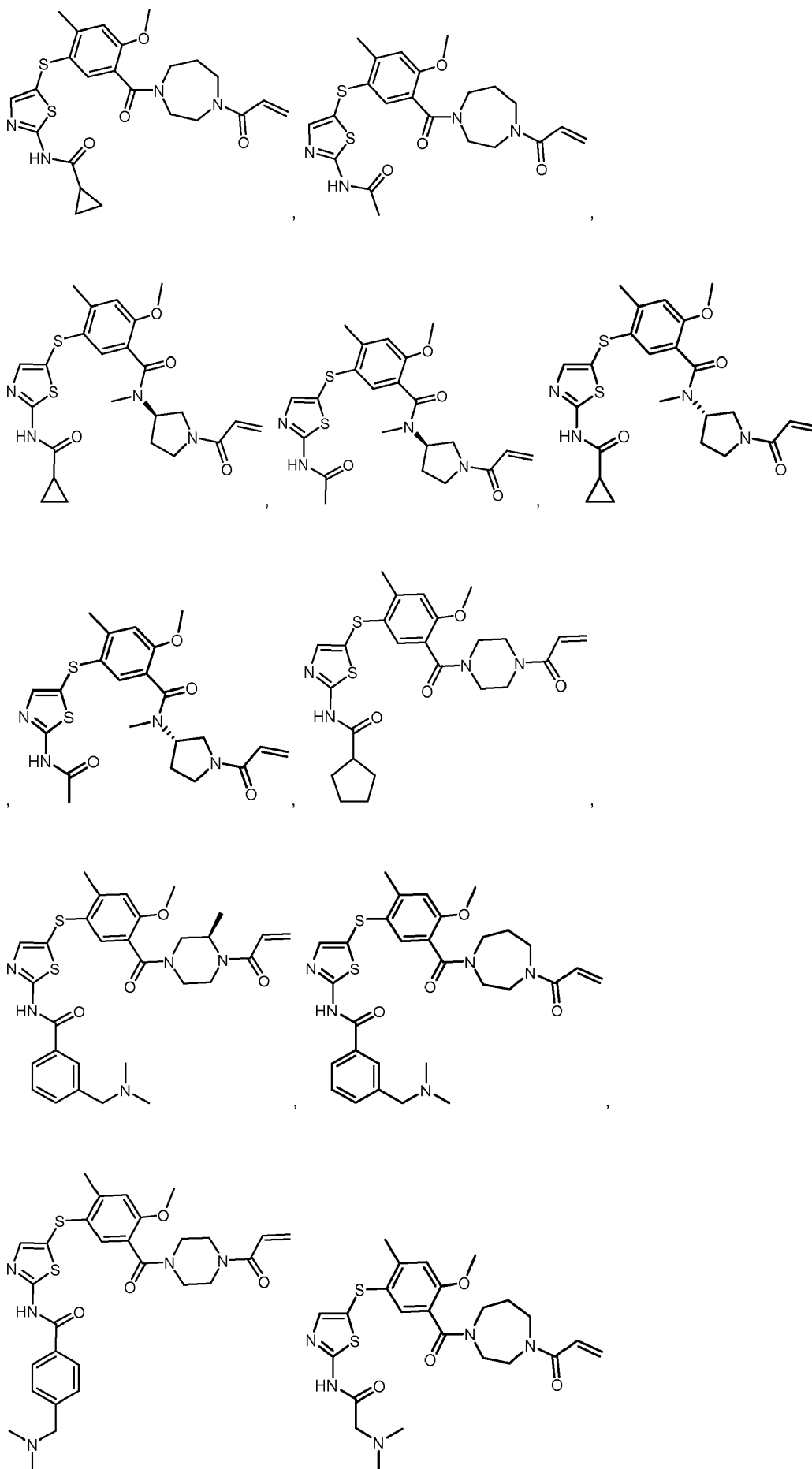
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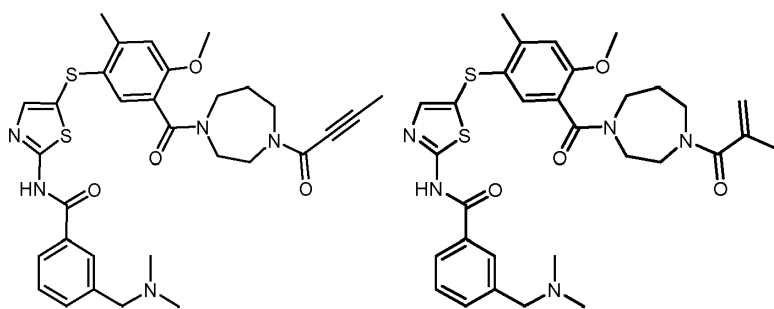
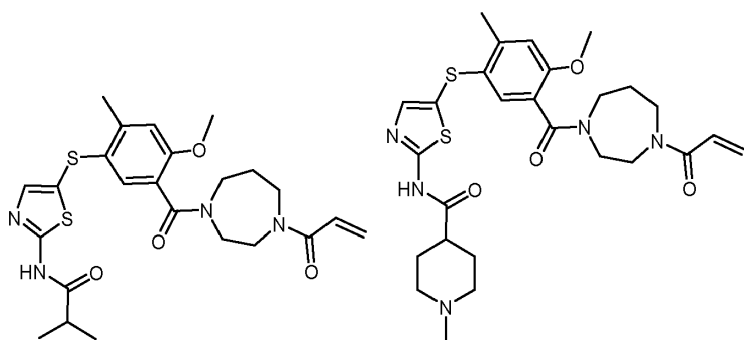


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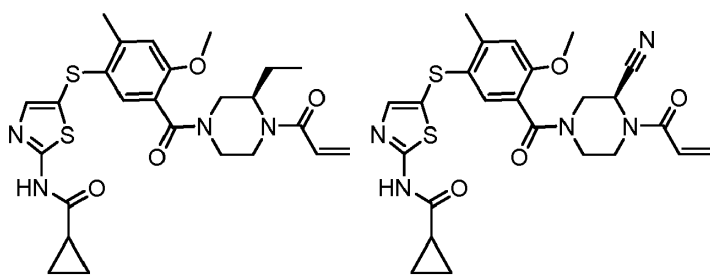
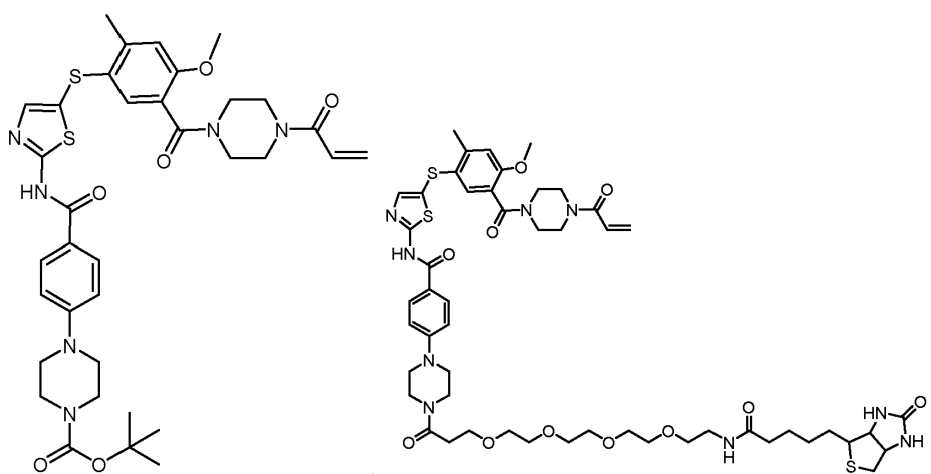
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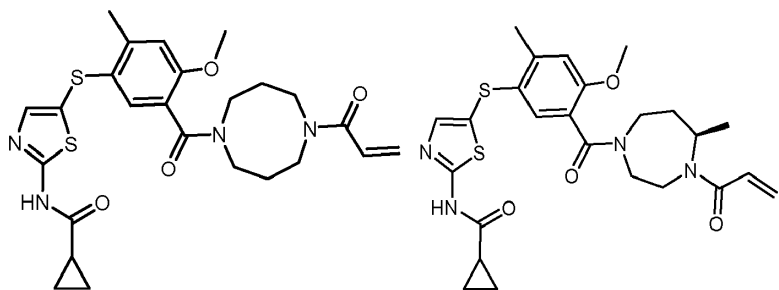
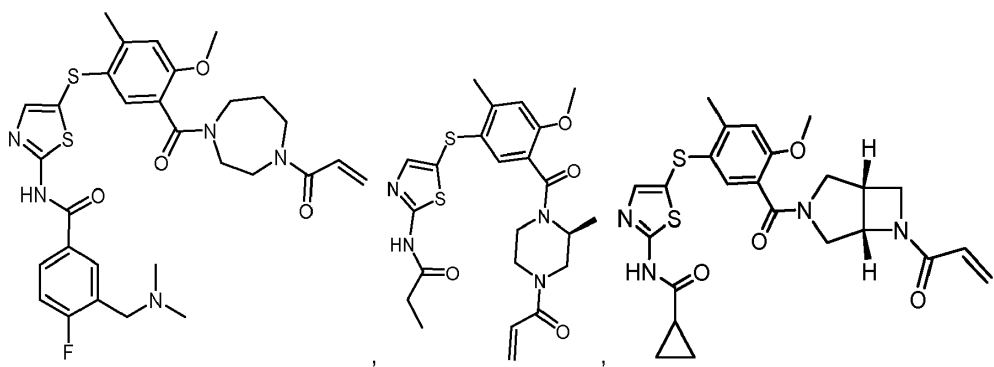
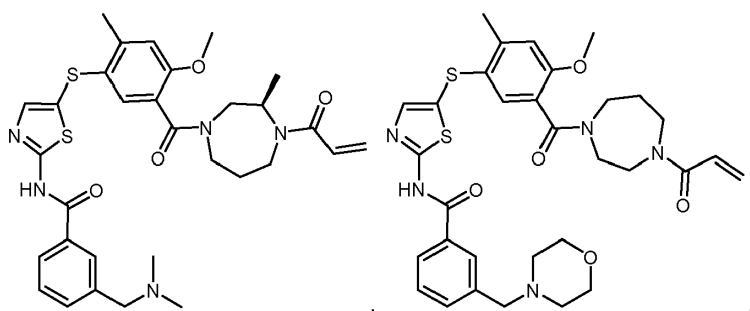
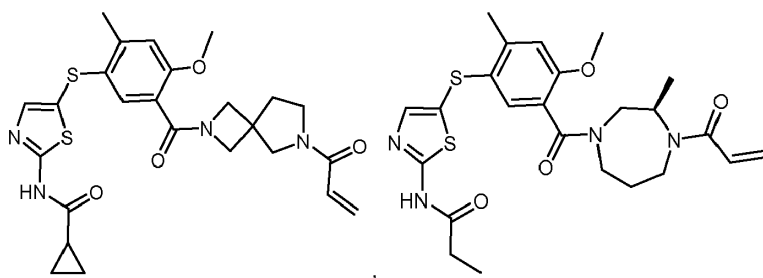
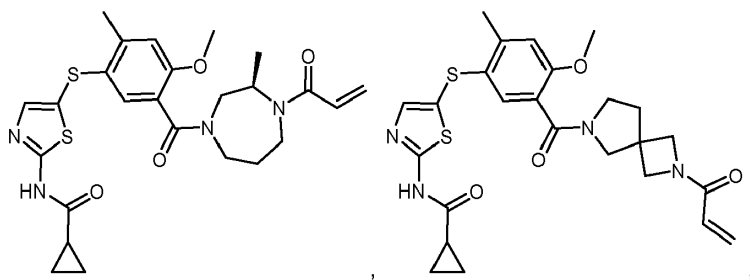


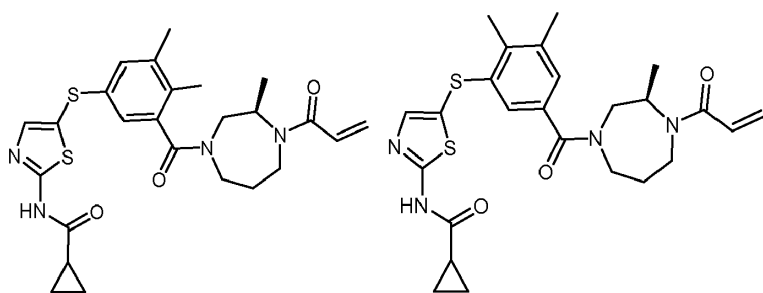


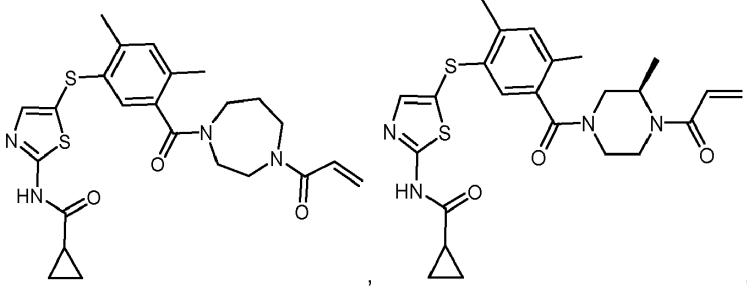
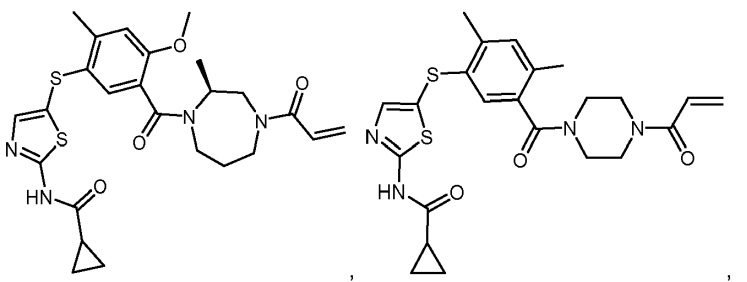
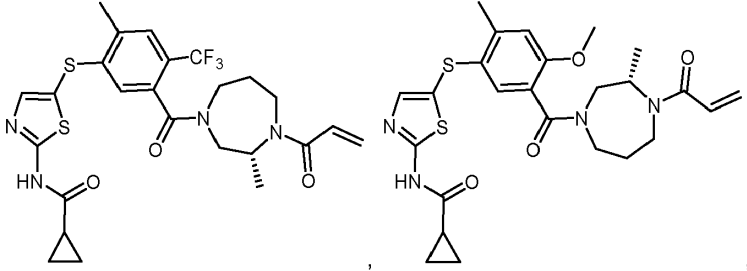
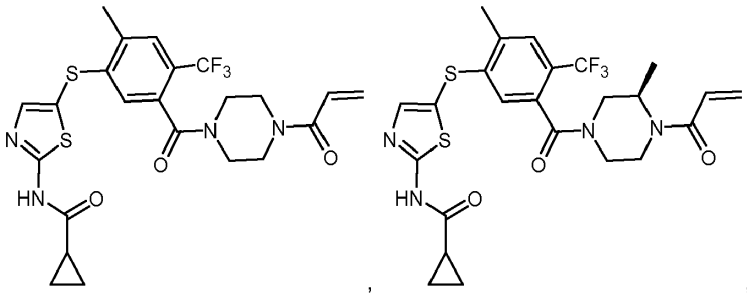
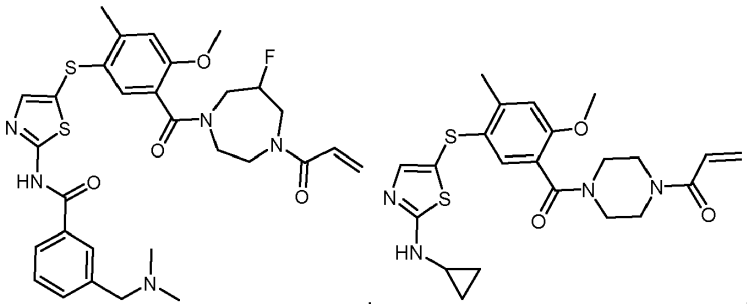


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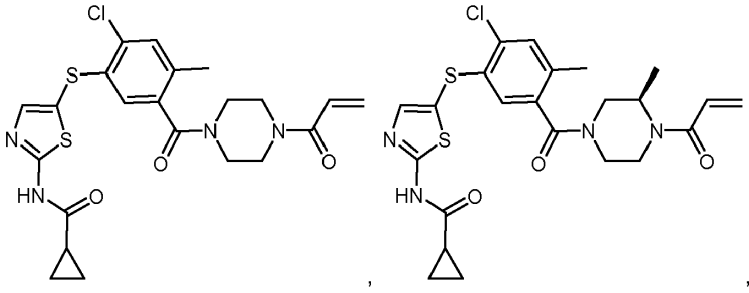
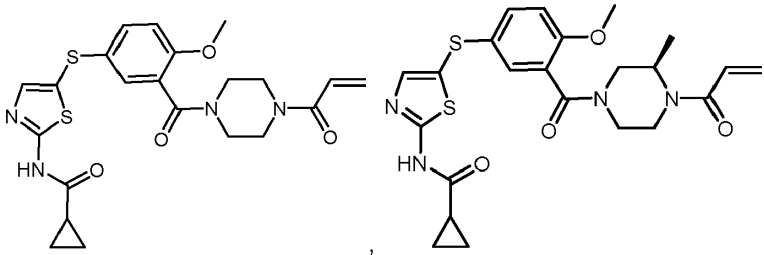
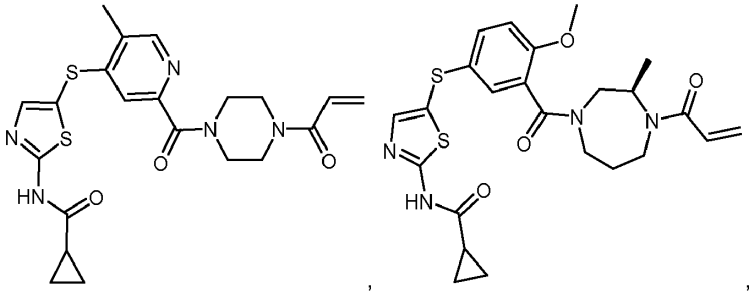




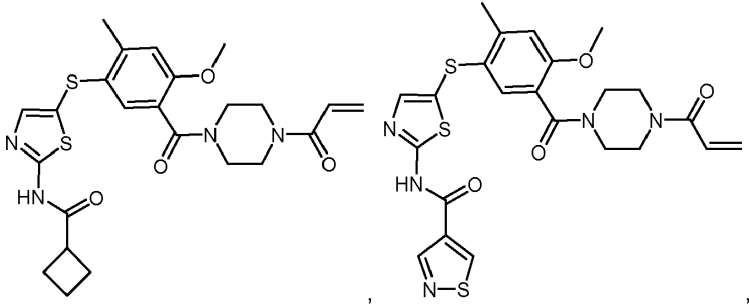
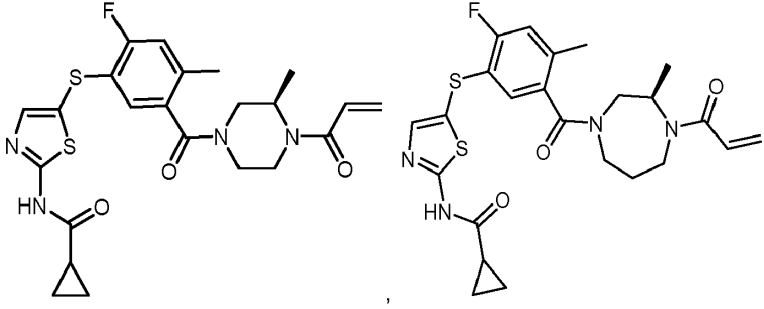


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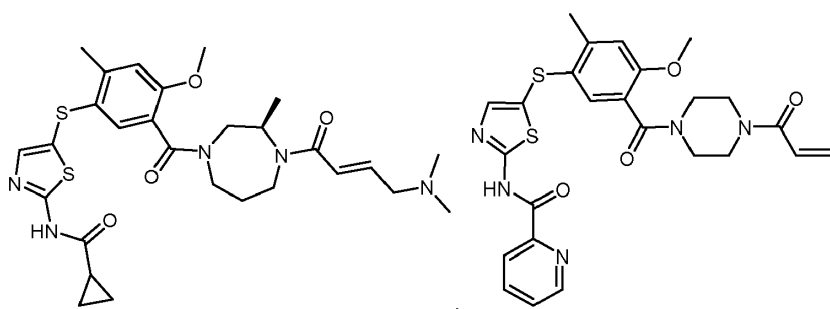
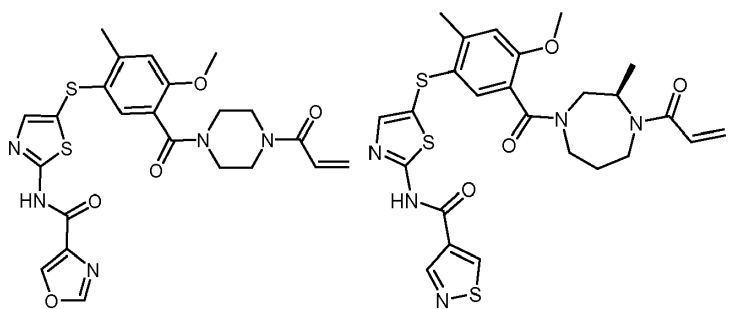
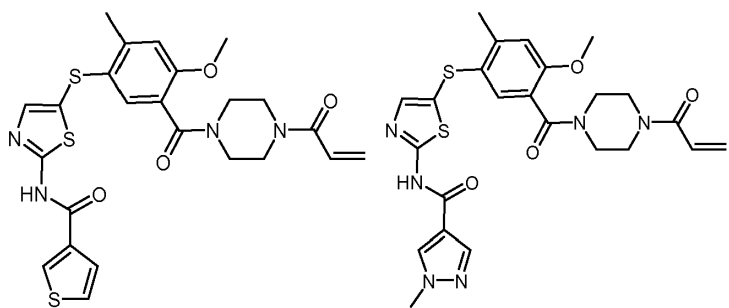
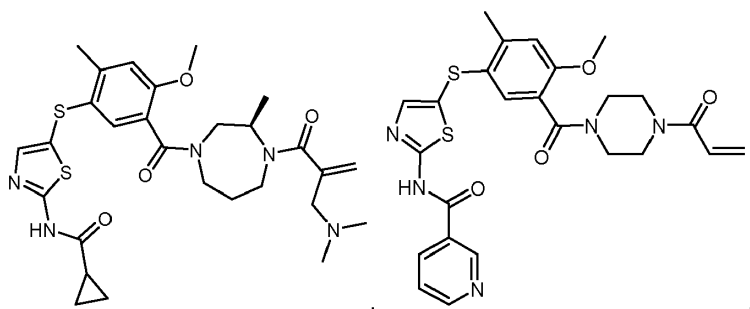
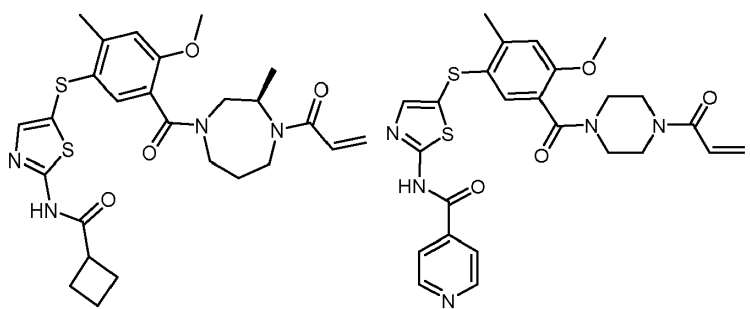
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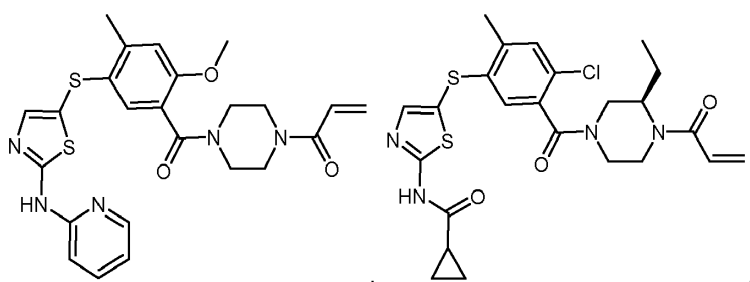
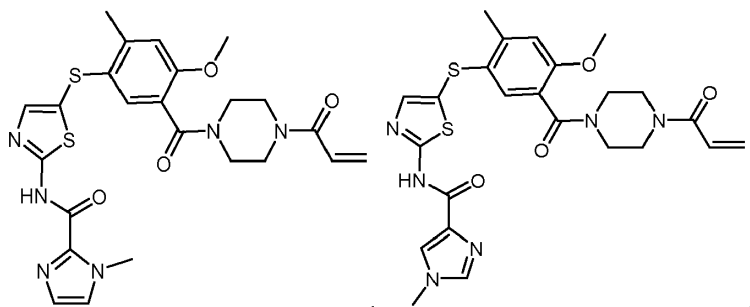


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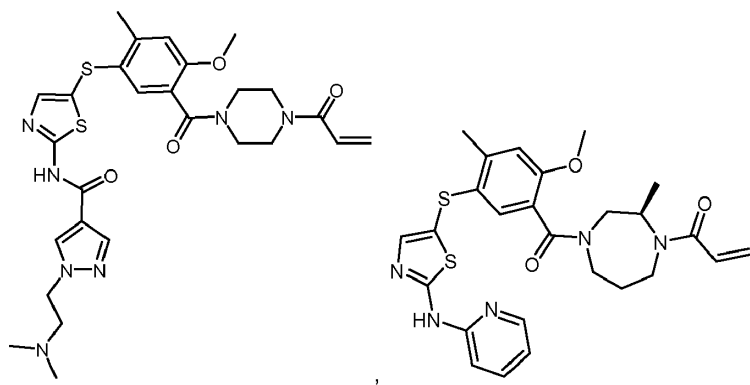
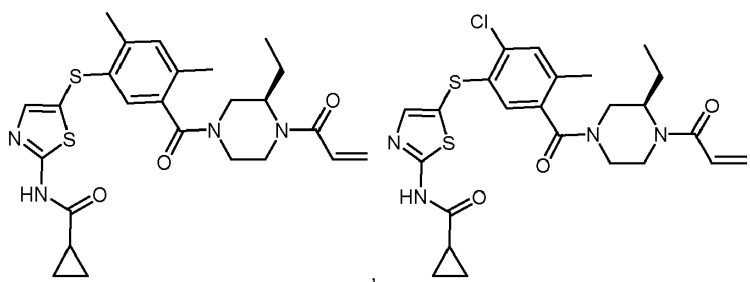


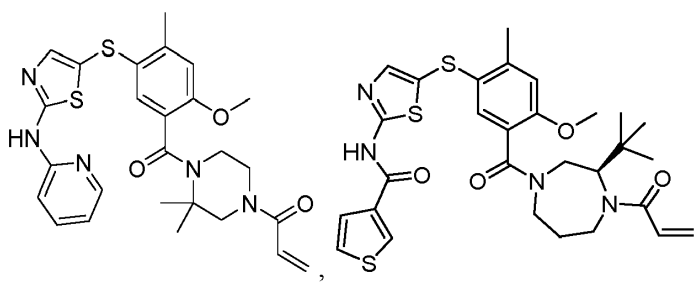
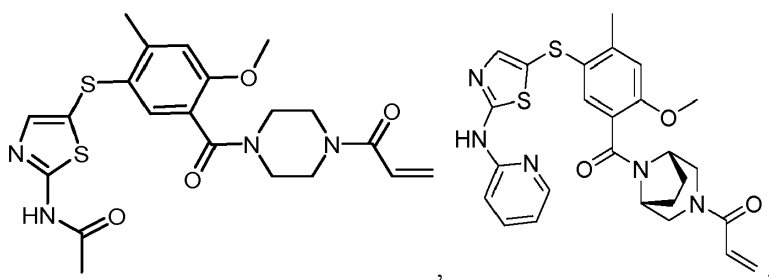
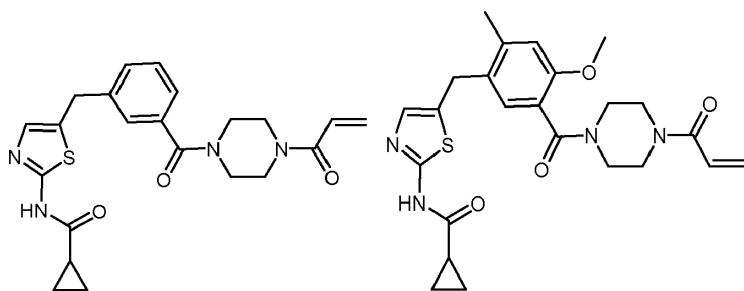
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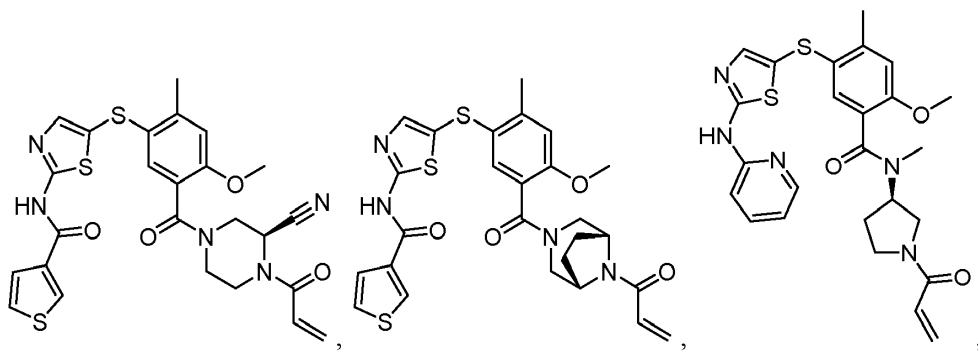
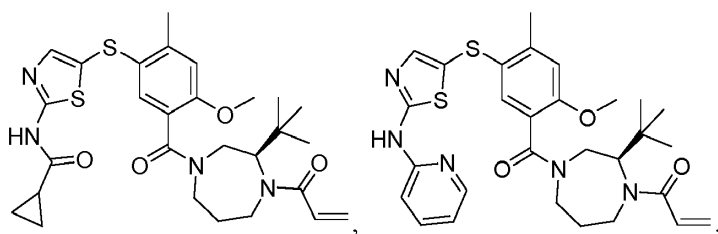


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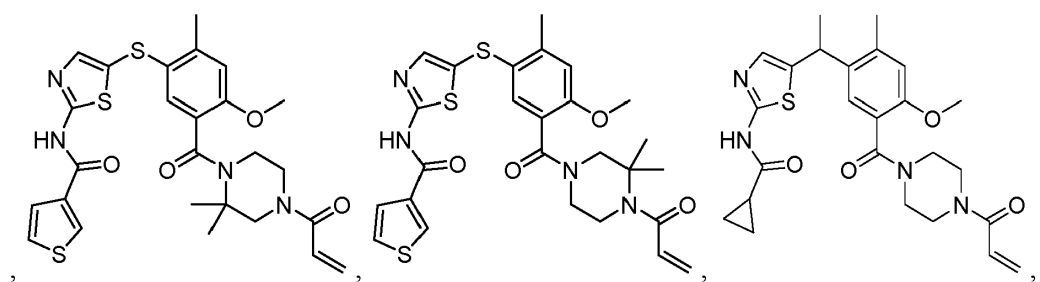
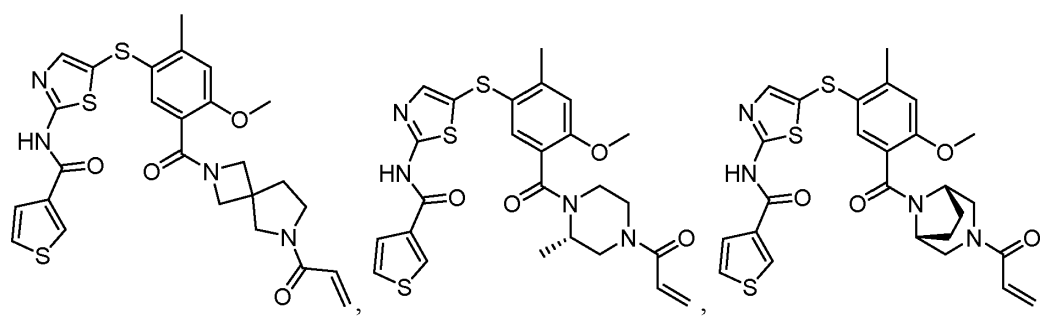
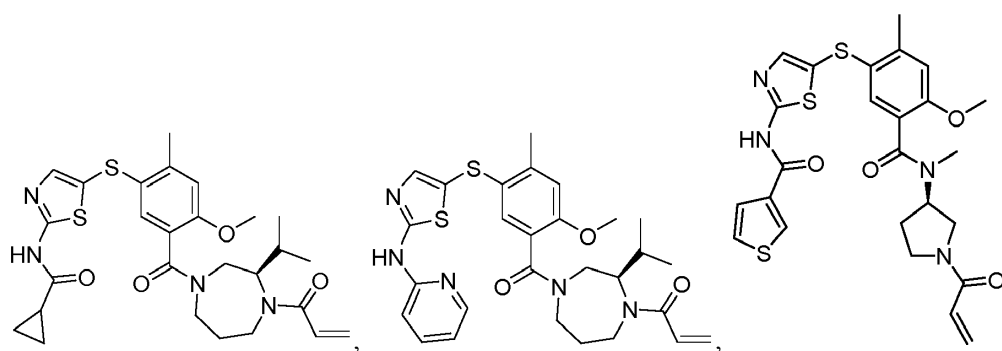
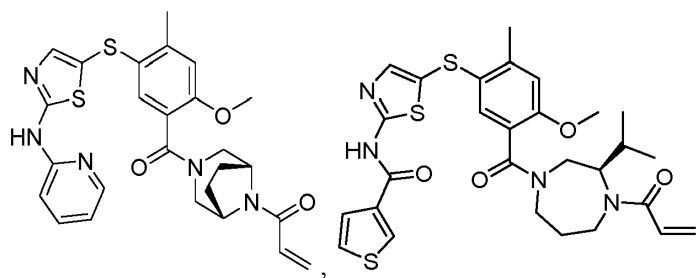
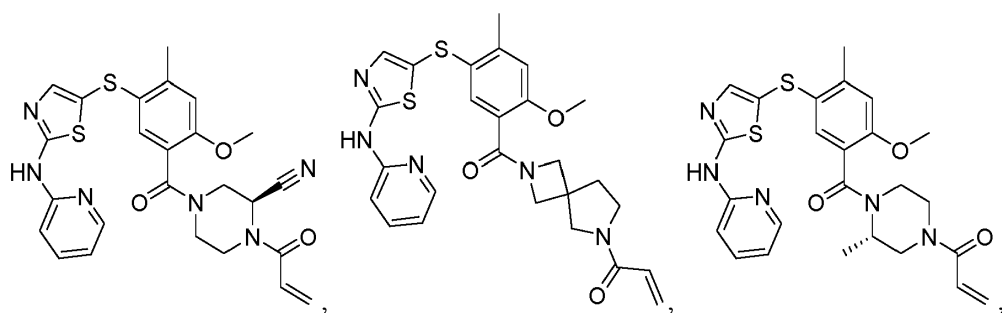




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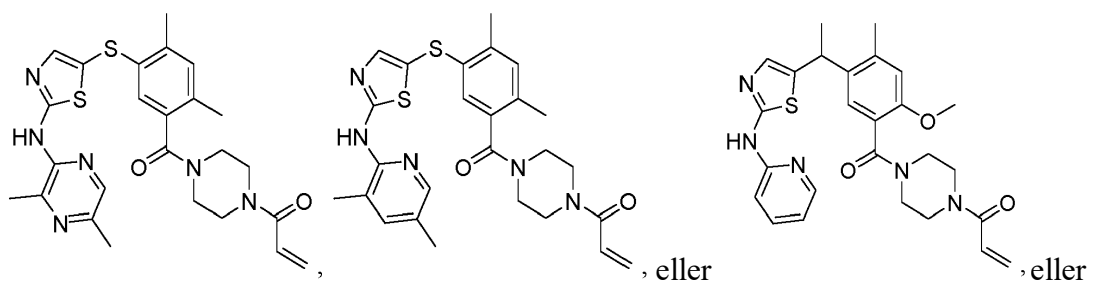


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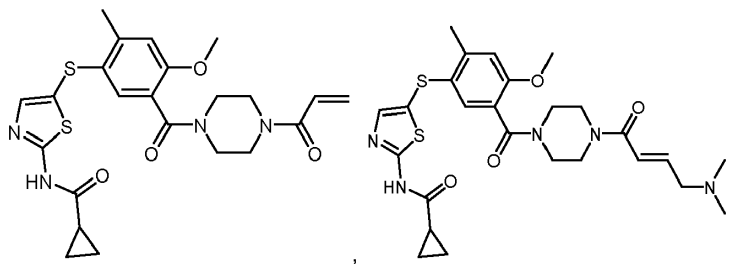
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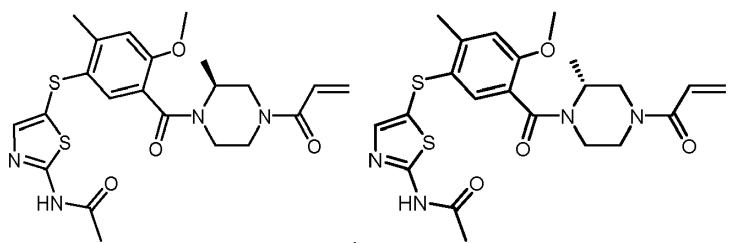
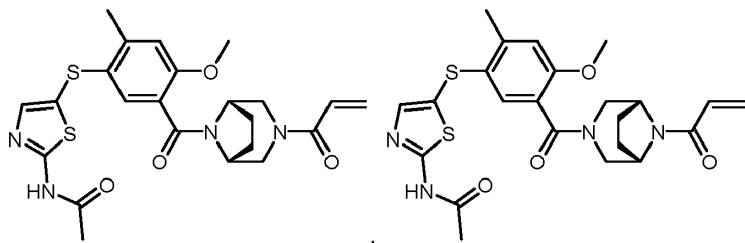
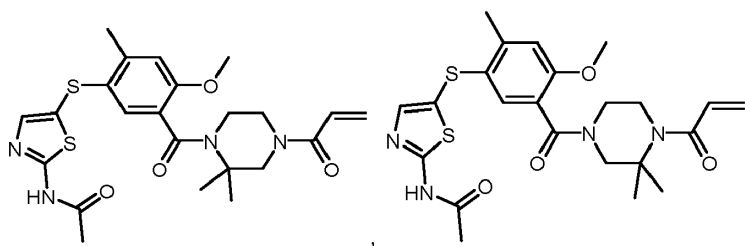


et farmasøytisk akseptabelt salt derav; eller

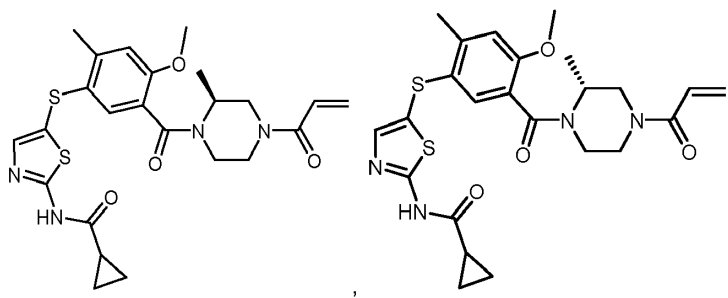
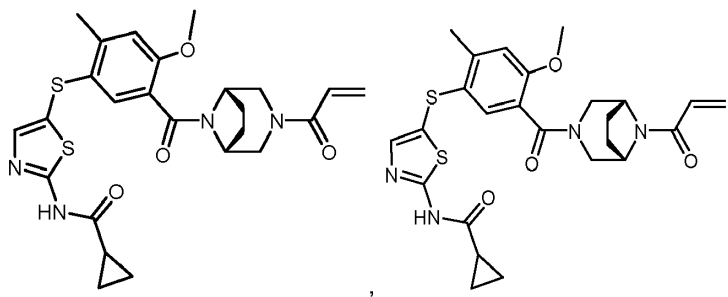
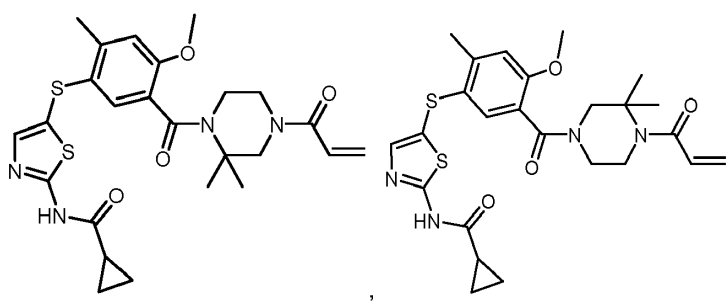
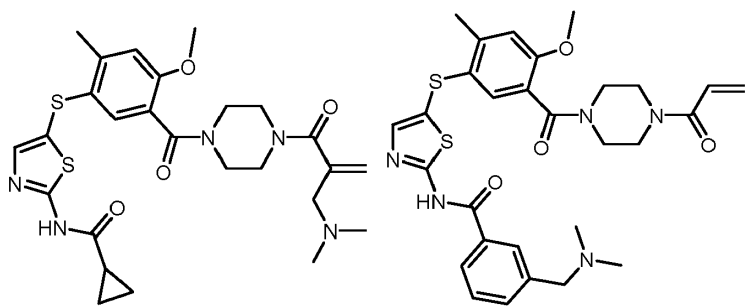
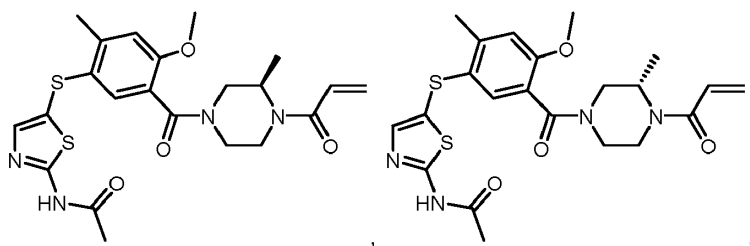
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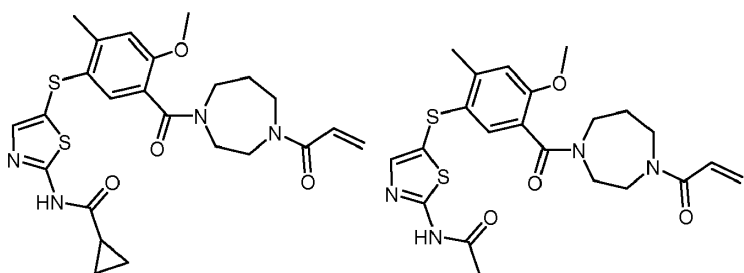
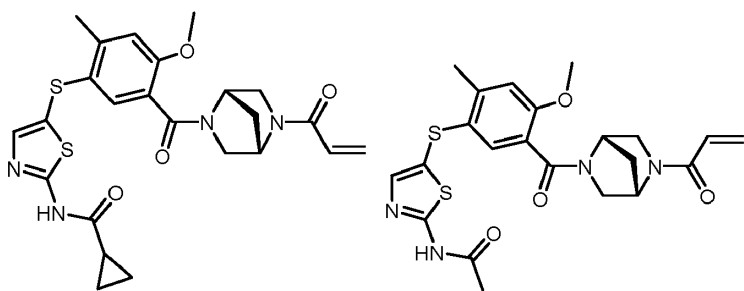
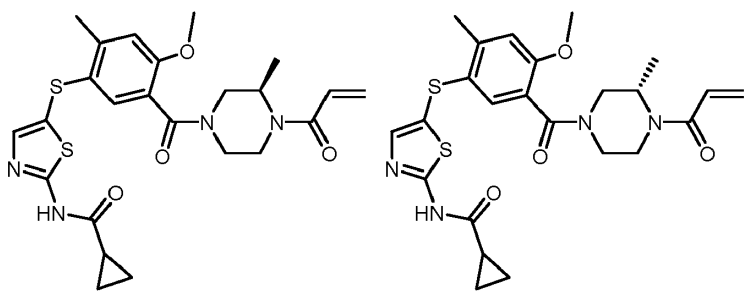


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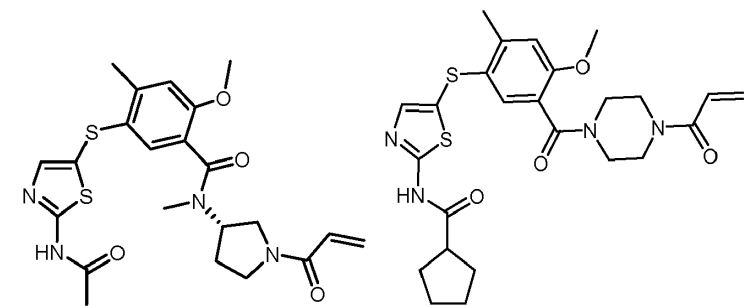
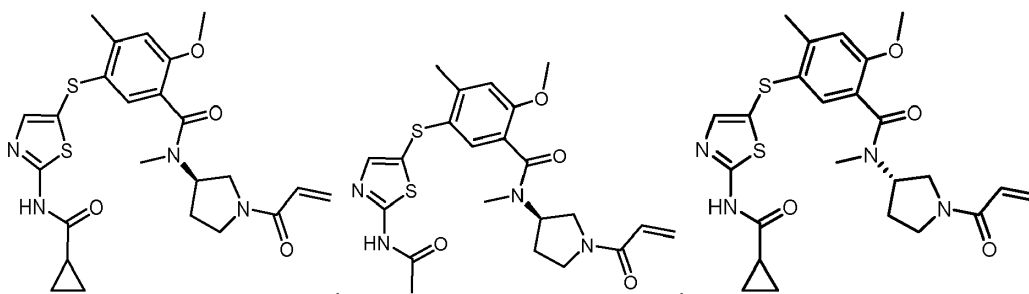


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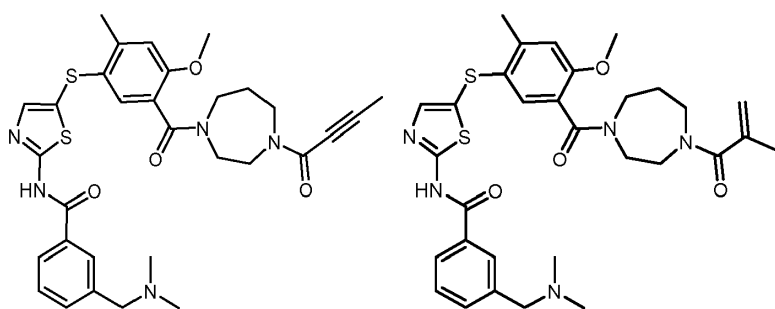
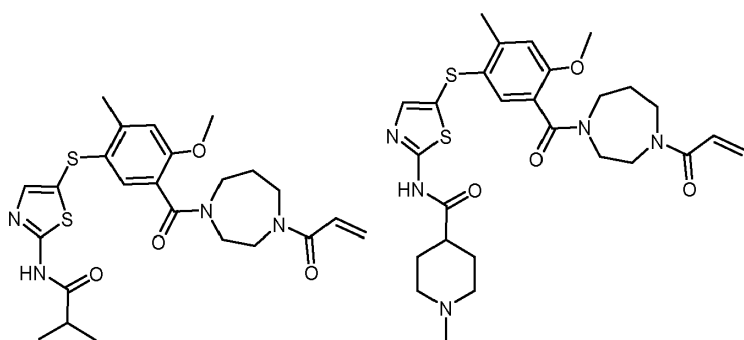
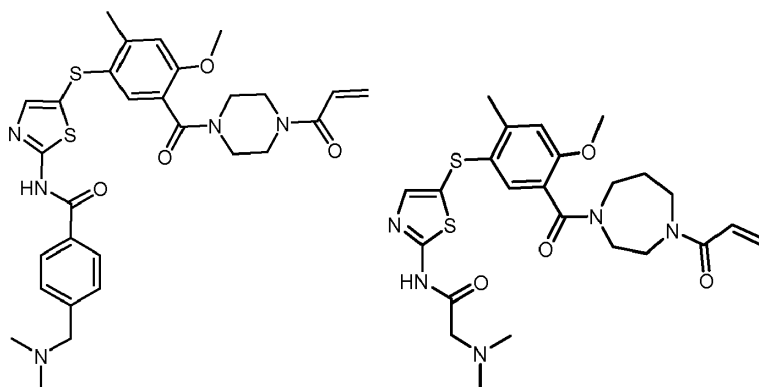
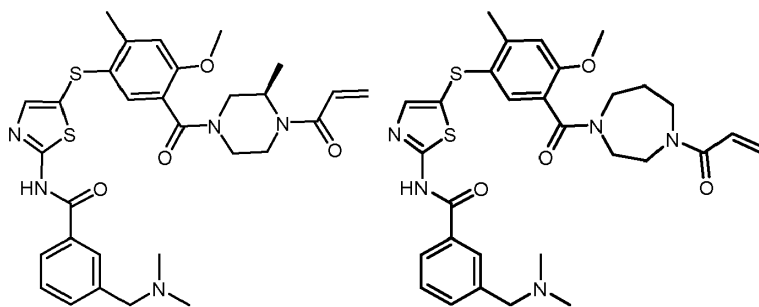


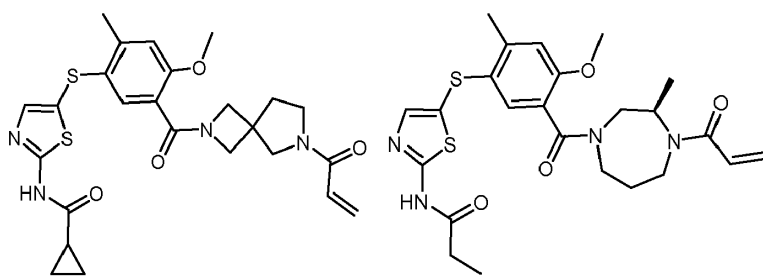
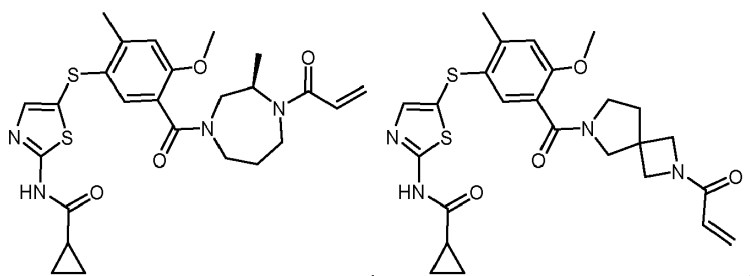
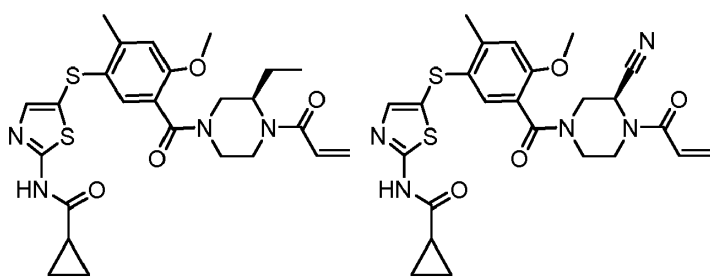
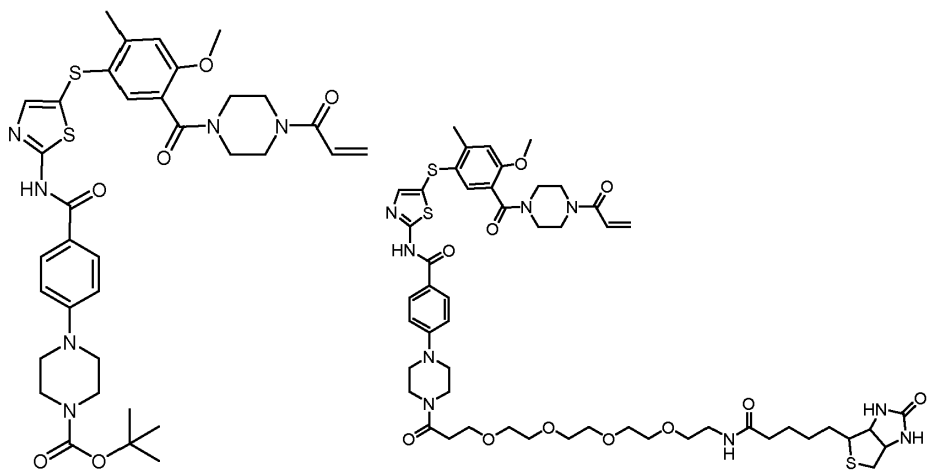


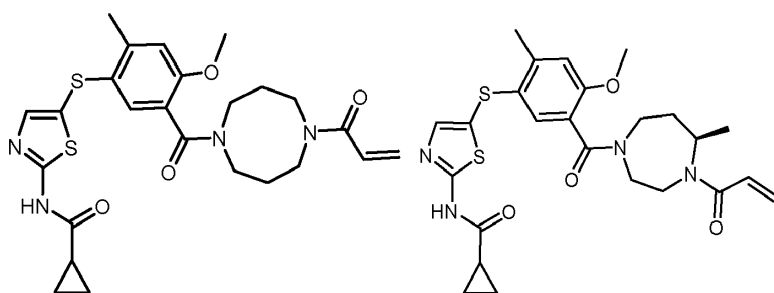
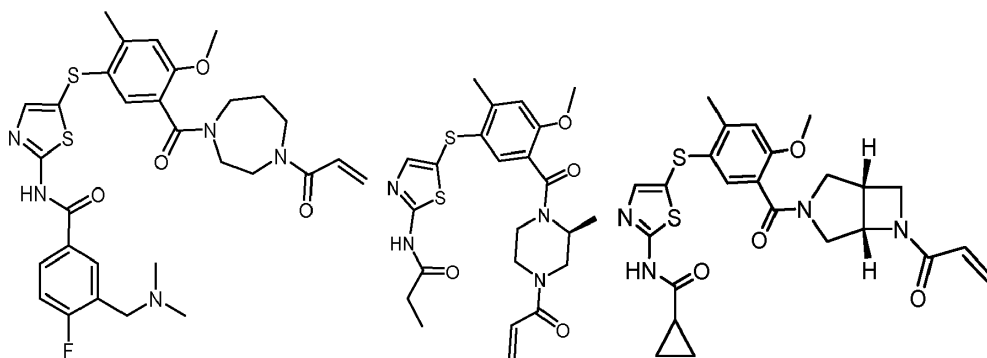
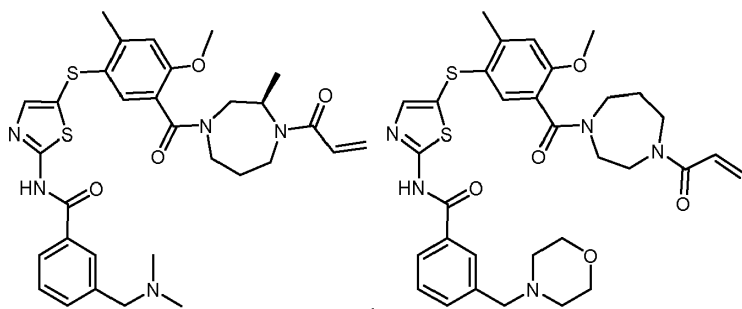
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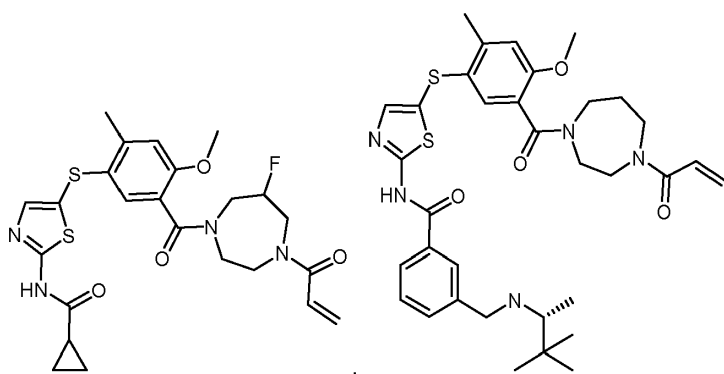
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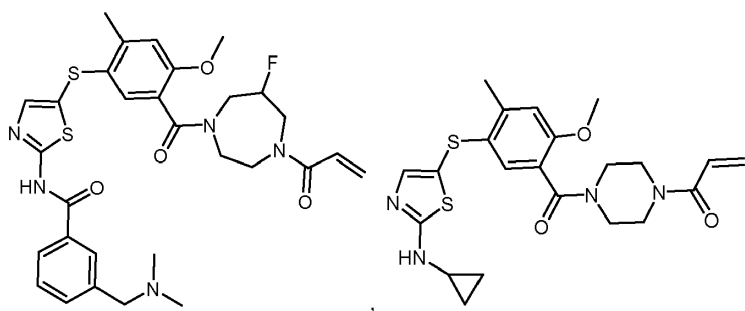
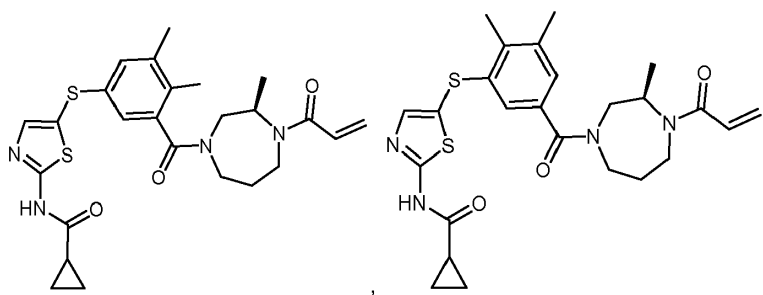
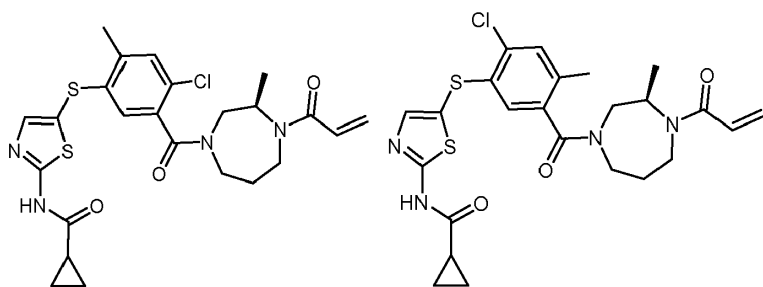
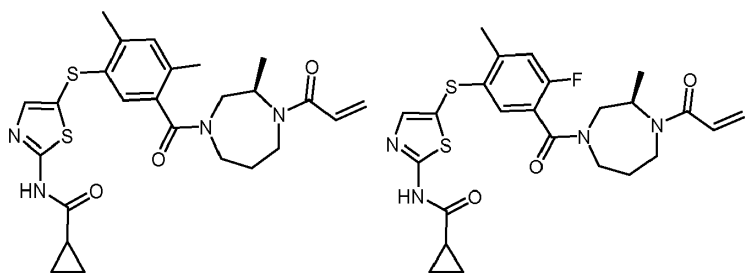
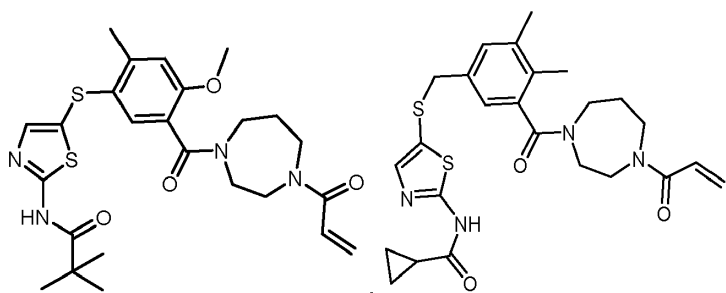


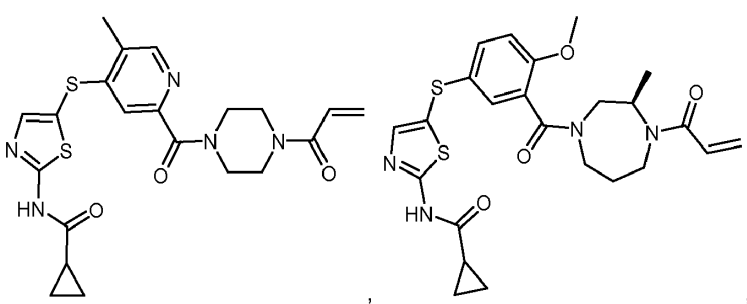
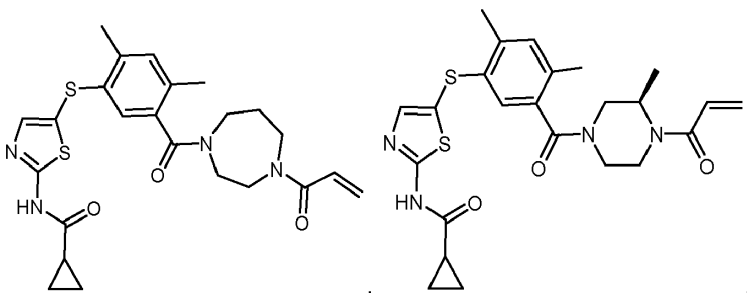
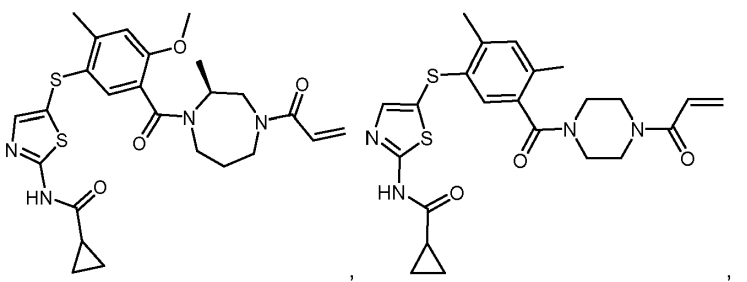
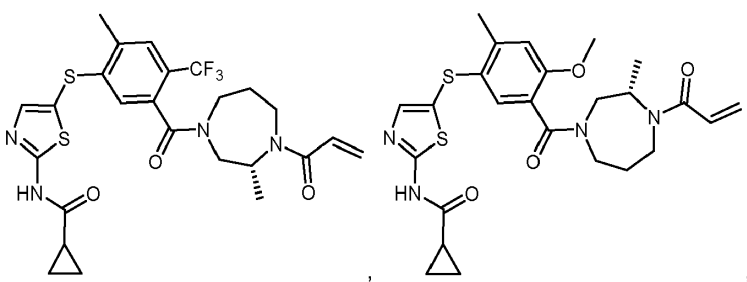
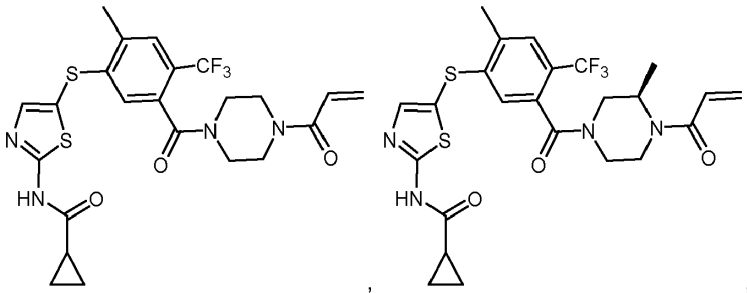




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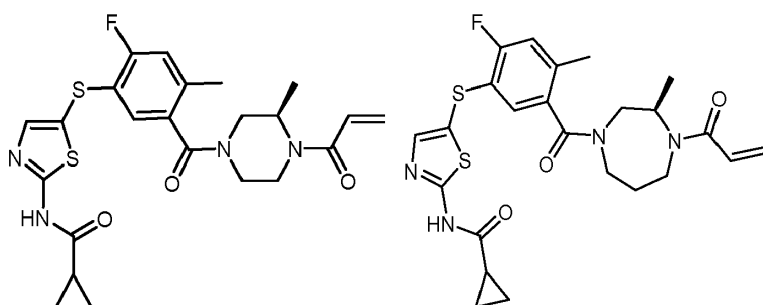
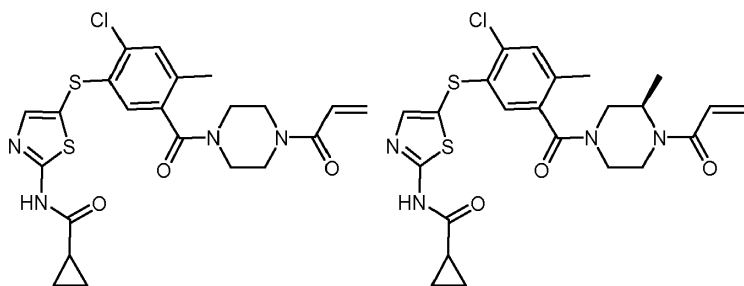
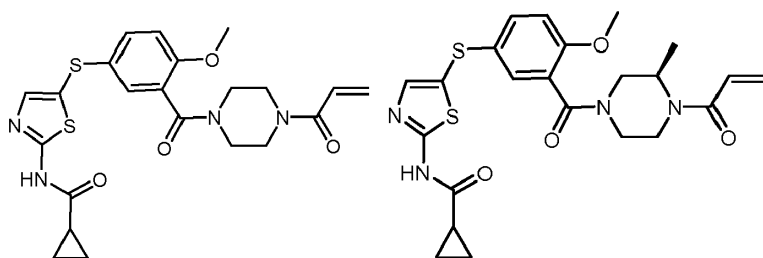




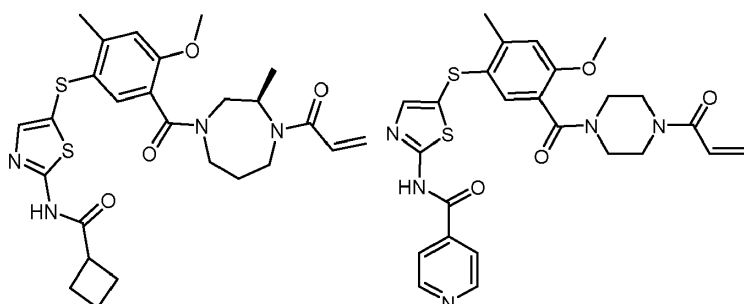
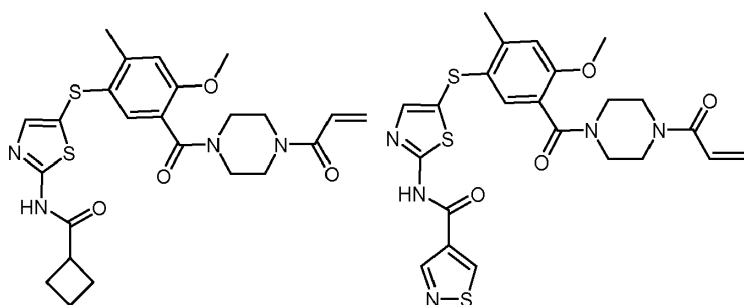


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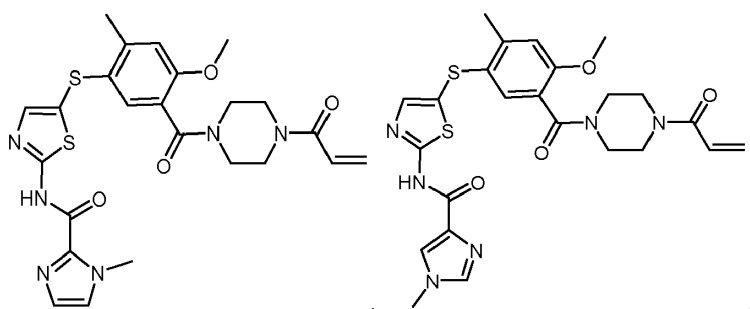
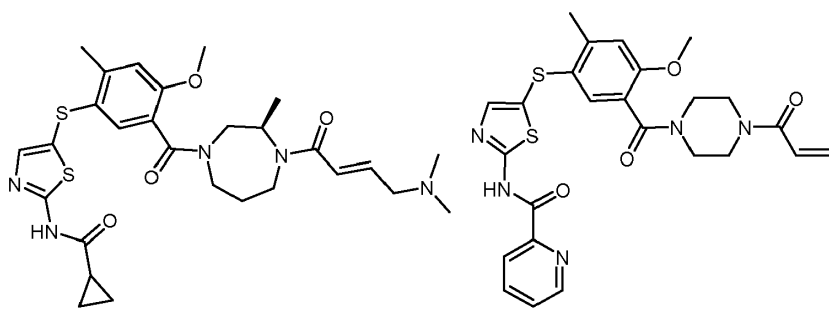
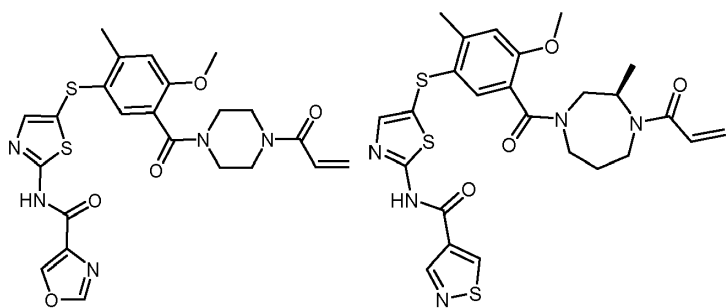
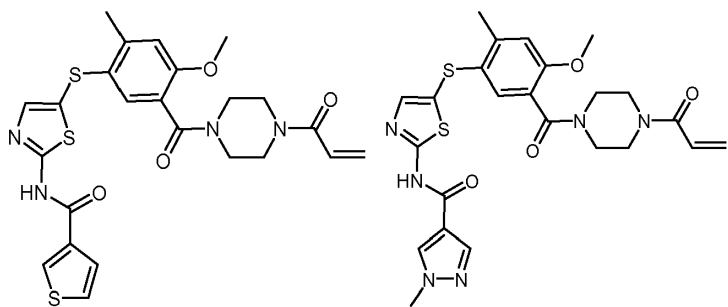
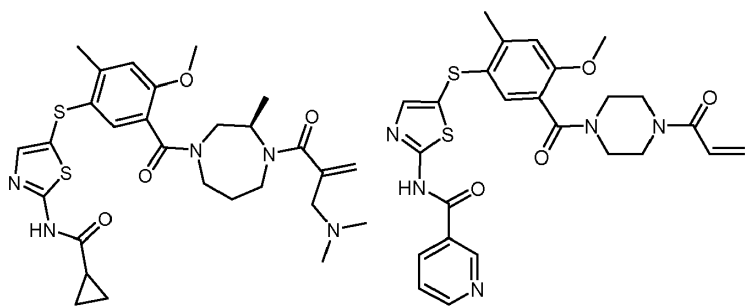
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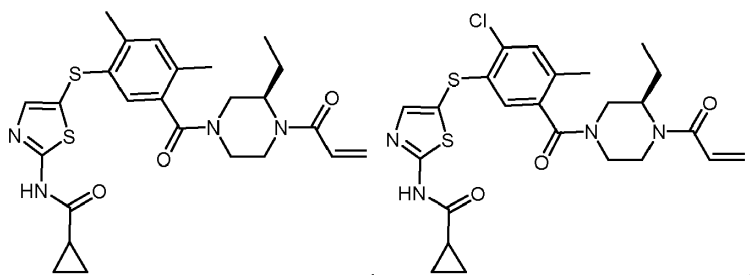
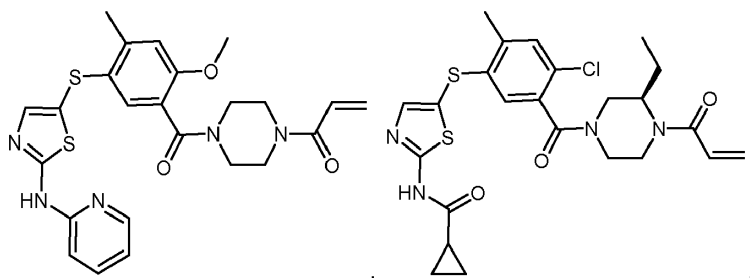


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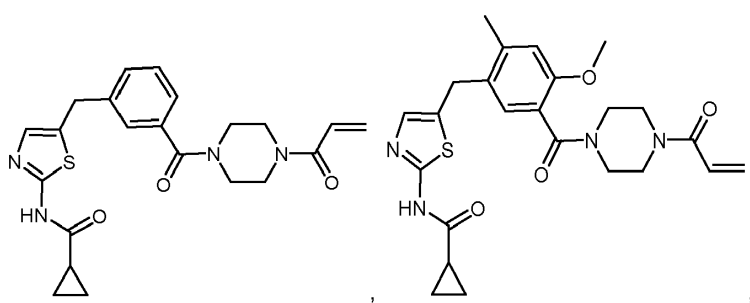
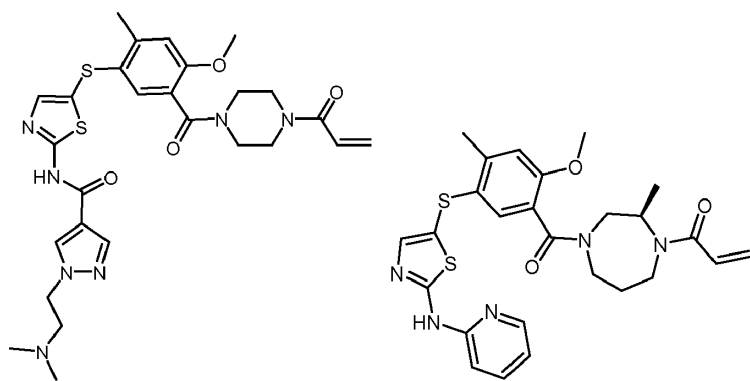


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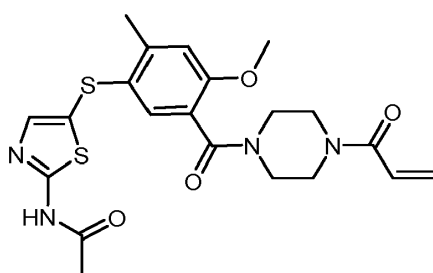




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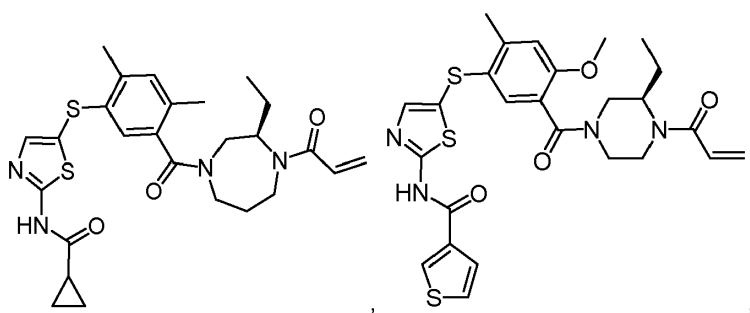
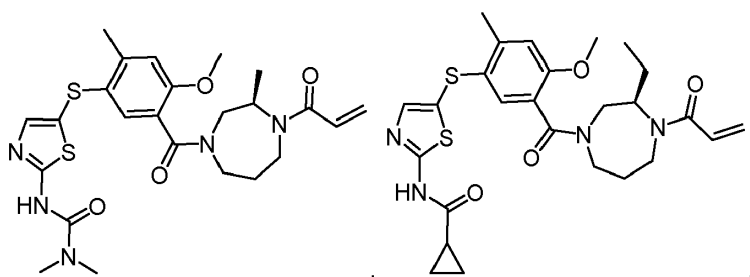
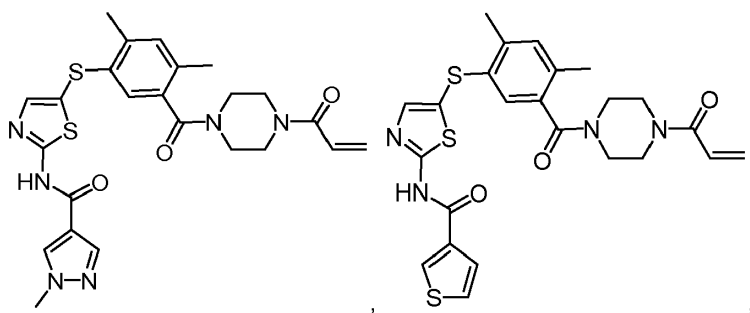
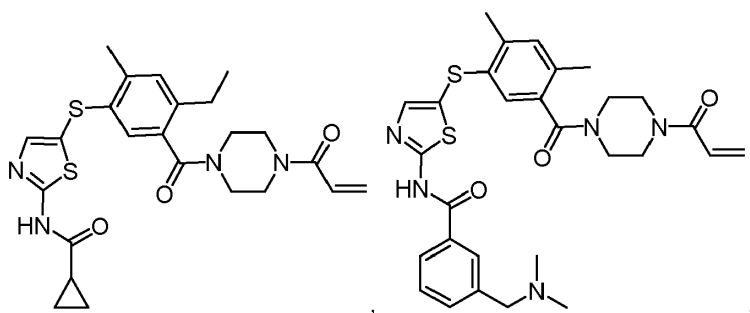
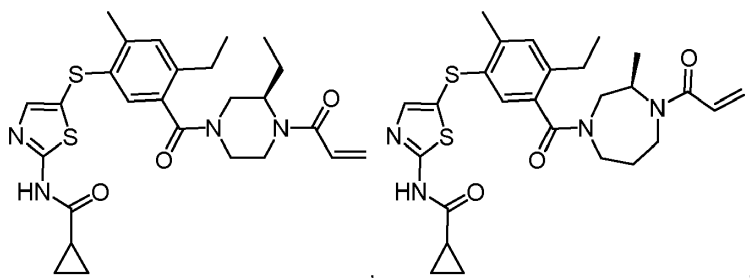
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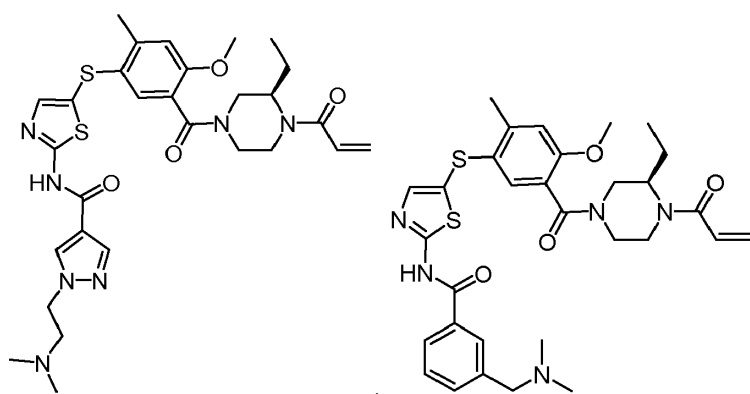
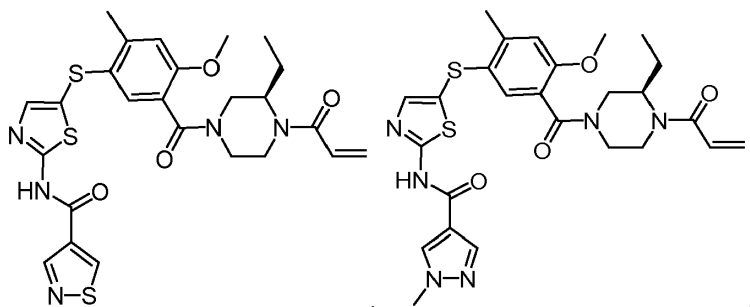


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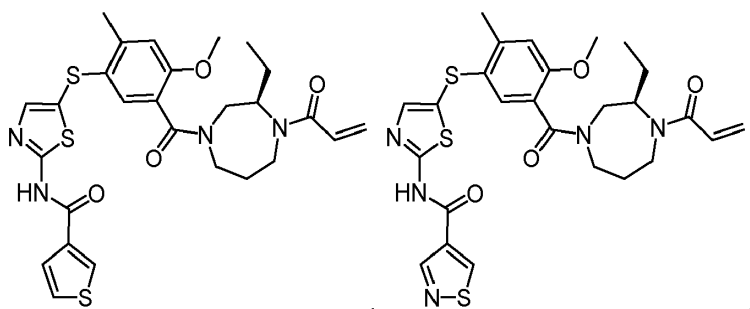
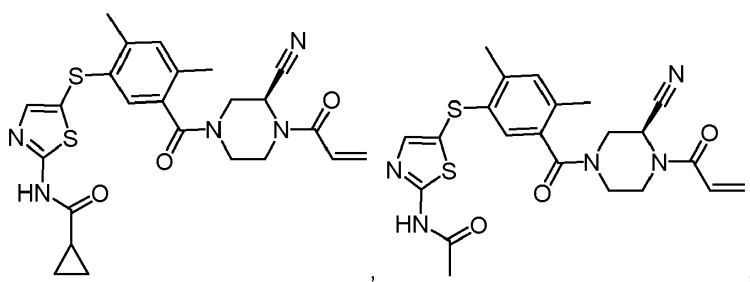
, eller et farmasøytisk akseptabelt salt derav; eller

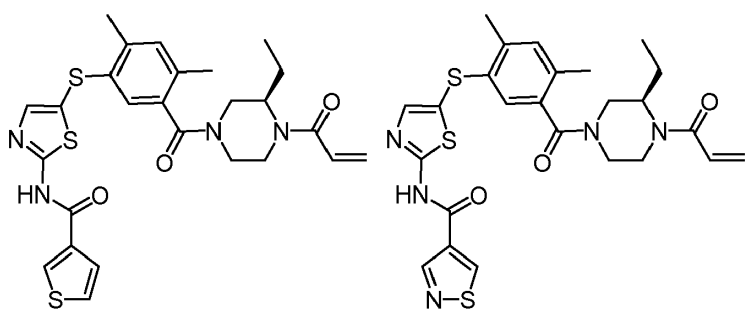
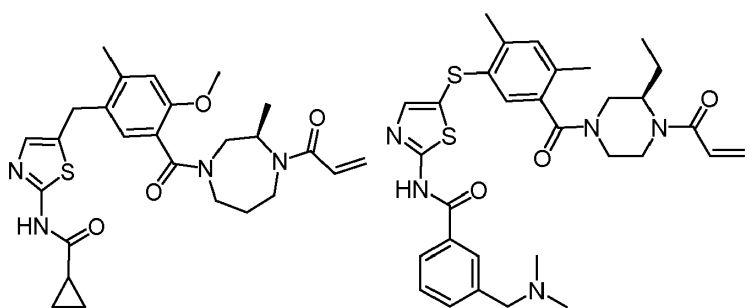
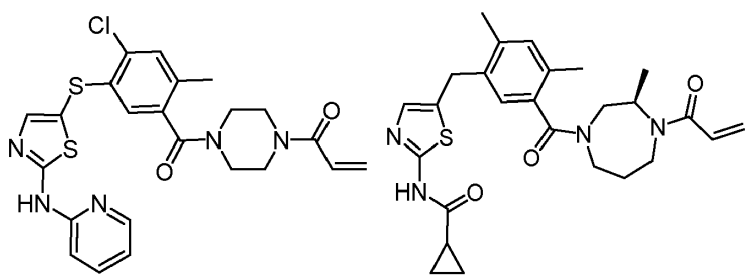
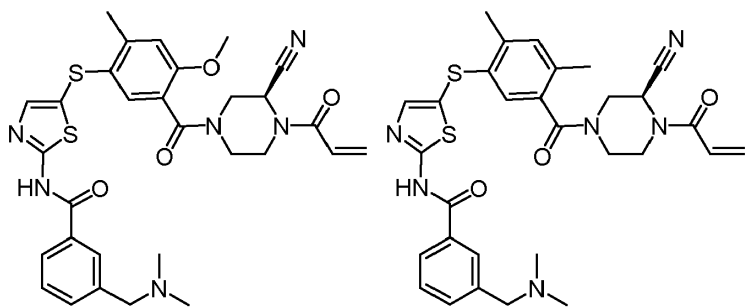
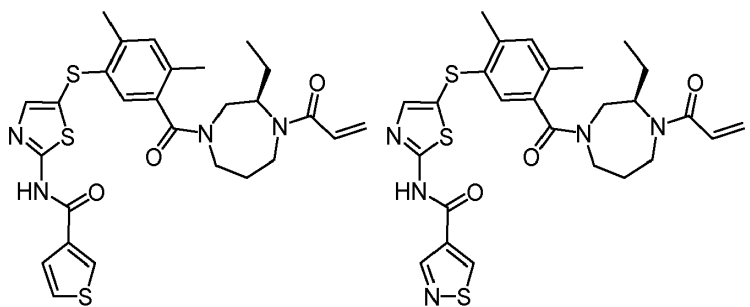
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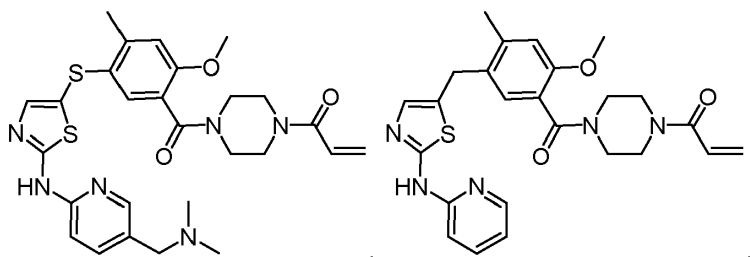
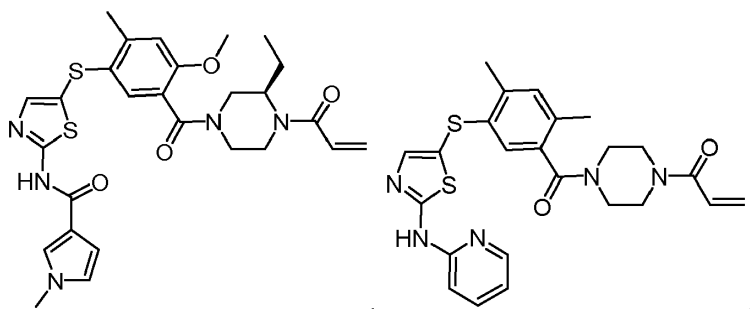




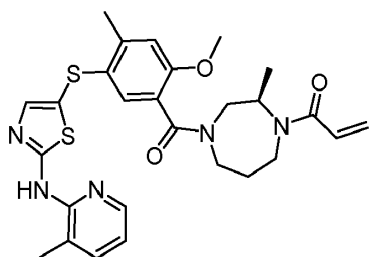
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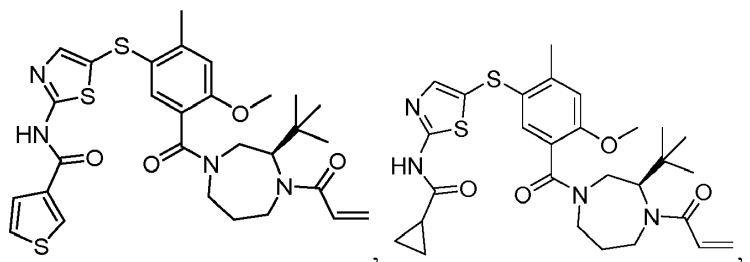
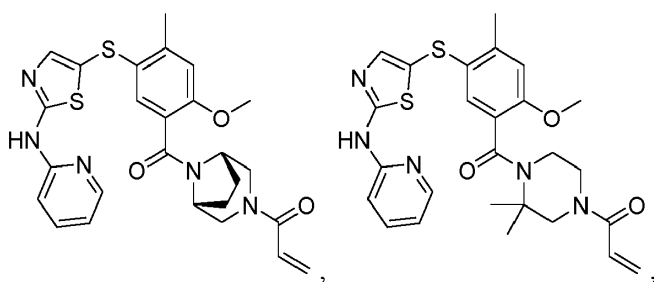
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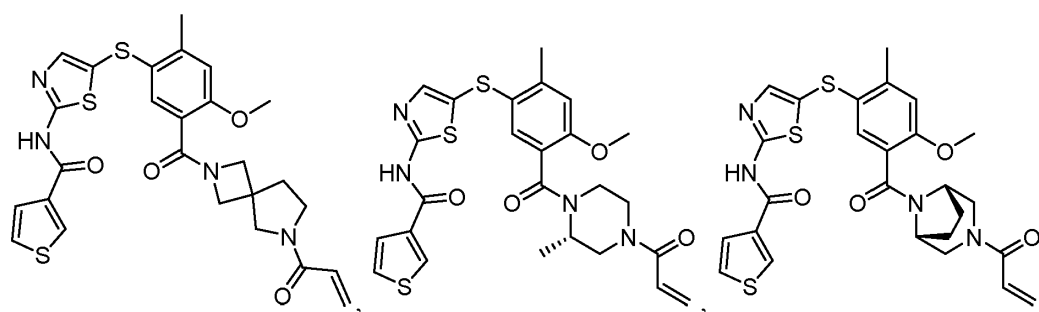
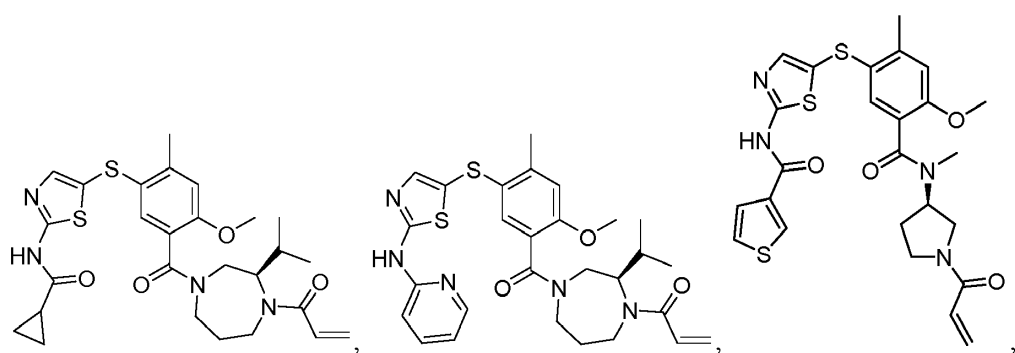
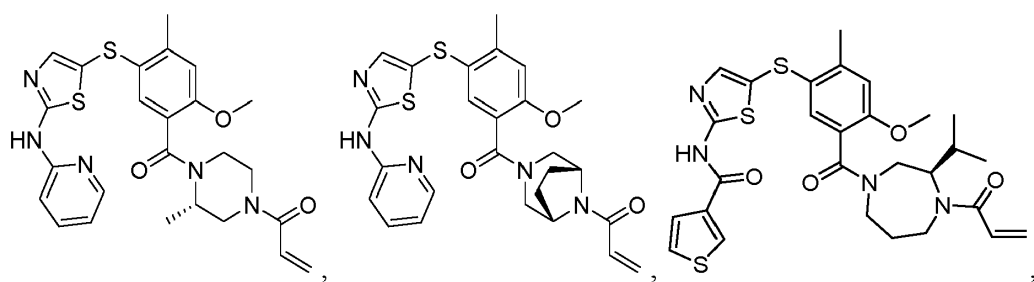
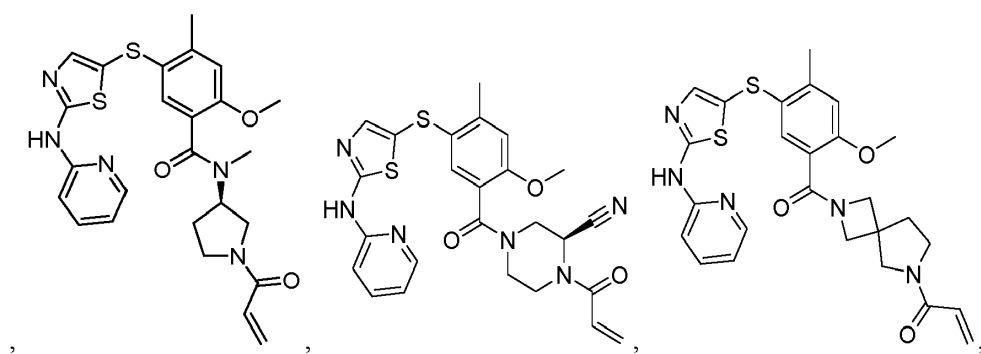
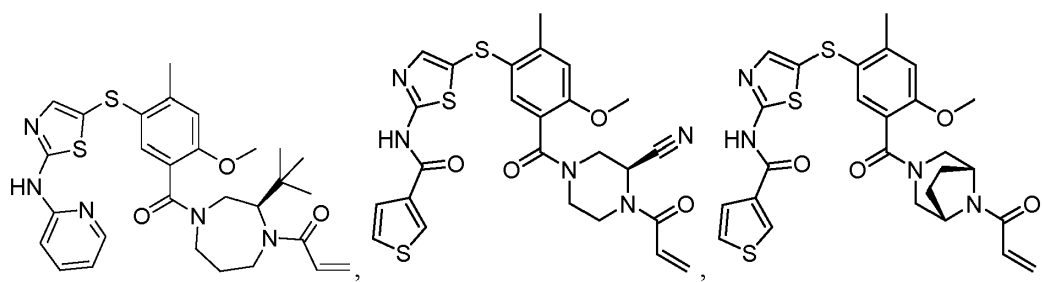
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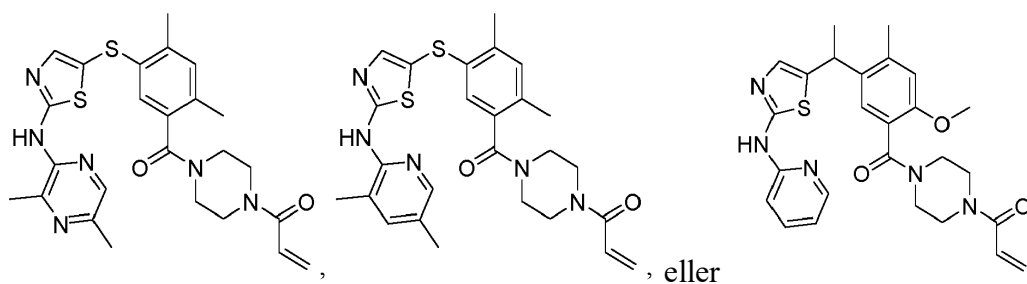
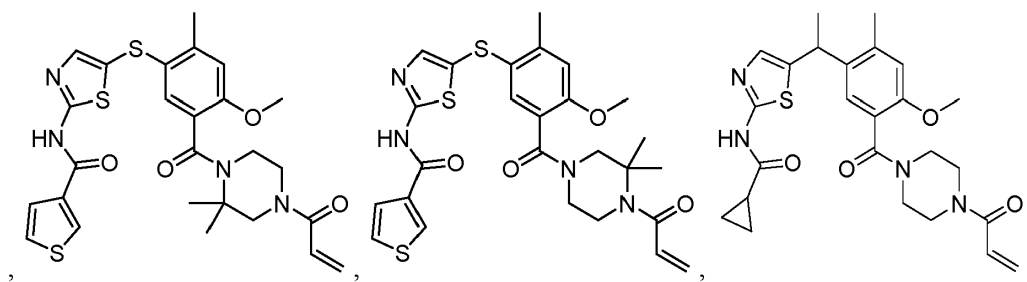
eller et farmasøytisk akseptabelt salt derav; eller

Alternativ D:



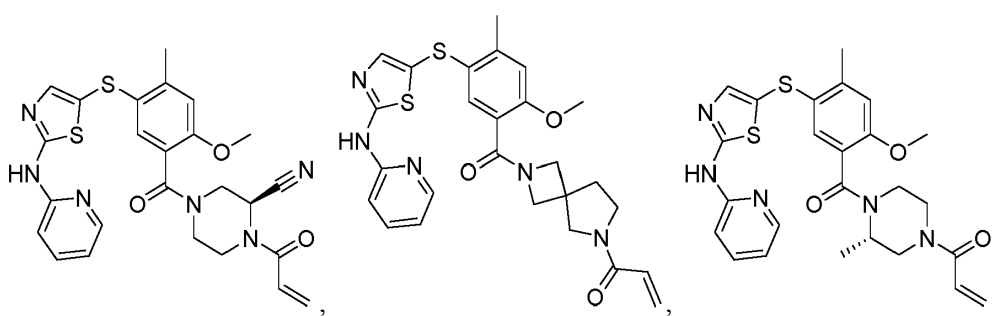
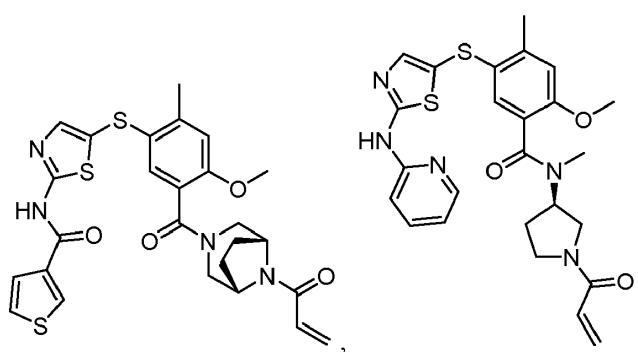
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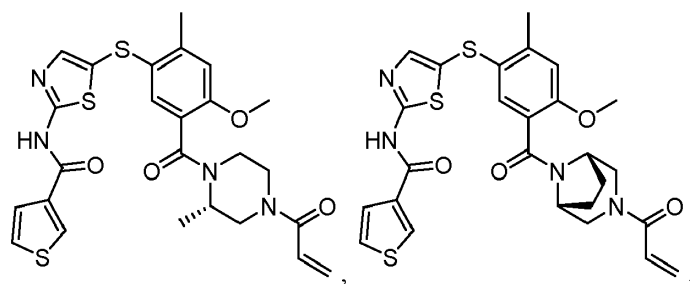
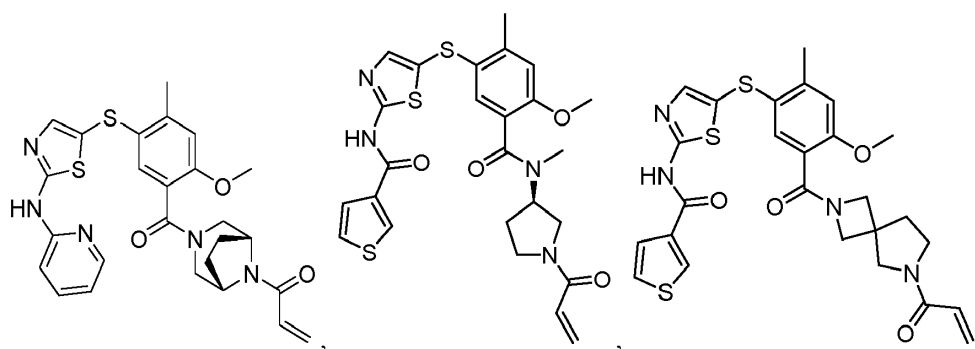




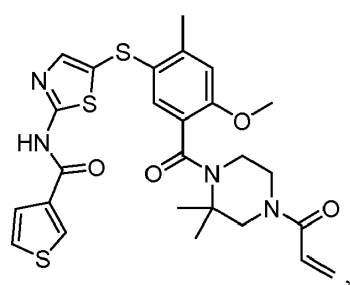
, eller

5 , eller et farmasøytisk akseptabelt salt derav; eller

Alternativ E:



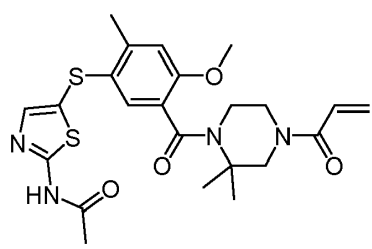
eller



5

eller et farmasøytisk akseptabelt salt derav; eller

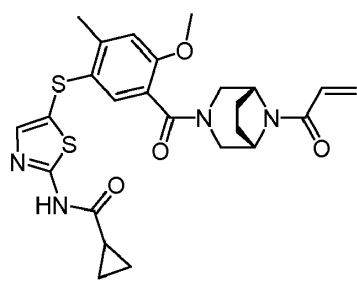
Alternativ F:



eller et farmasøytisk akseptabelt salt derav; eller

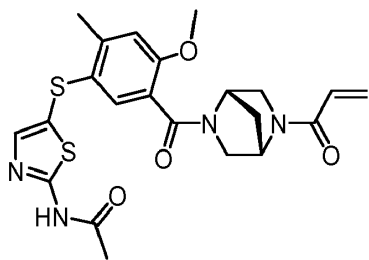
10

Alternativ G:



eller et farmasøytisk akseptabelt salt derav; eller

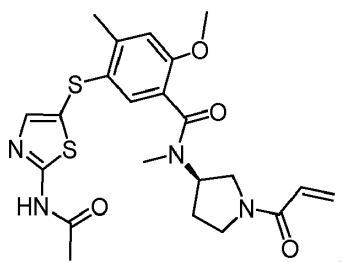
Alternativ H:



eller et farmasøytisk akseptabelt salt derav; eller

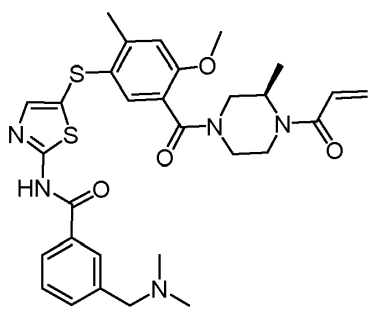
5

Alternativ I:



eller et farmasøytisk akseptabelt salt derav; eller

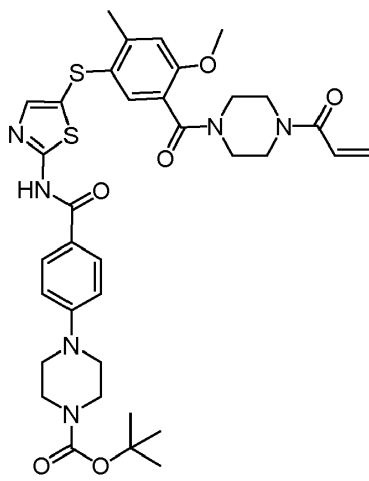
Alternativ J:



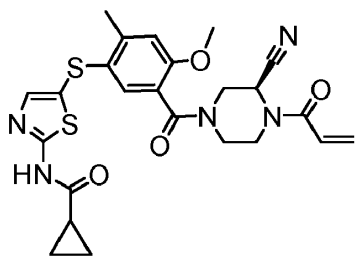
10

eller et farmasøytisk akseptabelt salt derav; eller

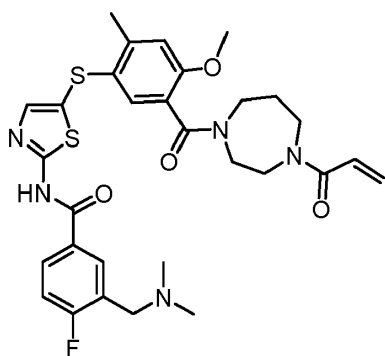
Alternativ K:



eller et farmasøytisk akseptabelt salt derav; eller

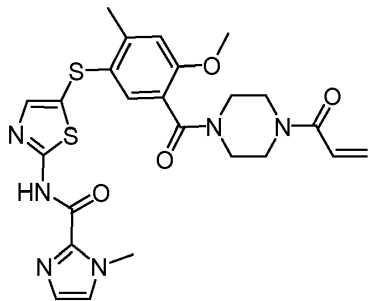
Alternativ L:

eller et farmasøytisk akseptabelt salt derav; eller

Alternativ M:

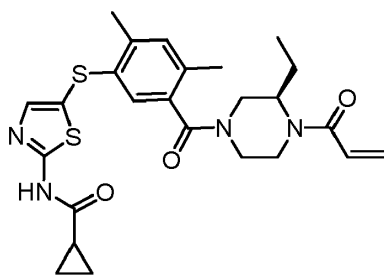
5

eller et farmasøytisk akseptabelt salt derav; eller

Alternativ N:

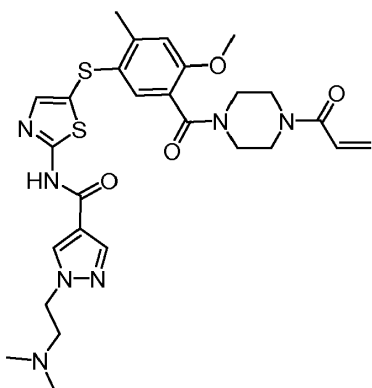
eller et farmasøytisk akseptabelt salt derav; eller

10

Alternativ O:

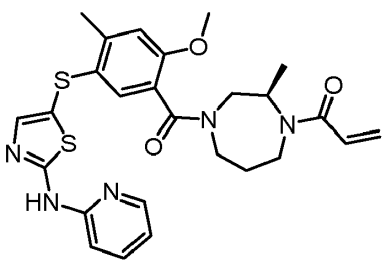
eller et farmasøytisk akseptabelt salt derav; eller

Alternativ P:



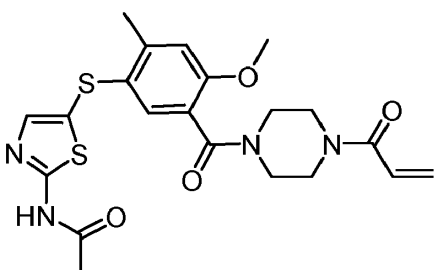
eller et farmasøytisk akseptabelt salt derav; eller

Alternativ Q:



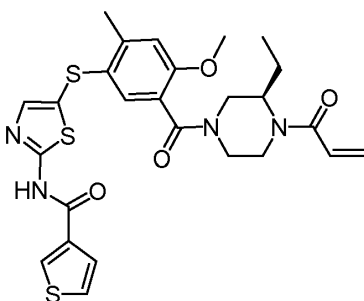
5 eller et farmasøytisk akseptabelt salt derav; eller

Alternativ R:



eller et farmasøytisk akseptabelt salt derav; eller

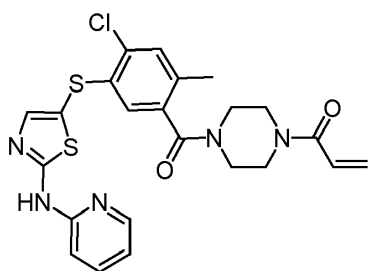
Alternativ S:



10

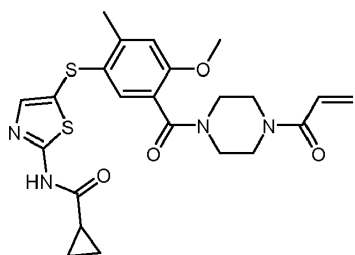
eller et farmasøytisk akseptabelt salt derav; eller

Alternativ T:



eller et farmasøytisk akseptabelt salt derav.

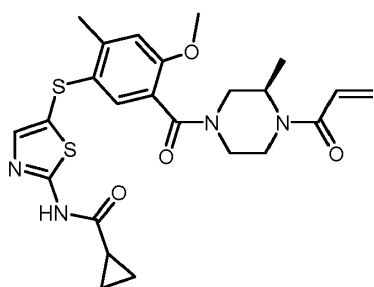
15. Forbindelsen ifølge krav 1, hvori forbindelsen er:



5

eller et farmasøytisk akseptabelt salt derav.

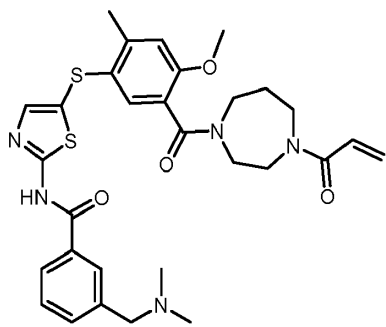
16. Forbindelsen ifølge krav 1, hvori forbindelsen er:



10

eller et farmasøytisk akseptabelt salt derav.

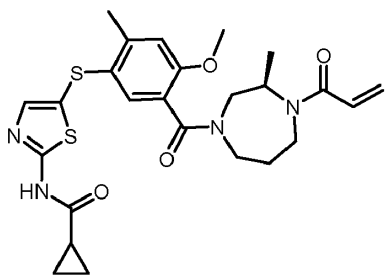
17. Forbindelsen ifølge krav 1, hvori forbindelsen er:



eller et farmasøytisk akseptabelt salt derav.

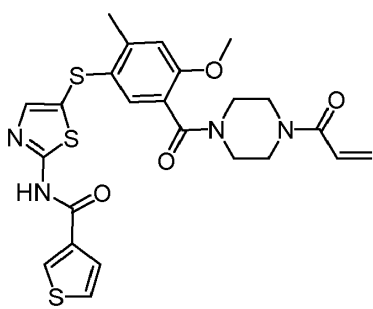
15

18. Forbindelsen ifølge krav 1, hvori forbindelsen er:



eller et farmasøytisk akseptabelt salt derav.

- 5 19. Forbindelsen ifølge krav 1, hvori forbindelsen er:



eller et farmasøytisk akseptabelt salt derav.

- 10 20. Farmasøytisk sammensetning omfattende forbindelsen ifølge et hvilket som helst av kravene 1 til 19 eller et farmasøytisk akseptabelt salt derav og en farmasøytisk akseptabel eksipiens.

- 15 21. Forbindelse ifølge et hvilket som helst av kravene 1 til 19 eller et farmasøytisk akseptabelt salt derav, for anvendelse i behandlingen av en kreft; eventuelt hvori:

Alternativ A:

kreften er T-cellelymfom; eller

Alternativ B:

kreften er T-cellelymfom, kutant T-cellelymfom eller perifert T-cellelymfom.