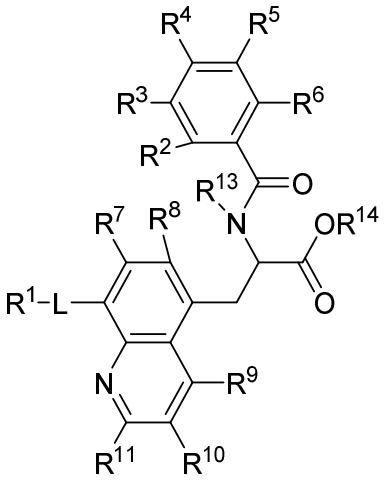
**What is claimed is:**

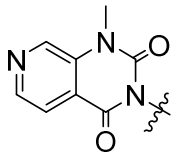
1. A compound of formula (II):



(II);

or a pharmaceutically acceptable salt thereof, wherein:

L is a bond;

$R^1$ is  substituted with one to two $R^a$; and wherein each $R^a$ is independently selected from $C_{1-4}$alkyl, $C_{1-4}$haloalkyl, $C_{1-4}$alkoxyl, and $C_{3-6}$cycloalkyl;

each $R^2$, $R^3$, $R^4$, $R^5$, and $R^6$ is independently selected from H, halo, cyano, hydroxyl, $C_{1-6}$alkyl, $C_{2-6}$alkenyl, $C_{2-6}$alkynyl, $C_{1-6}$alkoxyl, $C_{1-8}$haloalkyl, $C_{1-8}$haloalkoxyl, $-NR^{b1}R^{b2}$, $-R^{b3}S(O)_mR^{b4}$, $-S(O)_mR^{b4}$, $-NR^{b1}S(O)_nR^{b4}$, $-COOR^{b1}$, $-CONR^{b1}R^{b2}$, $-NR^{b1}COOR^{b2}$, $-NR^{b1}COR^{b4}$, $-R^{b3}NR^{b1}R^{b2}$, $-S(O)_nNR^{b1}R^{b2}$, $C_{3-12}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 3-12 membered heterocyclyl;

wherein each $C_{1-6}$alkyl, $C_{2-6}$alkenyl, $C_{2-6}$alkynyl, $C_{1-6}$alkoxyl, $C_{1-8}$haloalkyl, and $C_{1-8}$haloalkoxyl of $R^2$, $R^3$, $R^4$, $R^5$, and $R^6$ is unsubstituted or substituted with one to two $R^c$; wherein each $R^c$ is independently selected from azido, oxo, cyano, halo, hydroxyl, $-NR^{a1}R^{a2}$, $C_{1-4}$alkoxyl, $C_{3-8}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 4-6 membered heterocyclyl; wherein each $C_{3-8}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 4-6 membered heterocyclyl of $R^c$ is unsubstituted or substituted with one to three groups independently selected from halo, cyano, hydroxyl, $-NR^{a1}R^{a2}$, $C_{1-4}$alkyl, $C_{1-6}$haloalkyl, $C_{1-4}$alkoxyl, and $C_{3-6}$cycloalkyl;

wherein each $C_{6-10}$aryl, and 5-6 membered heteroaryl of $R^2$, $R^3$, $R^4$, $R^5$, and $R^6$ is independently unsubstituted or substituted with one to five $R^b$; and

wherein each $C_{3-12}$cycloalkyl, and 3-12 membered heterocyclyl of $R^2$, $R^3$, $R^4$, $R^5$, and $R^6$ is independently unsubstituted or substituted one to six groups independently selected from $=CR^{b1}R^{b2}$ and $R^b$;

wherein each $R^b$ is independently selected from azido, cyano, halo, hydroxyl, $-NR^{a1}R^{a2}$, $C_{1-6}$alkyl, $C_{1-8}$haloalkyl, $C_{1-6}$alkoxyl, $C_{3-6}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 4-6 membered heterocyclyl; wherein each $C_{3-6}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 4-6 membered heterocyclyl of $R^b$ is independently unsubstituted or substituted with one to three groups independently selected from halo, cyano, hydroxyl, $-NR^{a1}R^{a2}$, $C_{1-4}$alkyl, $C_{1-4}$haloalkyl, and $C_{1-4}$alkoxyl;

wherein each $R^{b1}$ and $R^{b2}$ is independently selected from H, $C_{1-8}$alkyl, $C_{1-8}$haloalkyl, $C_{3-8}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 3-8 membered heterocyclyl;

wherein each $C_{3-8}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 4-6 membered heterocyclyl of $R^{b1}$ and $R^{b2}$ is independently unsubstituted or substituted with one to three groups independently selected from halo, cyano, hydroxyl, $-NR^{a1}R^{a2}$, $C_{1-8}$alkyl, $C_{1-8}$haloalkyl, $C_{1-6}$alkoxyl, $C_{3-6}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 4-6 membered heterocyclyl; and

wherein each $C_{1-8}$alkyl and $C_{1-8}$haloalkyl of $R^{b1}$ and $R^{b2}$ is unsubstituted or substituted with one to two $R^{b5}$;

wherein $R^{b3}$ is $C_{1-4}$alkylene;

wherein $R^{b4}$ is selected from $C_{1-4}$alkyl, $C_{1-4}$haloalkyl, $C_{3-6}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 4-6 membered heterocyclyl; wherein each $C_{1-4}$alkyl, $C_{1-4}$haloalkyl, $C_{3-8}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and the 4-6 membered heterocyclyl of $R^{b4}$ is unsubstituted or substituted with one to three $R^{b6}$;

wherein each $R^{b5}$ is independently selected from cyano, hydroxyl, $C_{1-4}$alkoxyl, $C_{3-8}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 4-6 membered heterocyclyl; and each $C_{1-4}$alkoxyl, $C_{3-8}$cycloalkyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, and 4-6 membered heterocyclyl of $R^{b5}$ is unsubstituted or substituted with one to three groups independently selected from halo, cyano, hydroxyl, $-NR^{a1}R^{a2}$, $C_{1-4}$alkyl, $C_{1-4}$haloalkyl, $C_{1-4}$alkoxyl, and phenyl; and

wherein each $R^{b6}$ is independently selected from halo, cyano, $C_{1-4}$alkyl, $C_{1-4}$haloalkyl, $C_{1-4}$alkoxyl, $C_{3-6}$cycloalkyl, phenyl, 4-6 membered heterocyclyl, and 5-6 membered heteroaryl; wherein each $C_{3-6}$cycloalkyl, 4-6 membered heterocyclyl, and 5-6 membered heteroaryl of $R^{b6}$ is independently unsubstituted or substituted with

one to three groups independently selected from halo, cyano, -$NR^{a1}R^{a2}$, $C_{1-4}$alkyl, $C_{1-4}$haloalkyl, and $C_{1-4}$alkoxyl; or

$R^2$ and $R^3$, $R^3$ and $R^4$, or $R^5$ and $R^6$ together with the atoms to which they are attached form a $C_{6-10}$aryl, 5-6 membered heteroaryl, $C_{3-6}$cycloalkyl, or 5-6 membered heterocyclyl; wherein each $C_{6-10}$aryl, 5-6 membered heteroaryl, $C_{3-6}$cycloalkyl, and 5-6 membered heterocyclyl is independently unsubstituted or substituted with one to three groups independently selected from halo, cyano, -$NR^{a1}R^{a2}$, $C_{1-6}$alkyl, $C_{1-6}$alkoxyl, $C_{1-6}$haloalkyl, $C_{3-8}$cycloalkyl, 3-6 membered heterocyclyl, $C_{6-10}$aryl, 5-6 membered heteroaryl, $C_{1-4}$alkylene-$C_{3-8}$cycloalkyl, $C_{1-4}$alkylene-$C_{6-10}$aryl, and $C_{1-4}$alkylene-(5-6 membered heteroaryl);

each $R^7$, $R^8$, $R^9$, $R^{10}$, and $R^{11}$ is independently selected from H, halo, hydroxyl, cyano, $C_{1-6}$alkyl, $C_{1-6}$alkoxyl, $C_{1-6}$haloalkyl, $C_{1-6}$haloalkoxyl, and -$NR^{a1}R^{a2}$;

$R^{13}$ is selected from H, $C_{1-4}$alkyl, and $C_{1-4}$haloalkyl; and

$R^{14}$ is selected from H, $C_{1-6}$alkyl, -$C_{1-4}$alkylene-$NR^{a1}R^{a2}$, -$C_{1-4}$alkylene-C(O)$NR^{a1}R^{a2}$, -$C_{1-4}$alkylene-O-C(O)-$C_{1-4}$alkyl, -$C_{1-4}$alkylene-O-C(O)-O-$C_{1-4}$alkyl, -$C_{1-4}$alkylene-O-C(O)-$C_{1-4}$alkylene-$NR^{a1}R^{a2}$, -$C_{1-4}$alkylene-O-$C_{1-4}$alkyl, $C_{3-8}$cycloalkyl, -$C_{1-4}$alkylene-$C_{3-8}$cycloalkyl, 4-6 membered heterocyclyl, and -$C_{1-4}$alkylene-(4-6 membered heterocyclyl);

wherein each $C_{3-8}$cycloalkyl, -$C_{1-4}$alkylene-$C_{3-8}$cycloalkyl, 4-6 membered heterocyclyl, and -$C_{1-4}$alkylene-(4-6 membered heterocyclyl) of $R^{14}$ is unsubstituted or substituted with one to three groups independently selected from halo, $C_{1-4}$alkyl, $C_{1-4}$alkoxyl, and $C_{1-4}$haloalkyl; or

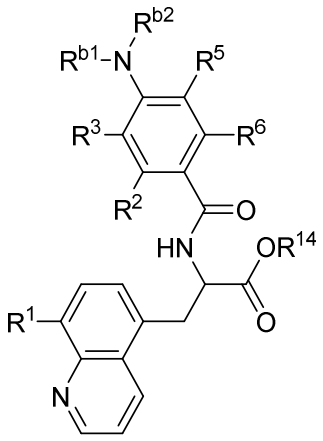
$R^{14}$ together with the N that attaches to $R^{13}$ forms a 5 membered heterocyclyl; wherein the 5 membered heterocyclyl is unsubstituted or substituted with one to two groups independently selected from halo, $C_{1-6}$alkyl, $C_{1-6}$alkoxyl, $C_{1-6}$haloalkyl, and $C_{6-10}$aryl; wherein the $C_{6-10}$aryl is unsubstituted or substituted with one to three groups independently selected from halo, $C_{1-6}$alkyl, $C_{1-6}$alkoxyl, $C_{1-6}$haloalkyl;

each $R^{a1}$ and $R^{a2}$ is independently selected from H, $C_{1-6}$alkyl, and $C_{1-6}$haloalkyl;

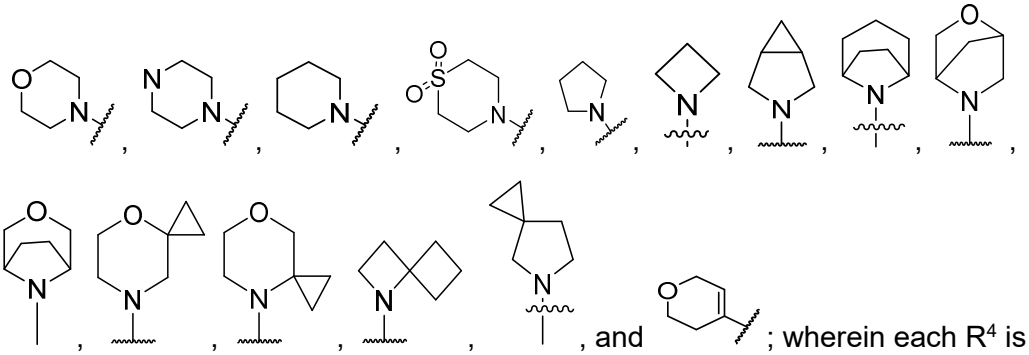
m is selected from 0, 1, and 2; and

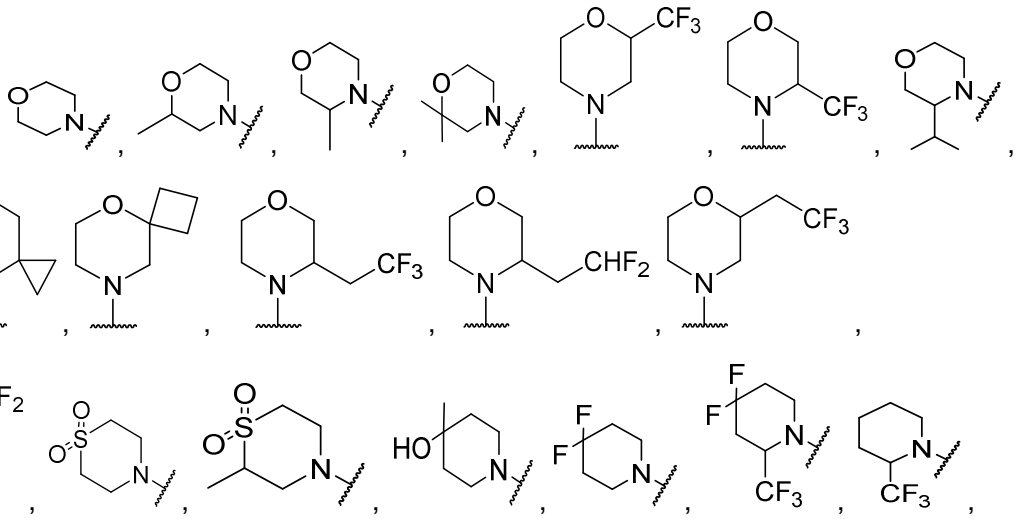
n is selected from 1, and 2.

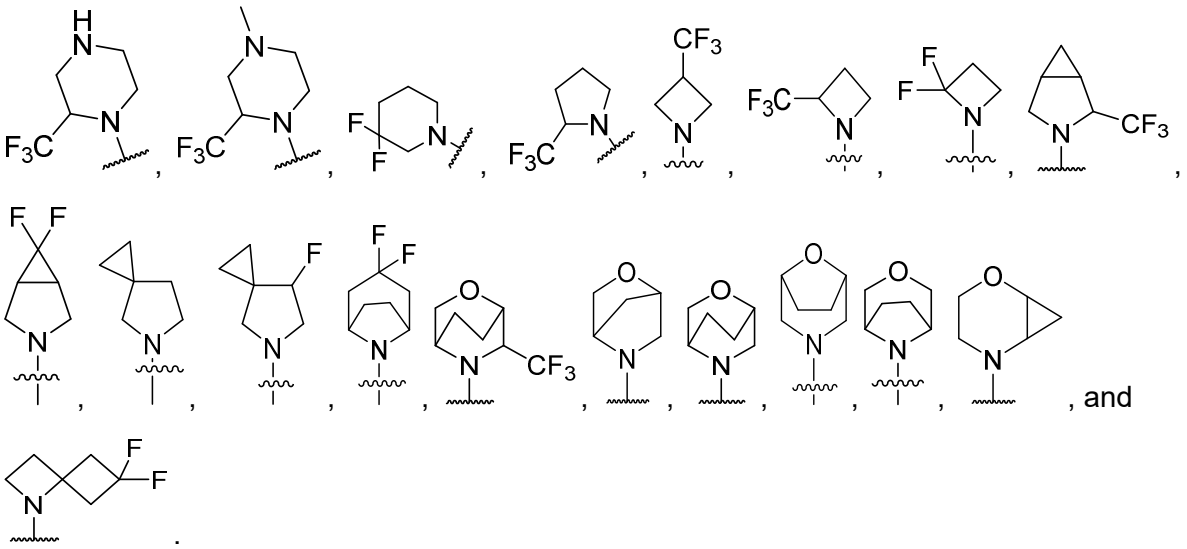
2. The compound of claim 1, or a pharmaceutically acceptable salt thereof, wherein the compound is of formula (IIa):

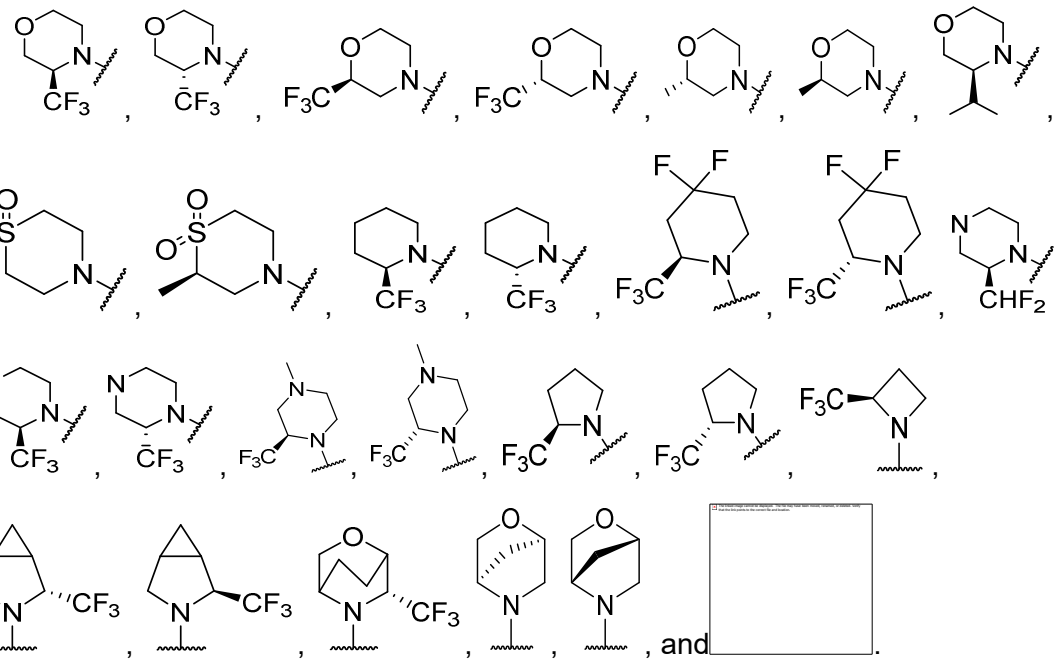


(IIa).

3. The compound of claim 1, or a pharmaceutically acceptable salt thereof, wherein $R^4$ is selected from , and ; wherein each $R^4$ is optionally substituted with one to three groups independently selected from F, Cl, cyano, hydroxyl, $NH_2$, $-CH_3$, $-CH(CH_3)_2$, $-CH_2F$, $-CHF_2$, $-CF_3$, $-CH_2CH_2F$, $-CH_2CHF_2$, $-CH_2CF_3$, $C_{3-6}$cycloalkyl, and $-CH_2C_{3-6}$cycloalkyl.

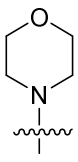
4. The compound of claim 1, or a pharmaceutically acceptable salt thereof, wherein $R^4$ is selected from 



5. The compound of claim 1, or a pharmaceutically acceptable salt thereof, wherein $R^4$ is selected from 

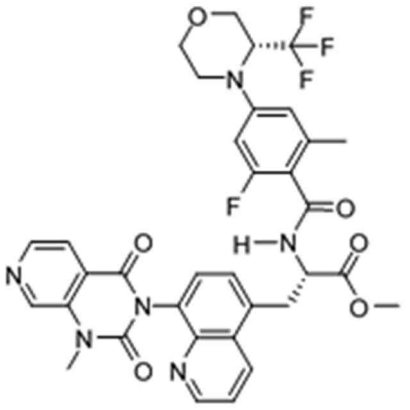
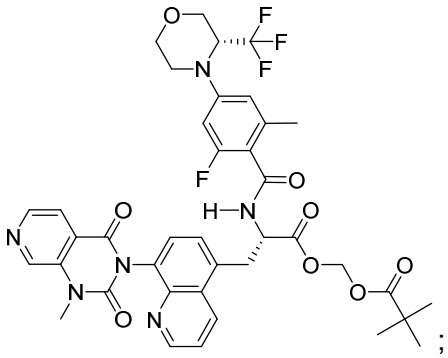
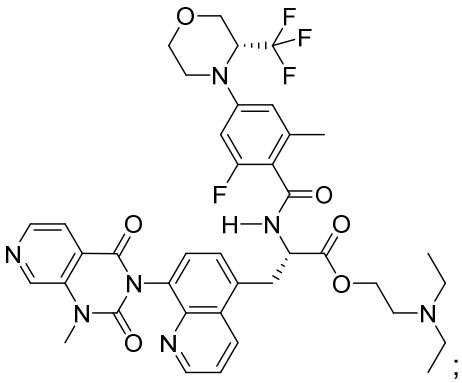
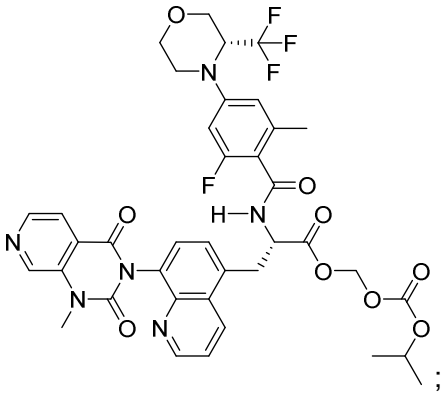
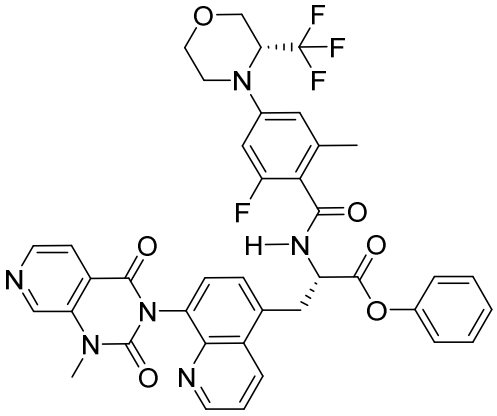
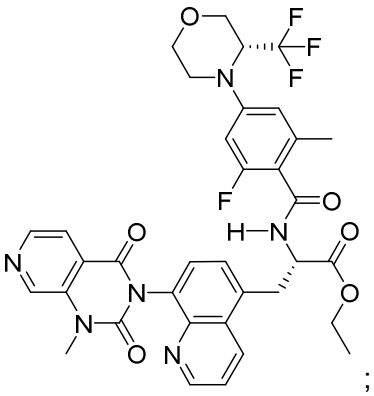
6. The compound of any one of claims 1 to 5, or a pharmaceutically acceptable salt thereof; wherein $R^3$ and $R^5$ are H.

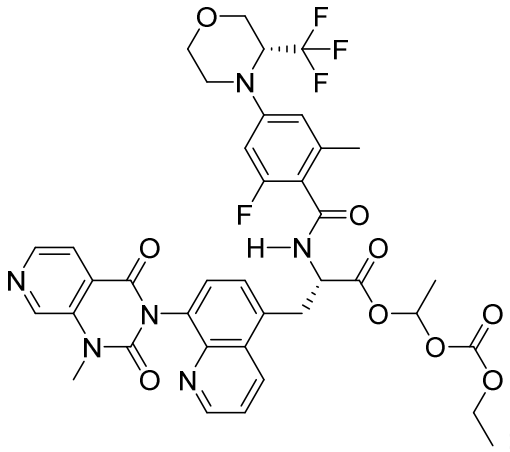
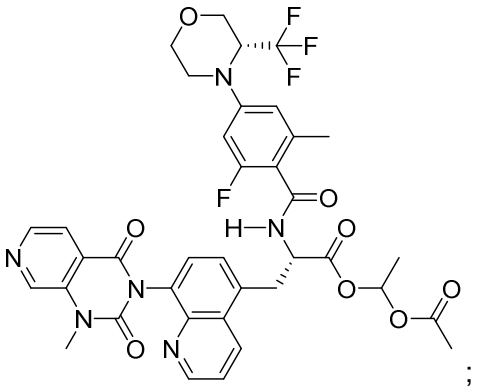
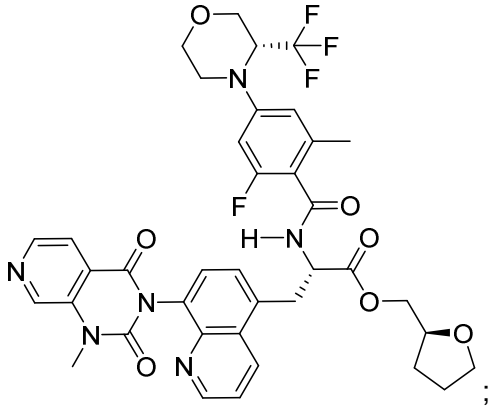
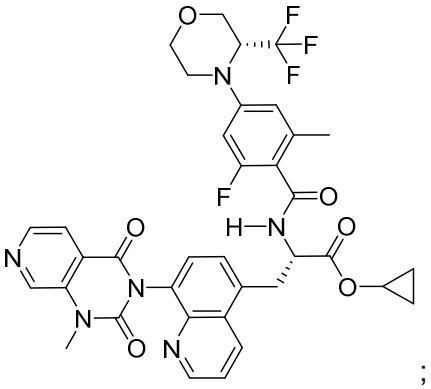
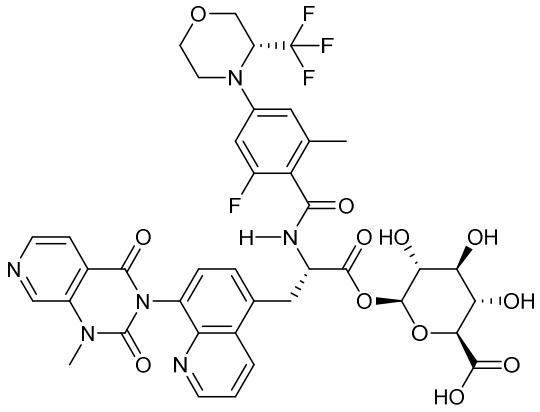
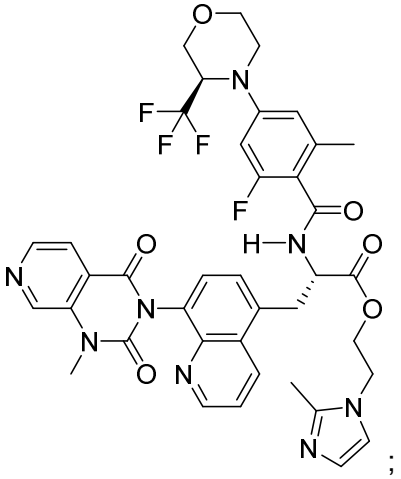
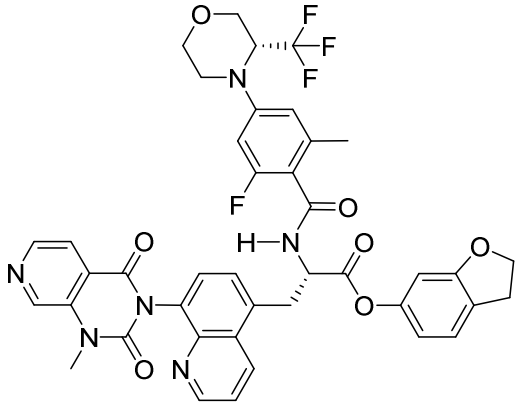
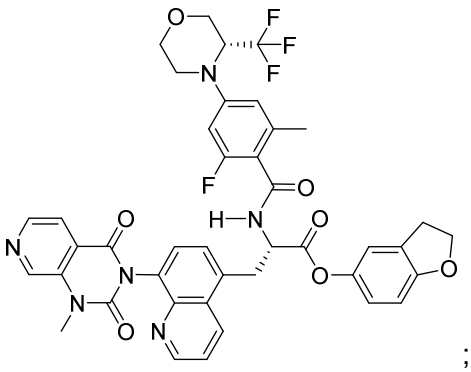
7. The compound of any one of claims 1 to 6, or a pharmaceutically acceptable salt thereof, wherein each $R^2$ and $R^6$ is independently selected from F and $-CH_3$.

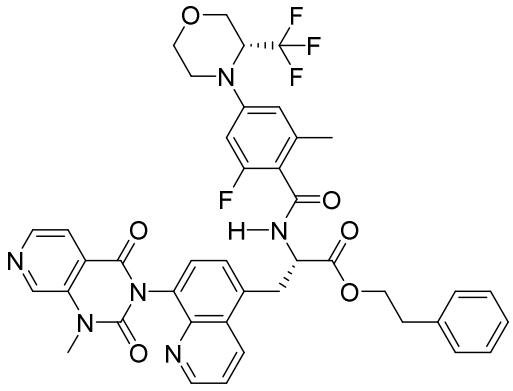
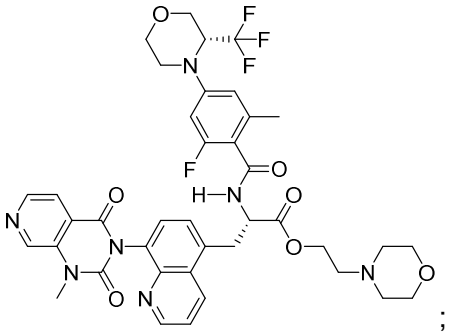
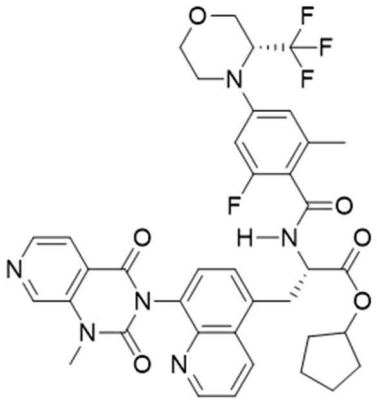
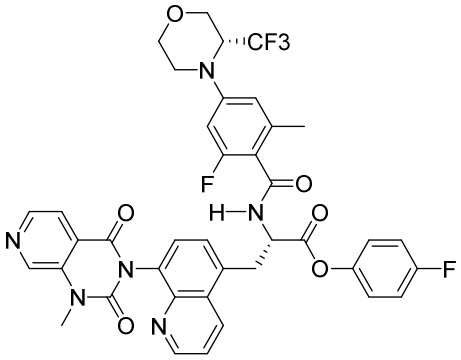
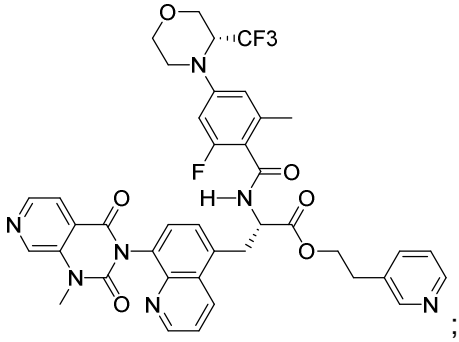
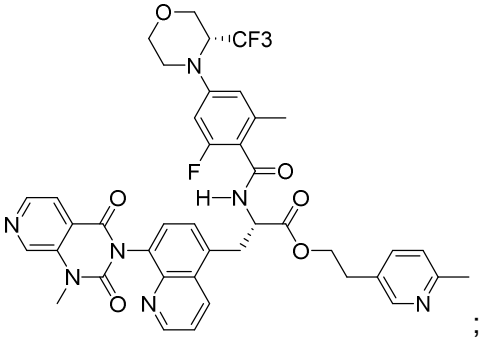
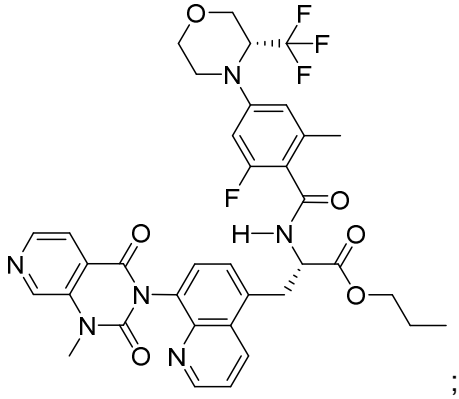
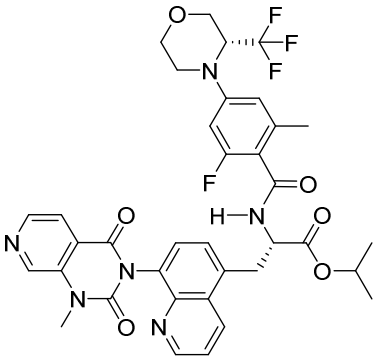
8. The compound of any one of claims 1, 3, 6 or 7, or a pharmaceutically acceptable salt thereof, wherein $R^4$ is  substituted with $-CF_3$.

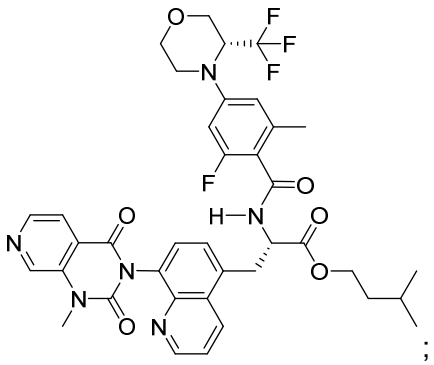
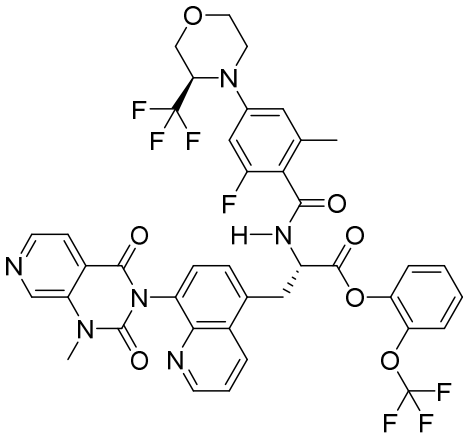
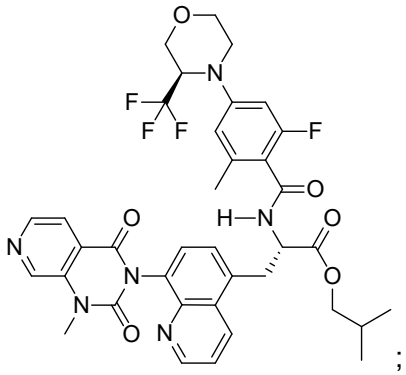
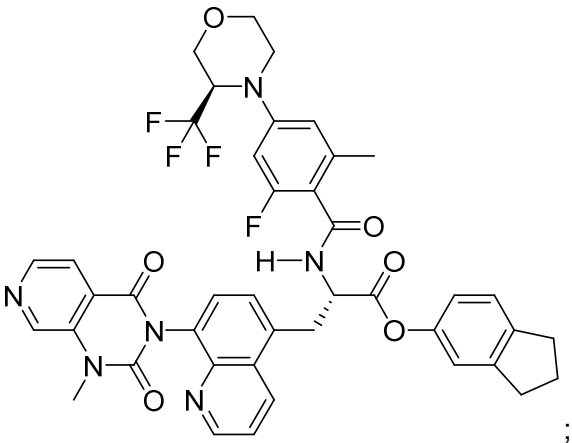
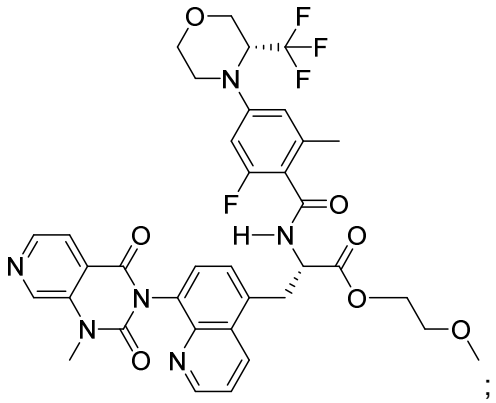
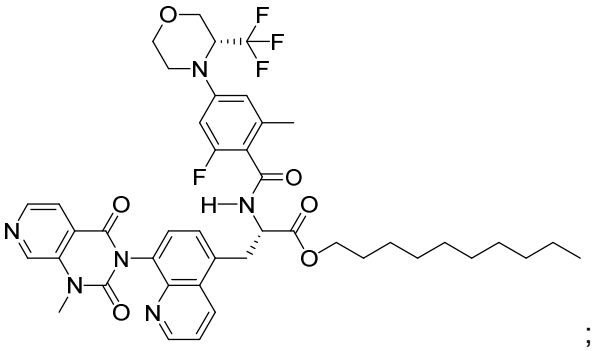
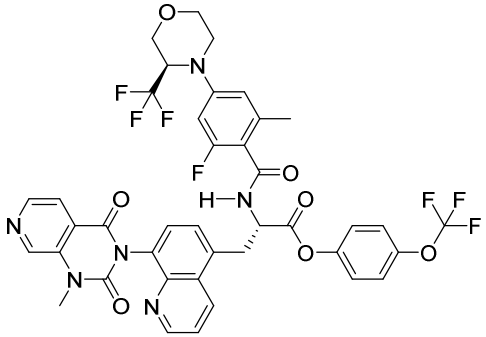
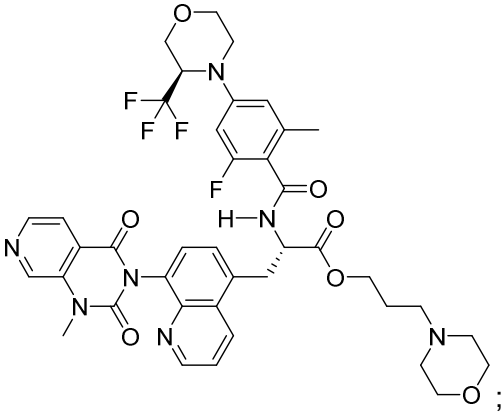
9. The compound of any one of claims 1 to 8, or a pharmaceutically acceptable salt thereof, wherein $R^{14}$ is selected from H, methyl, and ethyl.

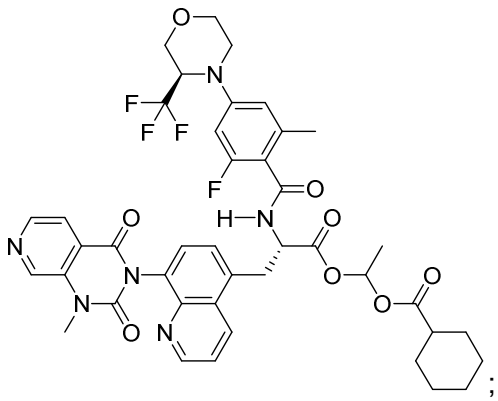
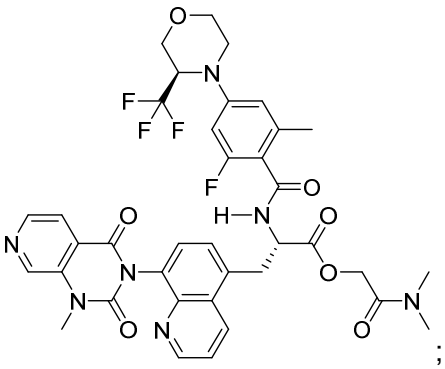
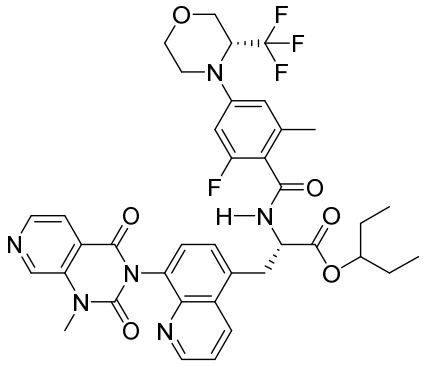
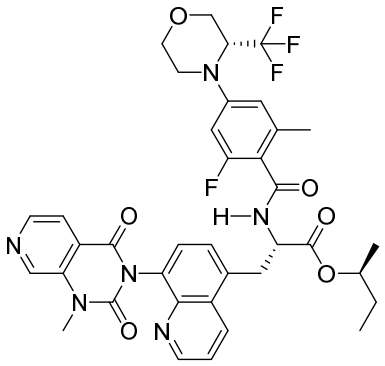
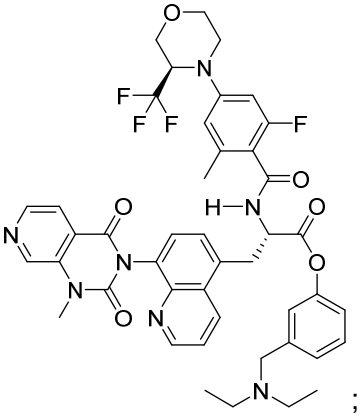
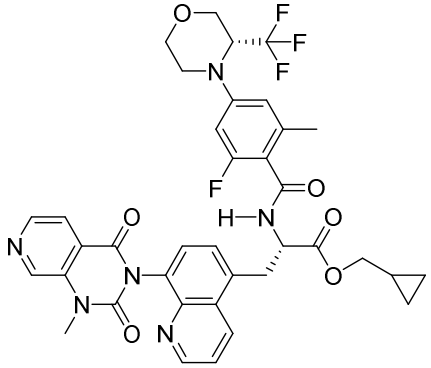
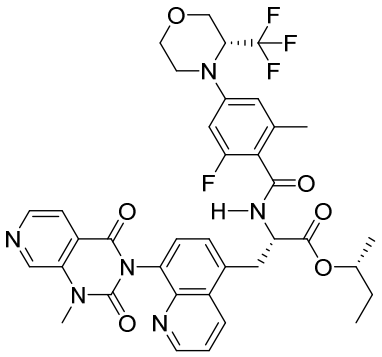
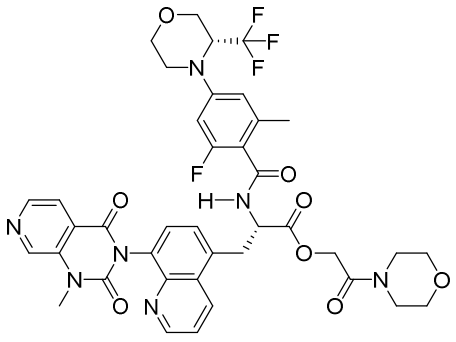
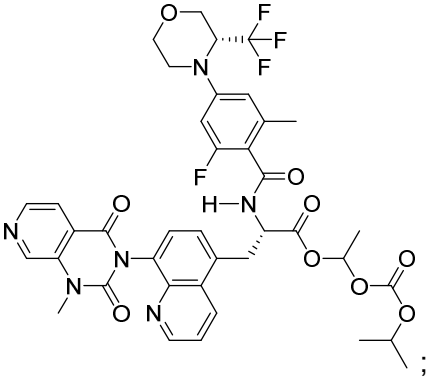
10. A compound, or a pharmaceutically acceptable salt thereof, wherein the compound is selected from:

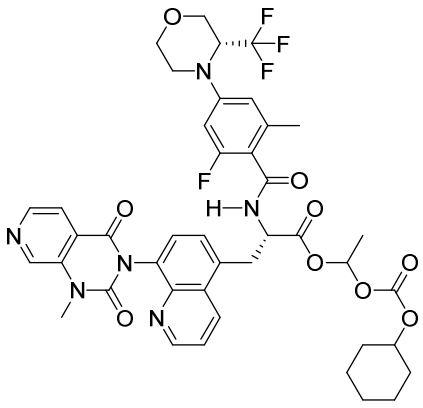
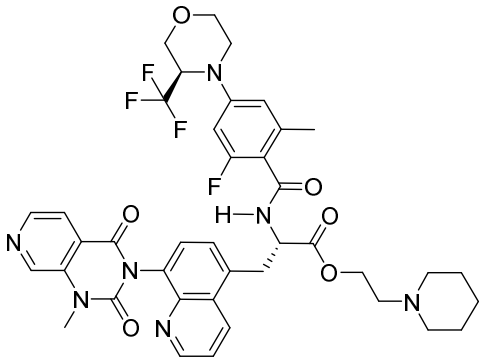
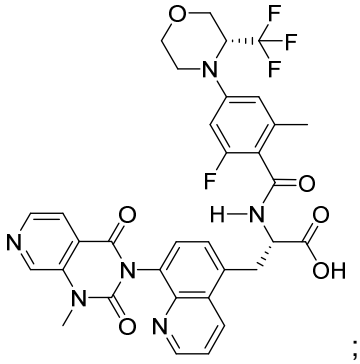
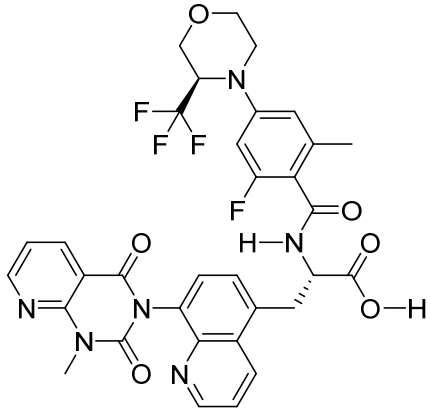
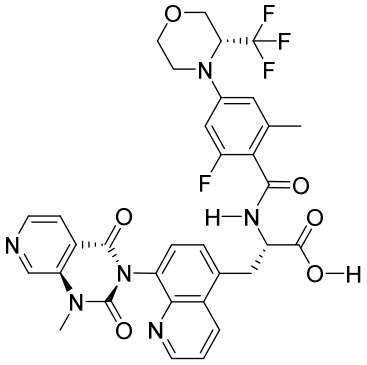
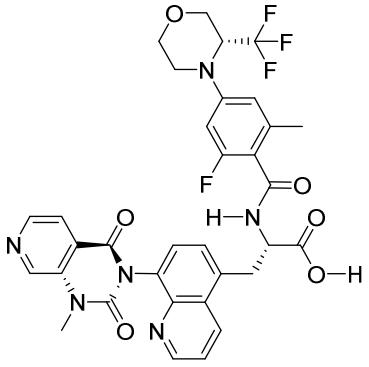
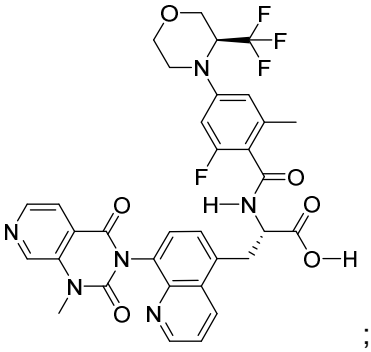
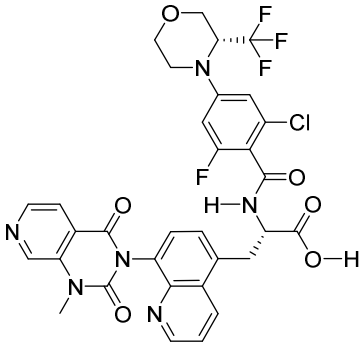
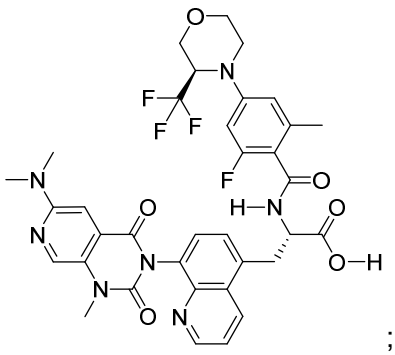
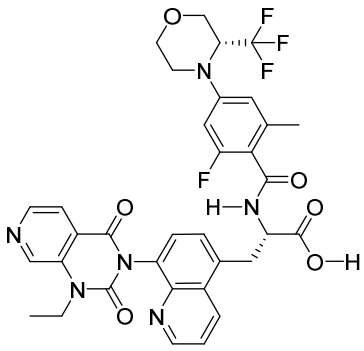
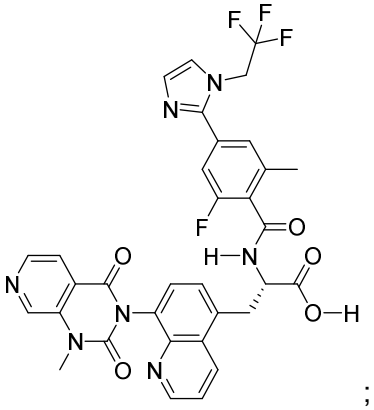

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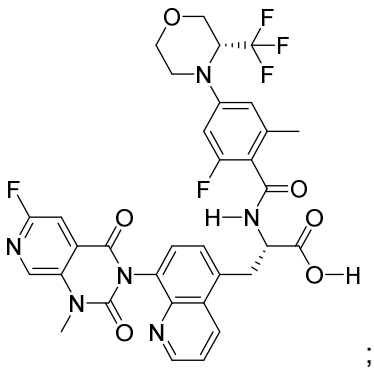
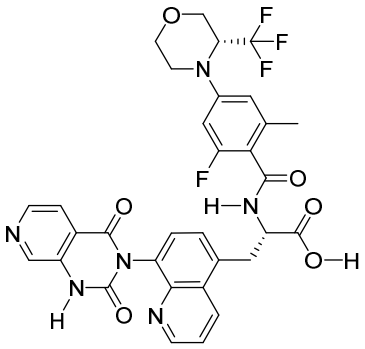
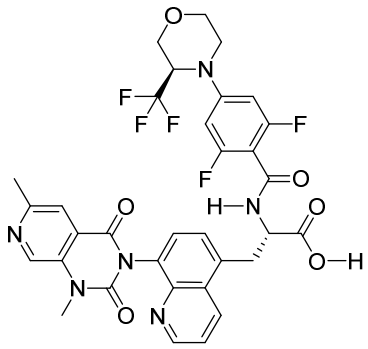
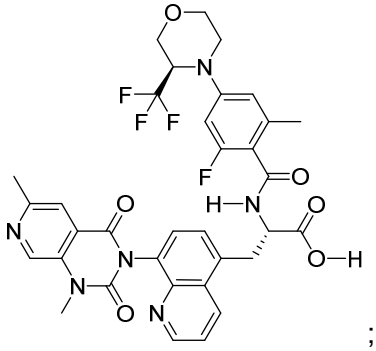
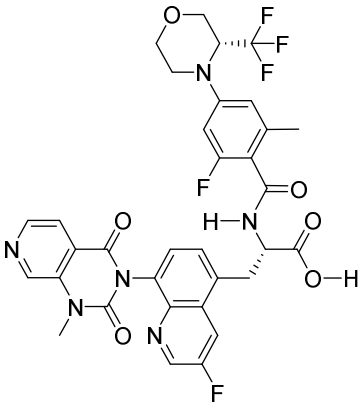
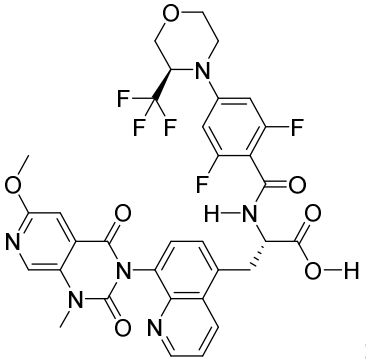
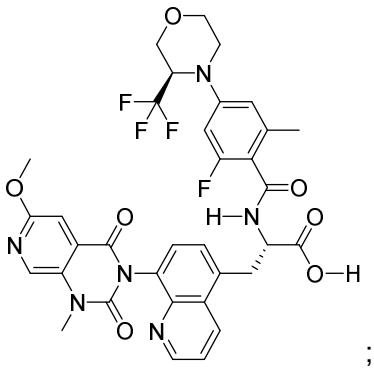
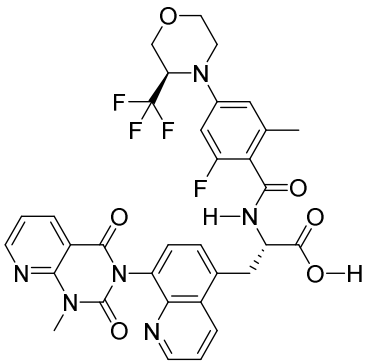
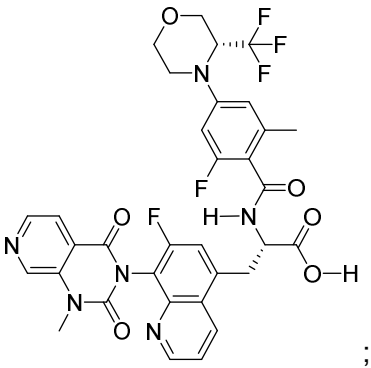
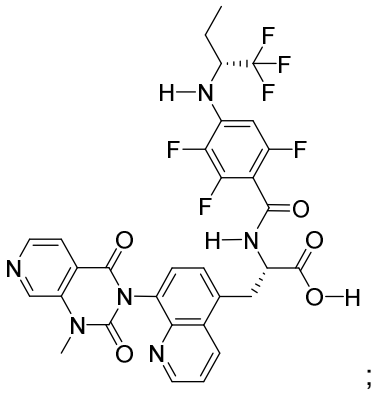
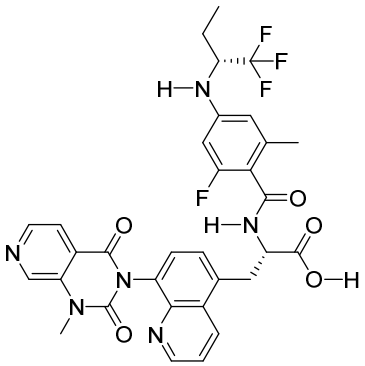
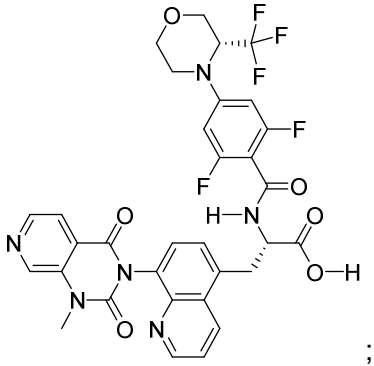

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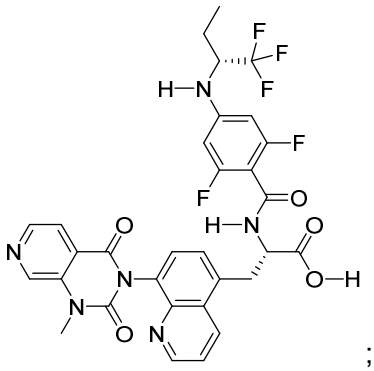
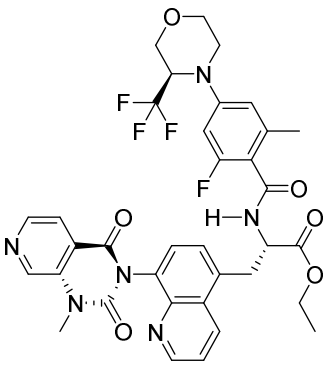
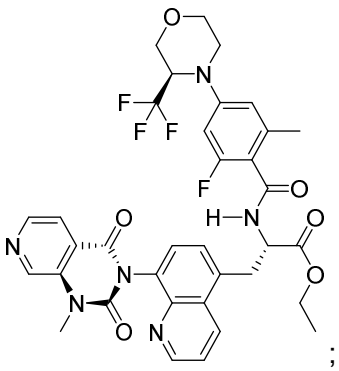
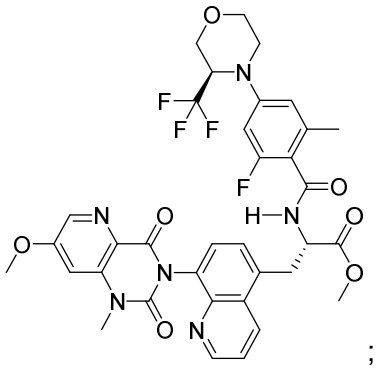
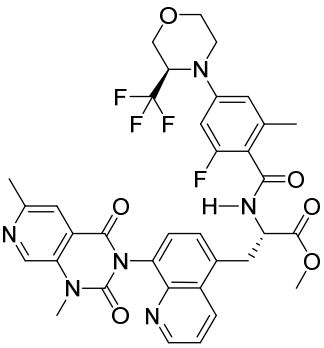
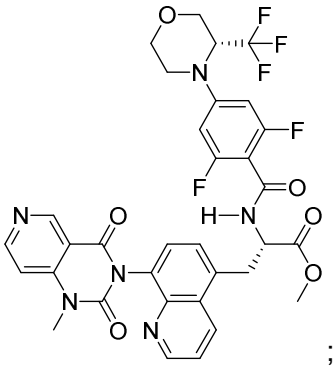
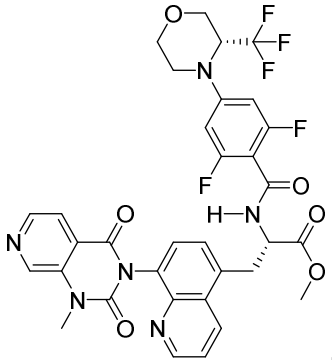

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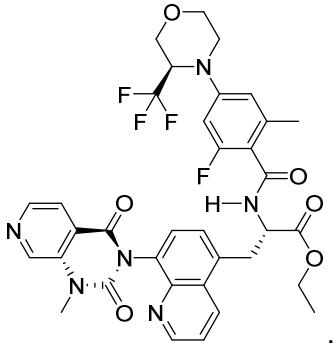

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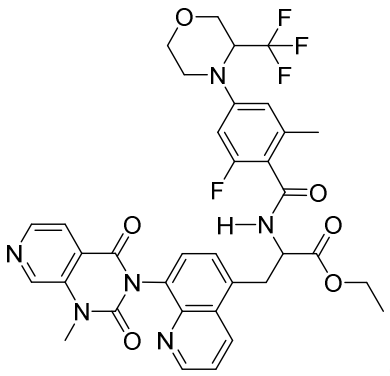

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11. The compound of claim 1, or a pharmaceutically acceptable salt thereof, wherein the compound is:

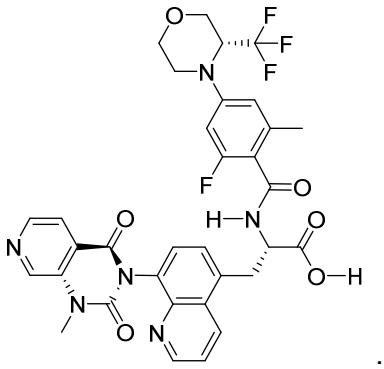

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12. The compound of claim 1 or a pharmaceutically acceptable salt thereof, wherein the compound is:



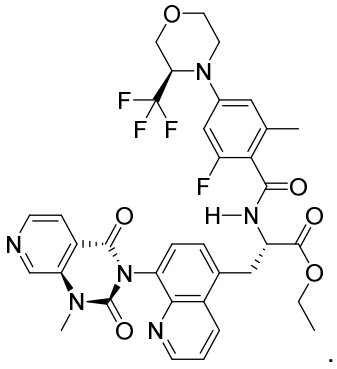
.

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



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



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