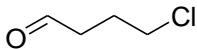
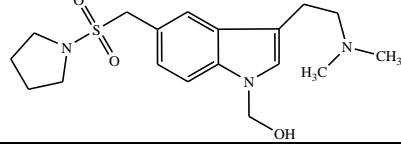
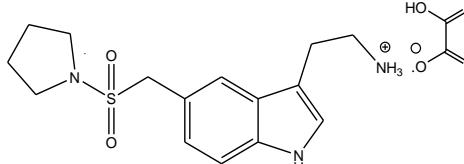
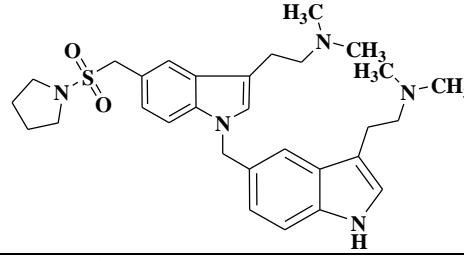
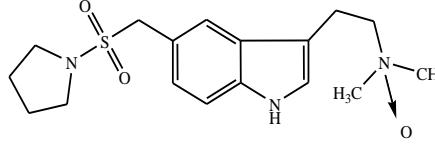
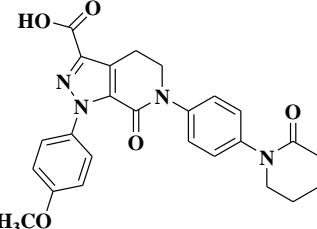
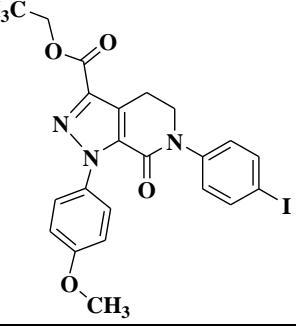
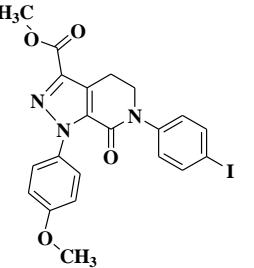
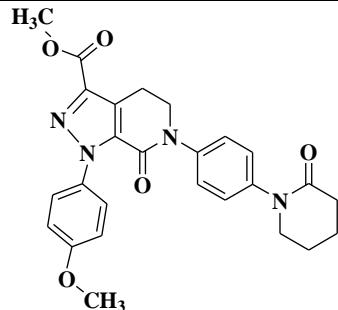
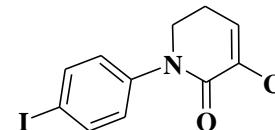
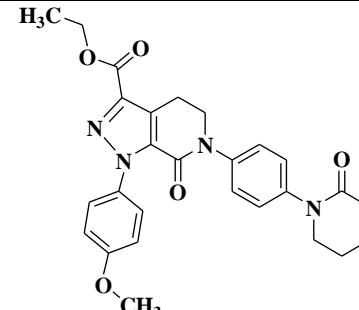
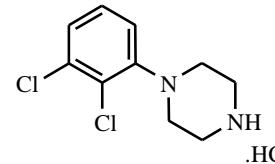
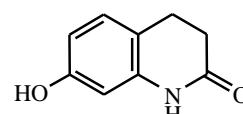
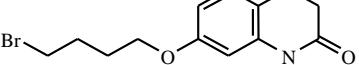
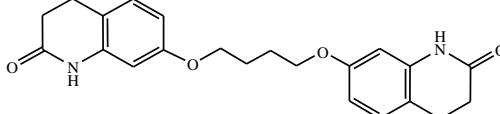
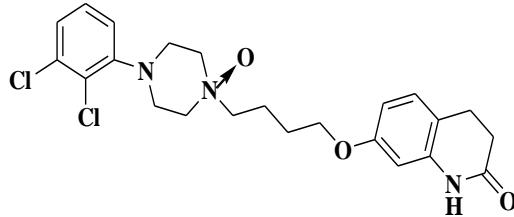


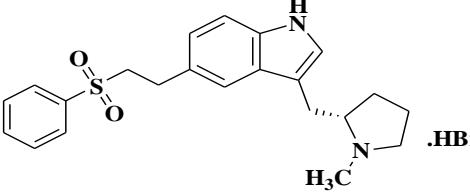
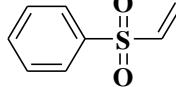
## LIST OF ANALYTICAL STANDARDS

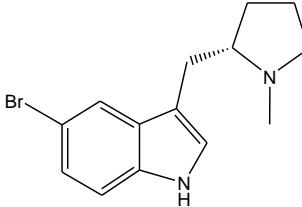
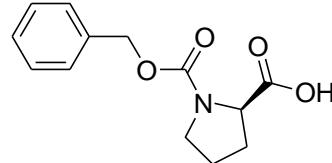
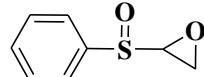
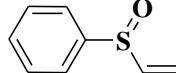
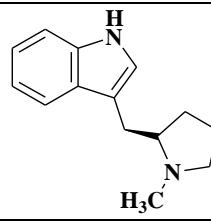
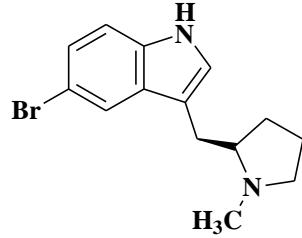
S. No.	Name of the sample	Chemical Name	Chemical Structure	CAS No.	Cost INR Rs./Mg
<b>ALMOTRIPTAN MALATE</b>					
1.	4-Chloro butyraldehyde (CBD-II)	4-Chloro butyraldehyde (CBD-II)		<b>6139-84-0</b>	
2.	Almotriptan impurity-A (1-Hydroxymethyl Almotriptan)	1-[[[3-[2-(Dimethyl amino) ethyl]-1-hydroxy methyl-indol-5-yl]methyl]sulfonyl] pyrrolidine		<b>1018676-02-2</b>	
3.	Almotriptan impurity-B (ATN-Tryptamine)	1-[[[3-[2-(Amino ethyl)]-1H-indol-5-yl] methyl]sulfonyl] pyrrolidine.oxalate		N/A	
4.	ALMO-N-Dimer Impurity	1-[[[3-[2-(Dimethyl amino) ethyl]-1-(N,N-dimethyl-2-(5-methyl-1H-indol-3-yl)ethanamine)-indol-5-yl]methyl]sulfonyl] pyrrolidine		N/A	
5.	Almotriptan impurity-C (ATN-N-Oxide)	1-[[[3-[2-(Dimethyl amino) ethyl]-1H-indol-5-yl] methyl]sulfonyl]pyrrolidine-N-oxide		N/A	

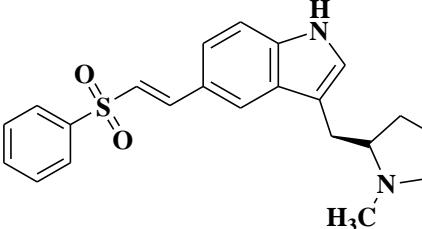
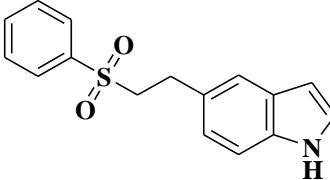
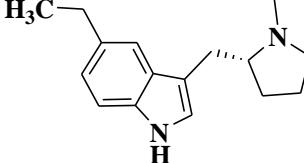
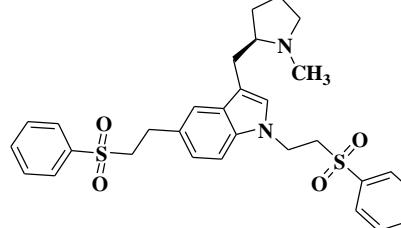
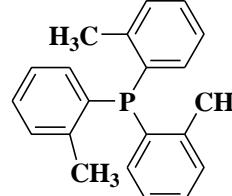
<b>APIXABAN</b>					
1.	APN-Impurity A/ Apixaban acid impurity	4,5,6,7-tetrahydro-1-(4-methoxyphenyl)-7-oxo-6-(4-(2-oxopiperidin-1-yl)phenyl)-1H-pyrazolo[3,4-c]pyridine-3-carboxylic acid		<b>503614-92-4</b>	
2.	APN-Impurity-H (APN-III)	ethyl 4,5,6,7-tetrahydro-6-(4-iodophenyl)-1-(4-methoxyphenyl)-7-oxo-1H-pyrazolo[3,4-c]pyridine-3-carboxylate			
3.	APN-III Methyl ester/ APN-Impurity-D	methyl 4,5,6,7-tetrahydro-6-(4-iodophenyl)-1-(4-methoxyphenyl)-7-oxo-1H-pyrazolo[3,4-c]pyridine-3-carboxylate			

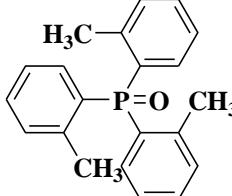
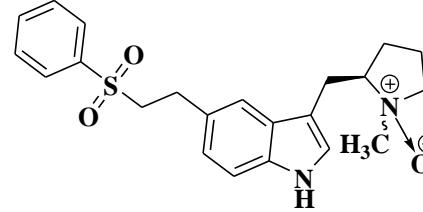
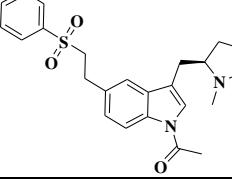
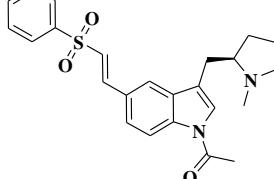
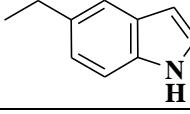
4.	<b>APN-V Methyl ester/ APN-Impurity-K</b>	methyl 4,5,6,7-tetrahydro-1-(4-methoxyphenyl)-7-oxo-6-(4-(2-oxopiperidin-1-yl)phenyl)-1H-pyrazolo[3,4-c]pyridine-3-carboxylate			
5.	<b>APN-Impurity-G (APN-II)</b>	3-chloro-5,6-dihydro-1-(4-iodophenyl)pyridin-2(1H)-one			
6.	<b>APN-Impurity-J (APN-IV)</b>	ethyl 4,5,6,7-tetrahydro-1-(4-methoxyphenyl)-7-oxo-6-(4-(2-oxopiperidin-1-yl)phenyl)-1H-pyrazolo[3,4-c]pyridine-3-carboxylate			
<b>ARIPIPRAZOLE</b>					
7.	DCP-II. HCl	1-(2,3-Dichloro-phenyl)-piperazine hydrochloride		<b>119532-26-2</b>	
8.	APZ-II	7-Hydroxy-3,4-dihydro-1H-quinolin-2-one (or) 3,4-Dihydro-7-hydroxy carbostyryl		<b>22246-18-0</b>	

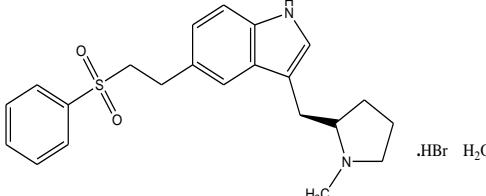
9.	APZ-III	7-(4-Bromo-butoxy)-3,4-dihydro-1H-quinolin-2-one		<b>129722-34-5</b>	
10.	APZ-III/Bis Impurity	1,4-Bis [3,4-dihydro-2(1H)-quinolinone-7-oxy]butane		N/A	
11.	Aripiprazole related compound-F	4-(2,3-Dichlorophenyl)-1-[4-(2-oxo-1,2,3,4-tetrahydro quinolin-7-yloxy)butyl] piperazine-1-oxide (or) 7-[4-[4-(2,3-Dichlorophenyl)-1-oxidopiperazin-1-yl]butoxy]-3,4-dihydroquinolin-2(1H)-one		N/A	
12.	1,4-Dibromo butane	1,4-Dibromo butane		N/A	

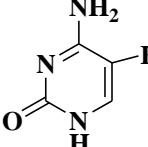
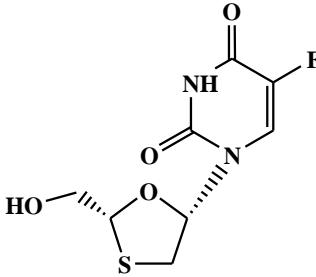
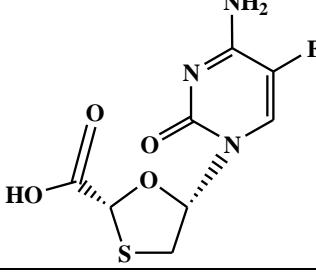
<b>ELETRIPTAN HYDROBROMIDE MONOHYDRATE</b>					
1.	Eletriptan hydrobromide (S-Isomer)	3-[((S)-1-Methylpyrrolidin-2-yl)methyl]-5-[2-(phenyl sulfonyl) ethyl]-1H-indole hydro bromide		<b>1252673-29-2</b>	
2.	ETNSC-II	Phenyl vinyl sulfone (or) 1-(Vinyl sulfonyl)benzene		<b>5535-48-8</b>	

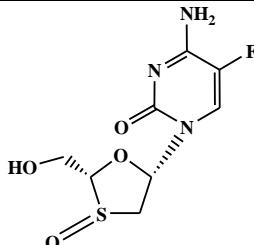
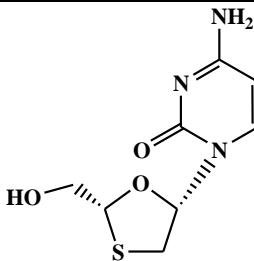
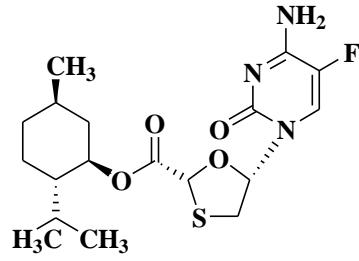
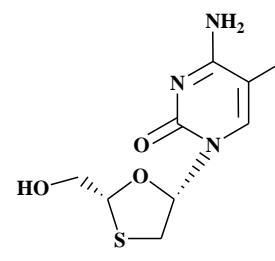
3.	ETN-IIA_(S)-isomer (ETNL-IIA)	5-bromo-3-((S)-1-methylpyrrolidin-2-yl)methyl)-1H-indole		N/A	
4.	ETN-I	(R)-1-((benzyloxy)carbonyl)pyrrolidine-2-carboxylic acid		<b>6404-31-5</b>	
5.	2-(Phenylsulfamoyl)Oxirane	2-(Phenylsulfinyl)Oxirane		N/A	
6.	(Vinylsulfinyl)benzene  (or)  1-(Vinylsulfinyl)benzene	(Vinylsulfinyl)benzene  (or)  1-(Vinylsulfinyl)benzene		N/A	
7.	Eletriptan impurity-A	3-[((R)-1-Methylpyrrolidin-2-yl)methyl]-1H-indole		N/A	
8.	Eletriptan impurity-B	5-Bromo-3-[((R)-1-methylpyrrolidin-2-yl)methyl]-1H-indole		<b>143322-57-0</b>	

9.	Eletriptan impurity-C	3-[((R)-1-Methylpyrrolidin-2-yl)methyl]-5-[(E)-2-(phenyl sulfonyl) vinyl]-1H-indole		N/A	
10.	Eletriptan impurity-D	5-[2-(Phenylsulfonyl)ethyl]-1H-indole		N/A	
11.	Eletriptan impurity-F Or <b>de-phenyl sulfonyl body</b>	5-Ethyl-3-((R)-1-methylpyrrolidin-2-yl)methyl)-1H-indole		<b>209682-64-4</b>	
12.	Eletriptan impurity-G	3-((R)-1-Methylpyrrolidin-2-yl)methyl)-1,5-bis(2-(phenyl sulfonyl)ethyl)-1H-indole		N/A	
13.	Eletriptan impurity-J	Tri(2-methylphenyl)-phosphine (or) Tri-o-tolylphosphine		<b>6163-58-2</b>	

14.	Eletriptan impurity-K	Tri-o-tolylphosphoryloxide		N/A	
15.	Eletriptan impurity-L	3-(((R)-1-Methyl-N-Oxide-pyrrolidin-2-yl)methyl)-5-(2-(phenylsulfonyl)ethyl)-1H-indole		N/A	
16.	<b>Acetyl adduct</b>	1-(3-(((R)-1-methylpyrrolidin-2-yl)methyl)-5-(2-(phenylsulfonyl)ethyl)-1H-indol-1-yl)ethanone		N/A	
17.	<b>Acetyl dehydro Adduct</b>	1-(3-(((R)-1-methylpyrrolidin-2-yl)methyl)-5-((E)-2-(phenylsulfonyl)vinyl)-1H-indol-1-yl)ethanone		<b>188113-71-5</b>	
18.	<b>5-Ethyl-1H-indole</b>	5-Ethyl-1H-indole		<b>68742-28-9</b>	

19.	Eletriptan hydrobromide monohydrate	3-[((R)1-Methylpyrrolidin-2-yl)methyl]-5-[2-(phenyl sulfonyl)ethyl]-1H-indole hydro bromide monohydrate		N/A	
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<b>EMTRICITABINE</b>					
1.	5-Fluorocytosine	5-Fluorocytosine (or) 4-Amino-5-fluoro-2(1H)-pyrimidinone		2022-85-7	
2.	Emtricitabine 5-fluoro uracil impurity	5-Fluoro-1-[(2R,5S)-2-(hydroxymethyl)-1,3-oxathiolan-5-yl]uracil		145986-11-4	
3.	Emtricitabine acid impurity	Cis-5-[4-Amino-5-fluoro-2-oxopyrimidin-1(2H)-yl]-1,3-oxathiolane-2-carboxylic acid		N/A	

4.	Emtricitabine sulfoxides impurity	5-Fluoro-1-[(2R,3RS,5S)-2-(hydroxy methyl)-3-oxo-1,3-oxathiolan-5-yl]cytosine		N/A	
5.	Emtricitabine Lamivudine impurity	(-)-1-[(2R,5S)-2-(Hydroxy methyl)-1,3-oxathiolan-5-yl]cytosine (or) 2(1H)-Pyrimidinone, 4-amino-1-[2-(hydroxyl methyl)-1,3-oxathiolan-5-yl]-(2R-cis)-		N/A	
6.	Emtricitabine methyl ester (ECB-I)	L-Methyl 5-(2R,5S)-[4-amino-5-fluoro-2-oxopyrimidin-1(2H)-yl]-1,3-oxathiolane-2-carboxylate		<b>147126-75-8</b>	
7.	Emtricitabine	4-Amino-5-fluoro-1-[(2R,5S)-2-(hydroxymethyl)-1,3-oxathiolan-5-yl]-2(1H)-pyrimidinone (or) 5-Fluoro-1-[(2R,5S)-2-(hydroxyl methyl)-1,3-oxathiolan-5-yl] cytosine		<b>143491-57-0</b>	

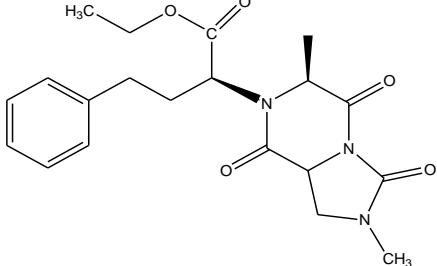
8.	Hydroxy compound (ECB-SM)	(2R,5R)-5-Hydroxy-[1,3]oxathio lane-2-carboxylic acid (1R,2S, 5R)-menth ylester (ECB-SM)		<b>147126-62-3</b>	
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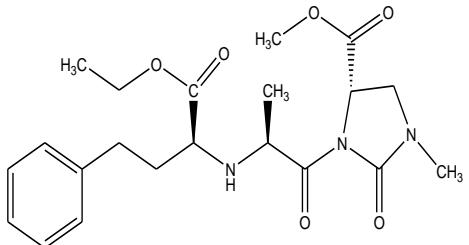
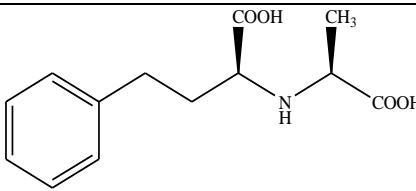
<b>FAMOTIDINE</b>					
9.	Famotidine impurity-A	3-[[[2-[(Diaminomethylene) amino]thiazol-4-yl] methyl] sulfanyl] propanimidamide Dihydrochloride		<b>88061-72-7</b>	
10.	Famotidine impurity-B	3,5-Bis[2-[[[2-[(diaminomethylidene) amino] thiazol-4-yl] methyl] sulfanyl] ethyl]-4H-1,2,4,6-thatriazine 1,1-dioxide		<b>89268-62-2</b>	
11.	Famotidine impurity-C	3-[[[2-[(Diaminomethylidene) amino] thiazol-4-yl] methyl] sulfanyl]-N-sulfamoylpropanamide		N/A	
12.	Famotidine impurity-D	3-[[[2-[(Diaminomethylidene) amino] thiazol-4-yl] methyl] sulfanyl]propamide		<b>76824-16-3</b>	
13.	Famotidine impurity-E	2,2'-[Disulfanediylbis(methylenethiazole-4,2-diyl)]diguanidine		<b>129083-44-9</b>	

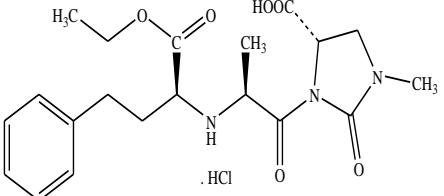
14.	Famotidine impurity-F	3-[[[2-(Diaminomethylidene)amino]thiazol-4-yl] methyl] sulfanyl] propanoic acid		<b>107880-74-0</b>	
15.	Famotidine impurity-G	N-Cyano-3-[[[2-(diaminomethylidene)amino]thiazol-4-yl] methyl] sulfanyl]propanemide		N/A	

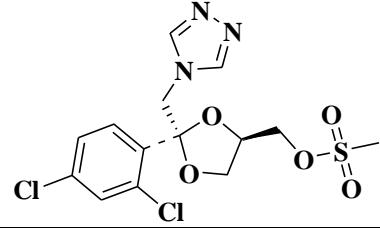
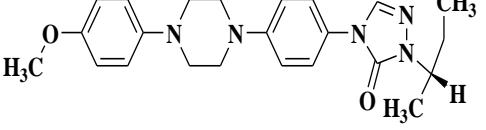
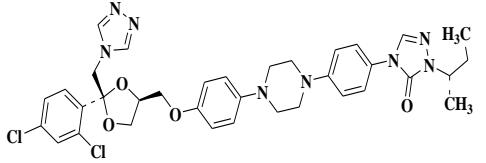
<b>GABAPENTIN</b>					
1.	GBT-SM	1,1-cyclohexanediamic acid monoamide		<b>99189-60-3</b>	
2.	1,1-Cyclohexanediamic acid	1,1-Cyclohexanediamic acid		<b>4355-11-7</b>	

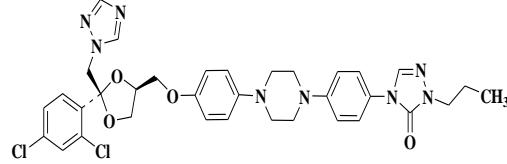
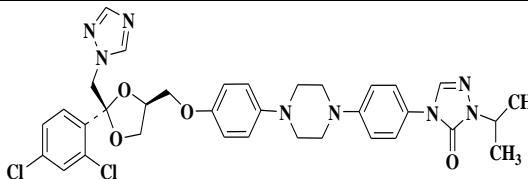
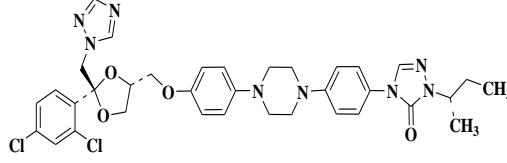
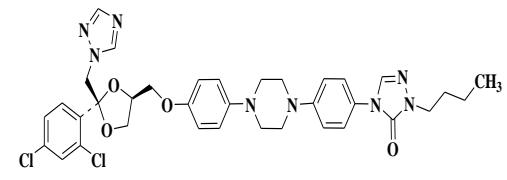
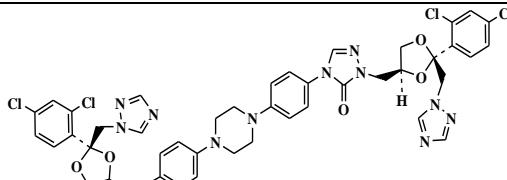
**IMIDAPRIL HYDROCHLORIDE**

1.	Imidapril impurity-A	(S)-Ethyl 2-((6S)-hexahydro-2,6-dimethyl-3,5,8-trioxoimidazo[1,5-a]pyrazin-7(1H)-yl)-4-phenyl butanoate		N/A	
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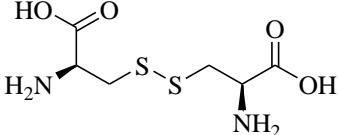
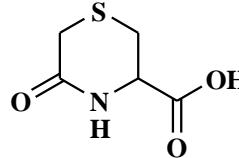
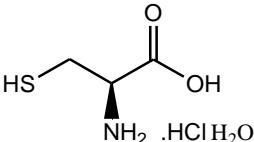
2.	Imidapril impurity-B	(4S)-3-[(2S)-2-[N-[(1S)-1-(Ethoxy carbonyl)-3-phenylpropyl] amino] propionyl]-1-methyl-2-oxoimidazolidine-4-methylester		N/A	
3.	Imidapril impurity-C	(S)-2-((S)-1-Carboxyethylamino)-4-phenylbutanoic acid		N/A	

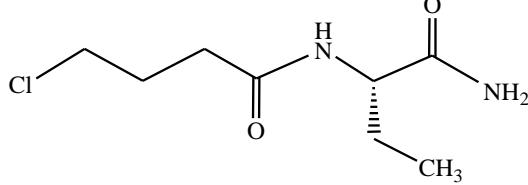
4.	Imidapril hydrochloride	(4S)-3-[(2S)-2-[N-[(1S)-1-(Ethoxycarbonyl)-3-phenylpropyl] amino]propionyl]-1-methyl-2-oxoimidazolidine-4-carboxylic acid hydrochloride		89396-94-1	
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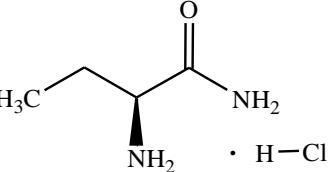
<b>ITRACONAZOLE</b>					
1.	Trans isomer of IT-IB	Trans-(2-(2,4-dichloro-phenyl)-2(1H-1,2,4-triazole-1-yl methyl)-1,3-dioxolan-4-yl) methyl methane sulphonate		N/A	
2.	Itraconazole impurity-A	4-[4-[4-(4-Methoxyphenyl)piperazin-1-yl]phenyl]-2-[(1RS)-1-methylpropyl]-2,4-dihydro-3H-1,2,4-triazol-3-one		252964-68-4	
3.	Itraconazole impurity-B	4-[4-[4-[4-[[Cis-2-(2,4-Dichlorophenyl)-2-(4H-1,2,4-triazol-4-ylmethyl)-1,3-dioxolan-4-yl]methoxy]phenyl]piperazin-1-yl]phenyl]-2-[(1 RS)-1-methylpropyl]-2,4-dihydro-3H-1,2,4-triazol-3-one		854372-77-3	

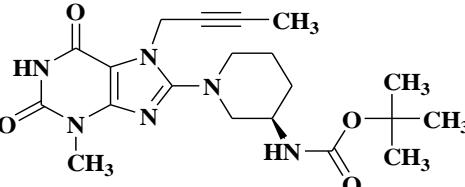
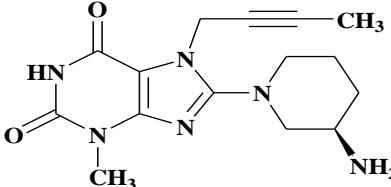
4.	Itraconazole impurity-C	4-[4-[4-[4-[[Cis-2-(2,4-dichlorophenyl)-2-(1H-1,2,4-triazol-1-ylmethyl)-1,3-dioxo lan-4-yl]methoxy] phenyl] piperazin-1-yl] phenyl]-2-propyl-2,4-dihydro-3H-1,2,4-triazol-3-one		<b>74855-91-7</b>	
5.	Itraconazole impurity-D	4-[4-[4-[4-[[Cis-2-(2,4-dichlorophenyl)-2-(1H-1,2,4-triazol-1-ylmethyl)-1,3-dioxo lan-4-yl]methoxy] phenyl] piperazin-1-yl] phenyl]-2-(1-methyl ethyl)-2,4-dihydro-3H-1,2,4-triazol-3-one		<b>89848-49-7</b>	
6.	Itraconazole impurity-E	4-[4-[4-[4-[[Trans-2-(2,4-dichlorophenyl)-2-(1H-1,2,4-triazol-1-ylmethyl)-1,3-dioxolan-4-yl]methoxy]phenyl] piperazin-1-yl]phenyl]-2-[(1RS)-1-methylpropyl]-2,4-dihydro-3H-1,2,4-triazol-3-one		<b>84604-65-9</b>	
7.	Itraconazole impurity-F	2-Butyl-4-[4-[4-[4-[[cis-2-(2,4-dichlorophenyl)-2-(1H-1,2,4-triazol-1-ylmethyl)-1,3-dioxolan-4-yl]methoxy]phenyl] piperazin-1-yl] phenyl]-2,4-dihydro-3H-1,2,4-triazol-3-one		<b>89848-51-1</b>	
8.	Itraconazole impurity-G	4-[4-[4-[4-[[Cis-2-(2,4-Dichlorophenyl)-2-(1H-1,2,4-triazol-1-ylmethyl)-1,3-dioxolan-4-yl]methoxy]phenyl] piperazin-1-yl]phenyl]-2-[[cis-2-(2,4-		<b>N/A</b>	

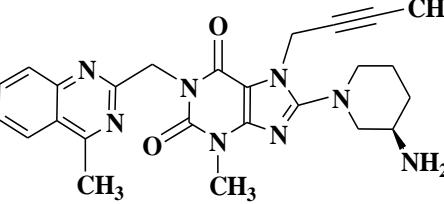
		dichlorophenyl)-2-(1H-1,2,4-triazol-1-ylmethyl)-1,3-dioxolan-4-yl]methyl-2,4-dihydro-3H-1,2,4-triazol-3-one			
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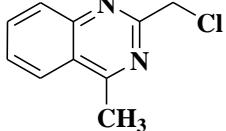
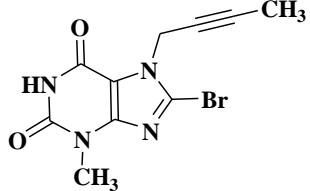
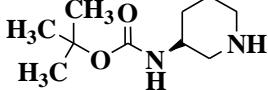
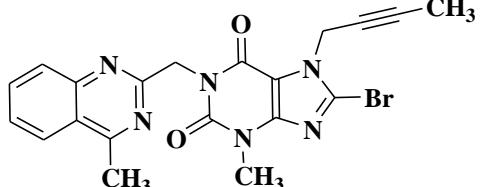
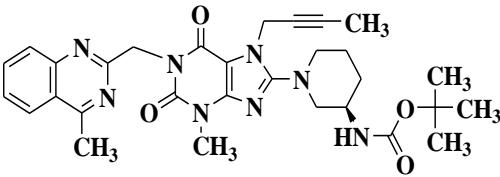
<b>L-CARBO CYSTEINE</b>					
1.	L-Cystine	( <i>R,R</i> )-3,3'-Dithiobis(2-aminopropionicacid)		<b>56-89-3</b>	
2.	TOCA impurity	5-oxothiomorpholine-3-carboxylic acid		<b>14226-97-2</b>	
3.	L-Cysteine HCl H <sub>2</sub> O	L-Cysteine hydrochloride monohydrate		<b>7048-04-6</b>	

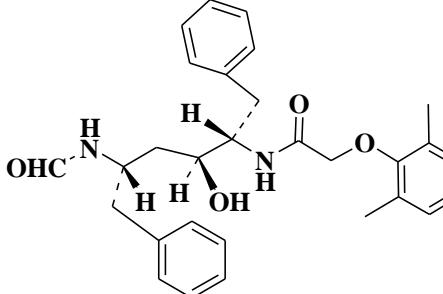
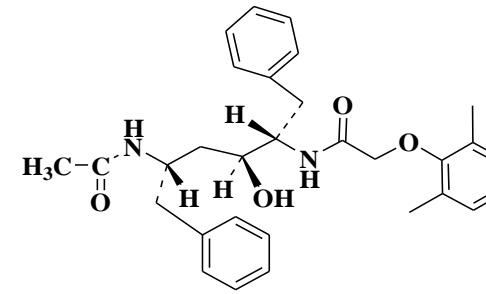
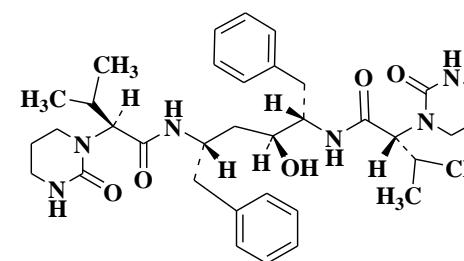
<b>LEVETIRACETAM</b>					
1.	Levetiracetam related compound-A	(S)-N-(1-Amino-1-oxobutan-2-yl)-4-chlorobutanamide		<b>102767-31-7</b>	

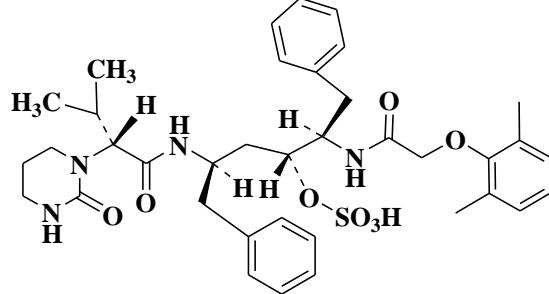
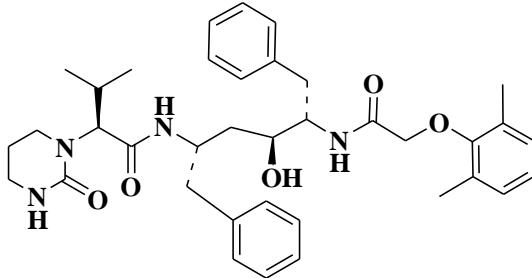
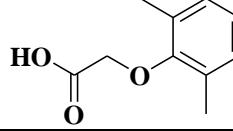
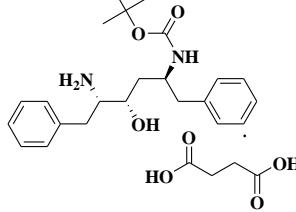
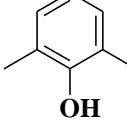
2.	Levetiracetam related compound-B	(S)-2-Aminobutanamide hydrochloride	 <p>Chemical structure: (S)-2-Aminobutanamide hydrochloride. It consists of a four-carbon chain with an amide group (-NH<sub>2</sub>-CO-) at one end and an amino group (-NH<sub>2</sub>) at the other. A vertical wedge bond is shown at the second carbon from the amide group, indicating stereochemistry.</p>	7682-20-4	
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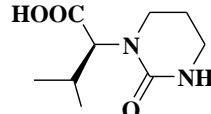
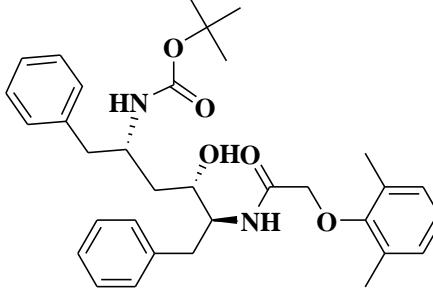
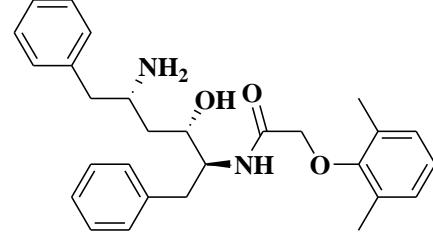
<b>LINAGLIPTIN</b>					
1.	Linagliptin impurity-A	Tert-butyl-1-(7(but-2ynyl)-2, 3, 6, 7-tetrahydro-3-methyl-2,6-dioxo-1H-purin-8-yl)piperidine-3-carbamate.	 <p>Chemical structure: Tert-butyl-1-(7(but-2ynyl)-2, 3, 6, 7-tetrahydro-3-methyl-2,6-dioxo-1H-purin-8-yl)piperidine-3-carbamate. It features a purine ring system with a tert-butyl group at position 1 and a piperidine-3-carbamoyl group at position 7.</p>	N/A	
2.	Linagliptin impurity-B	8-(3-Aminopiperidin-1-yl)-7-(but-2-ynyl)-3-methyl-1H-purine-2,6(3H,7H)-dione.	 <p>Chemical structure: 8-(3-Aminopiperidin-1-yl)-7-(but-2-ynyl)-3-methyl-1H-purine-2,6(3H,7H)-dione. It features a purine ring system with an amino group at position 8 and a but-2-ynyl group at position 7.</p>	N/A	

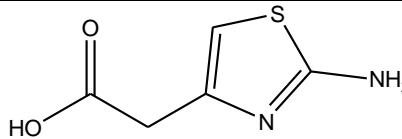
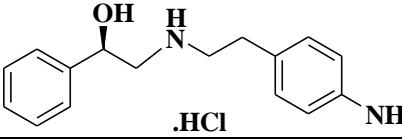
3.	Linagliptin	1-[(4-Methyl-quinazolin-2-yl)methyl]-3-methyl-7-(2-butyn-1-yl)-8-(3-(R)-amino-piperidin-1-yl)-xanthine (Or) 8-[(3R)-3-Aminopiperidin-1-yl]-7-(but-2-yn-1-yl)-3-methyl-1-[(4-methyl quinazolin-2-yl)methyl]-3,7-dihydro-1H-purine-2,6-dione	 <p>Chemical structure: Linagliptin. It is a complex molecule containing a purine ring system linked to a quinazoline ring via its 2-position. It has a methyl group at the 4-position of the quinazoline ring, a but-2-ynyl group at the 7-position of the purine ring, and an amino group at the 3-position of the purine ring.</p>	668270-12-0	
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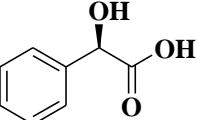
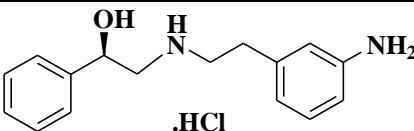
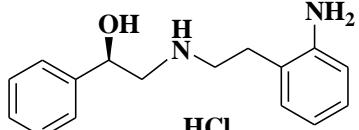
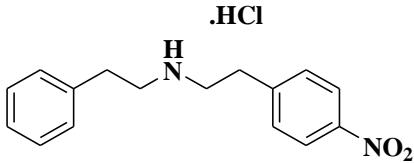
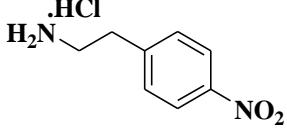
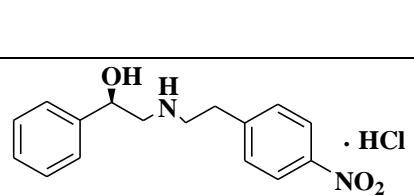
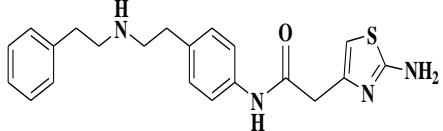
4.	2-(Chloromethyl)-4-methyl quiazoline	2-(Chloromethyl)-4-methyl quiazoline		<b>109113-72-6</b>	
5.	8-Bromo-7-(2-butynyl)-3,7-dihydro-3-methyl-1H-purine 2,6-dione	8-Bromo-7-(2-butynyl)-3,7-dihydro-3-methyl-1H-purine 2,6-dione		<b>666816-98-4</b>	
6.	( R ) -3-(Boc-amino) piperidine	( R ) -3-(Boc- amino) piperidine		<b>309956-78-3</b>	
7.	LGP-I	1-[(4-Methyl-quinazolin-2-yl)methyl]-3-methyl-7-(2-butyn-1-yl)-8-bromoxanthine		<b>853029-57-9</b>	
8.	N-BOC-Linagliptin (LGP-II)	tert-butyl (R)-1-(7-(but-2-ynyl)-2,3,6,7-tetrahydro-3-methyl-1-((4-methylquinazolin-2-yl)methyl)-2,6-dioxo-1H-purin-8-yl)piperidin-3-ylcarbamate		<b>668273-75-4</b>	

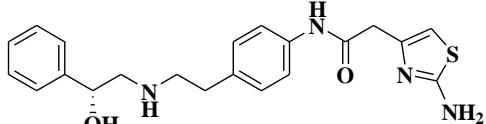
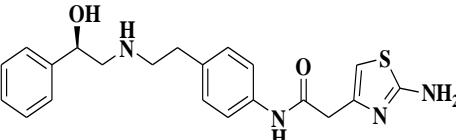
<b>LOPINAVIR</b>					
1.	Lopinavir N-formyl phenoxyacetamide	2-(2,6-Dimethylphenoxy)-N-[(2S, 3S, 5S)-5-formamido-3-hydroxy-1,6-diphenylhexan-2-yl]acetamide	 <p>Chemical structure of Lopinavir N-formyl phenoxyacetamide: A complex organic molecule featuring a central hexane-like core with multiple substituents. It includes two phenyl groups, a formyl group (-CHO), a hydroxyl group (-OH), and an acetamide group (-CONHAc). Stereochemistry is indicated at several points.</p>	N/A	
2.	Lopinavir N-acetyl phenoxy acetamide	N-[(2S, 3S, 5S)-5-Acetamido-3-hydroxy-1,6-diphenylhexan-2-yl]-2-(2,6-dimethylphenoxy) acetamide	 <p>Chemical structure of Lopinavir N-acetyl phenoxy acetamide: Similar to the first structure, but the formyl group is replaced by an acetyl group (-CH3CO-).</p>	N/A	
3.	Lopinavir divalinate	(2R)-N-[(1S,2S,4S)-1-benzyl-2-hydroxy-4-[[2S)-3-methyl-2-[2-oxotetrahydropyrimidin-1(2H)-yl]butanoyl]amino]-5-phenylpentyl]-3-methyl-2-[2-oxotetrahydropyrimidin-1(2H)-yl]butanamide	 <p>Chemical structure of Lopinavir divalinate: A highly complex molecule containing a pyrimidine ring, a piperazine ring, and various amide, ester, and hydroxyl groups, representing a prodrug form of lopinavir.</p>	N/A	

4.	Sulfolopinavir	(1R,3R)-1-[(1R)-1-[[2-(2,6-dimethylphenoxy)acetyl]amino]-2-phenylethyl]-3-[[2(R)-3-methyl-2-[2-oxotetrahydro pyrimidin-1(2H)-yl] butanoyl] amino]-4-phenylbutyl hydrogen sulfate		N/A	
5.	Lopinavir	[1S-[1R*(R*), 3R*, 4R*]]-N-[4[[2,6-Dimethylphenoxy) acetyl]amino]-3-hydroxy-5-phenyl-1-(phenylmethyl)pentyl]-tetrahydro- $\alpha$ -(1-methylethyl)-2-oxo-1(2H)-pyrimidineacetamide		<b>192725-17-0</b>	
6.	2,6 Dimethyl phenoxy acetic acid	2,6-Dimethylphenoxy acetic acid		<b>13335-71-2</b>	
7.	LRV-II (Succinate salt)	(2S,3S,5S)-2-Amino-3-hydroxy-5-(t-butyloxycarbonylamino)-1,6-diphenylhexane succinate salt		<b>169870-02-4</b>	
8.	2,6-Xylenol	2,6-Xylenol		<b>576-26-1</b>	

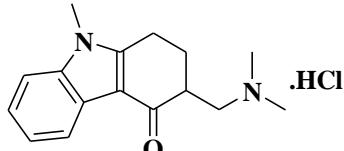
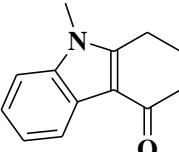
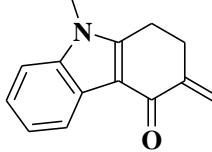
9.	THBA	2S-(1-Tetrahydro-pyrimid-2-onyl)-3-methyl butanoic acid		<b>192725-50-1</b>	
10.	LRV-III	(2S,3S,5S)-2-(2,6-dimethylphe noxyacetyl amino-3-hydroxy-5-(t butyloxycarbonyl amino)-1,6- diphenylhexane		<b>192725-45-4</b>	
11.	LRV-IV	(2S,3S,5S)-2-(2,6-Dimethyl phenoxy acetyl)amino-3-hydro xy- 5-amino-1,6-diphenylhexane		<b>192725-49-8</b>	

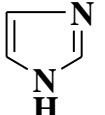
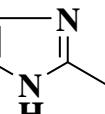
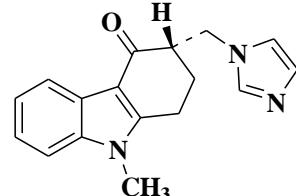
<b>MIRABEGRON</b>					
1.	2-(2-Amino thiazole-4-yl) acetic acid	2-(2-Amino thiazole-4-yl) acetic acid		<b>29676-71-9</b>	
2.	MBG-INT -III	(R)-2-(4-aminophenethylamino)-1- phenylethanol Hydrochloride.		<b>521284-22-0</b>	

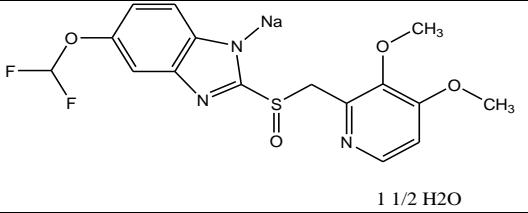
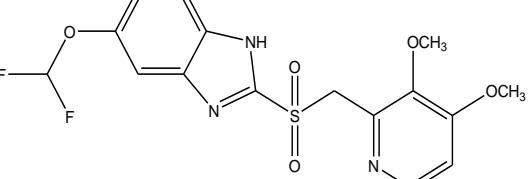
3.	R-Mandellic acid	(R)-2-Hydroxy-2-phenylacetic acid		<b>611-71-2</b>	
4.	MBG-INT-III (3-NITRO)	(R)-2-(3-aminophenethylamino)-1-phenylethanol hydrochloride		N/A	
5.	MBG-INT-III (2-NITRO)	(R)-2-(2-aminophenethylamino)-1-phenylethanol hydrochloride		N/A	
6.	MBG-INT -II DESOXY	2-(4-Nitrophenyl)-N-phenethyl ethanamine hydrochloride.		N/A	
7.	MBG-INT-I(SM-II)	4-Nitrophenyl ethylamine.Hydrochloride		<b>29968-78-3</b>	
8.	MBG-INT-II	(R)-2-(4-nitrophenethylamino)-1-phenylethanol hydrochloride		<b>521284-21-9</b>	
9.	Mirabergon Desoxy Impurity	2-(2-Aminothiazol-4-yl)-N-(4-(2-(phenethylamino)ethyl) phenyl) acetamide		N/A	

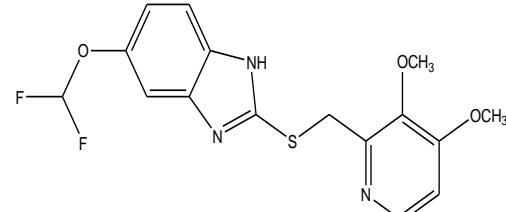
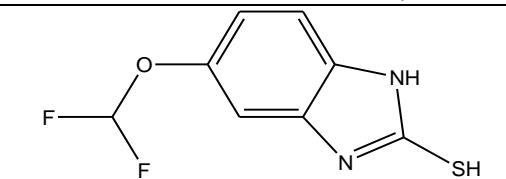
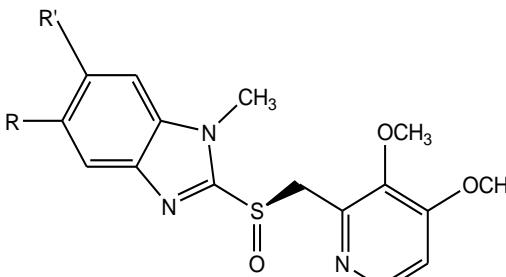
10.	Mirabergon (S-Isomer)	2-(2-amino-1,3-thiazol-4-yl)-N-[4-(2-[(2S)-2-hydroxy-2-phenylethyl] amino) ethyl] phenyl] acetamide.		N/A	
11.	Mirabegron	2-(2-Amino-1,3-thiazol-4-yl)-N-[4-(2-[(2R)-2-hydroxy-2-phenylethyl] amino)ethyl]phenyl]acetamide (or) 2-Amino-N-[4-[2-[(2R)-2-hydroxy-2-phenylethyl] amino]ethyl] phenyl]-4-thiazoleacetamide		<b>223673-61-8</b>	

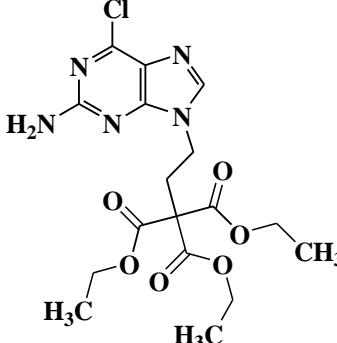
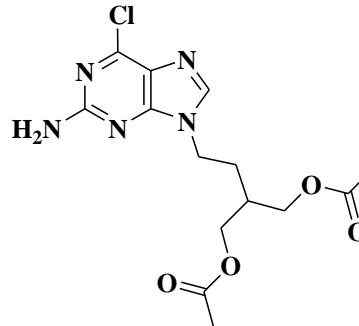
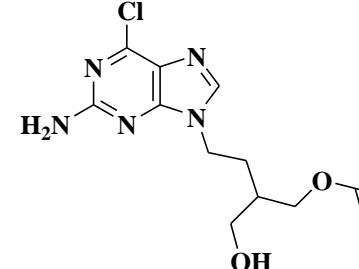
### ONDANSETRON HYDROCHLORIDE

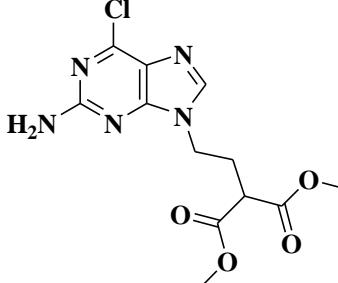
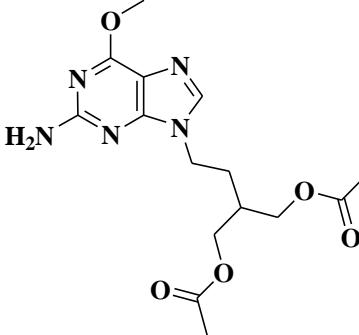
1.	Ondansetron related compound-A	[3-[(Dimethylamino)methyl]-1,2,3,9-tetrahydro-9-methyl-4H-carbazol-4-one] Hydrochloride		<b>119812-29-2</b>	
2.	Ondansetron related compound-C	1,2,3,9-tetrahydro-9-methyl-4H-carbazol-4-one		<b>27387-31-1</b>	
3.	Ondansetron related compound-D	2,3-dihydro-9-methyl-3-methylene-1H-carbazol-4(9H)-one		<b>99614-64-9</b>	

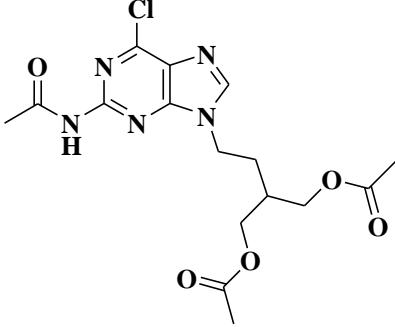
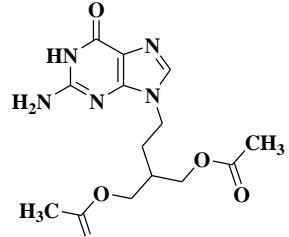
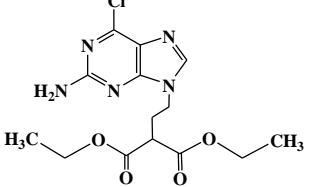
4.	Ondansetron related compound-E	1H-Imidazole		288-32-4	
5.	Ondansetron related compound-F	2-Methyl-1H-imidazole		693-98-1	
6.	Ondansetron impurity-G	(3RS)-3-[(1H-Imidazol-1-yl)methyl]-9-methyl-1,2,3,9-tetrahydro-4H-carbazol-4-one		99614-03-6	

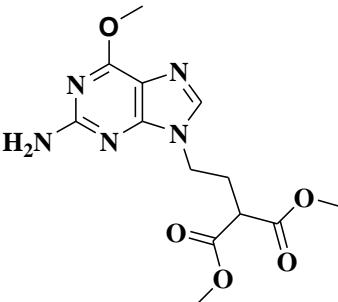
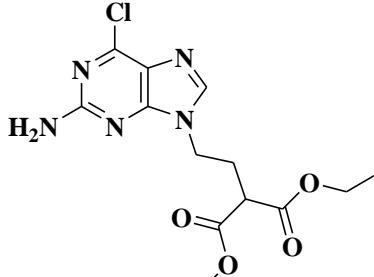
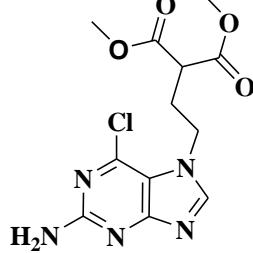
<b>PANTOPRAZOLE SODIUM SESQUIHYDRATE</b>					
1.	Pantoprazole	Sodium 5-(difluoro methoxy)-2-[(RS)-[(3,4-dimethoxypyridin-2-yl)methyl] sulfinyl]benimidazol-1-idesesqui hydrate	 1 1/2 H2O	164579-32-2	
2.	Pantoprazole impurity-A	5-(Difluoro methoxy)-2-[[3,4-dimethoxypyridin-2-yl)methyl] sulfonyl]-1H-benzimidazole		127780-16-9	

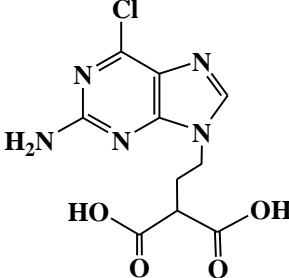
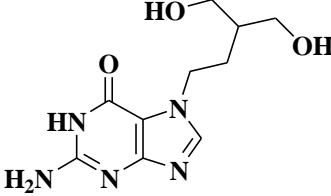
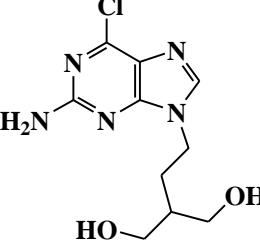
3.	Pantoprazole impurity-B	5-(Difluoro methoxy)-2-[(3,4-dime thoxypyridin-2-yl)methyl] sulfanyl]-1H-benzimidazole		<b>102625-64-9</b>	
4.	Pantoprazole impurity-C	5-Diflouromethoxy-1H-benzimidazole-2-thiol		<b>97963-62-7</b>	
5.	Pantoprazole impurity-D & F	R = OCHF <sub>2</sub> , R' = H; 5-(Difluoro methoxy)-2-[(RS)-[(3,4-dimethoxy pyridin-2-yl)methyl] sulfinyl]-1-meth yl-1H-benzimidazole (Impurity-D) & R = H, R' = OCHF <sub>2</sub> ; 6-(Difluoromethoxy)-2-[(RS)-[(3,4-dimethoxypyridin-2-yl)methyl]sulfinyl]-1-methyl-1H-benzimidazole (Impurity-F)		<b>624742-53-6</b>	

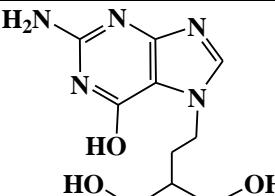
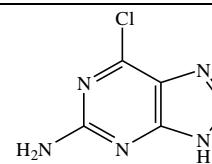
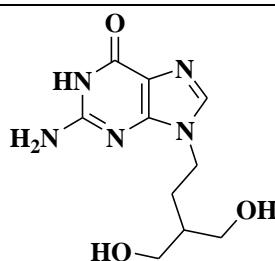
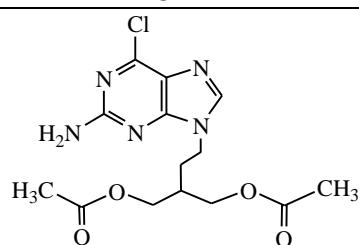
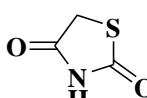
<b>PENCICLOVIR</b>					
1.	<b>PEN-IA</b>	2-Amino-6-chloro-9-(ethyl-2-bis-(carboethoxy)butanoate-4-yl)purine	 <p>The chemical structure shows a purine ring system with a chlorine atom at position 6 and an amino group at position 2. A long side chain is attached at position 9, consisting of an ethyl group, a carbonyl group, and two ester groups linked to a butanoate moiety.</p>	N/A	
2.	<b>BRL 46979/ PEN-II</b>	2-[2-(2-amino-6-chloro-9H-purin-9-yl)ethyl]-1,3-propanediol-1,3-diacetate	 <p>The chemical structure shows a purine ring system with a chlorine atom at position 6 and an amino group at position 2. A propyl side chain is attached at position 9, which further branches into a diacetate group and a hydroxyl group.</p>	N/A	
3.	<b>BRL-048420</b>	2-Amino-6-Chloro-9-(methyl-2-acetoxy-3-hydroxymethyl)butyl-1-ylpurine (Or)4-(2-amino-6-chloro-9H-purin-9-yl)-2-(hydroxy methyl)butyl acetate	 <p>The chemical structure shows a purine ring system with a chlorine atom at position 6 and an amino group at position 2. A propyl side chain is attached at position 9, which branches into a hydroxyl group and an acetate group.</p>	N/A	

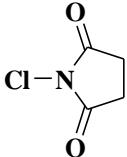
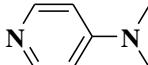
4.	<b>PEN-I</b>	2-Amino-6-Chloro-9-(methyl-2-carbomethoxybutanoate-1-yl)purine	 <p>Chemical structure of 2-Amino-6-Chloro-9-(methyl-2-carbomethoxybutanoate-1-yl)purine: A purine ring system with an amino group at position 2 and a chlorine atom at position 6. At position 9, there is a methylene group attached to a carbomethoxybutanoate side chain. The side chain consists of a four-carbon chain ending in a carbomethoxy group (-CH<sub>2</sub>-CH<sub>2</sub>-COOMe).</p>	N/A	
5.	<b>SB-224875</b>	9-(4-acetoxy-3-acetoxymethylbut-1-yl)-2-amino-6-methoxypurine	 <p>Chemical structure of 9-(4-acetoxy-3-acetoxymethylbut-1-yl)-2-amino-6-methoxypurine: A purine ring system with an amino group at position 2 and a methoxy group at position 6. At position 9, there is a methylene group attached to a side chain consisting of a four-carbon chain with two acetoxymethyl groups (-CH<sub>2</sub>-CH<sub>2</sub>-O-C(=O)-CH<sub>3</sub>) attached to the same carbon atom.</p>	N/A	

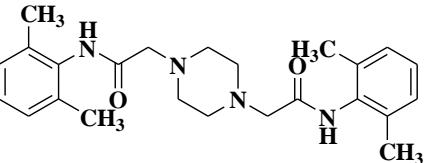
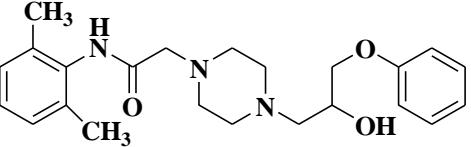
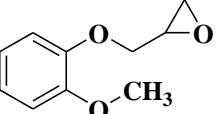
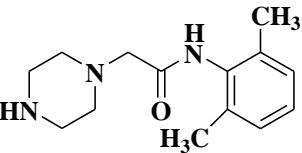
6.	<b>SB-216483</b>	2-Acetylaminio-6-chloro-9-(4-acetoxyethylbut-1-yl)purine		N/A	
7.	<b>PEN-Di-O-acetyl</b>	2-(2-(2-amino-1,6-dihydro-6-oxopurin-9-yl)ethyl)-1,3-diacetoxypropane Or Di-O-Acetoxy Penciclovir		<b>97845-72-2</b>	
8.	<b>PEN-I Diethyl ester</b>	diethyl 2-(2-(2-amino-6-chloro-9H-purin-9-yl)ethyl)malonate		N/A	

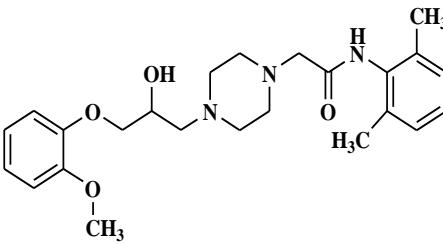
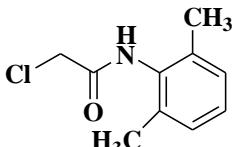
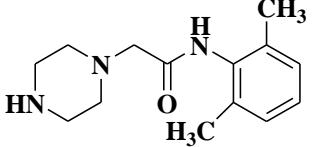
9.	<b>PEN-I Methoxy Impurity</b>	2-Amino-6-methoxy-9-(methyl-2-carbomethoxybutanoate-4-yl)purine	 <p>Chemical structure of 2-Amino-6-methoxy-9-(methyl-2-carbomethoxybutanoate-4-yl)purine: A purine ring system with an amino group at position 2 and a methoxy group at position 6. At position 9, there is a methyl-2-carbomethoxybutanoate side chain attached via an amide linkage.</p>	<b>131266-11-0</b>	
10.	<b>ACP methyl ethyl ester Or PEN-I Un symmetrical ester</b>	2-Amino-6-Chloro-9-(ethyl-2-carbomethoxybutanoate-4-yl)purine	 <p>Chemical structure of 2-Amino-6-Chloro-9-(ethyl-2-carbomethoxybutanoate-4-yl)purine: A purine ring system with an amino group at position 2 and a chlorine atom at position 6. At position 9, there is an ethyl-2-carbomethoxybutanoate side chain attached via an amide linkage.</p>	N/A	
11.	<b>PEN-I-7-Isomer</b>	Dimethyl 2-(2-(2-amino-6-chloro-7H-purin-7-yl)ethyl)malonate	 <p>Chemical structure of Dimethyl 2-(2-(2-amino-6-chloro-7H-purin-7-yl)ethyl)malonate: A purine ring system with an amino group at position 2 and a chlorine atom at position 6. At position 7, there is a dimethyl malonate side chain attached via an ester linkage.</p>	N/A	

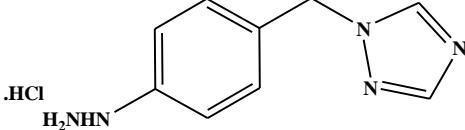
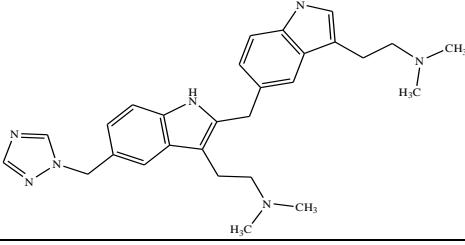
12.	<b>PEN-I-Acid</b>	2-(2-(2-amino-6-chloro-9H-purin-9-yl)ethyl)malonic acid	 <p>Chemical structure: A purine ring system with a chlorine atom at position 6 and an amino group at position 2. Attached to the 9-position is an ethyl group (-CH<sub>2</sub>CH<sub>3</sub>) which is further substituted with a malonic acid group (-CH<sub>2</sub>COOH).</p>	N/A	
13.	<b>PEN-III-N-7-Isomer</b>	7-(4-hydroxy-3-(hydroxymethyl)butyl)guanine	 <p>Chemical structure: A purine ring system with an amino group at position 2 and a chlorine atom at position 6. Attached to the 7-position is a butyl group (-CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>) which has hydroxyl groups (-OH) at positions 4 and 5.</p>	<b>127205-22-5</b>	
14.	<b>PEN-IIA</b>	2-Amino-6-chloro-9-(2-hydroxy methylbutanol-4-yl)purine	 <p>Chemical structure: A purine ring system with an amino group at position 2 and a chlorine atom at position 6. Attached to the 9-position is a butyl group (-CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>) which has a hydroxyl group (-OH) at position 2.</p>	<b>172529-94-1</b>	

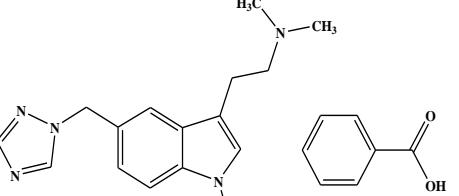
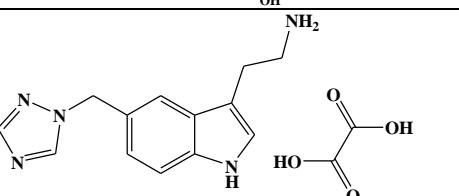
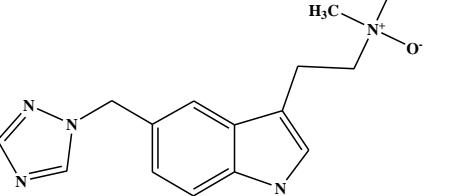
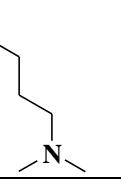
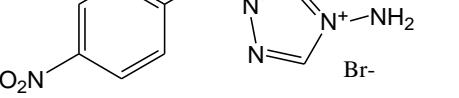
15.	<b>BRL46970</b>	7-(4-Hydroxy-3-hydroxymethylbut-1-yl)guanine	 <p>Chemical structure of 7-(4-Hydroxy-3-hydroxymethylbut-1-yl)guanine: A purine ring system with an amino group at position 6 and a guanine side chain at position 7. The guanine side chain consists of a four-carbon chain with hydroxyl groups at positions 3 and 4, and a terminal hydroxyl group.</p>	N/A	
16.	2-Amino-6-chloropurine (ACP)	2-Amino-6-chloropurine (ACP)	 <p>Chemical structure of 2-Amino-6-chloropurine (ACP): A purine ring system with an amino group at position 2 and a chlorine atom at position 6.</p>	<b>10310-21-1</b>	
17.	PEN-III	9-(4-hydroxy-3-(hydroxymethyl)butyl)guanine	 <p>Chemical structure of 9-(4-hydroxy-3-(hydroxymethyl)butyl)guanine: A purine ring system with an amino group at position 6 and a guanine side chain at position 9. The guanine side chain consists of a four-carbon chain with hydroxyl groups at positions 3 and 4, and a terminal hydroxyl group.</p>	<b>39809-25-1</b>	
18.	PEN-II	2-[2-(2-Amino-6-chloro-9H-purin-9-yl)ethyl]-1,3-propanediol-1,3-diacetate (PEN-II)	 <p>Chemical structure of 2-[2-(2-Amino-6-chloro-9H-purin-9-yl)ethyl]-1,3-propanediol-1,3-diacetate (PEN-II): A purine ring system with an amino group at position 2 and a chloropurine side chain at position 9. The side chain is further substituted with a propylene group, which is linked to a 1,3-propanediol-1,3-diacetate group.</p>	<b>97845-60-8</b>	
19.	2,4-Thiazolidinedione	2,4-Thiazolidinedione	 <p>Chemical structure of 2,4-Thiazolidinedione: A thiazolidinedione ring system.</p>	<b>2295-31-0</b>	

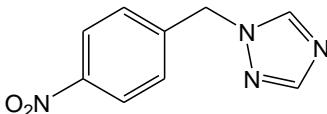
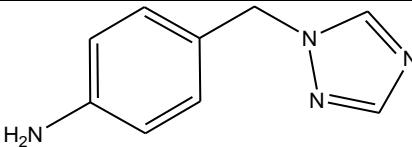
20.	NCS	N-Chlorosuccinimide		<b>128-09-6</b>	
21.	DMAP	4-Dimethylaminopyridine		<b>1122-58-3</b>	

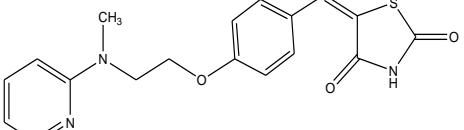
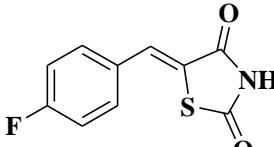
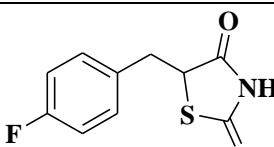
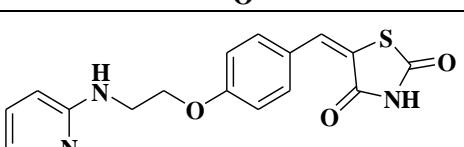
<b>RANOLAZINE</b>					
1.	RNZ-Bis piperazine acetamide impurity	2,2'-(Piperazine-1,4-diyl) bis[N-(2,6-dimethylphenyl)] acetamide		<b>380204-72-8</b>	
2.	RNZ-Desmethoxy impurity	N-(2,6-Dimethylphenyl)-2-[4-(2-hydroxy-3-phenoxypropyl)piperazin-1-yl]acetamide		<b>755711-09-2</b>	
3.	RNZ-INT	2-[(2-Methoxyphenoxy) methyl] Oxirane		<b>2210-74-4</b>	
4.	RNZ-II	N-(2, 6-Dimethylphenyl)-2-(piperazin-1-yl) acetamide		<b>5294-61-1</b>	

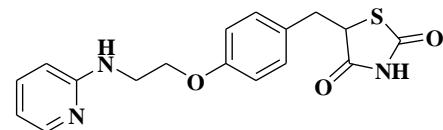
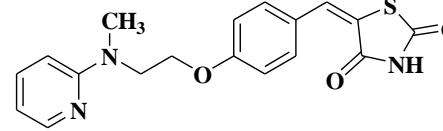
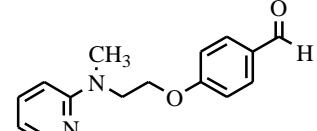
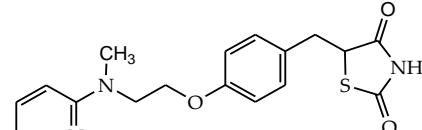
5.	Ranolazine	1-Piperazineacetamide, N-(2,6-dimethylphenyl)-4-[2-hydroxy-3-(2-methoxyphenoxy)propyl]-± (or) (±)-N-(2,6-Dimethylphenyl)-4-[2-hydroxy-3-(2-methoxy phenoxy)propyl]-piperazin-1-ylacetamide		<b>95635-55-5</b>	
6.	RNZ-I	2-Chloro-N-(2,6-dimethylphenyl)acetamide		<b>1131-01-7</b>	
7.	RNZ-II	N-(2,6-dimethylphenyl)-2-(piperazin-1-yl) acetamide		<b>5294-61-1</b>	

<b>RIZATRIPTAN BENZOATE</b>					
1.	RTZ-III A	1-((4-((1H-1,2,4-triazol-1-yl)methyl)phenyl)hydrazine Hydrochloride		<b>212248-62-9</b>	
2.	Rizatriptan impurity-2,5-Dimer	3-[2-(N,N-Dimethylamino) ethyl-2-[[3-[2-(N,N-dimethylamino)ethyl]-1H-indol-5-yl]methyl]-1H-indol-5-yl]methyl-1H-1,2,4-triazole benzoate		<b>1135479-44-5</b>	

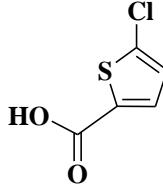
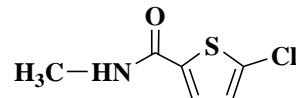
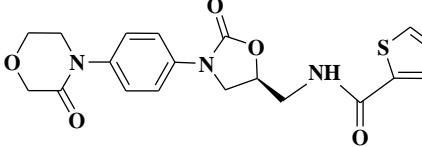
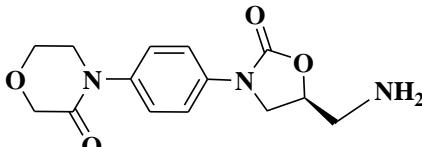
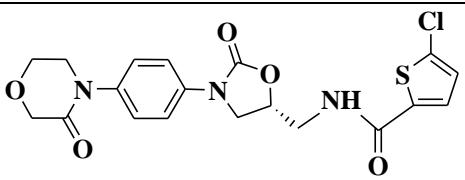
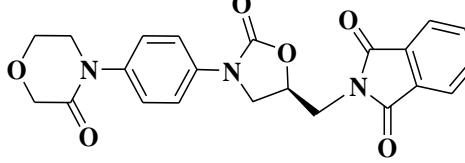
3.	Rizatriptan impurity-A	N,N-Dimethyl-5-((1H-1,2,4-triazol-1-yl)methyl)-1-hydroxymethyl-indol-3-yl)ethanamine.benzoate		<b>N/A</b>	
4.	Rizatriptan impurity-B	2-(5-((1H-1,2,4-Triazol-1-yl)methyl)-1H-indol-3-yl)ethanamine. oxalate		<b>N/A</b>	
5.	Rizatriptan impurity-C	2-(5-((1H-1,2,4-Triazol-1-yl)methyl)-1H-indol-3-yl)-N,N-dimethylethanamine oxide		<b>N/A</b>	
6.	4-(Dimethyl amine)butanal	4-(Dimethyl amine)butanal		<b>104459-70-3</b>	
7.	RTZ-I	4-Amino-1-(4-nitrobenzyl)-4H-1,2,4-triazolium bromide		<b>N/A</b>	

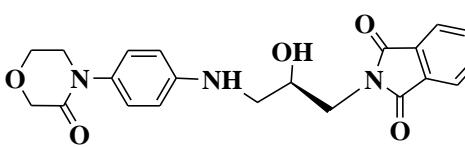
8.	RTZ-II	1H-1-[(4-Nitrophenyl)methyl]-[1,2,4]-triazole		<b>119192-09-5</b>	
9.	RTZ-III	1H-1-[(4-Aminophenyl) methyl]-1,2,4-triazole		<b>119192-10-8</b>	

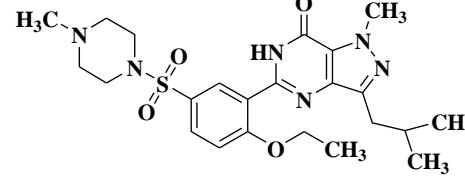
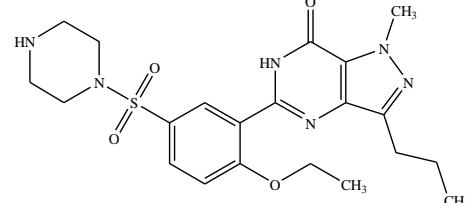
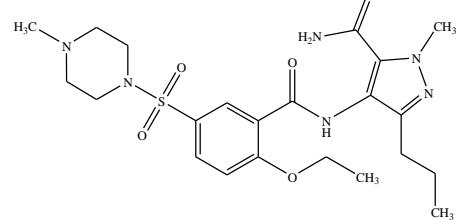
ROSIGLITAZONE MALATE					
1.	Rosiglitazone related compound-A	5-[4-[2-[N-Methyl-N-(2-pyridinyl)amino]ethoxy]benzylidene]thiazolidine-2,4-dione		<b>122320-74-5</b>	
2.	In-house Rosiglitazone impurity-A	(Z)-5-(4-Fluorobenzylidene)thiazolidine-2,4-dione		<b>291536-35-1</b>	
3.	In-house Rosiglitazone impurity-B	5-(4-Fluorobenzyl)thiazolidine-2,4-dione		<b>291536-42-0</b>	
4.	In-house Rosiglitazone impurity-C	(Z)-5-(4-(2-Pyridine-2-yl-amino)ethoxy)benzylidene)thiazolidine-2,4-dione		N/A	

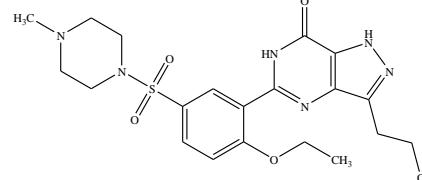
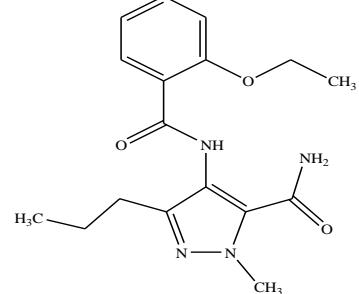
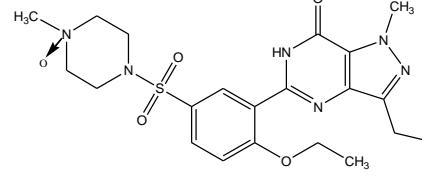
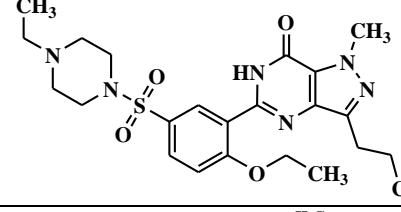
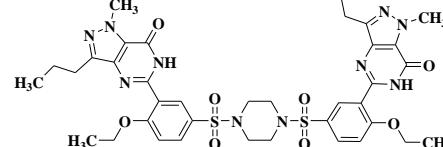
5.	In-house Rosiglitazone impurity-D	5-(4-(2-Pyridin-2-yl-amino)ethoxy)benzyl thiazolidine-2,4-dione		N/A	
6.	ROS-IB	5-[4-[2-[N-Methyl-N-(2-pyridinyl)amino] ethoxy] benzylidene] thiazolidine-2,4-dione		<b>122320-74-5</b>	
7.	ROS-IAB	4-[2-(Methyl-(2-pyridyl)amino)ethoxy] benzaldehyde		<b>122321-03-3</b>	
8.	ROS-II	5-[4-[2-(N-methyl-N-(2-pyridyl)amino)ethoxy]benzyl] thiazolidine-2,4-dione(ROS-II)		<b>122320-73-4</b>	

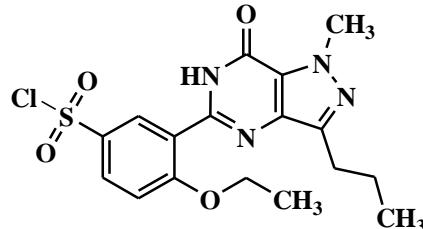
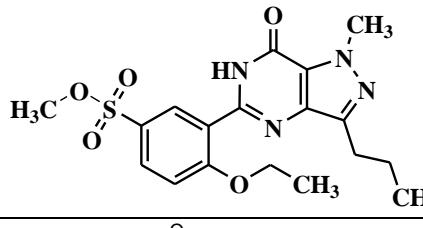
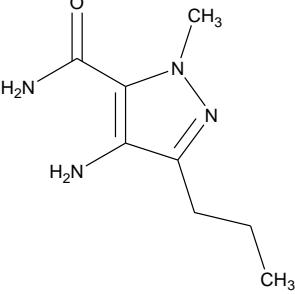
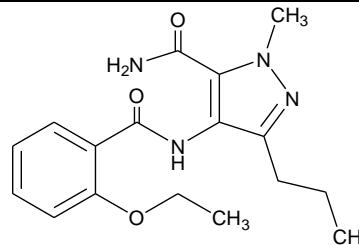
## RIVAROXABAN

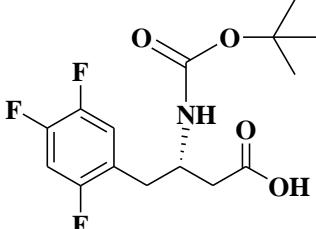
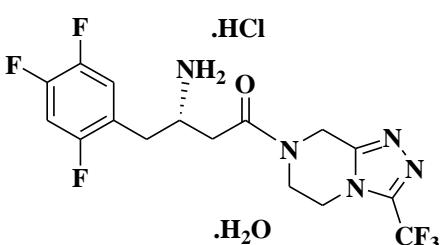
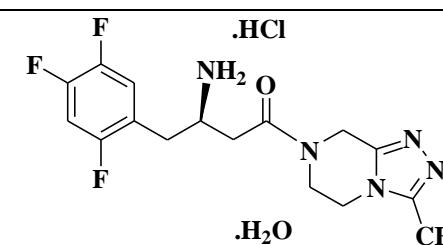
1.	Thiophene Acid impurity/ Impurity-A	5-Chlorothiophene-2-carboxylic acid		<b>24065-33-6</b>	
2.	Thiophene-N-methyl carboxamide/ Impurity-B	5-Chloro-N-methylthiophene-2-carboxamide		<b>31166-29-7</b>	
3.	Des Chloro Rivaroxaban / Impurity-C	N-((5S)-2-Oxo-3-[4-(3-oxo-4-morpholinyl)phenyl]-1,3-oxazolidin-5-yl)methyl)-2-thiophenecarboximide		<b>N/A</b>	
4.	Amino methyl impurity/ RBN-IIA	4-((S)-5-(aminomethyl)-2-oxooxazolidin-3-yl)phenyl)morpholin-3-one		<b>898543-06-1</b>	
5.	Rivaroxaban R-Isomer	5-ChloroN-((5R)-2-Oxo-3-[4-(3-oxo-4-morpholinyl)phenyl]-1,3-oxazolidin-5-yl)methyl)-2-thiophenecarboximide		<b>865479-71-6</b>	
6.	RBN-II	2-((5S)-2-Oxo-3-[4-(3-oxo-4-morpholinyl)phenyl]-1,3-oxazolidin-5-yl)methyl)-1H-isoindole-1,3(2H)-dione		<b>446292-08-6</b>	

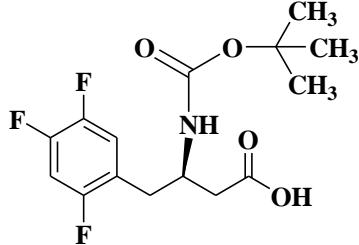
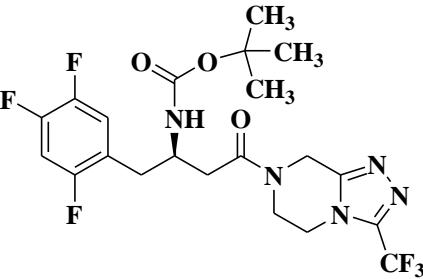
7.	RBN-I	2-((2R)-2-Hydroxy-3-{{[4-(3-oxo-4-morpholinyl) phenyl] amino} propyl}-1H-isoindole-1,3(2H)-dione		<b>446292-07-5</b>	
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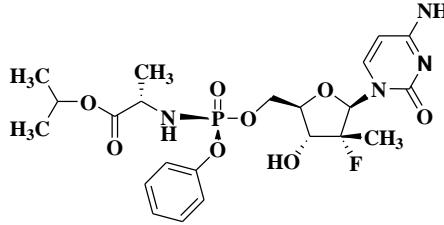
<b>SILDENAFIL CITRATE</b>					
1.	Sildenafil isobutyl impurity	5-[2-Ethoxy-5-[(4-methylpiperazin-1-yl)sulfonyl]phenyl]-1-methyl-3-(2-methylpropyl)-1,6-dihydro-7H-pyrazolo[4,3-d]pyrimidin-7-one		<b>1391053-95-4</b>	
2.	Sildenafil impurity-A	5-(2-Ethoxy-5-((piperazin-1-yl)sulfonyl)phenyl)-1-methyl-3-propyl-1H-pyrazolo[4,3-d]pyrimidin-7(6H)-one		<b>139755-82-1</b>	
3.	Sildenafil impurity-B	4-(2-Ethoxy-5-((4-methylpiperazin-1-yl)sulfonyl)benzamido)-1-methyl-3-propyl-1H-pyrazole-5-carboxamide		<b>N/A</b>	

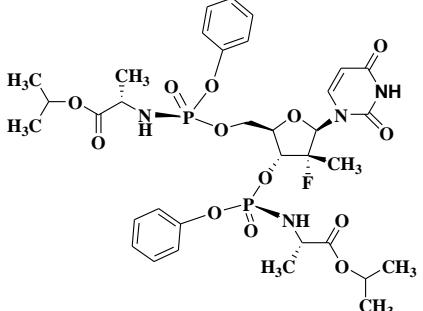
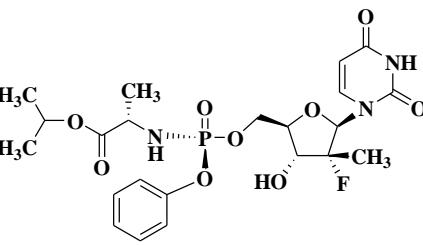
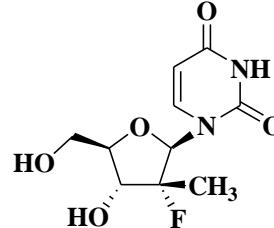
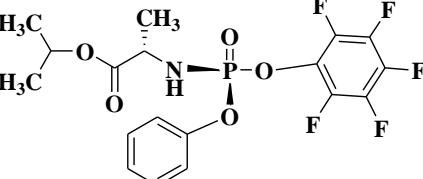
4.	Sildenafil impurity-C	5-(2-Ethoxy-5-((4-methylpiperazin-1-yl)sulfonyl)phenyl)-3-propyl-1H-pyrazolo[4,3-d] pyrimidin-7(6H)-one		N/A	
5.	Sildenafil impurity-D	4-(2-Ethoxybenzamido)-1-methyl-3-propyl-1H-pyrazole-5-carboxamide		N/A	
6.	Sildenafil N-oxide impurity	5-[2-Ethoxy-5-[(4-methyl-4-oxidopiperazin-1-yl)sulfonyl]phenyl]-1-methyl-3-propyl-1,6-dihydro-7H-pyrazolo[4,3-d]pyrimidin-7-one		<b>1094598-75-0</b>	
7.	Sildenafil N-Ethyl impurity	1-[[3-(4,7-Dihydro-1-methyl-7-oxo-3-propyl-1H-pyrazolo[4,3-d]-pyrimidin-5-yl)-4-ethoxyphenyl]sulfonyl]-4-ethylpiperazine		N/A	
8.	Sildenafil dimer impurity	Bis-1,4-[4-ethoxy-3-(6,7-dihydro-1-methyl-7-oxo-3-propyl-1H-pyrazolo [4,3-d] pyrimidin-5-yl)-benzenesulfonyl] piperazine		N/A	

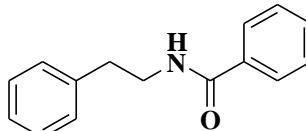
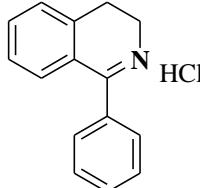
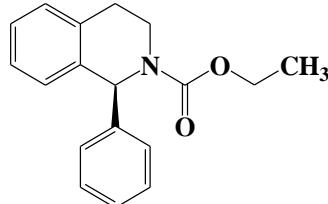
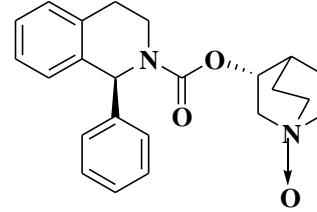
9.	Sildenafil chlorosulfonyl impurity	4-Ethoxy-3-(6,7-dihydro-1-methyl-7-oxo-3-propyl-1H-pyrazolo[4,3-d]pyrimidin-5-yl)benzene-1-sulfonyl chloride (or) 5-(5-Chlorosulphonyl-2-ethoxyphenyl)-1-methyl-3-n-propyl-1,6-dihydro-7H-pyrazolo [4,3-d] pyrimidine-7-one		<b>139756-22-2</b>	
10.	Sildenafil methyl sulfonate impurity	Methyl-4-ethoxy-3-(6,7-dihydro-1-methyl-7-oxo-3-propyl-1H-pyrazolo[4,3-d]pyrimidin-5-yl)benzenesulfonate		<b>N/A</b>	
11.	SLC-IIIA	1-Methyl-4-amino-3-propyl-1H-pyrazole-5-carboxamide		<b>139756-02-8</b>	
12.	SLC-IIIB	4-(2-Ethoxy-benzamido)-1-methyl-3-propyl-1H-pyrazole-5-carboxamide		<b>139756-03-9</b>	

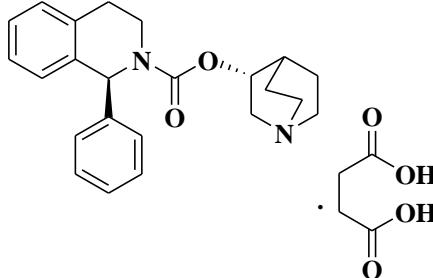
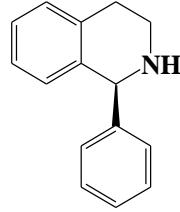
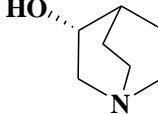
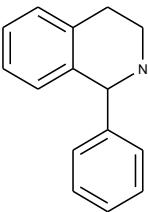
SITAGLIPTIN HYDROCHLORIDE					
1.	SGN-VII S-Isomer	Boc-(S)-3-Amino-4-(2,4,5-trifluorophenyl) butanoic acid	 <p>Detailed description: A chemical structure showing a 4-(2,4,5-trifluorophenyl)butanoic acid molecule substituted with a Boc group at the 3-position and an amino group at the 2-position. The carboxylic acid group is at the 4-position.</p>	<b>922178-94-7</b>	
2.	SGN-IX S-Isomer	(2S)-4-oxo-4-[3-(trifluoro methyl)-5,6-dihydro [1, 2, 4] triazolo [4,3-a]pyrazin-7(8H)-yl]-1-(2,4,5-trifluorophenyl) butan-2-amine hydrochloride monohydrate. (or) (3S)-3-Amino-1-[3-trifluorome thyl]-5,6-dihydro[1,2,4] triazolo[4,3-a]pyrazin-7(8H)-yl]-4-(2,4,5-trifluorophenyl) butan-1-one hydrochloride monohydrate.	 <p>Detailed description: A chemical structure showing a 4-(2,4,5-trifluorophenyl)butan-1-one molecule substituted with a 5,6-dihydro[1,2,4]triazolo[4,3-a]pyrazin-7(8H)-yl group at the 4-position and an amino group at the 2-position. It is shown as a monohydrate with HCl and H<sub>2</sub>O.</p>	N/A	
3.	Sitagliptin hydrochloride	(2R)-4-oxo-4-[3-(trifluoromethyl)-5,6-dihydro[1,2,4]triazolo [4,3-a]pyrazin-7(8H)-yl]-1-(2,4,5-trifluorophenyl)butan-2-amine Hydrochloride monohydrate.	 <p>Detailed description: A chemical structure showing a 4-(2,4,5-trifluorophenyl)butan-1-one molecule substituted with a 5,6-dihydro[1,2,4]triazolo[4,3-a]pyrazin-7(8H)-yl group at the 4-position and an amino group at the 2-position. It is shown as a monohydrate with HCl and H<sub>2</sub>O.</p>	<b>862156-92-1</b>	

4.	SGN-VII	Boc-(R)-3-Amino-4-(2,4,5-trifluorophenyl)butanoic acid	 <p>Chemical structure of Boc-(R)-3-Amino-4-(2,4,5-trifluorophenyl)butanoic acid: A benzene ring substituted with three fluorine atoms at positions 2, 4, and 5. It is connected via a methylene group to a chiral center (R). The chiral center is bonded to an amino group (NH), a tert-butyl carbamate group (O=C(=O)OC(CH<sub>3</sub>)<sub>3</sub>), and a carboxylic acid group (-COOH).</p>	486460-00-8	
5.	SGN-VIII	tert-Butyl (R)-4-(3-(trifluoromethyl)-5,6-dihydro-[1,2,4]triazolo [4,3-a]pyrazin-7(8H)-yl)-1-(2,4,5-trifluorophenyl)-4-oxobutan-2-ylcarbamate	 <p>Chemical structure of tert-Butyl (R)-4-(3-(trifluoromethyl)-5,6-dihydro-[1,2,4]triazolo [4,3-a]pyrazin-7(8H)-yl)-1-(2,4,5-trifluorophenyl)-4-oxobutan-2-ylcarbamate: A benzene ring substituted with three fluorine atoms at positions 2, 4, and 5. It is connected via a methylene group to a chiral center (R). The chiral center is bonded to an amino group (NH), a tert-butyl carbamate group (O=C(=O)OC(CH<sub>3</sub>)<sub>3</sub>), and a carbonyl group (-C(=O)-) which is further bonded to a 5,6-dihydro-[1,2,4]triazolo [4,3-a]pyrazin-7(8H)-yl group and a 2-oxobutan-2-yl group.</p>	N/A	

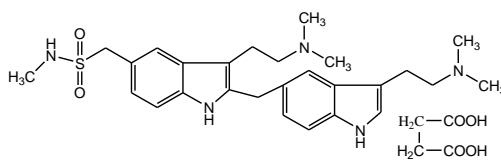
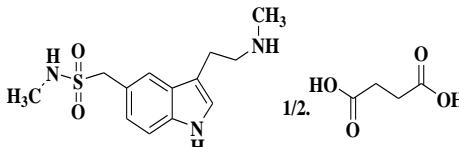
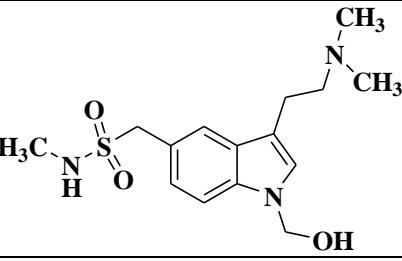
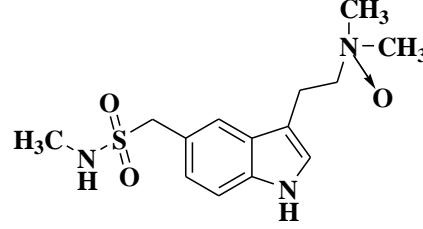
<b>SOFOSBUVIR</b>					
1.	Sofosbuvir amine impurity	Isopropyl ((S)-(((2R,3R,4R,5R)-5-(4-amino-2-oxopyrimidin-1(2H)-yl)-4-fluoro-3-hydroxy-4-methyltetrahydrofuran-2-yl)methoxy)(phenoxy)phosphoryl)-L-alaninate	 <p>Chemical structure of Sofosbuvir amine impurity: A complex molecule featuring a 2-aminopyrimidine ring fused to a 4-hydroxypiperidine-2-one ring. This is linked via a phosphorus atom to a 2-hydroxy-4-methyltetrahydrofuran-2-yl group and an isopropyl ((S)-((4R,5R)-5-(4-amino-2-oxopyrimidin-1(2H)-yl)-4-fluoro-3-hydroxy-4-methyltetrahydrofuran-2-yl)methoxy)phosphoryl)-L-alaninate group.</p>	N/A	

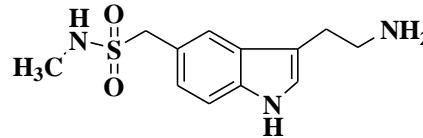
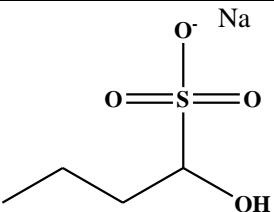
2.	Sofosbuvir dimer impurity	Isopropyl (((2R,3R,4R,5R)-5-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)-4-fluoro-2-(((S)-((S)-1-isopropoxy-1-oxopropan-2-yl)amino)(phenoxy)phosphoryl)oxy)methyl)-4-methyltetra hydrofuran-3-yl)oxy)(phenoxy)phosphoryl)-L-alaninate		1256490-29-5	
3.	Sofosbuvir	Isopropyl ((2S(P))-2-[[[(2R,3R,4R,5R)-5-(2,4-dioxopyrimidin-1-yl)-4-fluoro-3-hydroxy-4-methyl-tetrahydrofuran-2-yl]methoxy-phenoxy-phosphoryl]amino]propanoate		1190307-88-0	
4.	SBR SM-I	2'-Deoxy-2'-fluoro-2'-C-methylcytidine		863329-66-2	
5.	SBR SM-II OR Penta fluorophenyl analog	(S)-2-[-(S)-(2,3,4,5,6-Pentafluorophenoxy)-phenoxy-phosphorylamino]propionic acid isopropyl ester		1334513-02-8	

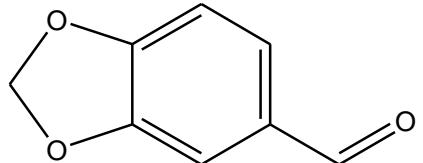
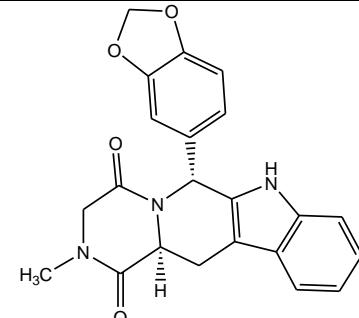
<b>SOLIFENACIN SUCCINATE</b>					
1.	Phenethylbenzamide (SFN-I)	N-Phenethylbenzamide		<b>3278-14-6</b>	
2.	Phenyldihydroisoquinoline (SFN-IIA)	1-Phenyl-3,4-dihydroisoquinoline hydrochloride		<b>52250-51-8</b>	
3.	Phenyltetrahydroisoquinoline ethyl carbamate (SFN-IVA)	Ethyl (S)-1-phenyl-3,4-dihydroisoquinoline-2(1H)-carboxylate		<b>180468-42-2</b>	
4.	Solifenacin N-oxide	(R)-3-{[(S)-1-Phenyl-1,2,3,4-tetrahydroisoquinoline-2-carbonyl]oxy}quinuclidine 1-oxide.		<b>180272-28-0</b>	

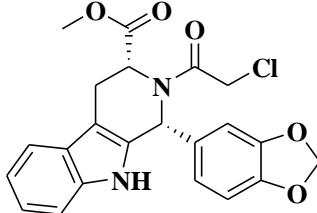
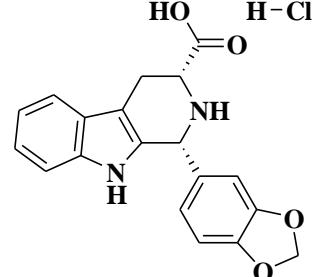
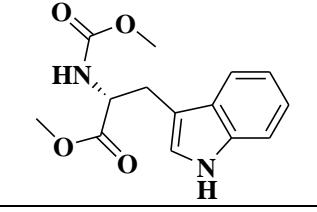
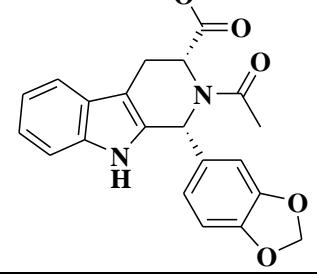
5.	Solifenacin succinate	Butanedioic acid, compound with (1S)-(3R)-1- azabicyclo[2.2.2]oct-3-yl 3,4-dihydro-1-phenyl-2(1H)-isoquinoline carboxylate (or) (R)-Quinuclidin-3-yl (S)-1-Phenyl-3,4-dihydroisoquinoline 2(1H)- carboxylate succinate salt		<b>242478-38-2</b>	
6.	<b>SFN-IV</b>	(S)-1-Phenyl-1,2,3,4-tetrahydroisoquinoline		<b>118864-75-8</b>	
7.	R-(-)-3Quinuclidinol	R-(-)-3Quinuclidinol		<b>25333-42-0</b>	
8.	<b>SFN-III</b>	1,2,3,4-tetrahydro-1-phenylisoquinoline		<b>22990-19-8</b>	

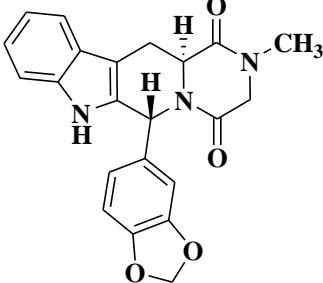
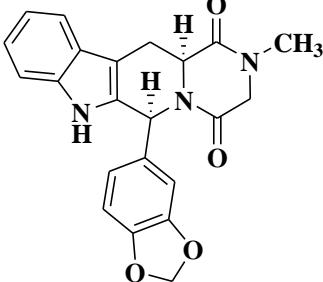
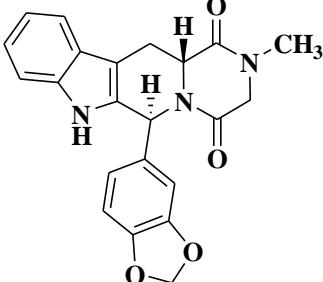
**SUMATRIPTAN SUCCINATE**

1.	Sumatriptan succinate related compound-A	[3-[2-(Dimethyl amino)ethyl]-2-[[3-[2-(dimethyl amino)ethyl]-1H-indol-5-yl]methyl]-1H-indole-5yl]-N-methylmethanesulphonamide succinate		N/A	
2.	Sumatriptan related compound-B	N-Methyl [3-[2-(methylamino)ethyl]-1H-indol-5-yl] methane sulphonamide. Hemi succinate (or) [3-[2-(Methyl amino)ethyl]-1H-indol-5-yl]-N-Methylmethane sulphonamide. Hemi succinate		N/A	
3.	Sumatriptan related compound-C	[3-[2-(Dimethyl amino)ethyl]-1-(hydroxy methyl)-1H-indol-5-yl]-N-methyl methanesulphonamide		<b>1797905-62-4</b>	
4.	Sumatriptan related compound-D	N,N-Dimethyl-2-[5-(methyl sulphonyl)methyl]-1H-indol-3-yl] Ethan amine-N-oxide (or) [3-[2-(Dimethylamino-N-oxide)ethyl]-1H-indol-5-yl]-N-methylmethanesulfonamide		N/A	

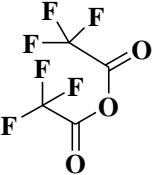
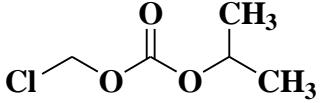
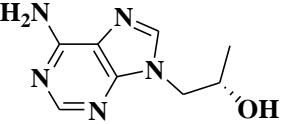
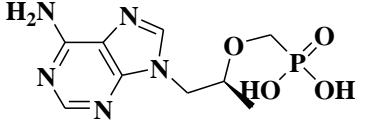
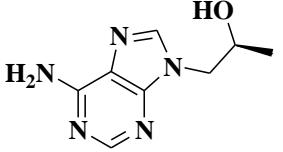
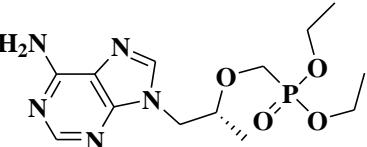
5.	Sumatriptan related compound-E	[3-(2-Aminoethyl)-1H-indol-5-yl]-N-methyl methanesulphonamide		<b>88919-22-6</b>	
6.	4-Chloro butyraldehyde sodium bisulphite	4-Chloro butyraldehyde sodium bisulphite		<b>54322-20-2</b>	

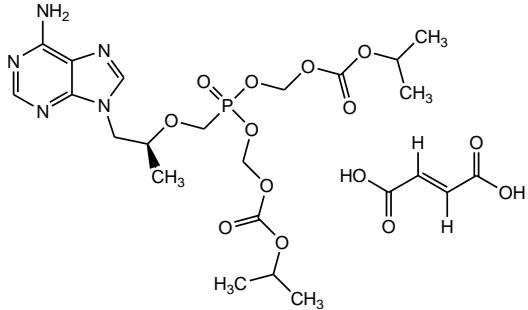
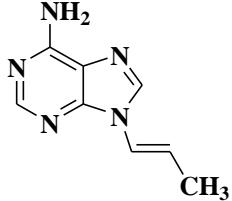
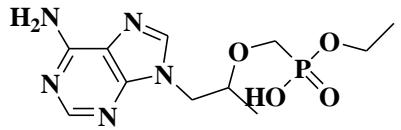
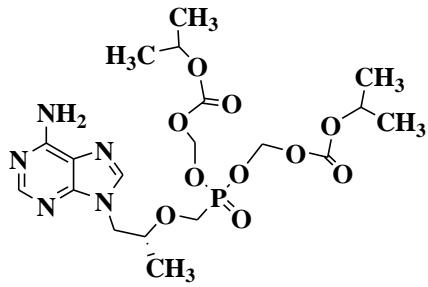
<b>TADALAFIL</b>					
1.	Tadalafil impurity-A	Benzo[d][1,3]dioxole-5-carbaldehyde		<b>120-57-0</b>	
2.	Tadalafil impurity-B	(6R-Cis)-6-(1,3-benzodioxol-5-yl)-2,3,6,7,12,12a-hexahydro-2-methylpyrazino[1',2':1,6]pyrido[3,4-b]indole-1,4-dione		N/A	

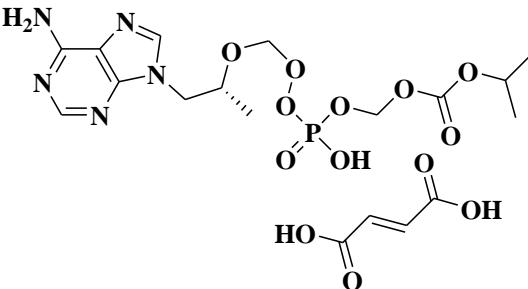
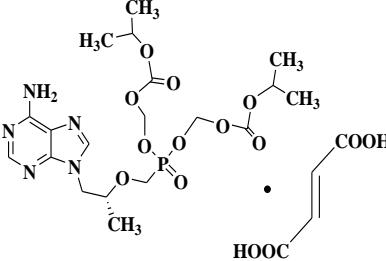
3.	Tadalafil impurity-C	(1R, 3R)-methyl-1-(benzo[d][1,3]dioxol-5-yl)-2-(2-chloroacetyl)-2,3,4,9-tetrahydro-1H-pyrido[3,4-b]indole-3-carboxylate		<b>171489-59-1</b>	
4.	Tadalafil impurity-1	(1R,3R)-1-(Benzo[d][1,3]dioxol-5-yl)-2,3,4,9-tetrahydro-1H-pyrido[3,4-b]indole-3-carboxylic acid hydrochloride		N/A	
5.	Tadalafil impurity-2	Methyl (R)-1-(methoxy carbonyl)-2-(1H-indol-3yl)ethyl carbamate		N/A	
6.	Tadalafil impurity-3	(1R,3R)-Methyl 2-acetyl-1-(benzo[d][1,3]dioxol-5yl)-2,3,4,9-tetrahydro-1H-pyrido[3,4-b]indole-3-carboxylate		N/A	

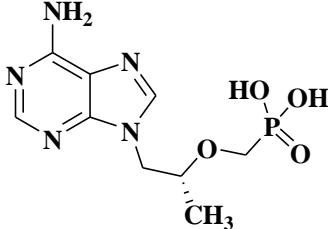
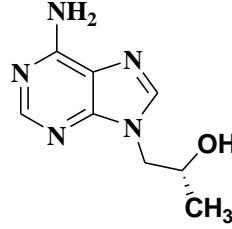
7.	Tadalafil diastereomer (6R, 12aS)	(6R, 12aS)-6-(1,3-Benzodioxol-5-yl)-2,3,6,7,12,12a-hexahydro-2-methyl-pyrazino[1',2':1,6]pyrido[3,4-b]indole-1,4-dione		<b>171596-27-3</b>	
8.	Tadalafil enantiomer (6S, 12aS)	(6S, 12aS)-6-(1,3-Benzodioxol-5-yl)-2,3,6,7,12,12a-hexahydro-2-methyl-pyrazino[1',2':1,6]pyrido[3,4-b]indole-1,4-dione		<b>629652-72-8</b>	
9.	Tadalafil diastereomer (6S, 12aR)	(6S, 12aR)-6-(1,3-Benzodioxol-5-yl)-2,3,6,7,12,12a-hexahydro-2-methyl-pyrazino [1',2':1,6]pyrido[3,4-b]indole-1,4-dione		<b>171596-28-4</b>	

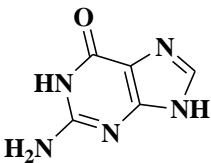
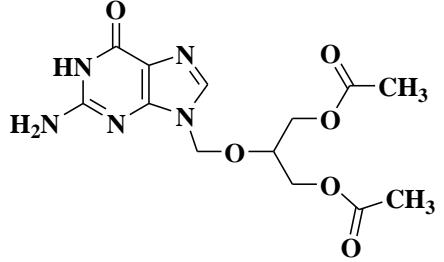
**TENOFOVIR DISOPROXIL FUMARATE**

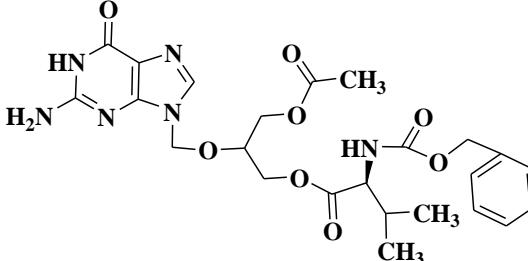
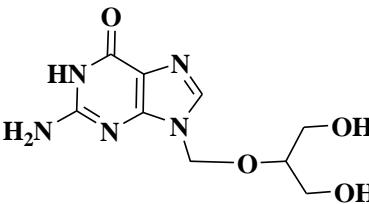
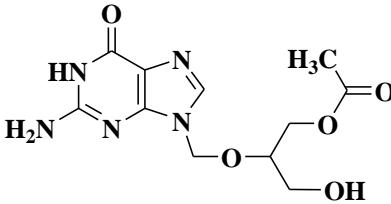
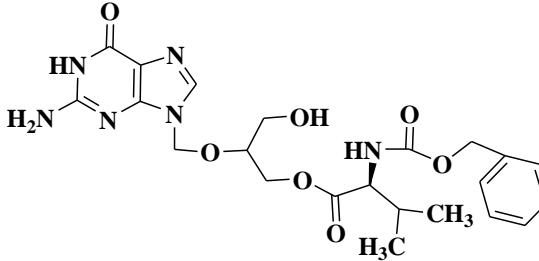
1.	Trifluoroacetic anhydride (TFAA)	Trifluoroacetic anhydride		<b>407-25-0</b>	
2.	Chloromethyl isopropyl carbonate (CMIC)	Chloromethylisopropyl carbonate		<b>35180-01-9</b>	
3.	TEN-I (ISOMER)	(S)-1-(6-Amino-9H-purin-9-yl)propan-2-ol		N/A	
4.	TEN-II (ISOMER)	Ethyl((S)-1-(6-amino-9H-purin-9-yl)propan-2-yloxy)methyl phosphonate		N/A	
5.	TEN-I/Regio isomer	(R)-1-(9H-purin-6-ylamino)propan-2-ol		N/A	
6.	TEN-IA	Diethyl ((R)-1-(6-amino-9H-purin-9-yl)propan-2-yloxy)methyl phosphonate		N/A	

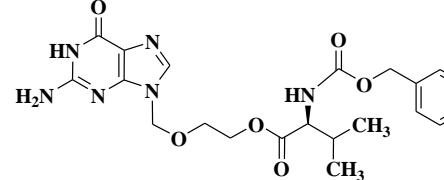
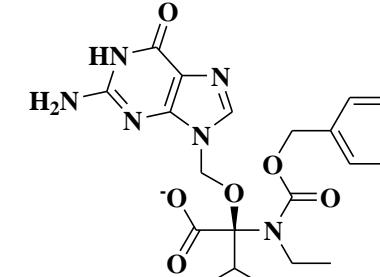
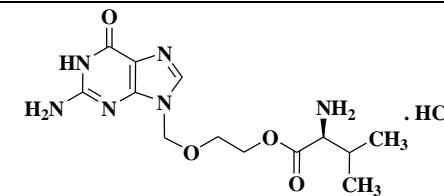
7.	Tenofovir disoproxil S-Isomer	[(1S)-2-(6-Amino-9H-purin-9-yl)- 1-methylethoxy]methyl phosphonic acid disoproxil fumarate		<b>1280130-08-6</b>	
8.	Tenofovir Related Compound -B	(E)-9-(Prop-1-enyl)-9H-adenine (or) (E)-9-(Prop-1-enyl)-9H-purin-6- amine		<b>4121-40-8</b>	
9.	Monoethyl ester	Ethyl hydrogen-(1-(6-amino-9H- purin-9yl)-propan-2-yloxy)methyl phosphonate		N/A	
10.	Tenofovir disoproxil	[(1R)-2(6-Amino-9H-purin-9-yl)- 1-methylethoxy]methyl phosphonic acid Disoproxil		<b>201341-05-1</b>	

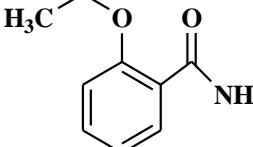
11.	Tenofovir isoproxil monoester	<p>(((R)-1-(6-Amino-9H-purin-9-yl)propan-2-yloxy)methylphosphonoyl) methyl isopropyl carbonate fumarate (or)</p> <p>({{(R)-1-(6-Amino-9H-purin-9-yl)propan-2-yloxy]methyl}(hydroxy) phosphoryloxy)methyl isopropyl carbonate fumarate</p>		N/A	
12.	Tenofovir disoproxil fumarate	<p>(R)-5-[[2-(6-Amino-9H-purin-9-yl)-1-methyl ethoxy]methyl]-2,4,6,8-tetraoxa-5-phosphanona nedioic acid, bis(1-methyl ethyl) ester, 5-oxide, (E)-2-butenedioate (1:1) (or)</p> <p>Bis(hydroxymethyl) [[(R)-2-(6-amino-9H-purin-9-yl)-1-methyl ethoxy] methyl]phosphonate, bis(isopropyl carbonate) (ester), fumarate (1:1) (or)</p> <p>9-[(R)-2-Bis[[isopropoxycarbonyloxy]methoxy] phosphinyl] methoxy]propyl]adenine fumarate (1:1)</p>		202138-50-9	

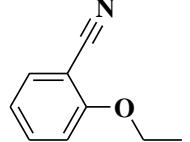
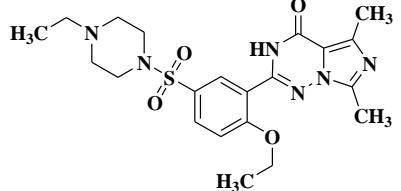
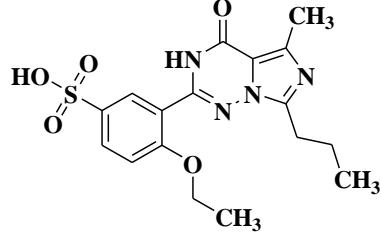
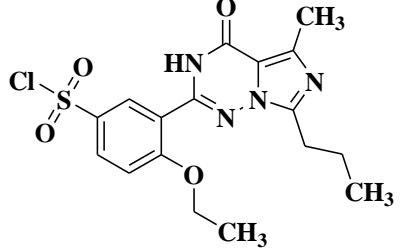
13.	TEN-II	((R)-1-(6-Amino-9H-purin-9-yl)propan-2-yloxy)methylphosphonic acid	 <p>Chemical structure: A purine ring system with an amino group at position 6 and a propan-2-yloxy group at position 9. The propan-2-yloxy group is connected via a phosphonate side chain where the chiral center is labeled R.</p>	<b>147127-20-6</b>	
14.	TEN-I	R-1-(6-Amino-purin-9-yl)-propan-2-ol	 <p>Chemical structure: A purine ring system with an amino group at position 6 and a propan-2-ol group at position 9. The propan-2-ol group is connected via a chiral center where the hydroxyl group is labeled R.</p>	<b>14047-28-0</b>	

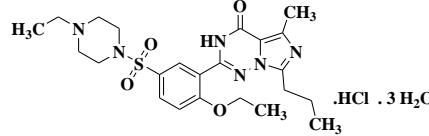
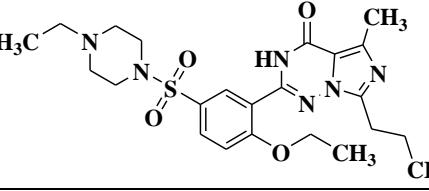
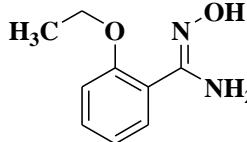
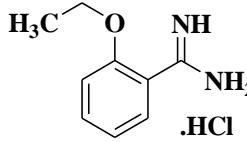
<b>VALGANCICLOVIR</b>					
1.	Guanine	Guanine	 <p>Chemical structure: A purine ring system with an amino group at position 6 and a carbonyl group at position 2.</p>	<b>73-40-5</b>	
2.	VGC-I	2-[(2-Amino-6-oxo-3,6-dihydro-9H-purin-9-yl)methoxy]-1,3-propanediyl diacetate	 <p>Chemical structure: A purine ring system with an amino group at position 6 and a methoxy group at position 2. The methoxy group is connected to a 1,3-propanediyl diacetate side chain.</p>	N/A	

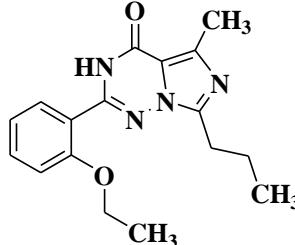
3.	VGC-III A	2-[(2-Amino-1,6-dihydro-6-oxo-1H-purin-9-yl)methoxy]-3-acetoxy propyl-N-(benzyloxy carbonyl)-L-valinate		N/A	
4.	VGC-II	2-Amino-1,9-dihydro-9-[[2-hydroxy-1-(hydroxymethyl)ethoxy]methyl]-6H-purin-6-one (or) 9-[(1,3-dihydroxy-2-propoxy)methyl]guanine		N/A	
5.	VGC-III	2-[(2-Amino-6-oxo-3,6-dihydro-9H-purin-9-yl)methoxy]-3-hydroxypropyl acetate		N/A	
6.	VGC-IV	2-[(2-Amino-6-oxo-1,6-dihydro-9H-purin-9-yl)methoxy]-3-hydroxypropyl N-[(benzyloxy)carbonyl]-L-valinate		N/A	

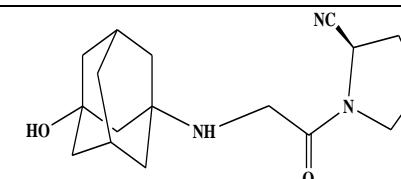
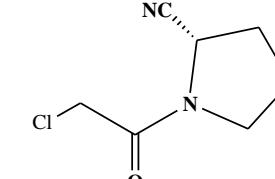
<b>VALACYCLOVIR HYDROCHLORIDE</b>					
1.	VAL-I	2-[(2-Amino-1,6-dihydro-6-oxo-9H-purin-yl)methoxy]ethyl-N-[(benzyloxy)carbonyl]-L-valinate		<b>124832-31-1</b>	
2.	VAL-I (D-Isomer)	2-[(2-Amino-1,6-dihydro-6-oxo-9H-purin-9-yl)methoxy]ethyl-N-[(benzyloxy)carbonyl]-D-valinate		N/A	
3.	Valacyclovir hydrochloride hydrate	L-Valine 2-[(2-amino-1,6-dihydro-6-oxo-9H-purin-9-yl) methoxy] ethyl ester Hydro chloride.		<b>124832-27-5</b>	

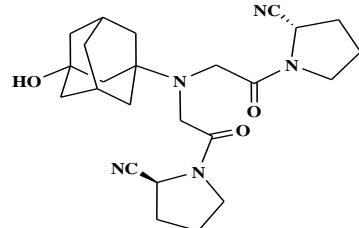
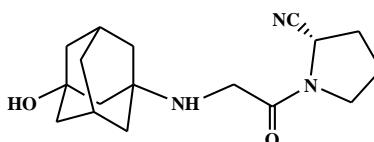
<b>VARDENAFIL HYDROCHLORIDE TRIHYDRATE</b>					
1.	2-Ethoxy benzamide	2-Ethoxybenzamide		<b>938-73-8</b>	

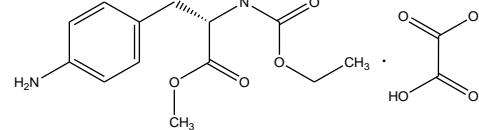
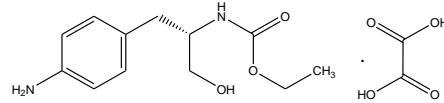
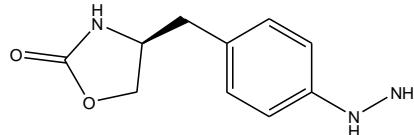
2.	2-Ethoxy benzonitrile	2-Ethoxy benzonitrile		<b>6609-57-0</b>	
3.	Vardenafil impurity-A	2-[2-Ethoxy-5-[(4-ethylpiperazine-1-yl)sulfonyl]phenyl]-5,7-dimethylimidazol[5,1-f][1,2,4]triazin-(3H)-one		N/A	
4.	Vardenafil impurity-B	4-Ethoxy-3-(5-methyl-4-oxo-7-propyl-3,4-dihydroimidazo[5,1-f][1,2,4]triazin-2-yl)benzenesulfonic acid		N/A	
5.	VDF-IV	2-[2-Ethoxy-5-[(4-ethylpiperezin-1-yl)sulfonyl] phenyl]-5-Methyl7propylimidazo[5, 1-f][Trizin-4(3H)-one hydrochloride trihydrate		N/A	

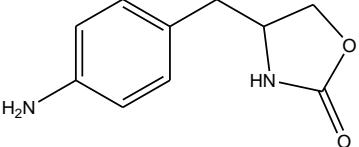
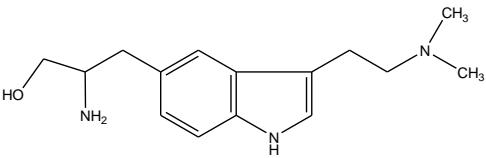
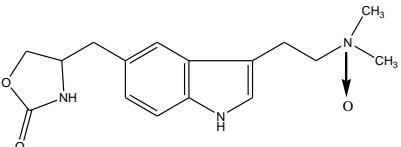
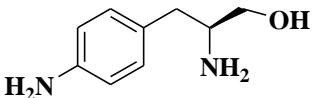
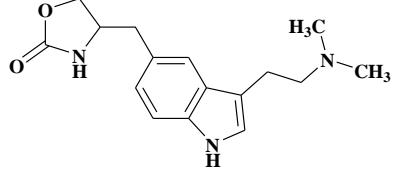
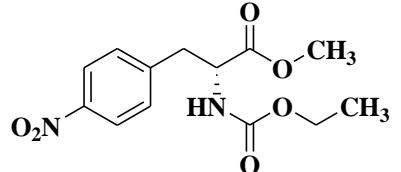
6.	Vardenafil hydrochloride trihydrate	2-[2-Ethoxy-5-[(4-ethylpiperazin-1-yl)sulfonyl]phenyl]-5-methyl-7-propylimidazo[5,1-f][1,2,4]triazin-4(3H)-one hydrochloride trihydrate (or) 1-[[3-(1,4-Dihydro-5-methyl-4-oxo-7-propylimidazo[5,1-f][1,2,4]triazin-2-yl)-4-ethoxyphenyl]sulfonyl]-4-ethyl-Piperazine hydrochloride trihydrate		<b>330808-88-3</b>	
7.	VDF-V	2-[2-Ethoxy-5-(4-ethyl-piperazin-1-sulphonyl)-phenyl]-5-methyl-7-propyl-3H-imadazo [5,1-f] [1,2,4]triazin-4-one		<b>224785-90-4</b>	
8.	VDF-I	2-Ethoxy-N'-hydroxybenzamidine		<b>879-57-2</b>	
9.	VDF-II	2-Ethoxybenzamidine hydrochloride		<b>18637-00-8</b>	

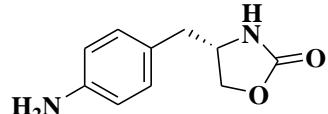
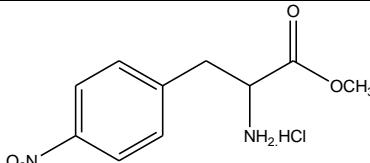
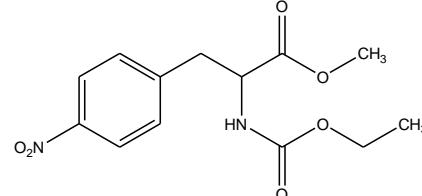
10.	VDF-III	2-(2-ethoxyphenyl)-5-methyl-7-propylimidazo[1,5-f][1,2,4]triazin-4(1H)-one		<b>224789-21-3</b>	
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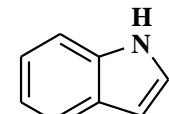
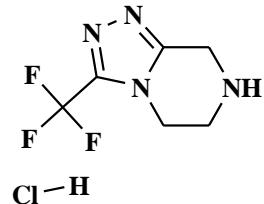
<b>VILDAGLIPTIN</b>					
1.	3-Amino-1-adamantanol	3-amino-1-adamantanol		<b>702-82-9</b>	
2.	Vildagliptin (R-Isomer)	(2R)-1-[(3-Hydroxy tricyclo[3.3.1.13,7]dec-1-yl)amino]acetyl]-2-pyrrolidinecarbonitrile. (or) 1-[[((3-Hydroxy-1-adamantyl) amino) acetyl]-2-cyano-(R)-pyrrolidine.		N/A	500
3.	VLG-III (R)-ISOMER	(R)-1-(2-chloroacetyl)pyrrolidine-2-carbonitrile		<b>565452-98-4</b>	

4.	Vildagliptin dimer impurity	(2S)-1-{{[2[(2S)-2-cyanopyrrolidin-1-yl]-2-oxyethyl][(3-hydroxytricyclo[3.3.1.13,7]dec-1-yl)amino]]acetyl}-pyrrolidine -2-carbonitrile		<b>1036959-23-5</b>	
5.	Vildagliptin	(2S)-1-[(3-Hydroxytricyclo[3.3.1.13,7]dec-1-yl)amino]acetyl]-2-pyrrolidinecarbonitrile (or) 1-[(3-Hydroxy-1-adamantyl)amino]acetyl]-2-cyano-(S)-pyrrolidine		<b>274901-16-5</b>	

<b>ZOLMITRIPTAN</b>					
1.	ZM-IIB	Methyl (S)-N-ethoxycarbonyl-4-aminophenylalaninate oxalate		<b>N/A</b>	
2.	ZM-IIC	(S)-N-Ethoxycarbonyl-4-aminophenylalaninol oxalate		<b>N/A</b>	
3.	ZM-III/N A	(S)-4-(4-Hydrazinobenzyl)-2-oxazolidinone		<b>171550-12-2</b>	

4.	Zolmitriptan impurity-A	(S)-4-(4-Aminobenzyl)-2-oxazolidinone		<b>152305-23-2</b>	
5.	Zolmitriptan impurity-B	2-Amino-3-[3-[2-(dimethylamino)ethyl]-1H-indol-5-yl]propanol		<b>139264-69-0</b>	
6.	Zolmitriptan impurity-C	(4S)-4-[[3-[2-(Dimethylamino)ethyl]-1H-indol-5-yl]methyl]-2-oxazolidinone N-oxide		<b>251451-30-6</b>	
7.	ZM-III/N alcohol impurity	(S)-2-Amino-3-(4-aminophenyl)propan-1-ol		<b>N/A</b>	
8.	Zolmitriptan (R-Isomer)	(4R)-4-[[3-[2-(Dimethylamino)ethyl]-1H-indol-5yl]methyl]-2-oxazolidinone		<b>N/A</b>	
9.	ZM-II/A (R-isomer)	Methyl (R)-N-ethoxycarbonyl-4-nitrophenylalaninate		<b>N/A</b>	

10.	ZM-III/N (R-isomer)	(R)-4-(4-Aminobenzyl)-2-oxazolidinone		<b>262857-90-9</b>	
11.	ZM-II	Methyl-4-nitrophenylalaninate hydrochloride		<b>N/A</b>	
12.	ZM-IIA	Methyl (S)-N-ethoxycarbonyl-4-nitrophenylalaninate		<b>N/A</b>	

13.	INDOLE	1H-Indole		<b>8047-67-4</b>	
14.	Triazolopyrazine	3-(Trifluoromethyl)-5,6,7,8-tetrahydro-[1,2,4]triazolo[4,3-a]pyrazine hydrochloride		<b>762240-92-6</b>	