International Nonproprietary Names for Pharmaceutical Substances (INN)

Recommended International Nonproprietary Names (Rec. INN): List 34

Notice is hereby given that, in accordance with paragraph 7 of the Procedure for the Selection of Recommended International Nonproprietary Names for Pharmaceutical Substances [Off. Rec. Wid Health Org., 1955, 60, 3 (Resolution EB15.R7); 1969, 173, 10 (Resolution EB43.R9)], the following names are selected as Recommended International Nonproprietary Names. The inclusion of a name in the lists of Recommended International Nonproprietary Names does not imply any recommendation of the use of the substance in medicine or pharmacy. Lists of Proposed (1-65) and Recommended (1-31) International Nonproprietary Names can be found in Cumulative List No. 8, 1992.

Dénominations communes internationales des Substances pharmaceutiques (DCI)

Dénominations communes internationales recommandées (DCI Rec): Liste 34

Il est notifié que, conformément aux dispositions du paragraphe 7 de la Procédure à suivre en vue du choix de Dénominations communes internationales recommandées pour les Substances pharmaceutiques (Actes off. Org. mond. Santé, 1955, 60, 3 (résolution EB15.R7); 1969, 173, 10 (résolution EB43.R9)] les dénominations ci-dessous sont mises à l'étude par l'Organisation mondiale de la Santé en tant que dénominations communes internationales proposées. L'inclusion d'une dénomination dans les listes de DCI proposées n'implique aucune recommandation en vue de l'utilisation de la substance correspondante en medecine ou en pharmacie. On trouvera d'autres listes de Dénominations communes internationales proposées (1-65) et recommandées

(1-31) dans la Liste récapitulative No. 8, 1992.

Denominaciones Comunes Internacionales para las Sustancias Farmacéuticas (DCI)

Denominaciones Comunes Internacionales Recomendadas (DCI Rec.): Lista 34

De conformidad con lo que dispone el párrafo 7 del Procedimiento de Selección de Denominaciones Comunes Internacionales Recomendadas para las Sustancias Farmacéuticas [Act. Of. Mund. Salud, 1955, 60, 3 (Resolución EB15.R7); 1969, 173, 10 (Resolución EB43.R9)], se comunica por el presente anuncio que las denominaciones que a continuación se expresan han sido seleccionadas como Denominaciones Comunes Internacionales Recomendadas. La inclusión de una denominación en las listas de las Denominaciones Comunes Recomendadas no supone recomendación alguna en favor del empleo de la sustancia respectiva en medicina o en farmacia,

Las listas de Denominaciones Comunes Internacionales Propuestas (1-65) y Recomendadas (1-31) se encuentran reunidas en Cumulative List No. 8, 1992.

Recommended INN

(Latin, English, French, Spanish)

DCI Recommandée DCI Recomendada

Chemical name or description and Molecular formula

Nom chimique ou description et Formule brute Nombre químico o descripción y Fórmula empírica

abclximabum

immunoglobulin G (human-mouse monoclonal c7E3 clone p7E3V_HhC_{v4} Fab abciximab

fragment anti-human glycoprotein llb/llla receptor), disulfide with human-mouse

monoclonal c7E3 clone p7E3V, hC, light chain

immunoglobuline G (fragment Fab de l'anticorps monoclonal homme-souris c7E3 abciximab

clone p7E3V_HhC_x anti-récepteur de la glycoprotéine IIb/IIIa humaine), ponts disulfure avec la chaîne légère de l'anticorps monoclonal homme-

souris c7E3 clone p7E3V_KhC_K

inmunoglobulina G (fragmento Fab del anticuerpo monoclonal hombre-ratón c7E3 abciximab

cion p7E3V_HhC_M antireceptor de la glicoproteina lib/illa humana), puentes disulturo con la cadena ligera dei anticuerpo monoclonal hombre-ratón c7E3

cion p7E3V_khC_k

acidum incadronicum

[(cycloheptylamino)methylene]diphosphonic acid incadronic acid

acide [(cycloheptylamino)méthylène]bisphosphonique acide incadronique

ácido [(cicloheptilamino)metilen]difosfónico acido incadrónico

C₈H₁₉NO₆P₂

adatanserinum

N-[2-[4-(2-pynmidinyl)-1-piperazinyl]ethyl]-1-adamantanecarboxamide adatanserin

N-[2-[4-(pyrimidin-2-yl)pipérazın-1-yl]éthyl]tricyclo[3.3.1.13,7]décane-1adatansérine

carboxamide

N-[2-[4-(2-pinmidinil)-1-piperazinil]etil]-1-adamantanocarboxamida adatansenna

C21H31N5O

adelmidrolum

N.N'-bis(2-hydroxyethyl)nonanediamide adelmidrol N.N'-bis(2-hydroxyéthyl)nonanediamide adelmidrol

N.N'-bis(2-hidroxietil)nonanodiamida adelmidrol

C13H26N2O4

afovirsenum

2'-deoxy-P-thiocytidylyl-(5'→3')-P-thiothymidylyl-(5'→3')-2'-deoxy-P-thioguanylylafovirsen $(5'\rightarrow 3')-2'-deoxy-P-thiocytidylyl-(5'\rightarrow 3')-P-thiothymidylyl-(5'\rightarrow 3')-2'-deoxy-P-thiothymidylyl-(5'\rightarrow 3')-2'-deoxy-P-thiothymidylyl-(5'\rightarrow 3')-2'-deoxy-P-thiothymidylyl-(5'\rightarrow 3')-2'-deoxy-P-thiothymidylyl-(5'\rightarrow 3')-2'-deoxy-P-thiothymidylyl-(5'\rightarrow 3')-P-thiothymidylyl-(5'\rightarrow 3')-2'-deoxy-P-thiothymidylyl-(5'\rightarrow 3')-P-thiothymidylyl-(5'\rightarrow 3')-2'-deoxy-P-thiothymidylyl-(5'\rightarrow 3')-P-thiothymidylyl-(5'\rightarrow 3')-2'-deoxy-P-thiothymidylyl-(5'\rightarrow 3')-P-thiothymidylyl-(5'\rightarrow 3')-2'-deoxy-P-thiothymidylyl-(5'\rightarrow 3')-P-thiothymidylyl-(5'\rightarrow 3')-P-thiothymidyl-(5'\rightarrow 3')-$

thiocytidylyl- $(5' \rightarrow 3')$ -2'-deoxy-P-thiocytidylyl- $(5' \rightarrow 3')$ -P-thiothymidylyl- $(5' \rightarrow 3')$ -P-thiothymidyl- $(5' \rightarrow$ thiothymidylyl- $(5'\rightarrow 3')$ -2'-deoxy-P-thiocytidylyl- $(5'\rightarrow 3')$ -P-thiothymidylyl- $(5'\rightarrow 3')$ -2'deoxy-P-thioadenylyl- $(5'\rightarrow 3')$ -2'-deoxy-P-thiocytidylyl- $(5'\rightarrow 3')$ -2'-deoxy-P $thio cytidylyl-(5'\rightarrow 3')-P-thio thymidylyl-(5'\rightarrow 3')-P-thio thymidylyl-(5'\rightarrow 3')-2'-deoxy-P-thio thymidylyl-(5'\rightarrow 3')-P-thio thymid$ thiocytidylyl- $(5'\rightarrow 3')$ -2'-deoxy-P-thioguanylyl- $(5'\rightarrow 3')$ -P-thiothymidylyl- $(5'\rightarrow 3')$ -

thymidine

afovirsen	2'-désoxy- P -thiocytidylyl- $(5'\rightarrow 3')$ - P -thiothymidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- P -thioguanylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- P -thiocytidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- P -thiothymidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- P -thiothymidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- $2'$ -thiothymidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- $2'$ -thiocytidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- $2'$ -désoxy- $2'$ -thiocytidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- $2'$ -thiocytidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- $2'$ -thiocytidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- $2'$ -thiothymidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- $2'$ -thiocytidylyl- $(5'\rightarrow 3')$ - $2'$ -désoxy- $2'$ -thiothymidylyl- $2'$ - $2'$ -désoxy- $2'$ -thiothymidylyl- $2'$ - $2'$ -désoxy- $2'$ -thiothymidylyl- $2'$ - $2'$ - $2'$ - $2'$ - $2'$ - $2'$ - $2'$ - $2'$
afovirseno	2'-deoxi- P -tiocitidili-(5' \rightarrow 3')- P -tiotimidili-(5' \rightarrow 3')-2'-deoxi- P -tioguanili-(5' \rightarrow 3')-2'-deoxi- P -tiocitidili-(5' \rightarrow 3')- P -tiotimidili-(5' \rightarrow 3')-2'-deoxi- P -tiocitidili-(5' \rightarrow 3')- P -tiotimidili-(5' \rightarrow 3')- P -tiotimidili-(5' \rightarrow 3')- P -tiotimidili-(5' \rightarrow 3')- P -tiocitidili-(5' \rightarrow 3')- P -tiotimidili-(5' \rightarrow 3')-timidine
	C ₁₉₂ H ₂₅₀ N ₅₇ O ₁₀₇ P ₁₉ S ₁₉
aglepristonum	
aglepristone	11β-[p-(dimethylamino)phenyl]-17β-hydroxy-17-[(Z)-propenyl]estra-4,9-dien-3-one
aglépristone	11β-[4-(diméthylamino)phényl]-17β-hydroxy-17-[(Z)-prop-1-ényl]estra-4,9-dién- 3-one
aglepristona	11 β -[ρ -(dimetilamino)fenil]-17 β -hidroxi-17-[(Z)-propenil]estra-4,9-dien-3-ona
	C ₂₉ H ₃₇ NO ₂
alnespironum	
alnespirone	$(+)\cdot(S)\cdot N\cdot[4\cdot[(5\cdot methoxy-3\cdot chromanyl)propylamino]butyl]-1,1\cdot cyclopentanediacetimide$
alnespirone	(+)-(S)-8-[4-[(5-méthoxy-3,4-dihydro-2 <i>H-c</i> hromén-3-yl)(propyl)amino]butyl]-8-azaspiro[4.5]décane-7,9-dione
atnespirona	(+)-(S)-N-[4-[(5-metoxi-3-cromanil)propilamino]butil]-1,1-ciclopentanodiacetimida
	C ₂₆ H ₃₆ N ₂ O ₄
alvirceptum sudotoxum	
alvircept sudotox	N ² -L-methionyl-1-178-antigen CD 4 (human clone pT4B protein moiety reduced) (178→248)-protein with 248-L-histidine-249-L-methionine-250-L-alanine-251-L-glutamic acid-248-613-exotoxin A (<i>Pseudomonas aeruginosa</i> reduced)
alvircept sudotox	N²-L-méthionyl-1-178-antigène CD 4 (partie protéique réduite de la substance issue du clone humain pT4B) (178→248′)-protéine avec la 248-L-histidine-249-L-méthionine-250-L-alanine-251-acide L-glutamique-248-613-exotoxine A (Pseudomonas aeruginosa) réduite
alvircept sudotox	N²-L-metionil-1-178-antigeno CD 4 (fracción proteica reducida de la sustancia obtenida del cion humano pT4B) (178→248′)-proteina con 248-L-histidina-249-L-metionina-250-L-alanina-251-L-ácido glutamico-248-613-exotoxina A

(Pseudomonas aeruginosa reducida)

C2500H4130N748OB12S10

arenidipinum

aranidipine (±)-acetonyl methyl 1,4-dihydro-2,6-dimethyl-4-(o-nitrophenyl)-3,5-pyridinedi=

carboxylate

aranidipine (RS)-2,6-diméthyl-4-(2-nitrophényl)-1,4-dihydropyridine-3,5-dicarboxylate de

méthyle et de 2-oxopropyle

aranidipino (±)-acetonil metil 1,4-dihidro-2,6-dimetil-4-(o-nitrofenil)-3,5-piridindicarboxılato

C₁₉H₂₀N₂O₇

arteflenum

artefiene (1.S.4.R.5.R.8.S)-4-[(Z)-2,4-bis(trifluoromethyl)styryl]-4,8-dimethyl-

2,3-dioxabicyclo[3.3.1]nonan-7-one

artéfiène (1S,4R,5R,8S)-4-[(Z)-2-[2,4-bis(trifiuorométhyl)phényl]-4,8-diméthyl-

2,3-dioxabicyclo[3.3.1]nonan-7-one

artefieno (1S,4R,5R,8S)-4-[(Z)-2,4-bis(trifluorometil)estiril]-4,8-dimetil-

2,3-dioxabiciclo[3.3.1]nonan-7-ona

C19H18F6O3

atevirdinum

atevirdine 1-[3-(ethylamino)-2-pyridyl]-4-[(5-methoxyındol-2-yl)carbonyl]piperazine

atévirdine 1-[3-(éthylamino)pyridin-2-yl]-4-[(5-méthoxy-1*H*-indol-2-yl)carbonyl]pipérazine

atevirdına 1-[3-(etilamino)-2-pırıdil]-4-[(5-metoxiindol-2-il)carbonil]piperazina

C21H25N5O2

azelnidipinum

azelnidípine 3-[1-(diphenylmethy/)-3-azetidínyl] 5-isopropyl (±)-2-amino-1,4-dihydro-6-methyl-

4-(m-nitrophenyl)-3,5-pyridinedicarboxylate

azelnidipine (RS)-2-amino-6-méthyl-4-(3-nitrophényl)-1,4-dıhydropyridine-3,5-dicarboxylate de

3-[1-(diphénylméthyl)azétidtn-3-yle] et de 5-(1-méthyléthyle)

azelnidipıno 3-[1-(dıfenilmetil)-3-azetidinil] 5-isopropil (±)-2-amino-1,4-dihidro-6-metil-4-(m-

nitrofenil)-3,5-piridindicarboxilato

C33H34N4O6

batimastatum

batimastat (2S,3R)-5-methyl-3-[[(αS)- α -(methylcarbamoyl)phenethyl]carbamoyl]-2-

[(2-thienylthio)methyl)hexanohydroxamic acid

batımastat (2S,3R)·N'-hydroxy-N⁴-[(S)-1-[(méthylamino)carbonyl]-2-phényléthyl]-

3-(2-méthylpropyl)-2-[(2-thiénylthio)méthyl]butanediamide

batimastat ácido (2S,3R)-5-metil-3- $[(\alpha S)$ - α -(metilcarbamoil)fenetil]carbamoil]-2-

[(2-tieniltio)metil]hexanohidroxamico

C23H31N3O4S2

beciparcilum

beciparcil p-[(5-thio- β -p-xylopyranosyl)thio]benzonitrile beciparcil 4-[(5-thio- β -p-xylopyranosyl)thio]benzonitrile

beciparcilo p-[(5-tio-β-p-xilopiranosii)tio]benzonitrilo

C₁₂H₁₃NO₃S₂

besipirdinum

1-(propyi-4-pyndylamino)indole besipirdine

(1 H-indol-1-yl)(propyl)(pyridin-4-yl)amıne bésipirdine

1-(propil-4-piridilamino)indol besipirdina

C16H17N3

biapenemum

6-[[(4R,5S,6S)-2-carboxy-6-[(1R)-1-hydroxyethyl]-4-methyl-7-oxo-1-aza= biapenem

bicyclo[3.2.0]hept-2-en-3-yl]thio]-6,7-dihydro-5H-pyrazolo[1,2-a]-s-triazol-4-ium

hydroxide, inner salt

6-[[(4R,5S,6S)-2-carboxylato-6-[(1R)-1-hydroxyéthyl]-4-méthyl-7-oxo-1-aza= biapénem

bicyclo[3.2.0]hept-2-en-3-yl]thio]-6,7-dihydro-5H-pyrazolo[1,2-a][1,2,4]tnazol-4-ium

6-[[(4R,5S,6S)-2-carboxi-6-[(1R)-1-hidroxietil]-4-metil-7-oxo-1-azabiciclo= biapenem

[3.2.0]hept-2-en-3-il]tio]-6,7-dihidro-5H-pirazolo[1,2-a]-s-triazol-4-io hidroxido, sal

interna

C15H18N4O4S

bicatutamidum

(\pm)-4'-cyano- α , α , α -trifluoro-3-[(p-fluorophenyl)sulfonyl]-2-methyl-m-lactotoluidide bicalutamide

(RS)-N-[4-cyano-3-(trifluorométhyl)phényl]-3-[(4-fluorophényl)sulfonyl]-2-hydroxybicalutamide

2-méthylpropanamide

(\pm)-4'-ciano- α , α , α -trifiuoro-3-[(p-fluorofenil)sulfonil]-2-metil-m-lactotoluidida bicalutamida

C18H14F4N2O4S

bosentanum

p-tert-butyl-N-[6-(2-hydroxyethoxy)-5-(o-methoxyphenoxy)-2-(2-pyrimidinyl)bosentan

4-pyrimidinyl]benzenesulfonamide

4-(1,1-diméthyléthyl)-N-[6-(2-hydroxyéthoxy)-5-(2-méthoxyphénoxy)bosentan

2-(pyrimidin-2-yl)pyrimidin-4-yl)benzenesulfonamide

p-terc-butil-N-[6-(2-hidroxietoxi)-5-(p-metoxifenoxi)-2-(2-pirimidinil)bosentano

4-pinmidinil]bencensulfonamida

C27H29N5O6S

candocuronii lodidum

17a,17a-dimethyl-3β-(1-methylpyrrolidinio)-17a-azonia-D-homoandrost-5-ene candocuronium jodide

diiodide

diiodure de 17a,17a-dıméthyl-3β-(1-méthylpyrrolıdınio)-17a-azonıaiodure de candocuronium

D-homoandrost-5-ène

17a,17a-dimetil-3β-(1-metilpirrolidinio)-17a-azonia-D-homoandrost-5-eno diioduro ioduro de candocuronio

C26H46I2N2

capromabum

immunoglobulin G 1 (mouse monoclonal 7E11-C5.3 anti-human prostatic capromab

carcinoma ceil), disulfide with mouse monocional 7E11-C5.3 light chain, dimer

immunoglobuline G 1 (anticorps monoclonal de souris 7E11-C5.3 anti-cellules de capromab

carcinome prostatique humain), dimère du disulture avec la chaîne légère de

l'anticorps monoclonal de souris 7E11-5.3

capromab

inmunoglobulina G1 (anticuerpo monoclonal 7E11-C5.3 de ratón anticélulas de carcinoma prostático humano), puentes disulfuro con la cadena ligera del

anticuerpo monocional 7E11-C5.3 de ratón, dimero

carvotrolinum

carvotroline 8-fluoro-2,3,4,5-tetrahydro-2-[2-(4-pyridyl)ethyl]-1*H*-pyrido[4,3-*b*] indole

carvotroline 8-fluoro-2-[2-(pyridin-4-yl)éthyl]-2,3,4,5-tétrahydro-1 H-pyrido[4,3-b]indole

carvotrolina 8-fluoro-2,3,4,5-tetrahidro-2-[2-(4-pindil)etil]-1*H*-pindo(4,3-b]indol

C₁₈H₁₈FN₃

cedefingolum

cedefingol N-[(1S,2S)-2-hydroxy-1-(hydroxymethyl)heptadecyl]acetamide

cédéfingol N-[(15,25)-2-hydroxy-1-(hydroxyméthyl)heptadécyl]acétamide

cedefingol N-[(15,25)-2-hidroxi-1-(hidroximetil)heptadecii]acetamida

C20H41NO3

cetcapenum

cefcapene (6R,7R)-7-[(Z)-2-(2-amino-4-thiazolyl)-2-pentenamido]-3-(hydroxymethyl)-8-oxo-

5-thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylic acid, carbamate (ester)

cefcapène acide (+)-(6R,7R)-7-[[(Z)-2-(2-aminothiazol-4-yf)pent-2-énoyt]amino]-

3-[[(aminocarbonyl)oxy]méthyl]-8-oxo-5-thia-1-azabicyclo[4.2.0]oct-2-ène-2-

carboxylique

cefcapeno ácido (+)-hidroximetii (6*R*,7*R*)-7-[(*Z*)-2-(2-amino-4-tiazolii)-2-pentenamido]-3-

(hidroximetil)-8-oxo-5-tia-1-azabicicio[4.2.0]oct-2-en-2-carboxílico

C17H19N5O6S2

certoparinum natricum

certoparin sodium

Sodium salt of depolymerized heparin obtained by isoamyl nitrite degradation of heparin from pork intestinal mucosa; the majority of the components have a 2-O-sulfo- α -L-idopyranosuronic acid structure at the non reducing end and a 6-O-sulfo-2,5-anhydro-p-mannose structure at the reducing end of their chain, the average relative molecular mass is 5000 to 7000; at least 70 per cent less than 10 000; the

degree of sulfatation is 2 to 2,5 per disaccharidic unit.

certoparine sodique

sel de sodium d'héparine dépolymérisée obtenue par fragmentation au moyen de nitrite d'isoamyle d'héparine de muqueuse intestinale de porc. La majorité des composants présentent une structure acide 2-O-sulfo-α-L-idopyranosuronique à l'extrémité non réductrice et une structure 6-O-sulfo-2,5-anhydro-p-mannose à l'extrémité réductrice de leur chaîne. La masse moléculaire relative moyenne est de 5000 à 7000, 70 pour cent au moins des composants ayant une masse moléculaire relative inférieure à 10 000. Le degré de sulfatation est de 2 à 2,5 par

unité disaccharidique.

certoparina sódica

Sal sódica de la heparina despolimerizada obtenida por fragmentación con nitrito de isoamilo de la heparina de la mucosa intestinal del cerdo; la mayoría de los compuestos tienen una estructura de ácido 2-*O*-sulfo-α-L-idopirano-surónico en el extremo no reductor y una estructura de 6-*O*-sulfo-2,5-anhidro-p-manitol en el extremo reductor de la cadena; la masa molecular relativa media es 5000 a 7000, al menos el 70% es menor de 10 000; el grado de sulfatación es de 2 a 2,5 por

unidad de disacárido.

cinalukastum

cinalukast

3'-[(E)-2-(4-cyclobutyl-2-thiazolyl)vinyl]-2,2-diethylsuccinanilic acid acide (E)-4-[[3-[2-(4-cyclobutylthiazol-2-yl)éthényl]phényl]amino]-2,2-diéthylcinalukast

4-oxobutanoïque

ácido 3'-[(E)-2-(4-ciclobutil-2-tiazolil)vinil]-2,2-dietilsuccinanilico cinalukast

C23H28N2O3S

ciprokirenum

 (αS) -N-[(1 S,2R,3S)-1-(cyclohexylmethyl)-3-cyclopropyl-2,3-dihydroxypropyl]- α ciprokiren

 $[(\alpha S)-\alpha \cdot [[1-methyl-1-(morpholinocarbonyl)ethyl]sulfonyl]methyl]hydrocinnam=$

amido)imidazole-4-propionamide

(S)-N-[(1S,2R,3S)-1-(cyclohexylméthyl)-3-cyclopropyl-2,3-dihydroxypropyl]-2ciprokirène

[[(5)-2-[[[1-methyl-1-[(morpholin-4-yl)carbonyl]éthyl]sulfonyl]méthyl]-3-

phénylpropanoyl]amino]-3-(1H·midazol-4-yl)propanamide

 (αS) - $\mathcal{N}[(1S,2R,3S)$ -1-(ciclohexilmetil)-3-ciclopropil-2,3-dihidroxipropil]- α - $[(\alpha S)$ ciprokireno

α-[[[1-metil-1-(morfolinocarbonil)etil]sulfonil]metil]hidrocinnamamido]imidazol-4-

propionamida

C37H55N5OaS

dapabutanum

(±)-3-[[3-(dodecylamino)propyllamino]butyric acid dapabutan

acide (RS)-3-[[3-(dodécylamino)propyl]amino]butanoïque dapabutan

ácido (±)-3-[[3-(dodecilamino)propil]amino]butírico dapabutano

C19H40N2O2

darqlitazonum

 (\pm) -5-[p-[3-(5-methyl-2-phenyl-4-oxazolyl)propionyl]benzyl]-2,4-thiazolidinedione darolitazone

(RS)-5-[4-[3-(5-méthyl-2-phényloxazol-4-yl)propanoyl]benzyl]thiazolidine-2,4-dione darglitazone

 (\pm) -5-[ρ -[3-(5-metil-2-fenil-4-oxazolil)propionil]bencil]-2,4-tiazolidindiona dargiitazona

C23H20N2O4S

darifenacinum

 $(S)\text{-}1\text{-}[2\text{-}(2,3\text{-}dihydro\text{-}5\text{-}benzofuranyl})\text{ethyl}]\text{-}\alpha, \alpha\text{-}diphenyl\text{-}3\text{-}pyrrolidineacetamide}$ darifenacin

(S)-2-[1-[2-(2,3-dihydrobenzofurɛn-5-yl)éthyl]pyrrolidin-3-yl]-2,2-diphénylacétamide darifénacine

(S)-1-[2-(2,3-dihidro-5-benzofuranii)etii]- α , α -difenii-3-pirrolidinacetamida darifenacina

C28H30N2O2

destrudinum

63-desulfohirudin (Hirudo medicinalis isoform HV1) desirudin

63-désulfohirudine (Hirudo medicinalis, isoform HV1) déstrudine

63-desulfohirudina (isoforma HV1 de Hirudo medicinalis) desirudina

C287H440N80O110S6

desmeninolum

desmeninol (±)-2-hydroxy-4-(methylthio)butyric acid

desméninol acide (RS)-2-hydroxy-4-(méthylthio)butanoïque

desmeninol ácido (±)-2-hidroxi-4-(metiltio)butírico

C5H10O3S

detumomabum

detumomab immunoglobulin (mouse monoclonal SPECIFID anti-human B lymphoma cell)

disulfide with mouse monoclonal SPECIFID light chain, dimer

détumomab immunoglobuline (anticorps monoclonal de souris SPECIFID anticellules de

lymphome B humain), dimère du disulfure avec la chaîne légère de l'anticorps

monoclonal de souris SPECIFID

detumomab inmunoglobulina (anticuerpo monoclonal SPECIFID de ratón anticélulas de

lintoma B humano), puentes disulfuro con la cadena ligera del anticuerpo

monoclonal SPECIFID de ratón, dimero

dexketoprofenum

dexketoprofen (+)-(S)-m-benzoylhydratropic acid

dexkétoprofène acide (+)-(S)-2-(3-benzoylphényl)propanoique

dexketoprofeno ácido (+)-(S)-m-benzoılhidratropico

C₁₆H₁₄O₃

dornasum alfa

dornase alfa deoxyribonuclease (human clone 18-1 protein moiety)

domase alfa désoxyribonucléase (partie protéique de la substance issue du clone humain 18-1)

domasa alfa desoxirribonucleasa (clon humano 18-1 fracción proteica)

C₁₃₂₁H₁₉₉₅N₃₃₉O₃₉₆S₉

edobacomabum

edobacomab immunoglobulin M (mouse monoclonal XMMEN-0E5 anti-endotoxin), disulfide

with mouse monoclonal XMMEN-0F5 light chain, pentameric dimer

édobacomab immunoglobuline M monoclonale de souris XMMEN-0E5 dirigée contre le

domaine lipidique A d'endotoxines de bactéries gram-négatives

edobacomab inmunoglobulina M monoclonal de ratón XMMEN-0E5 anti-endotoxina, unida

mediante enlace disulfuro con la cadena ligera del anticuerpo monoclonal de

ration XMMEN-OE5, dimero pentamérico

elopiprazolum

elopiprazole 1-(7-benzofuranyl)-4-[[5-(p-fluorophenyl)pyrrol-2-yl]methyl]piperazine

élopiprazole 1-(benzofuran-7-yl)-4-[[5-(4-fluorophényl)-1*H*-pyrrol-2-yl]met: ¿[pipérazine

elopiprazol 1-(7-benzofuranii)-4-[[5-(p-fluorofenii)pirrol-2-ii]metii]piperazina

C₂₃H₂₂FN₃O

emideltidum

emideltide L-tryptophyl-L-alanylglycylglycyl-L-α-aspartyl-L-alanyl-L-serylglycyl-L-glutamic acid

émideltide L-tryptophyi-L-alanyl-glycyl-glycyl-μα-aspartyl-L-alanyl-L-séryl-glycyl-

acide L-glutamique

emideltida ácido ι-triptofil-ι-alanılglicilglicil-ι-α-aspartil-ι-alanıl-ι-serilglicil-ι-glutamico

C35H48N10O15

enlimomabum

enlimomab immunoglobulin G 2a (mouse monoclonal BI-RR-1 anti-human-antigen CD 54),

disulfide with mouse monoclonal BI-RR-1 light chain, dimer

enlimomab immunoglobuline G 2a (anticorps monoclonal de souris BI-RR-1 anti-antigène

CD 54 humain), dimère du disulfure avec la chaîne légère de l'anticorps mono-

clonal de souris BI-RR-1

entimomab inmunoglobulina G2a (anticuerpo monoclonal BI-RR-1 de ratón anti-antigeno

CD 54 humano), puentes disulturo con la cadena ligera del anticuerpo mono-

cional BI-RR-1 de ratón

epristeridum

epristeride 17β-(tert-butylcarbamoyl)androsta-3,5-diene-3-carboxylic acid

epristende acide 17β-[[(1,1-diméthyléthyl)amino]carbonyl]androsta-3,5-diène-3-carboxylique

epristerida ácido 17β-(terc-butılcarbamoil)androsta-3,5-dien-3-carboxílico

C25H37NO3

fananserinum

fananserin 2-[3-[4-(p-fluorophenyl)-1-piperazinyl]propyl]-2H-naphth[1,8-cd]isothiazole

1.1-dioxide

fanansérine 2-[3-[4-(4-fluorophényl)pipérazin-1-yl]propyl]-2H-naphto[1,8-cd]isothiazole

1,1-dioxyde

tanansenna 2-[3-[4-(p-fluorofenil)-1-piperazinil)]propil]-2H-naft[1,8-co]isotiazol 1,1-dióxido

C23H24FN3O2S

ferpifosatum natricum

ferpifosate sodium. hexasodium tris[(4,5-dihydroxy-6-methyl-3-pyridinemethanol 3-phosphato)(3-)-

 O^3 , O^5]ferrate(6-)

ferpifosate sodique tris[[(4-hydroxy-6-méthyl-5-olatopyridin-3-yl)méthanol 3-phosphato](3)-

O³,O³,O⁵]ferrate(6-) d'hexasodium

ferpifosato sodico tris[(4,5-dihidroxi-6-metil-3-piridinometanol 3-fosfato)(3-)-O3,O3,O5]ferrate(6-) de

hexasodio

C₂₁H₂₁FeNa₆N₃O₁₈P₃

fosopaminum

fosopamine 4-[2-(methylamino)ethyl]pyrocatechol 1-(dihydrogen phosphate)

fosopamine dihydrogénophosphate de 2-hydroxy-4-[2-(méthylamino)éthyl]phényle

fosopamina 1-dihidrogeno fosfato de 4-[2-(metilamino)etil]pirocatecol

C₉H₁₄NO₅P

geclosporinum

geclosporin cyclo[[(2S,3R,4R,6E)-3-hydroxy-4-methyl-2-(methylamino)-6-octenoyi]-L-

norvalyl-N-methylglycyl-N-methyl-L-leucyl-L-valyl-N-methyl-L-leucyl-L-alanyl-D-

alanyl-N-methyl-L-leucyl-N-methyl-L-leucyl-N-methyl-L-valyl]

géclosporine cyclo[-[(6E)-(2S,3R,4R)-3-hydroxy-4-méthyl-2-(méthylamino)oct-6-énoyl]-L-

norvalyi-(N-méthyigyicyl)-(N-méthyi-L-leucyl)-L-valyi-(N-méthyi-L-leucyl)-L-alanyi-o-alanyi-(N-méthyi-L-leucyl)-(N-méthyi-L-leucyl)-(N-méthyi-L-valyl)-)

geclosporina ciclo[[(2S,3R,4R,6E)-3-hidroxi-4-metil-2-(metilamino)-6-octenoil]-L-norvalil-N-

metilglicil-//-metil-L-leucil-L-valil-//-metil-L-leucil-L-alanil-p-alanil-//-metil-L-leucil-//-

metil-L-leucil-N-metil-L-valil)

C63H113N11O12

glenvastatinum

glenvastatin (4R,6S)-6-[(E)-2-[4-(p-fluorophenyl)-2-isopropyl-6-phenyl-3-pyridyl]vinyl]tetra=

hydro-4-hydroxy-2H-pyran-2-one

glenvastatine (4R,6S)-6-[(E)-2-[4-(4-fluorophényl)-2-(1-méthyléthyl)-6-phénylpyridin-

3-yl]éthényl]-4-hydroxytétrahydro-2H-pyran-2-one

glenvastatina (4R,6S)-6-[(E)-2-[4-(p-fluorofenil)-2-isopropil-6-fenil-3-piridil]vinil]tetrahidro-

4-hidroxi-2H-piran-2-ona

C27H26FNO3

Icometasonii enbutas

icometasone enbutate 9-chloro-11β,17,21-trihydroxy-16α-methylpregna-1,4-diene-3,20-dione

17-butyrate 21-acetate

icométasone enbutate 21-acétate 17-butanoate de 9-chloro-11β,17,21-trihydroxy-16α-méthylprégna-

1,4-diène-3,20-dione

enbutato de icometasona 9-cloro-11β,17,21-trihidroxi-16α-metilpregna-1,4-dieno-3,20-diona 17-butirato

21-acetato

C26H37CIO7

iganidipinum

iganidipine (±)-3-(4-allyl-1-piperazinyl)-2,2-dimethylpropyl methyl 1,4-dihydro-2,6-dimethyl-

4-(m-nitrophenyl)-3,5-pyridinedicarboxylate

iganidipine (RS)-2,6-diméthyl-4-(3-nitrophényl)-1,4-dihydropyridine-3,5-dicarboxylate de

2,2-diméthyl-3-[4-(prop-2-ényl)piperazin-1-yl]propyle et de méthyle

iganidipino (±)-3-(4-alil-1-piperazinil)-2,2-dimetilpropil metil 1,4-dihidro-2,6-dimetil-4-

(m-nitrofenil)-3,5-piridindicarboxilato

C28H38N4O6

Hepcimidum

ilepcimide 1-[(E)-3,4-(methylenedioxy)cinnamoyl]piperidine

ileparmide 1-[(E)-3-(1,3-benzodioxof-5-yf)prop-2-énoyf)pipéridine

ilepcimida 1-{(E)-3,4-(metilendioxi)cinnamoil]piperidina

C₁₅H₁₇NO₃

72.

lliparcllum

iliparcil 4-ethyl-7-[(5-thio-β-p-xylopyranosyl)oxy)coumarin

iliparcil 4-éthyl-7-[(5-thio-β-b-xylopyranosyl)oxy]-2H-chromén-2-one

ıliparcilo 4-etil-7-[(5-tio-β-o-xilopiranosil)oxijcumarina

C16H18O6S

ilonidapum

ilonidap 6-chloro-5-fluoro-3-[(Z)-α-hydroxy-2-thenylidene]-2-oxo-1-indolinecariboxamide

ilonidap (Z)-6-chloro-5-fluoro-3-[hydroxy(2-thiényl)méthylène]-2-oxo-2,3-dihydrò-1 H-

indole-1-carboxamide

ilonidap 6-cloro-5-fluoro-3-[(Z)-α-hidroxi-2-tienilidene]-2-oxo-1-indolincarboxamida

C₁₄H₈CIFN₂O₃S

iloperidonum

iloperidone 4'-[3-[4-(6-fluoro-1,2-benzisoxazol-3-yl)piperidino]propoxy]-3'-methoxyaceto=

phenone

ilopéridone 1-[4-[[3-[4-(6-fluoro-1,2-benzisoxazol-3-yl)pipéridin-1-yl]propyl]oxy]-3-méthoxy=

phényl]éthanone

iloperidona 4'-[3-[4-(6-fluoro-1,2-bencisoxazol-3-il)piperidino]propoxi]-3'-metoxiacetofenona

C24H27FN2O4

imitrodastum

imitrodast 4,5-dihydro-2-(imidazol-1-ylmethyl)benzo[b]thiophene-6-carboxylic acid

imitrodast acide 2-[(1H-imidazol-1-yl)méthyl]-4,5-dihydrobenzo[b]thìophène-6-carboxylique

imitrodast ácido 4,5-dihidro-2-(imidazol-1-ilmetil)benzo[b]tiofeno-6-carboxílico

C₁₃H₁₂N₂O₂S

iralukastum

iralukast 7-[[(1S,2E,4Z)-9-(4-acetyl-3-hydroxy-2-propylphanoxy)-1-[(αF)- α -hydroxy-

m-(trifluoromethyl)benzyl]-2,4-nonadienyl]thio]-4-oxo-4H-1-benzopyran-2-

carboxylic acid

iralukast acide 7-[[(2E,4Z)-(1S)-9-(4-acétyl-3-hydroxy-2-propylphénoxy)-1-[(R)-hydroxy=

[3-(trifluorométhyl)phényl]methyl]ncna-2,4-diényl]thio]-4-oxo-4H-chromène-2-

carboxylique

iralukast ácido 7-[[(1S,2E,4Z)-9-(4-acetil-3-hidroxi-2-propilfenoxi)-1-[(αR)- α -hidroxi-

m-(trifluorometil)bencil]-2,4-nonadienil]tio]-4-oxo-4H-1-benzopiran-2-carboxílico

C38H37F3O8S

laflunimusum

laflunimus (Z)-α-cyano-α⁴',α⁴',α⁴'-trifluoro-β-hydroxycyclopropaneacrylo-3',4'-xylidide'

laflunimus (Z)-2-cyano-3-cyclopropyl-3-hydroxy-N-[3-méthyl-4-(trifluorométhyl)phényl]prop-

2-énamide

laflunimus (Z)-α-ciano-α⁴',α⁴'-trifluoro-β-hidroxiciclopropanacrilo-3',4'-xylidida

C₁₅H₁₃F₃N₂O₂

Recommended INN: List 34

latutidinum

lafutidine (±)-2-(furfurylsulfinyl)-N-[(Z)-4-[[4-(piperidinomethyl)-2-pyridyl]oxy]-2-butenyl]=

acetamide

 $|afutudine (\pm)-2-[(2-furylméthyl)sulfinyl]-N-[(Z)-4-[[4-(pipéridin-1-ylméthyl)pyridin-2-yl]oxy]=$

but-2-ényl]acétamide

lafutidina (±)-2-(furfurilsulfinit)-N-[(Z)-4-[[4-(piperidinometil)-2-piridif]oxi]-2-butenil]acetamıda

C22H29N3O4S

laurcetii bromidum

laurcetium bromide (carboxymethyl)dodecyldimethylammonium bromide, ethyl ester

bromure de laurcétium bromure de dodécyl[(éthoxycarbonyl)méthyl]diméthylammonium

bromuro de (aurcetio ester etílico del bromuro de (carboximetil)dodecildimetilamonio

C₁₈H₃₈BrNO₂

lecimibidum

lecimibide 3-{2,4-difluorophenyl}-1-[5-[(4,5-diphenylimidazol-2-yl)thio]pentyl]-1-heptylurea

lécimibide 3-(2,4-difluorophényl)-1-[5-[(4,5-diphényl-1*H*-imidazol-2-yl)thio]pentyl]-1-

heptylurée

lecimibida 3-(2,4-difluorofenil)-1-[5-[(4,5-difenilimidazol-2-il)tio]pentil]-1-heptilurea

C34H40F2N4OS

ledismasum

ledismase superoxide dismutase (human copper-zinc subunit), cyclic (57→146)-disulfide,

dimer

lédismase superoxyde dismutase humaine (dimère de deux sous-unités comportant chacune

un ion cuivre et un ion zinc et une liaison (57→146)-disulfure cyclique

ledismasa superoxido dismutasa (subunidad cobre-zinc humana), disulfuro cíclico

(57→146), dimero

C679H1083N203O224S4

lemildipinum

lemildipine 3-isopropyl 5-methyl (±)-4-(2,3-dichlorophenyl)-1,4-dihydro-2-(hydroxymethyl)-

6-methyl-3,5-pyridinedicarboxylate, carbamate (ester)

lémildipine (RS)-2-[[(aminocarbonyl)oxy]méthyl]-4-(2,3-dichlorophényl)-6-méthyl-1,4-dihydro=

pyridine-3,5-dicarboxylate de 5-méthyle et de 3-(1-méthyléthyle)

lemildipino carbamato de 3-isopropil 5-metil (±)-4-(2,3-dıclorofenil)-1,4-dihidro-2-

(hidroximetil)-6-metil-3,5-piridindicarboxilato

C20H22Cl2N2O6

lemoxinolum

lemoxinol α -(4,6-dichloro-*m*-tolyl)oxy- ω -hydroxypoly(oxyethylene)

Each *lemoxinol* name is followed by a number indicating the approximate number of oxyethylene groups present, e.g. *lemoxinol* 5, and the individual chemical names may contain a specific numerical syllable for the same purpose.

lémoxinol α -(2,4-dichloro-5-méthylphényl)- ω -hydroxypoly(oxyéthylène)

Chaque lémoxinol est suivi d'un nombre indiquant le nombre approximatif de groupe oxyéthylène présents (lémoxinol 5) et les noms chimiques individuels

peuvent contenir une syllabe numérique ayant la même signification.

lemoxino! α-[(4,6-dicloro-m-tolii)oxi]-ω-hidroxi poli(oxietileno)

Cada denominación de lemoxinol va seguida de un número que indica el número aproximado de grupos oxietileno presentes; p.ej., lemoxinol 5; la denominación química individual puede contener una sílaba numérica específica, con el mismo fin.

C7H6OCI2(C2H4O)7

lercanidipinum

lercanidipine (±)-2-[(3,3-diphenylpropyl)methylamino]-1,1-dimethylethyl methyl

1,4-dihydro-2,6-dimethyl-4-(m-nitrophenyl)-3,5-pyidinedicarboxytate

lercanidipine (RS)-2,6-dirnéthyl-4-(3-nitrophényl)-1,4-dihydropyridine-3,5-dicarboxylate de

2-[(3,3-diphénylpropyl)(méthyl)amino]-1,1-diméthyléthyle et de méthyle

1.4-dihidro-2,6-dimetil-4-(m-nitrofenil)-3,5-piridinodicarboxilato de lercanidipino

(±)-2-[(3,3-difenilpropil)metilamino]-1,1-dimetiletil metilo

C36H41N3Q6

lerisetronum

lerisetron 1-benzyl-2-(1-piperazinyl)benzımidazole

lérisétron 1-benzyl-2-(pipérazin-1-yl)-1H-benzimidazole

lerisetron 1-bencil-2-(1-piperazinil)bencimidazol

C₁₈H₂₀N₄

letrozolum

letrozole 4.4'-(1H-1,2,4-triazol-1-y/methylene)dibenzonitrile

létrozole 4,4'-[(1H-1,2,4-triazol-1-yl)méthylène]dibenzonitrile

letrozol 4,4'-(1H-1,2,4-triazol-1-ilmetilen)dibenzonitrilo

C17H11N5

lexipafantum

lexipafant \mathcal{N} -methyl- \mathcal{N} -[α -(2-methyl-1 \mathcal{H} -imidazo[4,5-c]pyridin-1-yl)-p-tolyl]sulfonyl]-

L-leucine, ethyl ester

lexipafant (S)-4-méthyl-2-[(méthyl)[4-[[(2-méthylimidazo[4,5-c]pyridin-1-yl)méthyl]]=

sulfonyl]amino]pentanoate d'éthyle

lexipafant ester etílico de la N-metil-N- $[\alpha$ -(2-metil-1 H-imidazo[4,5-c]piridin-1-il)-p-

tolil]sulfonil]-L-leucina

C23H30N4O4S

limazocicum

limazocic (-)-(R)-hexahydro-7,7-dimethyl-6-oxo-1,2,5-dithiazocine-4-carboxylic acid

Imazocio acide (-)-(R)-7,7-diméthyl-6-oxohexahydro-1,2,5-dithiazocine-4-carboxylique

limazocico ácido (--)-(R)-hexahidro-7,7-dimetil-6-oxo-1,2,5-ditiazocina-4-carboxílico

C₈H₁₃NO₃S₂

linotrobanum

linotroban [[5-(2-benzenesulfonamidoethyl)-2-thienyl]oxy]acetic acid

linotroban acide 2-[[5-[2-[(phénylsulfonyl)amino]éthyl]-2-thiényl]acyjacétique

linotroban ácido [[5-(2-bencensulfonamidoetil)-2-tienil]oxi]acetico

C14H15NO5S2

lopobutanum

lopobutan (±)-3-[[3-(dodecyloxy)propyl]amino]butyric acid

lopobutan acide (RS)-3-[[3-(dodécyloxy)propyl]amino]butanoïque

lopobutano ácido (±)-3-[[3-(dodeciloxi)propil]amino]butírico

C19H39NO3

łoviridum

loviride (±)-2-(6-acetyl-m-toluidino)-2-(2,6-dichlorophenyl)acetamide

loviride (RS)-2-[(2-acétyl-5-méthylphényl)amıno]-2-(2,6-dichlorophényl)acétamide

lovirida (±)-2-(6-acetil-m-toluidino)-2-(2,6-diclorofenii)acetamida

C17H16Cl2N2O2

lubeluzolum

lubeluzole (+)-(S)-4-(2-benzothiazolylmethylamino)- α -[(3,4-diffuorophenoxy)methyl]-

1-piperidineethanol

lubéluzole (+)-(S)-1-[4-[(benzothiazol-2-yl)(méthyl)amino]pipéridin-1-yl]-

3-(3,4-difluorophénoxy)propan-2-ol

lubeluzol (+)-(S)-4-(2-benzotiazolilmetilamino)- α -[(3,4-difluorofenoxi)metil]-

1-piperidinietanol

C22H25F2N3O2S

lurosetronum

lurosetron 6-fluoro-2,3,4,5-tetrahydro-5-methyl-2-[(5-methylimidazol-4-yl)methyl]-1*H*-

pyrido[4,3-b]indol-1-one

lurosétron 6-fluoro-5-méthyl-2-[(5-méthyl-1/+imidazol-4-yl)méthyl]-2,3,4,5-tétrahydro-1/+

pyrido[4,3-b]indol-1-one

lurosetron 6-fluoro-2,3,4,5-tetrahidro-5-metil-2-[(5-metilimidazol-4-il)metil]-1*H*-pindo=

[4,3-b]indol-1-ona

C17H17FN4O

merafloxacinum

merafloxacin (±)-1-ethyl-7-[3-[(ethylamino)methyl]-1-pyrrolidinyl]-6,8-difluoro-1,4-dihydro-4-oxo-

3-quinolinecarboxylic acid

mérafloxacine acide (RS)-1-éthyl-7-[3-[(éthylamino)méthyl]pyrrolidin-1-yl]-6,8-dıfluoro-4-oxo-

1,4-dihydroquinoléine-3-carboxylique

merafloxacino ácido (±)-1-etil-7-[3-[(etilamino)metil]-1-pirrolidinil]-6,8-difluoro-1,4-dihidro-4-oxo-

3-quinolincarboxilico

C₁₉H₂₃F₂N₃O₃

mofarotenun

mofarotene 4-[2-[p-[(E)-2-(5,6,7,8-tetrahydro-5,5,8,8-tetramethyl-2-naphthyl)propenyi]=

phenoxy)ethyl)morpholine

mofarotène 4-[2-[4-[(E)-2-(5,5,8,8-tétraméthyl-5,6,7,8-tétrahydronaphtalén-2-yl)prop-1-ényl]=

phénoxy]éthyl]morpholine

mofaroteno 4-[2-[p-(E)-2-(5,6,7,8-tetrahidro-5,5,8,8-tetrametil-2-naftil)propenil](enoxi]etil=

morfolina

C29H39NO2

mofegilinum

mofegiline (E)-2-(fluoromethylene)-4-(p-fluorophenyl)butylamine mofégiline (E)-3-fluoro-2-[2-(4-fluorophényl)éthyl]prop-2-énylamine

mofegilina (E)-2-(fluorometileno)-4-(p-fluorofenil)butilamina

C₁₁H₁₃F₂N

naratriptanum

naratriptan N-methyl-3-(1-methyl-4-piperidyl)indole-5-ethanesulfonamide

naratriptan N-méthyl-2-[3-(1-méthylpipéridin-4-yl)indol-5-yl]éthanesulfonamide

naratrıptan N-metil-3-(1-metil-4-pipericil)indol-5-etanosulfonamida

C₁₇H₂₅N₃O₂S

nedaplatinum

nedaplatin cis-diammine(glycolato-O1,O2)platinum

nédaplatine cis-diammine[2-hydroxyacétato(2-)-O1,O2]platine

nedaplatino cis-diamina(glicolato-O1,O2)platino

C₂H₈N₂O₃Pt

nupafantum

nupafant $\mathcal{N}[(S)-1-(ethoxymethyl)-3-methylbutyl]-N-methyl-<math>\alpha$ -(2-methyl-1H-

imidazo[4,5-c]pyridin-1-yl)-p-toluenesulfonamide

nupafant \(\mathcal{N}\-[(S)-1-(\(\delta\)toxym\(\delta\))-3-m\(\delta\)tylbutyl}-\(N\-\delta\)tyl-4-[(2-m\(\delta\)tyl-1\(H\)

imidazo[4,5-c]pyridin-1-yl)méth/d]benzènesulfonamide

nupafant $N[(S)-1-(etoximetil)-3-metil-butil]-N-metil-\alpha-(2-metil-1)$ imidazo[4,5-c]piridin-

1-ii)-p-toluensulfonamida

C23H32N4O3S

olprinonum

olprinone 1,2-dihydro-5-imidazo[1,2-a]pyridin-6-yl-6-methyl-2-oxonicotinonitrile

olprinone 5-(imidazo[1,2-a]pyridin-6-yl)-6-méthyl-2-oxo-1,2-dihydropyridine-3-carbonitrile

ofprinona 1,2-dihidro-5-imidazo[1,2-a]piridin-6-il-6-metil-2-oxonicotinonitrilo

C14H10N4O

ormeloxifenum

omeloxifene (±)-1-[2-[p-(trans-7-methoxy-2,2-dimethyl-3-phenyl-4-chromanyl)phenoxy]ethyl]=

pyrrolidine

orméloxifène (±)-1-[2-[4-(trans-7-méthoxy-2,2-diméthyl-3-phényl-3,4-dihydro-2*H*-chromén-4-

yl)phenoxy]éthyl]pyrrolidine

ormeloxifeno (±)-1-[2-[p-(trans-7-metoxi-2,2-dimetil-3-fenil-4-cromanil)fenoxi]etil]pirrolidina

C₃₀H₃₅NO₃

oxeclosporinum

oxeclosporin cyclo[[(2S,3R,4R,6E)-3-hydroxy-4-methyl-2-(methylamino)-6-octenov[]-L-

2-aminobutyryl-N-methylglycyl-N-methyl-L-leucyl-L-valyl-N-methyl-L-leucyl-L-alanyl-O-(2-hydroxyethyl)-o-seryl-N-methyl-L-leucyl-N-methyl-N-methyl-L-leucyl-N-methyl-N-methyl-L-leucyl-N-methyl-N

L-valyi]

oxéclosporine cyclo[-[(6E)-(2S,3F,4F)-3-hydroxy-4-méthyl-2-(méthylamino)oct-6-énoyl]-L-

2-aminobutyryl-(N-méthylglycyl)-(N-méthyl-L-leucyl)-L-valyl-(N-méthyl-L-leucyl)-L-alanyl-[O-(2-hydroxyéthyl)-p-séryl]-(N-méthyl-L-leucyl)-(N-méthyl-L-leucyl)-

(N-méthyl-L-valyl)-]

oxeclosporina ciclo[[(2S,3R,4R,6E)-3-hidroxi-4-metil-2-(metilamino)-6-octenoil]-L-

2-aminobutiril-N-metilglicil-N-metil-L-leucil-L-valil-N-metil-L-leucil-L-alanil-O-(2-

hidroxietil)-D-seril-N-metil-L-leucil-N-metil-L-leucil-N-metil-L-valil)

C64H115N11O14

pamicogrelum

pamicogrel ethyl 2-[4,5-bis(p-methoxyphenyl)-2-thiazolyl]pyrrole-1-acetate

pamicogrel 2-[2-[4,5-bis(4-méthoxyphényl)thiazol-2-yl]-1H-pyrrol-1-yl]acétate d'éthyle

pamicogrel etil 2-[4,5-bis(p-metoxifenil)-2-tiazolil]pirrole-1-acetato

C25H24N2O4S

patamostatum

patamostat p-[(2-succinimidoethyl)thio]phenyl p-guanidinobenzoate

patamostat 4-guanidinobenzoate de 4-[[2-(2,5-dioxopyrrolidin-1-yl)éthyl]thio}phényle

patamostat p-guanidinobenzoato de p-[(2-succinimidoetil)tio]fenil

C20H20N4O4S

pazinaclonum

pazinaclone (±)-8-[[2-(7-chloro-1,8-naphthyridin-2-yl)-3-oxo-1-isoindolinyl]acetyl]-1,4-dioxa-

B-azaspiro[4.5]decane

pazinacione (RS)-8-[2-[2-(7-chloro-1,8-naphtyridin-2-yl)-3-oxo-2,3-dihydro-1H-iso-indol-

1-yl]acétyl]-1,4-dioxa-8-azaspiro[4.5]décane

pazinaciona (±)-8-[[2-(7-cloro-1,8-naftindin-2-il)-3-oxo-1-isoindolinif]acetil]-1,4-dioxa-

8-azaspiro[4.5]decano

C25H23CIN4O4

pimagedinum

pimagedine aminoguanidine pimagédine aminoguanidine pimagedina aminoguanidina

CH₆N₄

pobilukastum

pobilukast (2S,3R)-3-[(2-carboxyethyl)thio]-3-[o-(8-phenyloctyl)phenyl]lactic acid

pobilukast acide (2*S*,3*R*)-3-[(2-carboxyéthyl)thio]-2-hydroxy-3-[2-(8-phényloctyl)phényl]=

propanoïque

pobilukast ácido (2S,3R)-3-[(2-carboxietil)tio]-3-[o-(8-feniloctil)fenil]lactico

C₂₆H₃₄O₅S

polixetonii chloridum

polizetonium chloride poly[oxyethylene(dimethyliminio)ethylene(dimethyliminio)ethylene dichloride]

chlorure de polixétonium poly[dichlorure d'oxyéthylène(diméthyliminio)éthylène[diméthyliminio)éthylène]

cloruro de polixetonio poli[dicloruro de oxietileno(dimetiliminio)etileno[dimetiliminio)etileno[

(C₁₀H₂₄Cl₂N₂O)_n

rabeprazolum

rabeprazole 2-[[[4-(3-methoxypropoxy)-3-methyl-2-pyridyl]methyl]sulfinyl]benzimidazole

rabéprazole 2-[[[4-[(3-méthoxypropyl)oxy]-3-méthylpyridin-2-yl]méthyl]sulfinyl]-1 H-benzimi=

dazole

rabeprazoł 2-[[[4-(3-metoxipropoxi)-3-metil-2-piridil]metil]sulfinii]benzimidazol

C18H21N3O3S

ramosetronum

ramosetron (-)-(R)-1-methylindol-3-yl 4,5,6,7-tetrahydro-5-benzimidazolyl ketone

ramosétron (-)-(A)-(1-méthyl-1-H-indol-3-yl)(4,5,6,7-tétrahydro-1-H-benzimidazol-5-yl)=

méthanone

ramosetron (--)-(R)-1-metilindol-3-il 4,5,6,7-tetrahidro-5-bencimidazolil cetona

C₁₇H₁₇N₃O

rasagillnum

rasagiline (R)-N-2-propynyl-1-indanamine

rasagiline [(R)-2,3-dihydro-1H-indén-1-yl](prop-2-ynyl)amine

rasagılina (R)-N-2-propinil-1-indanamina

C₁₂H₁₃N

reteplasum

reteplase 173-L-serine-174-L-tyrosine-175-L-glutamine-173-527-plasminogen activator

(human tissue-type)

rétéplase 173-t-sérine-174-t-tyrosine-175-t-glutamine-173-527-activateur du

plasminogène (type tissulaire humain)

reteplasa 173-L-serina-174-L-tirosina-175-L-glutamina-173-527-activador del

plasminogeno (tipo tisular humano)

C₁₇₃₆H₂₆₅₃N₄₉₉O₅₂₂S₂₂

ricasetronum

ricasetron 3,3-dimethyl-N- $1\alpha H$,5 αH -tropan- 3α -yl-1-indolinecarboxamide

ricasétron 3,3-diméthyl-*N*-[(1*R*,3*r*,5*S*)-8-méthyl-8-azabicyclo[3.2.1]oct-3-yl]-2,3-dihydro-

1H-indole-1-carboxamide

ricasetron 3,3-dimetil-N-1αH,5αH-tropan-3α-il-1-indolinacarboxamida

C₁₉H₂₇N₃O

safingolum

safingol (2S,3S)-2-amino-1,3-octadecanediol safingol (2S,3S)-2-aminooctadecane-1,3-diol safingol (2S,3S)-2-amino-1,3-octadecanodiol

C18H39NO2

sameridinum

sameridine N-ethyl-1-hexyl-N-methyl-4-phenylisonipecotamide

saméridine N-éthyl-1-hexyl-N-méthyl-4-phénylpipéridine-4-carboxamide

C₂₁H₃₄N₂O

saquinavirum

saquinavir $(S)-N-[(\alpha S)-\alpha-[(1H)-2-[(3S,4aS,8aS)-3-(tert-butylcarbamoyl)octahydro-2(1H)-$

isoquinolyl]-1-hydroxyethyl]phenethyl]-2-quinaldamido succinamide

saquinavir (2S)-N'-[(1S,2R)-1-benzyl-3-[(3S,4aS,8aS)-3-[[(1,1-diméthyléthyl)amino]=

carbonyl]octahydro-isoquinoléin-2(1H)-yl]-2-hydroxypropyl]-2-[[(quinoléin-2-yl)=

carbonyl]amino]butanediamide

saquinavir (S)-N-[(α S)- α -[(1R)-2-[(3S,4aS,8aS)-3-(terc-butilcarbamoil)octahidro-2(1H)-

isoquinolil]-1-hidroxietil]fenetil]-2-quinaldamida succinamida

C38H50N6O5

selfotelum

selfotel cis-4-(phosphonomethyl)pipecolic acid

selfotei acide cis-4-(phosphonométhyl)pipéridine-2-carboxylique

selfotel ácido cis-4-(fosfonometil)pipecolico

C7H14NO5P

seratrodastum

seratrodast (±)-2,4,5-trimethyl-3,6-dioxo-ζ-phenyl-1,4-cyclohexadiene-1-heptanoic acid

sératrodast acide (RS)-7-phényl-7-(2,4,5-triméthyl-3,6-dioxocyclohexa-1,4-dién-1-yl)=

heptanoïque

seratrodast ácido (±)-2,4,5-trimetil-3,6-dioxo-ζ-fenil-1,4-ciclohexadieno-1-heptanoico

C22H26O4

5,5-bis(4-pyridylmethyl)-5*H*-cyclopenta[2,1-*b*:3,4-*b*']dipyridine 5,5-bis[(pyridin-4-yl)methyl]-5*H*-cyclopenta[2,1-*b*:3,4-*b*']dipyridine

5,5-bis(4-piridilmetil)-5H-ciclopenta[2,1-b:3,4-b']dipiridina

C₂₃H₁₈N₄

sirolimusum

sirolimus (3*S*,6*R*,7*E*,9*R*,10*R*,12*R*,14*S*,15*E*,17*E*,19*E*,21*S*,23*S*,26*R*,27*R*,34a*S*)-9,10,12,13.

14,21,22,23,24,25,26,27,32,33,34,34a-hexadecahydro-9,27-dihydroxy-3-[(1*R*)-2-[(1*S*,3*R*,4*R*)-4-hydroxy-3-methoxycyclohexyl]-1-methylethyl]-10,21-dimethoxy-6,8,12,14,20,26-hexamethyl-23,27-epoxy-3*H*-pyrido[2,1-c][1,4]oxaazacyclohen=

triacontine-1,5,11,28,29(4H,6H,31H)-pentone

sirolimus (7E,15E,17E,19E)-(3S,6R,9R,10R,12R,14S,21S,23S,26R,27R,34aS)-9,27-

dihydroxy-3-[(1*R*)-2-[(1*S*,3*R*,4*R*)-4-hydroxy-3-méthoxycyclohexyl]-1-méthyléthyl]-10,21-diméthoxy-6,8,12,14,20,26-hexaméthyl-9,10,12,13,14,21,22,23,24,25,26,27,32,33,34,34a-hexadécahydro-23,27-époxy-3*H*-pyrido[2,1-c][1,4]oxaaza=

cyclohentriacontène-1,5,11,28,29(4H,6H,31H)-pentone

sirolimus (3*S*,6*R*,7*E*,9*R*,10*R*,12*R*,14*S*,15*E*,17*E*,19*E*,21*S*,23*S*,26*R*,27*R*,34a*S*)-9,10,12,13, 14,21,22,23,24,25,26,27,32,33,34,34a-hexadecahidro-9,27-dihidroxi-3-[(1*R*)-2-

[(1*S*,3*R*,4*R*)-4-hidroxi-3-metoxiciclohexil]-1-metiletil]-10,21-dimetoxi-6,8,12,14, 20,26-hexametil-23,27-epoxi-3*H*-pirido[2,1-*c*][1,4]oxaazaciclohentriacontina-

1,5,11,28,29(4*H*,6*H*,31*H*)-pentona

C₅₁H₇₉NO₁₃

somatosalmum

somatosalm somatotropin (Oncorhyncus mykiss clone ptGH-II isoform II reduced)

somatosalm somatotropine (isoforme II réduite issue du clone de Oncorhyncus mykiss ptGH-II)

somatosalm somatotropina (isoforma II reducida del clon ptGHII de Oncorhyncus mykiss)

C952H1524N266O290S8

spiroglumidum

spiroglumide (R)-γ-(3,5-dichlorobenzamido)-δ-oxo-8-azaspiro[4.5]decane-8-valeric acid

spiroglumide acide (F)-5-(8-azaspiro[4.5]déc-8-yl)-4-[(3,5-dichlorobenzoyl)amino]-5-

oxopentanoïque

espiroglumida ácido (R)-γ-(3,5-diclorobenzamido)-δ-oxo-8-azaspiro[4.5]decan-8-valérico

C21H26Cl2N2O4

sprodiamidum

sprodiamide aqua[N,N-bis[2-[(carboxymethyl)[(methylcarbamoyl)methyl]=

glycinato(3-)]dysprosium, hydrate

sprodiamide aqua[N,N-bis[2-[(carboxym-śɪnyl)[[(méthylamino)carbonyl]méthyl]amino]éthyl]=

glycinato(3-)]dysprosium hydraté

esprodiamida aquo[N,N-bis[2-[(carboximetil)[(metilcarbamoil)metil]amino]etil]=

glicinato(3-)]disprosio, hidrato

C16H2IIDyN5O9. xH2O

suritozolum

suritozole 3-(*m*-fluorophenyl)-1,4-dimethyl-Δ²-1,2,4-triazoline-5-thione

suritozole 5-(3-fluorophényl)-2,4-diméthyl-2,4-dihydro-3/+1,2,4-triazole-3-thione

suritozol 3-(*m*-fluorofenil)-1,4-dimetil-Δ²-1,2,4-triazolina-5-tiona

C₁₀H₁₀FN₃S

technetium (99mTc) furlfosminum

technetium (99mTc) furifosmin

(OC-6-13)-[[4,4'-[ethylenebis(nitrilomethylidyne)]bis[dihydro-2,2,5,5-tetramethyl-3(2H)-furanonato]](2-)-N,N',O3,O3']bis[tris(3-methoxypropyl)phosphine-P][99mTc]=

technetium(1+) chloride

technétium (99mTc) furifosmine

(OC-6-13)-chlorure de [[4,4'-[éthylènebis(nitrilométhylidyne)]bis= [2,2,5,5-tétraméthyldihydrofuran-3(2H)-onato]](2-)-N,N',O3,O3']bis=

[tris(3-methoxypropyl)phosphine-P][99mTc]technetium(1+)

furifosmina de technetium (99mTc)

cloruro de (OC-6-13)-[[4,4'-[etilenbis(nitrilometilidina)]bis[dihidro-2,2,5,5tetrametil-3(2H)-furanonato]](2-)-N,N',O3,O3]bis[tris(3-metoxipropil)fostina-

P][99mTc]tecnecio(1+)

C44H64CIN2O10P299mTc

telmisartanum

telmisartan

4'-[[4-methyl-6-(1-methyl-2-benzimidazolyl]-2-propyl-1-benzimidazolyl]methyl]-

2-biphenylcarboxylic acid

telmisartan

acide 4'-[[4-méthyl-6-(1-méthyl-1H-benzimidazol-2-yl)-2-propyl-1H-benzimidazol-

1-yl]méthyl]biphényl-2-carboxylique

telmisarian

ácido 4'-[[4-metil-6-(1-metil-2-bencimidazolil)-2-propil-1-bencimidazolil]metil]-

2-bifenilcarboxílico

C33H30N4O2

temoporfinum

temoporfin

3.3',3"',3"'-(7,8-dihydroporphyrin-5,10,15,20-tetrayl)tetraphenol

témoporfine

3,3',3"'-(7,8-dihydroporphyrine-5,10,15,20-tetrayl)tétraphénol

temoporfina

3.3',3",3"-(7,8-dihidroporfirin-5,10,15,20-tetrail)tetrafenol

C44H32N4O4

tolafentrinum

tolafentrine

(-)-4'-(cis-1,2,3,4,4a,10b-hexahydro-8,9-dimethoxy-2-methylbenzo[c][1,6]=

naphthyridin-6-yl)-p-toluenesulfonanilide

tolafentrine

(-)-N-[4-(cis-8,9-diméthoxy-2-méthyl-1,2,3,4,4a,10b-hexahvdrobenzolcl(1,6]=

naphtyridin-6-yl)phényl]-4-méthylbenzènesulfonamide

tolafentrina

(-)-4'-(cis-1,2,3,4,4a,10b-hexahidro-8,9-dimetoxi-2-metilbenzo[c][1,6]=

naftiridin-6-il)-p-toluenosulfonanilida

C28H31N3O4S

tradecamidum

tradecamide tradécamide 13-hydroxy-N,N-dimethyltridecanamide

tradecamida

13-hydroxy-N,N-diméthyltridécanamide

13-hidroxi-N,N-dimetiltridecanamida

C15H31NO2

ularitidum

ularitide

L-threonyl-L-alanyl-L-prolyl-L-arginyl-L-seryl-L-leucyl-L-arginyl-L-arginyl-L-seryl-L-seryl-L-cysteinyl-L-phenylalanylgiycylglycyl-L-arginyl-L-methionyl-L-aspartyl-

L-arginyl-L-isoleucylglycyl-L-alanyl-L-glutaminyl-L-serylglycyl-L-leucylglycyl-

L-cysteinyl-L-asparaginyl-L-seryl-L-phenylalanyl-L-arginyl-L-tyrosine cyclic (11→27)-

disulfide

ularitide (11→27)-disulfure cyclique de t-thréonyl-t-alanyl-t-prolyl-t-arginyl-t-séryl-

L-leucyl-L-arginyl-L-arginyl-L-séryl-L-séryl-L-cystéinyl-L-phénylatanyl-glycyl-glycyl-L-arginyl-L-méthionyl-L-aspartyl-L-arginyl-L-isoleucyl-glycyl-L-alanyl-L-glytaminyl-L-séryl-glycyl-L-leucyl-glycyl-L-cystéinyl-L-asparaginyl-L-séryl-L-phénylatanyl-

L-arginyl-L-tyrosine

ularitida c-treonil-c-alanil-c-prolil-c-arginil-c-seril-c-leucil-c-arginil-c-arginil-c-seril-c-seril-

L-cisteinil-L-fenilalanilglicilglicil-L-arginil-L-metionil-L-aspartıl-L-arginil-L-isoleu= cilglicil-L-alanil-L-glutaminil-L-serilglicil-L-leucilglicil-L-cisteinil-L-asparaginil-L-serilglicil-L-cisteinil-L-asparaginil-L-serilglicil-L-cisteinil-L-asparaginil-L-serilglicil-L-cisteinil-L-asparaginil-L-serilglicil-L-cisteinil-L-asparaginil-L-serilglicil-L-cisteinil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-asparaginil-L-serilglicil-L-serilglicil-L-asparaginil-L-serilglicil-L-serilglicil-L-asparaginil-L-serilglicil-L-serilglicil-L-asparaginil-L-serilglicil-L-s

L-fenilalanil-L-arginil-L-tirosina disulfuro cíclico (11→27)

C145H234N52O44S3

valaciclovirum

valaciclovir L-valine, ester with 9-[(2-hydroxyethoxy)methyl]guanine

valaciclovir (S)-2-amino-3-méthylbutanoate de 2-[(2-amino-6-oxo-1,6-dihydro-9-H-purin-9-yl)=

méthoxy]éthyle

valaciclovir éster de la L-valina, con 9-[(2-hidroxietoxi)metil]guanina

C13H20N6O4

vebufloxacinum

vebufloxacin (±)-9-fluoro-6,7-dihydro-5-methyl-8-(4-methyl-1-piperazinyl)-1-oxo-1H,5H

benzo[ii]quinolizine-2-carboxylic acid

vébufloxacine acide (RS)-9-fluoro-5-méthyl-8-(4-méthylpipérazin-1-yl)-1-oxo-6,7-dihydro 1H,5H-

benzo[ij]quinolizine-2-carboxylique

vebufloxacıno ácido (±)-9-fluoro-6,7-dihidro-5-metil-8-(4-metil-1-piperazinil)-1-oxo-1H,5H

benzo[ij]quinolizina-2-carboxílico

C19H29FN3O3

votumumabum

votumumab immunoglobulin G3 (human monoclonal 88-BV59 heavy chain anti-human

carcinoma-associated antigen), disulfide with human monoclonal 88-BV59

κ-chain, dimer

votumumab immunoglobuline G 3 (chaîne lourde de l'anticorps monoclonal humain 88-BV59

anti-antigène associé aux carcinomes humains), dimère du disulfure avec la

chaîne κ de l'anticorps monoclonal humain 88-BV59

votumumab inmunoglobulina G3 (cadena pesada del anticuerpo monoclonal 88-Bv59 humano

anti-antigeno asociado a los carcinomas humanos), puentes disulfuro con la

cadena x del anticuerpo monoclonal 88-BV59 humano, dimero

xanomelinum

xanomeline 3-[4-(hexyloxy)-1,2,5-thiadiazol-3-yl]-1,2,5,6-tetrahydro-1-methylpyridine

xanoméline 3-[4-(hexyloxy)-1,2,5-thiadiazol-3-yl]-1-méthyl-1,2,5,6-tétrahydropyridine

xanomelina 3-[4-(hexiloxi)-1,2,5-tiadiazol-3-il]-1,2,5,6-tetrahidro-1-metilpiridina

C14H23N3OS

zolasartanum

zolasartan 1-[[3-bromo-2-(o-1/H-tetrazol-5-yiphenyl)-5-benzofuranyl]methyl]-2-butyl-4-

chloroimidazole-5-carboxylic acid

zolasartan acide 1-[[3-bromo-2-[2-(1H-tétrazol-5-yl)phényl]benzofuran-5-yl]méthyl]-2-butyl-

4-chloro-1H-imidazole-5-carboxylique

zolasartan ácido 1-[[3-bromo-2-(o-1H-tetrazol-5-ilfenil)-5-benzofuranil]metil]-2-butil-4-

cloroimidazol-5-carboxílico

C24H20BrClN6O3

zolimomabum aritoxum

zolimomab aritox immunoglobulin G 1 (mouse monocional H65-RTA anti-human antigen CD 5

heavy chain), disulfide with mouse monoclonal H65-RTA light chain, dimer,

disulfide with ricin (castor-oil plant A-chain protein moiety)

zolimomab aritox immunotoxine obtenue par couplage, par une liaison disulfure, de l'immuno=

globuline G1 monoclonale de souris H65-RTA dirigée contre l'antigène de surface CD 5 des lymphocytes T humains et de la chaîne A de la ricine

zolimornab aritox inmunoglobulina G1 monocional de ratón H65-RTA anti(antigeno de

superficie CD5 de los linfocitos T humano), unida mediante enlace disulfuro con la cadena ligera de anticuerpo monoclonal de ratón H65-RTA, dimero, disulfuro

con la cadena A de la ricina

AMENDMENTS TO PREVIOUS LISTS

Supplement to WHO Chronicle Vol. 37, No. 6, 1983

Recommended International Nonproprietary Names (Rec. INN): List 23

p. 5 iloprostum

replace the chemical name by the following:

iloprost

(E)-(3aS,4R,5R,6aS)-hexahydro-5-hydroxy-4-[(E)-(3S,4RS)-3-hydroxy-4-

methyl-1-octen-6-ynyl]Δ2(1H),δ-pentalenevaleric acid

p. 6 mitindomidum

replace the chemical name by the following:

mitindomide

(1R*,25*,3R*,45*,5R*,65*,75*,8R*)-tricyclo[4,2,2,0^{2,5}]dec-9-ene-3,4,7,8-

tetracarboxylic 3,4:7,8-diimide

Supplement to WHO Chronicle Vol. 38, No. 6, 1984

Recommended International Nonproprietary Names (Rec. INN): List 24

p. 10 valproatum seminatricum

valproate semisodium

replace the chemical name and the molecular formula by the following:

sodium hydrogen bis(2-propylvalerate), oligomer

(C₁₆H₃₁NaO₄)_n

WHO Drug Information, Vol. 5, No. 3, 1991

Recommended International Nonproprietary Names (Rec. INN): List 31

p. 2 aprikalimum

replace the chemical name by the following:

aprikalım

(-)-(1R, 2R)-tetrahydro-N-methyl-2-(3-pyridyl)thio-2H-thiopyran-2-carboxamide

1-oxide

p. 6

gadodiamidum

replace the chemical name and the molecular formula by the following:

gadodiamide [N,N-bis[2-[(carboxymethyl)[methylcarbamoyl)methyl]amino]ethyl]glycinato=

(3-)gadolinium

C16H26GdN5O8

p. 6

gadoteridolum

replace the chemical name by the following:

gadoteridol

(±)-[10-(2-hydroxypropyl)-1,4,7,10-tetraazacyclododecane-1,4,7-triacetato=

[(3-)]gadolinium

MODIFICATIONS APPORTÉÉS AUX LISTES ANTÉRIEURES

Supplément à la Chronique OMS, Vol. 37, No. 6, 1983

Dénominations communes internationales recommandées (DCI Rec.): Liste 23

p. 6 mitindomidum

remplacer le nom chimique par:

mitindomide

(1R*,25*,3R*,45*,5R*,65*,75*,8R*)-tricyclo[4.2.2.0^{2,5}]déc-9-ène-3,4,7,8-

tétracarboxy-3,4:7,8-dumide

Supplément à la Chronique OMS, Vol. 38, No. 6, 1984

Dénominations communes internationales recommandées (DCI Rec.): Liste 24

p. 10 valproatum seminatricum

remplacer le nom chimique et la formule brute par:

valproate semisodique

oligomère du complexe d'acide 2-propylpetanoïque et de 2-propylpentanoate

de sodium (C₁₆H₃₁NaO₄),

Informations pharmaceutiques OMS, Vol. 5, No. 3, 1991

Dénominations communes internationales recommandées (DCI Rec.): Liste 31

p. 2 aprikalimum

remplacer le nom chimique par:

aprikalim

(-)-(1R, 2R)-N-methyl-2-(pyndin-3-yl)tétrahydro-2H-thiopyrane-2-

carbothioamide 1-oxyde

p. 6 gadodiamidum

gadodiamide

remplacer le nom chimique et la formule brute par.

[N,N-bis[2-

[(carboxyméthyl)[(méthylamino)carbonyl]méthyl]amino]éthyl]glycinato=

(3-)gadolinium C₁₆H₂₆GdN₅O₈

MODIFICACIONES A LAS LISTAS ANTERIORES

Suplemento de Crónica de la OMS, Vol. 37, No. 6, 1983

Denominaciones Comunes Internacionales Recomendadas (DCI Rec.): Lista 23

p. 6 mitindomidum

sustituir el nombre químico por lo siguiente:

mitindomida

(1R*,25*,3R*,45*,5R*,65*,75*,8R*)-triciclo[4.2.2.0^{2.5}]dec-9-eno-3,4,7,8-

tetracarboxílico 3,4:7,8-diimida

Suplemento de Crónica de la OMS, Vol. 38, No. 6, 1984

Denominaciones Comunes Internacionales Recomendadas (DCI Rec.): Lista 24

p. 10 valproatum seminatricum valproato semisódico sustituyase el nombre químico y la fórmula empirica por los siguientes:

bis(2-propilvalerato) de hidrogeno y sodio, oligómero

(C₁₆H₃₁NaO₄)_{/1}

Información Farmacéutica, de la OMS, Vol. 5, No. 3, 1991

Denominaciones Comunes Internacionales Recomendadas (DCI Rec.): Lista 31

p. 2 aprikalimum

sustituyase el nombre químico por lo siguiente:

aprikalim

(-)-(1R, 2R)-tetrahidro-N-metil-2-(3-piridil)tıo-2H-tiopiran-2-carboxamida 1-óxido

p. 6 gadodiamidum gadodiamida sustituir el nombre químico y la fórmula empírica por los siguientes:

[N,N-bis[2-[(carboximetil)[(metilcarbamoil)metil]amino]etil]glicinato= (3-) gadolinio

 $C_{16}H_{26}GdN_5O_8$