International Non-Proprietary Names for Pharmaceutical Preparations

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There is need to avoid the confusion which exists when different non-proprietary names come into use for the same medicinal substance. This multiplicity of names can be the source of difficulties in the daily work of the physician and of the pharmacist, as well as in therapeutic research throughout the world.

In many countries attempts have been made and progress achieved in obtaining the desired uniformity on a national level (examples: "Generic Names" of the Council on Pharmacy and Chemistry of the American Medical Association, "Approved Names" of the General Medical Council in the United Kingdom, "Nordiske Farmakopenaevn" of the Scandinavian Pharmacopoeia Council, "dénominations communes" in France, etc.) The World Health Organization was asked to co-ordinate these efforts at an international level, and special requests were also made to select international non-proprietary names for the drugs liable to produce addiction in order to facilitate their international control.

According to the following Procedure for the Selection of Recommended International Non-Proprietary Names for Pharmaceutical Preparations, WHO receives requests for the establishment of international non-proprietary names for new pharmaceutical preparations which can be used freely in all countries, and publishes lists of proposed international non-proprietary names and of recommended international non-proprietary names.

PROCEDURE FOR THE SELECTION OF RECOMMENDED INTERNATIONAL NON-PROPRIETARY NAMES FOR PHARMACEUTICAL PREPARATIONS ¹

The following procedure shall be followed by the World Health Organization in the selection of recommended international nonproprietary names for pharmaceutical preparations, in accordance with the World Health Assembly resolution WHA3 11:

- 1. Proposals for recommended international non-proprietary names shall be submitted to the World Health Organization on the form provided therefor.
- 2. Such proposals shall be submitted by the Director-General of the World Health Organization to the members of the Expert Advisory Panel on the International Pharmacopoeia and Pharmaceutical Preparations

¹The amended text reproduced here was adopted by the Executive Board in resolution EB15.R7 and supersedes that adopted by the Board at its twelfth session (see Official Records No. 49, Annex 6).

designated for this purpose, for consideration in accordance with the "General principles for guidance in devising International Non-proprietary Names", appended to this procedure. The name used by the person discovering or first developing and marketing a pharmaceutical preparation shall be accepted, unless there are compelling reasons to the contrary.

- 3. Subsequent to the examination provided for in article 2, the Director-General of the World Health Organization shall give notice that a proposed international non-proprietary name is being considered.
 - A. Such notice shall be given by publication in the Chronicle of the World Health Organization and by letter to Member States and to national pharmacopoeia commissions or other bodies designated by Member States.
 - (i) Notice may also be sent to specific persons known to be concerned with a name under consideration.
 - B. Such notice shall:
 - (i) set forth the name under consideration;
 - (ii) identify the person who submitted a proposal for naming the substance, if so requested by such person;
 - (iii) identify the substance for which a name is being considered;
 - (iv) set forth the time within which comments and objections will be received and the person and place to whom they should be directed:
 - (v) state the authority under which the World Health Organization is acting and refer to these rules of procedure.
 - C. In forwarding the notice, the Director-General of the World Health Organization shall request that Member States take such steps as are necessary to prevent the acquisition of proprietary rights in the proposed name during the period it is under consideration by the World Health Organization.

- 4. Comments on the proposed name may be forwarded by any person to the World Health Organization within four months of the date of publication, under article 3, of the name in the Chronicle of the World Health Organization.
- 5. A formal objection to a proposed name may be filed by any interested person within four months of the date of publication, under article 3, of the name in the Chronicle of the World Health Organization:
 - A. Such objection shall:
 - (i) identify the person objecting;
 - (ii) state his interest in the name;
 - (iii) set forth the reasons for his objection to the name proposed.
- 6. Where there is a formal objection under article 5, the World Health Organization may either reconsider the proposed name or use its good offices to attempt to obtain withdrawal of the objection. Without prejudice to the consideration by the World Health Organization of a substitute name or names, a name shall not be selected by the World Health Organization as a recommended international non-proprietary name while there exists a formal objection thereto filed under article 5 which has not been withdrawn
- 7. Where no objection has been filed under article 5, or all objections previously filed have been withdrawn, the Director-General of the World Health Organization shall give notice in accordance with subsection A of article 3 that the name has been selected by the World Health Organization as a recommended international non-proprietary name.
- 8. In forwarding a recommended intentional non-proprietary name to Memoc. States under article 7, the Director-General of the World Health Organization shall:
 - A. request that it be recognized as the non-proprietary name for the substance; and
 - B. request that Member States take such steps as are necessary to prevent the acquisition of proprietary rights in the name, including prohibiting registration of the name as a trade-mark or trade-name.

GENERAL PRINCIPLES FOR GUIDANCE IN DEVISING INTERNATIONAL NON-PROPRIETARY NAMES

- 1. Names should, preferably, be free from any anatomical, physiological, pathological or therapeutic suggestion,
- 2. An attempt should first be made to form a name by the combination of syllables in such a way as to indicate the significant chemical groupings of the compound and/or its pharmacological classification. Preference should be given to the following syllables:

Latın	English	French	
inum	ine	ine	for alkaloids and organic bases
inum	ia	ine	for glycerides and neutral principles
olum	ol	ol	for alcohols and phenols (-OH group)
alum	al	al	for aldehydes
onum	one	оле	for ketones and other substances containing the CO group
enum	ene	ène	for unsaturated hydrocarbons
าถนต	ane	ane	for saturated hydrocarbons
cainum	caine	caine	for local anaesthetics
mer	mer	mer	for mercurial compounds
sulfonum	sulfone	sulfone	for sulfone derivatives
quinum	quine	quine	for antimalarial substances containing a quinoline group
crinum	crine	crine	for antimalarial substances containing an acridine group
sulfa	sulfa	sulfa	for derivatives of sulfanilamide
dionum	dione	dione	for anti-epileptics derived from oxazolidinedione
toinum	toin	toine	for anti-epileptics derived from hydantoin
stigmmum	stigmine	stigmine	for anticholinesterases

- 3. Names should be distinctive in sound and spelling. They should not be inconveniently long and should not be liable to confusion with names already in use,
- 4. The addition of a terminal capital letter or number should be avoided as far as possible.
- 5. Names proposed by the person discovering or first developing and marketing a pharmaceutical preparation, or already officially adopted in any country, or used in the national pharmacopoeias, or in works of reference such as "New and Non-official Remedies", should receive preferential consideration.

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In accordance with article 3 of the Procedure for the Selection of Recommended International Non-Proprietary Names for Pharmaceutical Preparations, notice is hereby given that the following names are under consideration by the World Health Organization as Proposed International Non-Proprietary Names.

Comments on, or formal objections to, these proposed names may be forwarded by any person to the World Health Organization within four months from 1 February 1956.

The inclusion of a name in this list does not imply any recommendation for the use of the substance in medicine or pharmacy.

PROPOSED INTERNATIONAL NON-PROPRIETARY NAMES (Prop. I.N.N.): LIST 4 2

Proposed International Non-Proprietary Name (Lahn, English)

Chemical Name or Description

(Latin, English)
acetarsolum

3-acetamido-4-hydroxyphenylarsonic acid

acetarsol

acetylcholini chloridum acetylcholine chloride 2-acetoxyethyltrimethylammonium chloride

Other lists of proposed international non-proprietary names will be found in Chron. Wid Hith Org., 1953, 7, 297; 1954, 1, 216, 313. A list of recommended international non-proprietary names was published in Chron. Wid Hith Org., 1955, 9, 185.

Chemical Name or Description

acıdum ascorbicum

ascorbic acid

acidum folicum

folic acid

acidum iophenoicum iophenoic acid

acidum nicotinicum nicotinic acid

adipiodonum

adipiodone aethisteronum ethisterone

aethylis biscoumacetas ethyl biscoumacetate

amidopyrinum amidocyrine amınıtı əzolum

aminophyllinum aminophyllinum aminophylline

aminopterinum natricum aminopterin sodium

amphetaminum amphetamine androstanolonum androstanolone

arsthiaolum arsthiaol

atropint methoritras atropine methoritrate

barbitalum barbital

barbit dum natricum barbit d sodium

benzatropint methanesulfonas benzatropine methanesulfonate

bromazinum broma<u>nine</u>

buclizmum buclizme

butacamum butacaine

butalbitalum butalbital 3-oxo-L-gulofuranolactone (enolic form)

N-[4'-([2-amino-4-hydroxy-6-pteridyl]methyl]amino) benzoyl]-L

(+)glutamic acid

a-(3-hydroxy-2,4,6-triiodobenzyl)-butyric acid

pyridine-3-carboxylic acıd

adipic acid bis-(2,4,6-triiodo-3-carboxyanilide)

 17α -ethynyl- 17β -hydroxy-3-oxoandrostene-4

ethyl 4,4'-dihydroxy-3,3'-dicoumarinylacetate

2,3-dimethyl-4-dimethylamino-1-phenyl-5-pyrazolone

2-acetamido-5-nitrothiazole

mixture of theophylline and ethylenediamine

sodium 4-ammofolate

(±) 2-amino-1-phenylpropane

3a-hydroxy-17-oxoandrostane

2-(3'-acetamido-4'-hydroxyphenyl)-1,3-dithia-2-arsa-4-cyclopentyl-methanol

(\pm) 8-methyl-3-tropoyloxytropane nitrate

5,5-diethylbarbituric acid

sodium salt of 5,5-diethylbarbituric acid

tropine benzhydryl ether methanesulfonate

2-(4-bromophenyl-phenylmethoxy)-ethyldimethylamine

1-(4-chlorobenzhydryl)-4-(4-tert butylbenzyl)-piperazine

3-dibutylaminopropyl p-aminobenzoate

5-allyl-5-isobutylbarbituric acid

Chemical Name or Description

calciferolum calciferol

calcium saccharate

calcium D-glucarate

carbacholum carbachol

carbamoylcholine chloride

carbarsonum carbarsone 4-ureidophenylarsonic acid

carbimazolum carbimazole

2-ethoxycarbonylthio-1-methylimidazole

carbinoxaminum carbinoxamine 2-dimethylaminoethoxy-2-pyridyl-4-chlorophenylmethane

carzenidum carzenide 4-carboxybenzenesulfonamide

ceromacrogolum 1000 cetomacrogol 1000 polyethylene glycol 1000 monocetyl ether

chiniofonum chiniofon

mixture of four parts by weight of 7-iodo-8-hydroxyquinoli ie 5-sulfonic acid and one part of sodium bicarbonate

N-(2,4-dichlorobenzyl)-N-(dichloroacetyl)-ethanolamine

chlorbetamidum chlorbetamide

(3-chloromercuri-2-methoxypropyl)-urea

chlormerodrinum chlormerodrine

I, I, 1-trichloro-2-methylpropanoI-2

chlorobutanolum chlorobutanol

6-chloro-3-hydroxytoluene

chlorocresolum chlorocresol chloroquinum

7-chloro-4-(4'-diethylamino-1'-methylbutylamino)-quinoline

chloroquin chlorphenaminum chlorphenamine

1-(p-chlorophenyl)-1-(2-pyridyl)-3-dîmethylaminopropane

chlortetracyclinum chlortetracycline

10-chloro-1-dimethylamino-1,4,6,11,12,13,14,18-octahydro-2,5,711, 14-pentahydroxy-4,6-dioxo-11-methylnaphthacene-3-carbonamide

cholini chloridum Sholine chloride

2-hydroxyethyltrimethylammonium chloride

conessinum conessine an alkaloid obtained from the seeds of Holarrhena antidysenterica L.

cycriminum cycrimine

1-phenyl-1-cyclopentyl-3-piperidino-1-propanol

desoxycortonum desoxycortone

21-hydroxy-3,20-dioxopregnene-4

dibrompropamidinum dibrompropamidine

1,3-bis-(4-amidino-2-bromophenoxy)-propane

dichloroxylenolum dichloroxylenol 2,4-dichloro-3,5-dimethylphenol

Chemical Name or Description

diethyistilboestrolum diethylstilboestrol

trans-3,4-bis-(4-hydroxyphenyl)-hexene-3

digoxinum digoxin glycoside obtained from the leaves of Digitalis lanata Ehrh.

dihexyverinum dihexyverine 2-piperidinoethyl 1-cyclohexylcyclohexanecarboxylate

dihydralazınum dıhydralazine

dimazolum

1,4-dihydrazinophthalazine

dimazole

2-dimethylamino-6-(2-diethylaminoethoxy)-benzothiazole

diphemanili methylsulfas diphemanil methylsulfate N, N-dimethyl-4-piperidylidene-diphenylmethane methylsulfate

dithranolum dithranol ectylarea ectylarea 1,8,9-anthratriol

2-ethylcrotonylurea

edrophonii chloridum edrophonium chloride ethyldimethyl-3-hydroxyphenylammonium chloride

ergometrinum ergometrine

an alkaloid obtained from ergot

ergotaminum ergotamine

an alkaloid obtained from ergot

erythromycinum erythromycin

an antibiotic produced by a strain of Streptoinyces erythreus

ethaverinum

6,7-diethoxy-1-(3',4'-diethoxybenzyl)-isoquinoline

gitaliirim amorphum gitalin amorphous a glycosidal principle of Digitalis purpurea L.

glycerolum glycerol

propanetriol

hepatimum hepatim hexylcalnum hexylcalne

sodium salt of a complex organic acid having the characteristic property of delaying the clotting of blood

I-cyclohexylamino-2-propyl benzoate

hydroxystilbamidinum hydroxystilbamidine

1-(4-amidino-2-hydroxyphenyl)-2-(4-amidinophenyl)-ethene

isophanum insulinum isophane insulin a sterile suspension of insulin with the isophanic equivalent of protamine and zinc chloride

lanatosidum C lanatoside C glycoside obtained from the leaves of Digitalis lanata Ehrh.

laudexii methylsulfas laudexium methylsulfate decamethylene-w-bis-[1-(3',4'-dimethoxybenzyl)-1,2,3,4-tefranydro-6,7-dimethoxy-2,2-dimethyl:soquinolinium_sulfate]

levorphanolum* levorphanol

levothyroxinum natricum levothyroxine sodium

lobelinum lobeline rucanthonum lucanthone

macrogolum, 400 macrogol 400

macrogol 1000 macrogol 1000

macrosolum 4000 macrogol 4000

macrogoli laurate 600 macrogoli laurate 600 macrogoli oleas 600 macrogoli oleate 600

macrogoli stearas 600 macrogol stearate 600

macrogoli stearas 1000 macrogol stearate 1000

meclozine melarsoprolam melarsoprol

mepacrinum mepacrine meprykainum meprykaine

mercumatilinum natricum

mersalylum mersalyl - _ methioninum methionine

methoxaminum methoxamine

methylcellulosum methylcellulose

methylpentynolum methylpentynol Chemical Name or Description

(-) 3-hydroxy-N-methylmorphinan

L-β-[(4-hydroxy-3,5-diiodophenoxy)-3,5-diiodophenyl]-alanıne

L-2-(2-hydroxy-2-phenylethyl)-I-methyl-6-phenacyIpiperidine

1-(2-diethylaminoethylamino)-4-methylthiaxanthone

polyethylene glycol 400

polyethylene glycol 1000

polyethylene glycol 4000

mono ester of lauric acid and polyethylene glycol 600

mono ester of oleic acid and polyethylene glycol 600

mono ester of steams acid and polyethylene glycol 600

mono ester of stearic acid and polyethylene glycol 1000

1-(4-chlorobenzhydryl)-4-(3-methylbenzyl)-piperazine

2-[4-(4,6-diamino-2-s-triazinylamino)-phenyl]-4-hydroxymethyl-1,3, 2-dithiaarsenolidine

3-chloro-9-(4'-diethylamino-1'-methylbutylamino)-7-methoxyacridine

2-methyl-2-propylamino-propyl benzoate

sodium 8-(2'-methoxy-3'-hydroxymercuripropyl)-coumarin-3carboxylate (sodium mercumallylate) and theophylline

sodium salt of 2-[(3-hydroxymercuri-2-methoxy-propyl)-carbamoyl]phenoxyacetic acid

(a) 2-amino-4-methylthro-butyric acid

2-amino-1-(2,5-dimethoxyphenyl)-propanol-1

cellulose methyl ether containing about 30 per cent. w/w of

methoxyl

3-methy/pentyn-1-ol-3

^{*}This name is to replace "leverphanum" ("leverphanum"), which had been proposed for the same substance (see Chron. Wid Hith Org., 7, 311)

Chemical Name or Description

methyltestosteronum methyltestosterone

178-hydroxy-17a-methyl-3-oxoandrostene-4

natrii acetrizoas sodium acetrizoate

sodium 3-acetamido-2,4,6-triiodobenzoate

natrii amidotrizoas sodium amidotrizoate

sodium 3,5-diacetamido-2,4,6-triiodobenzoate

neoarsphenaminum neoarsphenamine

sodium 3,3'-diamino-4,4'-dihydroxyarsenobenzene-N-methylene-

neostigmini bromidum

sulfoxylate dimethylcarbamic ester of 3-hydroxy-phenyltrimethylammonium

neostigmine bromide nicethamidum

bromide pyridine-3-carboxylic acid diethylamide

nikethamide nicotinamidum nicotinamide

pyridine-3-carboxylic acid amide

nortestosteroni cypionas nortestosterone cypionate 17β-(3-cyclopentylpropionoxy)-3-oxoestrene-4

oestradioli benzoas oestradiol benzoate 3-benzoyloxy-17β-hydroxyoestratriene-1,3,5 (10)

qestradiolum

 $3,17\beta$ -dihydroxyoestratriene-1,3,5 (10)

pestradiol destronum destrone

3-hydroxy-17-oxoestratriene-1,3,5 (10)

pamaquinum pamaquin

8-(4-diethylamino-1-methylbutylamino)-6-methoxyquinoline salt of 2,2'-dihydroxy-1,1'-dinaphthylmethane-3,3'-dicarboxylic acid

pentaquinum pentaquin

8-(5-isopropylamino-pentylamino)-6-methoxyquinoline

pentetrazolum pentetrazol

pentamethylene-1,5-tetrazole

pentolonium pentolonium pethidmum pethidine

1,5-(1,1'-dimethyl-2,2'-dipyrrolidyl)-pentane

phenacainum phenacaine

ethy! 1-methyl-4-phenyl-piperidyl-4-carboxylate

phenacetinum

N1, N2-bis-(4-ethoxyphenyl)-acetamidine

phenacetin phenazonum phenazone

acetyl-4-phenetidine

phenazopyridinum phenazopyridine

2,3-dimethyl-1-phenyl-5-pyrazolone

phenobarbitalum

2,6-diamino-3-phenylazopyridine

phenobarbital phenobarbitalum natricum phenobarbital sodium

5-ethyl-5-phenyl-barbitume acid

phenylhydraegyri boras phenylmercuric borate

sodium salt of 5-ethyl-5-phenyl-barbituric acid

equimolecular compound of phenylmercuric borate and phenylmercuric hydroxide

Chemical Name or Description

phenytoinum
phenytoin

pramocainum pramocaine

primidonum primidone

progesteronum progesterone proguanilum

proguanil
propoxycamum
propoxycame
quinisocainum
quinisocaine

reserpinum reserpine riboflavinum riboflavine secobarbitalum

secobarbital solasulfonum solasulfone

sorbimacrogoli oleas 100 sorbimacrogoli oleas 300 sorbimacrogoli oleas 300 sorbimacrogoli oleate 300 shibamidini isethionas stilbamidine isethionate succinylsulfathiazolum

succinylsulfathiazole sulfadiazinum sulfadiazine

suifadiazinum natricum suifadiazine sodium sulfadicramidum sulfadicramide sulfaguanidinum sulfaguanidine

sulfamerazinum sulfamerazine

sulfamerazinum natricum sulfamerazine sodium

sulfanılamidum sulfanılamide sulfaproxylinum sulfaproxyline 5,5-diphenylhydantoin

1-[3-(4-morpholino)-propoxy]-4-butoxybenzene

5-ethyl-5-phenyl-4,6-dioxo-hexahydropyrimidine

3,20-dioxopregnene-4

N1-4-chlorophenyl-N5-isopropylbiguanide

2-diethylaminoethyl 4-amino-2-propoxybenzoate

1-(2-dimethylaminoethoxy)-3-butylisoquinoline

alkaloid from the roots of various species of Rauwolfia

6,7-dimethyl-9-(D-1'-ribityl)iso-alloxazine

5-allyl-5-(1-methylbutyl)-barbituric acid

tetrasodium salt of 4,4'-bis-(3-phenyl-1,3-disulfopropylamino)-diphenylsulfone

mono ester of oleic acid and tripolyethyleneglycol 100-sorbitan

mono ester of oleic acid and tripolyethyleneglycol 300-sorbitan

1,2-bis-(4-amidinophenyl)-ethene di-(2-hydroxyethane)-sulfonate

2-(N4-3-carboxypropionyl-sulfanilamido)-thiazole

2-sulfanilamidopyrimidine

sedium derivative of 2-sulfanilamidopyrimidine

N¹-(3,3-dimethylacroyl)-sulfanilamide

N1-amidinosulfanilamide

2-sulfanilamido-4-methylpyrimidme

sodium derivative of 2-sulfanilamido-4-methylpyrimidine

4-aminobenzenesulfonamide

N1-(4-isopropoxybenzoyl)-sulfanılamıde

Chemical Name or Description

sulfarsphenaminum disodium 3,3'-diamino-4,4'-dihydroxyarsenobenzene-N,N'-bissulfarsphenamine methylenebisulfite sulfathiazolum 2-sulfanilamidothiazole sulfathiazole a sterile buffered suspension of the amorphous form of insulin with suspensio insulini cum zinco (amorphum) insulin zinc suspension (amornhous) zinc chloride suspensio insulmi cum zinco (crystala sterile buffered suspension of the crystalline form of insulin with lisatum) zine chloride insulin zinc suspension (crystalline) testosteronum 178-hydroxy-3-oxoandrostene testosterone tetrabarbitalom 5-ethyl-5-(1-ethylbutyl)-barbituric acid telrabarbital tetracainum 2-dimethylaminoethyl 4-butylaminobenzoate tetracaine tetracyclinum 1-dimethylamino-1,4,6,11,12,13,14,18-octahydro-2,5,7,11,14-pentatetracycline hydroxy-4,6-dioxo-11-methylnaphthacene-3-carbonamide thialbarbitalum 5-allyi-5-(2-cyclohexenyl)-2-thiobarbituric acid thialbarbual thiaminum 3-(4-amino-2-methyl-5-pyrimidinylmethyl)-4-methyl-5-(2-hydroxythiamine ethyl)-thiazolium chloride thiopentalum natricum mono-sodium salt of 5-ethyl-5-(I-methylbutyl)-2-thiobarbituric acid thiopental sodium thiotetrabarbitalum 5-ethyl-5-(1-ethylbutyl)-2-thiobarbituric acid thiotetrabarbital tolonii chloridum 3-amino-7-dimethylamino-2-methyl-phenazathionium chloride tolonium chloride tosylchloramidum natricum sodium toluene-4-sulfonylchloramide

tosylchloramide sodium

tretamınım tretamine

tricyclamoli chloridum tricyclamol chloride

tryparsamidum tryparsamide

viomycinum an antibiotic obtained from certain strains of Streptomyces puniceu. viomycin

2,4,6-tri-(ethyleneimino)-s-triazine

(±) 1-(3-cyclohexyl-3-hydroxy-3-phenylpropyl)-1-methylpyrrolidi-

nium chloride

sodium N-phenylglycylamide-4-arsonate

or Streptomyces floridae