
First published in the Government *Gazette*, Electronic Edition, on 30 April 2020 at 5 pm.

No. S 349

**MISUSE OF DRUGS ACT
(CHAPTER 185)**

**MISUSE OF DRUGS ACT
(AMENDMENT OF FIRST, THIRD AND
FIFTH SCHEDULES) ORDER 2020**

In exercise of the powers conferred by sections 58A and 59 of the Misuse of Drugs Act, the Minister for Home Affairs makes the following Order:

Citation and commencement

1. This Order is the Misuse of Drugs Act (Amendment of First, Third and Fifth Schedules) Order 2020 and comes into operation on 1 May 2020.

Amendment of Part I of First Schedule

2. Part I of the First Schedule to the Misuse of Drugs Act is amended —

(a) by deleting item (19) of paragraph 1 and substituting the following item:

“(19) 1-(4-Bromo-2,5-dimethoxyphenyl)propan-2-amine
(also known as
4-Bromo-2,5-dimethoxy- α -methylphenethylamine,
4-Bromo-2,5-dimethoxyamphetamine,
Brolamfetamine or DOB) and its bromo and
dimethoxy positional isomers in the phenyl ring”;

(b) by deleting item (44) of paragraph 1 and substituting the following item:

“(44) 1-(2,5-Dimethoxy-4-methylphenyl)propan-2-amine
(also known as 2,5-Dimethoxy-4,
 α -dimethylphenethylamine, 2-Amino-1-
(2,5-dimethoxy-4-methyl)phenylpropane or DOM)

and its methyl and dimethoxy positional isomers in the phenyl ring”;

- (c) by inserting, immediately after item (8) of paragraph 7, the following item:

“(8AAA) Crotonylfentanyl”;

- (d) by inserting, immediately after item (13) of paragraph 7, the following item:

“(14) Valeryl fentanyl”;

- (e) by deleting item (12A) of paragraph 16 and substituting the following item:

“(12A) 2-(Ethylamino)-1-phenylhexan-1-one (also known as N-Ethylhexedrone or Ethyl-hexedrone)”;

- (f) by inserting, immediately after item (24) of paragraph 16, the following item:

“(24A) 1-Phenyl-2-(pyrrolidin-1-yl)hexan-1-one (also known as Alpha-Pyrrolidinohexiophenone or Alpha-PHP or α -PHP)”;

- (g) by inserting, immediately after item (2) of paragraph 18, the following item:

“(2AAA) 6-Allyl-6-nor-lysergic acid diethylamide (also known as N-Allyl-nor-LSD or AL-LAD) and its acyclic secondary and tertiary amide structural isomers”;

- (h) by inserting, immediately after item (2A) of paragraph 18, the following item:

“(2AAB) N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-butyl-1H-indazole-3-carboxamide (also known as 2-[(1-Butyl-1H-indazol-3-yl)formamido]-3,3-dimethylbutanamide or ADB-BUTINACA) and its N-(1-amino-1-oxohexan-2-yl) isomers”;

- (i) by inserting, immediately after item (2CA) of paragraph 18, the following item:

“(2CB) N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-hydroxybutyl)-1H-indazole-3-carboxamide and its N-(1-amino-1-oxohexan-2-yl) isomers and their

respective hydroxy positional isomers in the butyl group”;

- (j) by deleting item (2G) of paragraph 18 and substituting the following item:

“(2G) N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (also known as N-[1-Amino-3-methyl-1-oxobutan-2-yl]-1-[(4-fluorophenyl)methyl]-1H-indazole-3-carboxamide or AB-FUBINACA) and its N-(1-amino-1-oxopentan-2-yl) and N-(1-amino-2-methyl-1-oxobutan-2-yl) isomers and their respective fluoro positional isomers in the phenyl ring”;

- (k) by inserting, immediately after item (6G) of paragraph 18, the following items:

“(6H) 1-Butanoyl-N,N-diethyllysergamide (also known as 1-Butanoyl-LSD or 1B-LSD) and its acyclic secondary and tertiary amide structural isomers

(6I) 2-(1-Butyl-1H-indazole-3-carboxamido)-3,3-dimethylbutanoic acid and its hexanoic acid isomers”;

- (l) by inserting, immediately after item (8) of paragraph 18, the following item:

“(8AAA) 1-(4-Chloro-2,5-dimethoxyphenyl)propan-2-amine (also known as 4-Chloro-2,5-dimethoxy- α -methylphenethylamine, 4-Chloro-2,5-dimethoxyamphetamine, DOC, 3C-C, 4-Cl-2,5-DMA or 4-Chloro-2,5-DMA) and its chloro and dimethoxy positional isomers in the phenyl ring”;

- (m) by inserting, immediately after item (14) of paragraph 18, the following item:

“(14A) 2-[1-(4,5-Dihydroxypentyl)-1H-indazole-3-carboxamido]-3,3-dimethylbutanoic acid and its hexanoic acid isomers and their respective dihydroxy positional isomers in the pentyl group”;

- (n) by inserting, immediately after item (19AAB) of paragraph 18, the following item: