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MISUSE OF DRUGS ACT
(CHAPTER 185)

MISUSE OF DRUGS (AMENDMENT)
REGULATIONS 2018

In exercise of the powers conferred by section 58(1) of the Misuse of Drugs Act, the Minister for Home Affairs makes the following Regulations:

Citation and commencement

1. These Regulations are the Misuse of Drugs (Amendment) Regulations 2018 and come into operation on 1 May 2018.

Amendment of Second Schedule

2. The Second Schedule to the Misuse of Drugs Regulations (Rg 1) is amended —

- (a) by deleting the item “Remifentanil” in paragraph 1;
- (b) by deleting the item “3-methylthiofentanyl” in paragraph 1;
- (c) by deleting the full-stop at the end of “4-Phenylpiperidine-4-carboxylic acid ethyl ester” in paragraph 1, and by inserting immediately thereafter the following item:

“6,7,8,14-Tetrahydro-4,5- α -epoxy-6-methoxy-17-methylmorphinan-3-ol (also known as 3-O-demethylthebaine or Oripavine).”;

- (d) by inserting, immediately after item (1) of paragraph 10, the following item:

“(1A) Acryloylfentanyl (also known as Acrylfentanyl)”;

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- (e) by inserting, immediately after item (9) of paragraph 10, the following items:
- “(9A) 4-Fluoroisobutyrfentanyl (also known as 4-FIBF or pFIBF)
 - (9B) Furanyl fentanyl”;
- (f) by inserting, immediately after item (11) of paragraph 10, the following items:
- “(11A) 3-Methylthiofentanyl
 - (11B) Ocfentanil”;
- (g) by inserting, immediately after item (12) of paragraph 10, the following item:
- “(12A) Remifentanil”;
- (h) by inserting, immediately after item (13) of paragraph 10, the following item:
- “(13A) Tetrahydrofuranyl fentanyl (also known as THF-F)”.

Amendment of Fourth Schedule

3. The Fourth Schedule to the Misuse of Drugs Regulations is amended —

- (a) by inserting, immediately after item (1) of paragraph 15, the following item:
- “(1AA) 1-(1,3-Benzodioxol-5-yl)-2-(dimethylamino)-1-butanone (also known as Dibutylone or bk-DMBDB)”;
- (b) by inserting, immediately after item (1A) of paragraph 15, the following items:
- “(1AB) 1-(1,3-Benzodioxol-5-yl)-2-(ethylamino)-1-butanone (also known as Eutylone)
 - (1AC) 1-(1,3-Benzodioxol-5-yl)-2-(ethylamino)pentan-1-one (also known as N-Ethylpentylone)”;

(c) by inserting, immediately after item (8A) of paragraph 15, the following item:

“(8AAA) 1-(4-Bromophenyl)-2-(ethylamino)propan-1-one
(also known as 4-Bromoethcathinone or 4-BEC)”;

(d) by inserting, immediately after item (8AA) of paragraph 15, the following item:

“(8AB) 1-(4-Chlorophenyl)-2-(ethylamino)propan-1-one
(also known as 4-Chloroethcathinone or 4-CEC)”;

(e) by inserting, immediately after item (10B) of paragraph 15, the following item:

“(10C) 2-(Ethylamino)-1-(4-methylphenyl)pentan-1-one
(also known as 4-Methyl- α -ethylaminopentiophenone
or 4-MEAPP)”;

(f) by inserting, immediately after item (12) of paragraph 15, the following item:

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

(g) by inserting, immediately after item (15) of paragraph 15, the following item:

“(15AA) 1-(4-Fluorophenyl)-2-(pyrrolidin-1-yl)hexan-1-one
(also known as 4-Fluoro-PHP)”;

(h) by inserting, immediately after item (1) of paragraph 17, the following item:

“(1AA) N-(1-Adamantyl)-1-(4-fluorobenzyl)-1H-indazole-3-
carboxamide (also known as FUB-APINACA or
FUB-AKB48) and its fluoro positional isomers in
the phenyl ring”;

(i) by inserting, immediately after item (2B) of paragraph 17, the following item:

“(2BA) N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-
fluorobenzyl)-1H-indazole-3-carboxamide (also
known as ADB-FUBINACA) and its fluoro
positional isomers in the phenyl ring”;

(j) by inserting, immediately after item (2D) of paragraph 17, the following item:

“(2DA) N-(1-Amino-3,3-dimethyl-1-oxobutan-2-yl)-1-pentyl-1H-indole-3-carboxamide (also known as ADBICA)”;

(k) by deleting items (2E), (2F), (2G), (2H) and (2I) of paragraph 17 and substituting the following items:

“(2E) N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-chloropentyl)-1H-indazole-3-carboxamide (also known as 5-Chloro-AB-PINACA) and its N-(1-amino-1-oxopentan-2-yl) and N-(1-amino-2-methyl-1-oxobutan-2-yl) isomers and their respective chloro positional isomers in the pentyl group

(2F) N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide (also known as AB-CHMINACA) and its N-(1-amino-1-oxopentan-2-yl) and N-(1-amino-2-methyl-1-oxobutan-2-yl) isomers

(2G) N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide (also known as 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

(2H) N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (also known as 5-Fluoro-AB-PINACA) 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 pentyl group

(2I) N-(1-Amino-3-methyl-1-oxobutan-2-yl)-1-pentyl-1H-indazole-3-carboxamide (also known as AB-PINACA) and its N-(1-amino-1-oxopentan-2-yl) and N-(1-amino-2-methyl-1-oxobutan-2-yl) isomers

(2IA) N-(1-Amino-1-oxo-3-phenylpropan-2-yl)-1-(cyclohexylmethyl)-1H-indazole-3-carboxamide (also known as PX-3 or APP-CHMINACA)”;