

**POISONS STANDARD AMENDMENT No. 2 OF 2011**

I, ANTHONY GILL, a delegate of the Secretary to the Department of Health and Ageing for the purposes of paragraph 52D(2)(a) of the *Therapeutic Goods Act 1989* (the Act) and acting in accordance with the Secretary’s power under that paragraph of the Act, hereby amend the Poisons Standard 2010 in the manner set out in Schedule 1.

The amendments to the Poisons Standard 2010 as set out in Schedule 1 commence on

8 July 2011.

*(signed by)*

ANTHONY GILL

Delegate of the Secretary to the Department of Health and Ageing

Dated this 5th day of July 2011

**Schedule 1-Amendments to the Poisons Standard 2010**

STANDARD

FOR THE

UNIFORM SCHEDULING

OF

MEDICINES AND POISONS

No. 1

AMENDMENT No. 3



Effective Date – 8 July 2011

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The amendments listed in this document are a result of a decision made by the Secretary of the Department of Health and Ageing or the Secretary’s delegate. The basis of these amendments can be found in the ‘Reasons for delegate’s final decisions’, which can be accessed from the scheduling website:

[www.tga.gov.au/industry/scheduling-decisions-final.htm](http://www.tga.gov.au/industry/scheduling-decisions-final.htm)

Further inquiries should be directed to:

The Secretary

Medicines and Poisons Scheduling Secretariat

Office of Chemical Safety (MDP 88)

GPO Box 9848

CANBERRA ACT 2601

Or by email: [SMP@health.gov.au](mailto:SMP@health.gov.au)

Media Liaison Unit

Australian Government Department of Health and Ageing

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# Part A – AMENDMENTS TO THE SUSMP No. 1

**Amendments to the Standard for the Uniform Scheduling of Medicines and Poisons**

The Secretary of the Department of Health and Ageing directs that the amendments below be applied to the Standard for the Uniform Scheduling of Medicines and Poisons No. 1 (SUSMP 1) and recommends that these amendments be adopted by the States and Territories with effect from 8 July 2011.

# Part 4 – The Schedules

**SCHEDULE 9 – NEW ENTRIES**

1-(5-FLUOROPENTYL)-3-(2-IODOBENZOYL)INDOLE \*(AM-694).

2-[(1R,3S)-3-HYDROXYCYCLOHEXYL]-5-(2-METHYLNONAN-2-YL)PHENOL \*(Cannabicyclohexanol or CP 47,497 C8 homologue).

2-[(1R,3S)-3-HYDROXYCYCLOHEXYL]- 5-(2-METHYLOCTAN-2-YL)PHENOL \*(CP 47,497).

2-(2-METHOXYPHENYL)-1-(1-PENTYLINDOL-3-YL)ETHANONE \*(JWH – 250).

(1-(2-MORPHOLIN-4-YLETHYL)INDOL-3-YL)-NAPHTHALEN-1-YLMETHANONE \*(JWH – 200).

NAPHTHALEN-1-YL-(1-BUTYLINDOL-3-YL)METHANONE \*(JWH – 073).

1-PENTYL-3-(4-METHYL-1-NAPHTHOYL)INDOLE \*(JWH – 122).

1-PENTYL-3-(1-NAPHTHOYL)INDOLE \*(JWH - 018).