| | 10/20/10 | REVISOR | SGS/DI | RD3986 | | |
|------|--|---------------------------|------------------------|------------|--|--|
| 1.1 | Board of Pharmacy | | | | | |
| 1.2 | Proposed Permanent Rules Relating to | Controlled Substar | ices | | | |
| 1.3 | 6800.4210 SCHEDULE I CONTROLL | LED SUBSTANCES | 5. | | | |
| 1.4 | Schedule I shall consist of the drugs an | nd other substances, l | by whatever official | name, | | |
| 1.5 | common or usual name, chemical name, o | or brand name design | ated, listed in this p | oart. | | |
| 1.6 | [For text of ite | ems A to F, see M.R. | 1 | | | |
| 1.7 | G. Cannabinoids. Unless specifically | y excepted or unless | listed in another | | | |
| 1.8 | schedule, any natural or synthetic materia | l, compound, mixtur | e, or preparation w | hich | | |
| 1.9 | contains any quantity of a substance that i | is a cannabinoid rece | ptor agonist, includ | ing but | | |
| 1.10 | not limited to the following substances an | d their analogs (inclu | uding homologues), | isomers | | |
| 1.11 | (whether optical, positional, or geometric) |), esters, ethers, salts, | , and salts of isomer | s, esters, | | |
| 1.12 | and ethers, whenever the existence of such isomers, esters, ethers, or salts is possible | | | | | |
| 1.13 | within the specific chemical designation: | | | | | |
| 1.14 | (1) <u>1-pentyl-2-methyl-3-(1-naphtl</u> | hoyl)indole (JWH-00 | <u>)7)</u> | | | |
| 1.15 | (2) (2-methyl-1-propyl-1H-indol- | 3-yl)-1-naphthaleny | Imethanone | | | |
| 1.16 | <u>(JWH-015)</u> | | | | | |
| 1.17 | (3) <u>1-pentyl-3-(1-naphthoyl)indol</u> | e (JWH-018) | | | | |
| 1.18 | (4) <u>1-hexyl-3-(naphthalen-1-oyl)i</u> | ndole (JWH-019) | | | | |
| 1.19 | (5) <u>1-butyl-3-(1-naphthoyl)indole</u> | e (JWH-073) | | | | |
| 1.20 | (6) <u>4-methoxynaphthalen-1-yl-(1-</u> | -pentylindol-3-yl)me | thanone (JWH-081) | <u>)</u> | | |
| 1.21 | (7) 4-methoxynaphthalen-1-yl-(1- | -pentyl-2-methylindo | ol-3-yl)methanone | | | |
| 1.22 | <u>(JWH-098)</u> | | | | | |
| 1.23 | (8) (1-(2-morpholin-4-ylethyl)ind | lol-3-yl)-naphthalen- | 1-ylmethanone | | | |
| 1.24 | <u>(JWH-200)</u> | | | | | |

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|------|---|-----------------------|-------------------|--------|--|--|
| 2.1 | (9) <u>7-methoxynaphthalen-1-yl-(1-</u> | pentylindol-3-yl)me | thanone (JWH-164 |) | | |
| 2.2 | (10) 2-(2-chlorophenyl)-1-(1-pent | ylindol-3-yl)ethanon | e (JWH-203) | | | |
| 2.3 | (11) 4-ethylnaphthalen-1-yl-(1-pe | ntylindol-3-yl)metha | none (JWH-210) | | | |
| 2.4 | (12) $2-(2-methoxyphenyl)-1-(1-pethoxyphenyl)$ | entylindol-3-yl)ethan | one (JWH-250) | | | |
| 2.5 | (13) 1-pentyl-3-(4-chloro-1-napht | hoyl)indole (JWH-3 | <u>98)</u> | | | |
| 2.6 | (14) (6aR,10aR)-9-(hydroxymeth | yl)-6, | | | | |
| 2.7 | 6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol) | | | | | |
| 2.8 | <u>(HU-210)</u> | | | | | |
| 2.9 | (15) (dexanabinol, (6aS,10aS)-9- | (hydroxymethyl)-6, | | | | |
| 2.10 | 6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol) | | | | | |
| 2.11 | <u>(HU-211)</u> | | | | | |
| 2.12 | (16) (R)-(+)-[2,3-dihydro-5-methy | yl-3-(4-morpholinylı | nethyl) | | | |
| 2.13 | pyrrolo(1,2,3-de)-1,4-benzoxazin-6-yl]-1- | naphthalenylmethan | one (WIN-55/212-2 | 2) | | |
| 2.14 | (17) 2 -[(1R,3S)-3-hydroxycycloho | exyl]-5-(2-methyloct | an-2-yl)phenol | | | |

2.15 <u>(CP47,497)</u>