Session of 2010

SENATE BILL No. 348

By Senators V. Schmidt, D. Schmidt and Petersen

1-11

| 0. | AN A | CT concerning controlled substances; relating to certain synt | hetio |
|----|---------|--|--------|
| 1 | can | nabinoids schedule I drugs; amending K.S.A. 2009 Supp. 65- | 4105 |
| 2 | | repealing the existing section. | |
| .3 | | 1 0 0 | |
| 4 | Be it e | nacted by the Legislature of the State of Kansas: | |
| 5 | | tion 1. K.S.A. 2009 Supp. 65-4105 is hereby amended to re | ad as |
| 6 | | s: 65-4105. (a) The controlled substances listed in this sectio | |
| 7 | include | ed in schedule I and the number set forth opposite each dr | ug o |
| .8 | | nce is the DEA controlled substances code which has been assi | |
| 9 | to it. | | |
| 20 | (b) | Any of the following opiates, including their isomers, e | sters |
| 21 | ethers, | , salts, and salts of isomers, esters and ethers, unless specif | ically |
| 22 | except | ed, whenever the existence of these isomers, esters, ethers and | l salt |
| 23 | is poss | ible within the specific chemical designation: | |
| 24 | (1) | Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl)-4- | |
| 25 | | piperidinyl]-N-phenylacetamide | 9815 |
| 26 | (2) | Acetylmethadol | 960 |
| 27 | (3) | Allylprodine | 9602 |
| 28 | (4) | Alphacetylmethadol | 9603 |
| 29 | | (except levo-alphacetylmethadol also known as levo-alpha- | |
| 80 | | acetylmethadol, levomethadyl acetate or LAAM) | |
| 31 | (5) | Alphameprodine | 9604 |
| 32 | | Alphamethadol | 9605 |
| 3 | (7) | Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4- | |
| 34 | | piperidyl] propionanilide; 1-(1-methyl-2-phenylethyl)-4- | |
| 35 | | (N-propanilido) piperidine) | 9814 |
| 86 | (8) | Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl) ethyl-4- | |
| 37 | | piperidinyl]-N-phenylpropanamide) | 9832 |
| 88 | | Benzethidine | 9606 |
| 89 | | Betacetylmethadol | 960 |
| 0 | (11) | $Beta-hydroxy fentanyl\ (N-[1-(2-hydroxy-2-phenethyl)-4-piperidinyl]-\\$ | |
| 1 | | N-phenylpropanamide | 9830 |
| 2 | (12) | Beta-hydroxy-3-methylfentanyl (other name: N-[1-(2-hydroxy-2- | |
| 13 | | phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide | 983 |
| | | | |

| 1 | (13) | Betameprodine |
|----|------|---|
| 2 | (14) | Betamethadol |
| 3 | (15) | Betaprodine |
| 4 | (16) | Clonitazene |
| 5 | (17) | Dextromoramide |
| 6 | (18) | Diampromide |
| 7 | (19) | Diethylthiambutene |
| 8 | (20) | Difenoxin |
| 9 | (21) | Dimenoxadol |
| 10 | (22) | Dimepheptanol |
| 11 | (23) | Dimethylthiambutene |
| 12 | (24) | Dioxaphetyl butyrate |
| 13 | (25) | Dipipanone |
| 14 | | Ethylmethylthiambutene |
| 15 | (27) | Etonitazene |
| 16 | (28) | Etoxeridine |
| 17 | (29) | Furethidine |
| 18 | (30) | Hydroxypethidine |
| 19 | (31) | Ketobemidone |
| 20 | (32) | Levomoramide |
| 21 | (33) | Levophenacylmorphan |
| 22 | (34) | 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]-N- |
| 23 | | phenylpropanamide) |
| 24 | (35) | 3-Methylthiofentanyl (N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]- |
| 25 | | N-phenylpropanamide) |
| 26 | (36) | Morpheridine |
| 27 | (37) | MPPP (1-methyl-4-phenyl-4-propionoxypiperidine) |
| 28 | (38) | Noracymethadol |
| 29 | | Norlevorphanol |
| 30 | (40) | Normethadone |
| 31 | (41) | Norpipanone |
| 32 | | Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4- |
| 33 | | piperidinyl] propanamide |
| 34 | (43) | PEPAP (1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine) |
| 35 | | Phenadoxone |
| 36 | | Phenampromide |
| 37 | | Phenomorphan |
| 38 | | Phenoperidine |
| 39 | | Piritramide |
| 40 | . , | Proheptazine |
| 41 | | Properidine |
| 42 | | Propiram |
| 13 | | Pagamaranida |

| 1 | (53) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]- | |
|----|---|---------|
| 2 | propanamide | 9835 |
| 3 | (54) Tilidine | 9750 |
| 4 | (55) Trimeperidine | 9646 |
| 5 | (c) Any of the following opium derivatives, their salts, isome | rs and |
| 6 | salts of isomers, unless specifically excepted, whenever the existe | nce o |
| 7 | these salts, isomers and salts of isomers is possible within the s | pecific |
| 8 | chemical designation: | - |
| 9 | (1) Acetorphine | 9319 |
| 10 | (2) Acetyldihydrocodeine | 905] |
| 11 | (3) Benzylmorphine | 9052 |
| 12 | (4) Codeine methylbromide | 9070 |
| 13 | (5) Codeine-N-Oxide | 9053 |
| 14 | (6) Cyprenorphine | 9054 |
| 15 | (7) Desomorphine | 9055 |
| 16 | (8) Dihydromorphine | 914 |
| 17 | (9) Drotebanol | 933 |
| 18 | (10) Etorphine (except hydrochloride salt) | 9050 |
| 19 | (11) Heroin | 9200 |
| 20 | (12) Hydromorphinol | 930 |
| 21 | (13) Methyldesorphine | 930 |
| 22 | (14) Methyldihydromorphine | 9304 |
| 23 | (15) Morphine methylbromide | 9305 |
| 24 | (16) Morphine methylsulfonate | 9300 |
| 25 | (17) Morphine-N-Oxide | 930 |
| 26 | (18) Myrophine | 930 |
| 27 | (19) Nicocodeine | 930 |
| 28 | (20) Nicomorphine | 931 |
| 29 | (21) Normorphine | 931 |
| 30 | (22) Pholcodine | 931 |
| 31 | (23) Thebacon | 931 |
| 32 | (d) Any material, compound, mixture or preparation which co | ontain |
| 33 | any quantity of the following hallucinogenic substances, their sale | |
| 34 | mers and salts of isomers, unless specifically excepted, whenever t | |
| 35 | istence of these salts, isomers and salts of isomers is possible with | |
| 36 | specific chemical designation: | |
| 37 | (1) 4-bromo-2,5-dimethoxy-amphetamine | 739 |
| 38 | Some trade or other names: 4-bromo-2,5-dimethoxy-alpha- | |
| 39 | methylphenethylamine; 4-bromo-2,5-DMA. | |
| 40 | (2) 2,5-dimethoxyamphetamine | 739 |
| 41 | Some trade or other names: 2,5-dimethoxy-alpha-methyl- | |
| 42 | phenethylamine; 2,5-DMA. | |
| 43 | (3) 4-methovyamphetamine | 741 |

| 1 | | Some trade or other names: 4-methoxy-alpha-methylphene-thylamine; |
|---|------|--|
| 2 | | paramethoxyamphetamine; PMA. |
| 3 | (4) | 5-methoxy-3,4-methylenedioxy-amphetamine |
| 4 | (5) | 4-methyl-2,5-dimethoxy-amphetamine |
| 5 | | Some trade or other names: 4-methyl-2,5-dimethoxy-alpha- |
| 6 | | methylphenethylamine; "DOM"; and "STP". |
| 7 | (6) | 3,4-methylenedioxy amphetamine |
| 8 | (7) | 3,4-methylenedioxymethamphetamine (MDMA) |
| 9 | (8) | 3,4-methylenedioxy-N-ethylamphetamine (also known as N-ethyl- |
| 0 | | alpha-methyl-3,4 (methylenedioxy) phenethylamine, N-ethyl MDA, |
| 1 | | MDE, and MDEA) |
| 2 | (9) | N-hydroxy-3,4-methylenedioxyamphetamine (also known as N-hydroxy-1,0) and the state of the sta |
| 3 | | alpha-methyl-3,4(methylenedioxy) phenethylamine, and N-hydroxy |
| 4 | | MDA) |
| 5 | (10) | 3,4,5-trimethoxy amphetamine |
| 3 | (11) | Bufotenine |
| 7 | | Some trade or other names: 3-(Beta-Dimethyl-aminoethyl)-5- |
| 8 | | hydroxyindole; 3-(2-dimethyl-aminoethyl)-5-indolol; N,N- |
| 9 | | dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine. |
| 0 | (12) | Diethyltryptamine |
| L | , , | Some trade or other names: N,N-Diethyltryptamine; DET. |
| 2 | (13) | Dimethyltryptamine |
| 3 | , , | Some trade or other names: DMT. |
| 1 | (14) | Ibogaine |
| 5 | , , | Some trade or other names: 7-Ethyl-6,6 Beta,7,8,9,10,12,13-octahydro- |
| 3 | | 2-methoxy-6,9-methano-5H-pyrido[1',2':1,2] azepino [5,4-b]indole; |
| 7 | | Tabernanthe iboga. |
| 3 | (15) | Lysergic acid diethylamide |
|) | | Marihuana |
|) | , , | Mescaline |
| L | , , | Parahexyl |
| 2 | (/ | Some trade or other names: 3-Hexyl-l-hydroxy-7,8,9,10-tetrahydro- |
| 3 | | 6,6,9-trimethyl-6H-dibenzo[b,d]pyran; Synhexyl. |
| 1 | (19) | Peyote |
| 5 | (/ | Meaning all parts of the plant presently classified botanically as Lopho- |
| 3 | | phora williamsii Lemaire, whether growing or not, the seeds thereof, |
| 7 | | any extract from any part of such plant, and every compound, manu- |
| 3 | | facture, salts, derivative, mixture or preparation of such plant, its seeds |
| 9 | | or extracts. |
| 0 | (20) | N-ethyl-3-piperidyl benzilate |
| 1 | | N-methyl-3-piperidyl benzilate |
| 2 | | Psilocybin |
| 2 | | Deiloom |

| 1 | (24) | Tetrahydrocannabinols | 7370 |
|----------|-------|--|------|
| 2 | | Synthetic equivalents of the substances contained in the plant, or in the | |
| 3 | | resinous extractives of Cannabis, sp. and/or synthetic substances, deriv- | |
| 4 | | atives, and their isomers with similar chemical structure and pharma- | |
| 5 | | cological activity such as the following: Delta 1 cis or trans tetrahydro- | |
| 6 | | cannabinol, and their optical isomers Delta 6 cis or trans | |
| 7 | | tetrahydrocannabinol, and their optical isomers Delta 3,4 cis or trans | |
| 8 | | tetrahydrocannabinol, and its optical isomers (Since nomenclature of | |
| 9 | | $these \ substances \ is \ not \ internationally \ standardized, \ compounds \ of \ these$ | |
| 10 | | structures, regardless of numerical designation of atomic positions cov- | |
| 11 | | ered.) | |
| 12 | (25) | Ethylamine analog of phencyclidine | 7455 |
| 13 | | Some trade or other names: N-ethyl-1-phenyl-cyclo-hexylamine; | |
| 14 | | (1-phenylcyclohexyl) ethylamine; N-(1-phenylcyclohexyl) ethylamine; | |
| 15 | | cyclohexamine; PCE. | |
| 16 | (26) | Pyrrolidine analog of phencyclidine | 7458 |
| 17 | | Some trade or other names: 1-(1-phenylcyclo-hexyl)-pyrrolidine; PCPy; | |
| 18 | | PHP. | |
| 19 | (27) | Thiophene analog of phencyclidine | 7470 |
| 20 | | Some trade or other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine; 2- $$ | |
| 21 | | thienylanalog of phencyclidine; TPCP; TCP. | |
| 22 | (28) | 1-[1-(2-thienyl)-cyclohexyl] pyrrolidine | 7473 |
| 23 | | Some other names: TCPy. | |
| 24 | (29) | 2,5-dimethoxy-4-ethylamphetamine | 7399 |
| 25 | | Some trade or other names: DOET. | |
| 26 | (30) | Salvia divinorum or salvinorum A; all parts of the plant presently clas- | |
| 27 | | sified botanically as salvia divinorum, whether growing or not, the seeds | |
| 28 | | thereof, any extract from any part of such plant, and every compound, | |
| 29 | | manufacture, salts, derivative, mixture or preparation of such plant, its | |
| 30 | | seeds or extracts. | |
| 31 | (31) | Datura stramonium, commonly known as gypsum weed or jimson weed; | |
| 32 | | all parts of the plant presently classified botanically as datura stramo- | |
| 33 | | nium, whether growing or not, the seeds thereof, any extract from any | |
| 34 | | part of such plant, and every compound, manufacture, salts, derivative, | |
| 35 | | mixture or preparation of such plant, its seeds or extracts. | |
| 36 | (32) | (6aR,10aR)-9-(hydroxymethyl)-6, 6-dimethyl-3-(2-methyloctan-2-yl)- | |
| 37 | | 6a,7,10,10a-tetrahydrobenzo[c]chromen-1-ol | 7370 |
| 38 | | Some trade or other names: HU-210. | |
| 39 | (33) | 1-Pentyl-3-(1-naphthoyl)indole | |
| 40 | y = | Some trade or other names: JWH-018. | |
| 41 | (34) | 1-Butyl-3-(1-naphthoyl)indole | |
| 42 42 | / | Some trade or other names: JWH-073. | 7402 |
| | /·>=\ | N. Is assertantes asserting a | 7409 |

| 1 | Some trade or other names: BZP. |
|----|---|
| 2 | (e) Any material, compound, mixture or preparation which contains |
| 3 | any quantity of the following substances having a depressant effect on the |
| 4 | central nervous system, including its salts, isomers, and salts of isomers |
| 5 | whenever the existence of such salts, isomers, and salts of isomers is |
| 6 | possible within the specific chemical designation: |
| 7 | (1) Mecloqualone |
| 8 | (2) Methaqualone |
| 9 | (3) Gamma hydroxybutyric acid |
| 10 | (f) Unless specifically excepted or unless listed in another schedule |
| 11 | any material, compound, mixture or preparation which contains any quan- |
| 12 | tity of the following substances having a stimulant effect on the central |
| 13 | nervous system, including its salts, isomers and salts of isomers: |
| 14 | (1) Fenethylline |
| 15 | (2) N-ethylamphetamine |
| 16 | (3) (+)cis-4-methylaminorex ((+)cis-4,5-dihydro-4-methyl-5-phenyl-2- |
| 17 | oxazolamine |
| 18 | (4) N,N-dimethylamphetamine (also known as N,N-alpha-trimethyl- |
| 19 | benzeneethanamine; N,N-alpha-trimethylphenethylamine) |
| 20 | (5) Cathinone (some other names: 2-amino-1-phenol-1-propanone, alpha- |
| 21 | amino propiophenone, 2-amino propiophenone and norphedrone) 1235 |
| 22 | (g) Any material, compound, mixture or preparation which contains |
| 23 | any quantity of the following substances: |
| 24 | (1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its |
| 25 | optical isomers, salts and salts of isomers |
| 26 | (2) N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide |
| 27 | (thenylfentanyl), its optical isomers, salts and salts of isomers |
| 28 | (3) Methcathinone (some other names: 2-methylamino-1-phenylpropan-1- |
| 29 | one: Ephedrone: Monomethylpropion: UR 1431, its salts, optical iso- |
| 30 | mers and salts of optical isomers) |
| 31 | (4) Aminorex (some other names: Aminoxaphen 2-amino-5-phenyl-2- |
| 32 | oxazoline or 4,5-dihydro-5-phenyl-2-oxazolamine, its salts, optical iso- |
| 33 | mers and salts of optical isomers) |
| 34 | (5) Alpha-ethyltryptamine, its optical isomers, salts and salts of isomers 7248 |
| 35 | Some other names: etryptamine, alpha-methyl-1H-indole-3- |
| 36 | ethanamine; 3-(2-aminobutyl) indole. |
| 37 | Sec. 2. K.S.A. 2009 Supp. 65-4105 is hereby repealed. |
| 38 | Sec. 3. This act shall take effect and be in force from and after its |
| 30 | nublication in the statute book Kanage register |