	CONTROLLED SUBSTANCES AMENDMENTS
	2024 GENERAL SESSION
	STATE OF UTAH
	Chief Sponsor: Jennifer Dailey-Provost
	Senate Sponsor: Evan J. Vickers
L	ONG TITLE
G	eneral Description:
	This bill adds gabapentin to the list of controlled substances.
H	ighlighted Provisions:
	This bill:
	<ul> <li>adds gabapentin to Schedule V of the list of controlled substances; and</li> </ul>
	<ul> <li>makes technical and conforming changes.</li> </ul>
M	Ioney Appropriated in this Bill:
	None
0	ther Special Clauses:
	None
U	tah Code Sections Affected:
A	MENDS:
	58-37-4, as last amended by Laws of Utah 2022, Chapter 165
Be	e it enacted by the Legislature of the state of Utah:
	Section 1. Section <b>58-37-4</b> is amended to read:
	58-37-4. Schedules of controlled substances Schedules I through V Findings
re	equired Specific substances included in schedules.
	(1) There are established five schedules of controlled substances known as Schedules I,
II,	, III, IV, and V which consist of substances listed in this section.

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<ul> <li>the official name, common or usual name, chemical name, or brand name designated:</li> <li>(a) Schedule I:</li> <li>(i) Unless specifically excepted or unless listed in another schedule, any of the</li> <li>following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and</li> <li>ethers, when the existence of the isomers, esters, ethers, and salts is possible within the specific</li> <li>chemical designation:</li> <li>(A) Acetyl-alpha-methylfentanyl</li> <li>(N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide);</li> <li>(B) Acetyl fentanyl: (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(C) Acetylmethadol;</li> <li>(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(E) Allylprodine;</li> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(M) Betacetylmethadol;</li> <li>(M) Betacetylmethadol;</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>phenethyl-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	28	(2) Schedules I, II, III, IV, and V consist of the following drugs or other substances by
<ul> <li>(i) Unless specifically excepted or unless listed in another schedule, any of the</li> <li>following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and</li> <li>ethers, when the existence of the isomers, esters, ethers, and salts is possible within the specific</li> <li>chemical designation:</li> <li>(A) Acctyl-alpha-methylfentanyl</li> <li>(N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide);</li> <li>(C) Acetyl fentanyl: (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(C) Acetyl-methadol;</li> <li>(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(E) Allylprodine;</li> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alphamethylfontanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthofentanyl (N-[1-(alpha-methyl-2-(2-thienyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthofentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthofentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthofentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(I) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	29	the official name, common or usual name, chemical name, or brand name designated:
32following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and33ethers, when the existence of the isomers, esters, ethers, and salts is possible within the specific34chemical designation:35(A) Acctyl-alpha-methylfentanyl36(N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide);37(B) Acctyl fentanyl: (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);38(C) Acctylmethadol;39(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacetamide);40(E) Allylprodine;41(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as42levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;43(G) Alphameprodine;44(H) Alphamethadol;45(I) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]46projonanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);47(J) Alpha-methylthiofentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]48piperidinyl]-N-phenylpropanamide);49(K) Benzylpiperazine;50(L) Benzethidine;51(M) Betacetylmethadol;52(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-53piperidinyl]-N-phenylpropanamide);54(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-55phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;56(P) Betameprodine;	30	(a) Schedule I:
<ul> <li>ethers, when the existence of the isomers, esters, ethers, and salts is possible within the specific</li> <li>chemical designation:</li> <li>(A) Acetyl-alpha-methylfentanyl</li> <li>(N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide);</li> <li>(B) Acetyl fentanyl: (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(C) Acetylmethadol;</li> <li>(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(E) Allylprodine;</li> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidyl]</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(L) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(O) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	31	(i) Unless specifically excepted or unless listed in another schedule, any of the
<ul> <li>chemical designation:</li> <li>(A) Acetyl-alpha-methylfentanyl</li> <li>(N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide);</li> <li>(B) Acetyl fentanyl: (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(C) Acetylmethadol;</li> <li>(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(E) Allylprodine;</li> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidyl]</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(I) Benzethidine;</li> <li>(I) Benzethidin</li></ul>	32	following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and
<ul> <li>(A) Acetyl-alpha-methylfentanyl</li> <li>(N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide);</li> <li>(B) Acetyl fentanyl: (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(C) Acetylmethadol;</li> <li>(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacetylamide);</li> <li>(E) Allylprodine;</li> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidyl]</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(I) Betaetylmethadol;</li> <li>(I) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	33	ethers, when the existence of the isomers, esters, ethers, and salts is possible within the specific
<ul> <li>(N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide);</li> <li>(B) Acetyl fentanyl: (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(C) Acetylmethadol;</li> <li>(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacrylamide);</li> <li>(E) Allylprodine;</li> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alphamethylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-piperidyl]</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(L) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(O) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	34	chemical designation:
<ul> <li>(B) Acetyl fentanyl: (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);</li> <li>(C) Acetylmethadol;</li> <li>(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacrylamide);</li> <li>(E) Allylprodine;</li> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alpharmethylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(I) Alpha-methylthofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(M) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	35	(A) Acetyl-alpha-methylfentanyl
<ul> <li>(C) Acetylmethadol;</li> <li>(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacrylamide);</li> <li>(E) Allylprodine;</li> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alphamethadol;</li> <li>(I) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-(methyl-2-(2-thienyl)ethyl-4-piperidyl]</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(L) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-piperidyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	36	(N-[1-(1-methyl-2-phenethyl)-4-piperidinyl]-N-phenylacetamide);
<ul> <li>(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacrylamide);</li> <li>(E) Allylprodine;</li> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alphamethadol;</li> <li>(I) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(I) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(I) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	37	(B) Acetyl fentanyl: (N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide);
<ul> <li>(E) Allylprodine;</li> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alphamethadol;</li> <li>(I) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(I) Betacetylmethadol;</li> <li>(N) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	38	(C) Acetylmethadol;
<ul> <li>(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as</li> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alphamethadol;</li> <li>(I) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(L) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	39	(D) Acryl fentanyl (N-(1-Phenethylpiperidin-4-yl)-N-phenylacrylamide);
<ul> <li>levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;</li> <li>(G) Alphameprodine;</li> <li>(H) Alphamethadol;</li> <li>(I) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(L) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	40	(E) Allylprodine;
<ul> <li>43 (G) Alphameprodine;</li> <li>44 (H) Alphamethadol;</li> <li>45 (I) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>46 propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>47 (J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>48 piperidinyl]-N-phenylpropanamide);</li> <li>49 (K) Benzylpiperazine;</li> <li>50 (L) Benzethidine;</li> <li>51 (M) Betacetylmethadol;</li> <li>52 (N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>53 piperidinyl]-N-phenylpropanamide);</li> <li>54 (O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>55 phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>56 (P) Betameprodine;</li> </ul>	41	(F) Alphacetylmethadol, except levo-alphacetylmethadol also known as
<ul> <li>(H) Alphamethadol;</li> <li>(I) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(L) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	42	levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM;
<ul> <li>(I) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]</li> <li>propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(L) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	43	(G) Alphameprodine;
<ul> <li>46 propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);</li> <li>47 (J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>48 piperidinyl]-N-phenylpropanamide);</li> <li>49 (K) Benzylpiperazine;</li> <li>50 (L) Benzethidine;</li> <li>51 (M) Betacetylmethadol;</li> <li>52 (N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>53 piperidinyl]-N-phenylpropanamide);</li> <li>54 (O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>55 phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>56 (P) Betameprodine;</li> </ul>	44	(H) Alphamethadol;
<ul> <li>(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(K) Benzylpiperazine;</li> <li>(L) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	45	(I) Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl)ethyl-4-piperidyl]
<ul> <li>48 piperidinyl]-N-phenylpropanamide);</li> <li>49 (K) Benzylpiperazine;</li> <li>50 (L) Benzethidine;</li> <li>51 (M) Betacetylmethadol;</li> <li>52 (N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>53 piperidinyl]-N-phenylpropanamide);</li> <li>54 (O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>55 phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>56 (P) Betameprodine;</li> </ul>	46	propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-propanilido) piperidine);
<ul> <li>49 (K) Benzylpiperazine;</li> <li>50 (L) Benzethidine;</li> <li>51 (M) Betacetylmethadol;</li> <li>52 (N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>53 piperidinyl]-N-phenylpropanamide);</li> <li>54 (O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>55 phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>56 (P) Betameprodine;</li> </ul>	47	(J) Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl)ethyl-4-
<ul> <li>(L) Benzethidine;</li> <li>(M) Betacetylmethadol;</li> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	48	piperidinyl]-N-phenylpropanamide);
<ul> <li>51 (M) Betacetylmethadol;</li> <li>52 (N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>53 piperidinyl]-N-phenylpropanamide);</li> <li>54 (O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>55 phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>56 (P) Betameprodine;</li> </ul>	49	(K) Benzylpiperazine;
<ul> <li>(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-</li> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	50	(L) Benzethidine;
<ul> <li>piperidinyl]-N-phenylpropanamide);</li> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	51	(M) Betacetylmethadol;
<ul> <li>(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-</li> <li>phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>(P) Betameprodine;</li> </ul>	52	(N) Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl)-4-
<ul> <li>55 phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;</li> <li>56 (P) Betameprodine;</li> </ul>	53	piperidinyl]-N-phenylpropanamide);
56 (P) Betameprodine;	54	(O) Beta-hydroxy-3-methylfentanyl, other name: N-[1-(2-hydroxy-2-
	55	phenethyl)-3-methyl-4-piperidinyl]-N-phenylpropanamide;
57 (O) Betamethadol:	56	(P) Betameprodine;
	57	(Q) Betamethadol;
58 (R) Betaprodine;	58	(R) Betaprodine;

59	(S) Butyryl fentanyl (N-(1-(2-phenylethyl)-4-piperidinyl)-N-phenylbutyramide);
60	(T) Clonitazene;
61	(U) Cyclopropyl fentanyl
62	(N-(1-Phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide);
63	(V) Dextromoramide;
64	(W) Diampromide;
65	(X) Diethylthiambutene;
66	(Y) Difenoxin;
67	(Z) Dimenoxadol;
68	(AA) Dimepheptanol;
69	(BB) Dimethylthiambutene;
70	(CC) Dioxaphetyl butyrate;
71	(DD) Dipipanone;
72	(EE) Ethylmethylthiambutene;
73	(FF) Etizolam
74	(1-Methyl-6-o-chlorophenyl-8-ethyl-4H-s-triazolo[3,4-c]thieno[2,3-e]1,4-diazepine);
75	(GG) Etonitazene;
76	(HH) Etoxeridine;
77	(II) Furanyl fentanyl (N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]
78	furan-2-carboxamide);
79	(JJ) Furethidine;
80	(KK) Hydroxypethidine;
81	(LL) Ketobemidone;
82	(MM) Levomoramide;
83	(NN) Levophenacylmorphan;
84	(OO) Methoxyacetyl fentanyl
85	(2-Methoxy-N-(1-phenylethylpiperidinyl-4-yl)-N-acetamide);
86	(PP) Morpheridine;
87	(QQ) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
88	(RR) Noracymethadol;
89	(SS) Norlevorphanol;

90	(TT) Normethadone;
91	(UU) Norpipanone;
92	(VV) Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4- piperidinyl]
93	propanamide);
94	(WW) Para-fluoroisobutyryl fentanyl
95	(N-(4-Fluorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide);
96	(XX) PEPAP (1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);
97	(YY) Phenadoxone;
98	(ZZ) Phenampromide;
99	(AAA) Phenomorphan;
100	(BBB) Phenoperidine;
101	(CCC) Piritramide;
102	(DDD) Proheptazine;
103	(EEE) Properidine;
104	(FFF) Propiram;
105	(GGG) Racemoramide;
106	(HHH) Tetrahydrofuran fentanyl
107	(N-(1-Phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-carboxamide);
108	(III) Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]- propanamide;
109	(JJJ) Tilidine;
110	(KKK) Trimeperidine;
111	(LLL) 3-methylfentanyl, including the optical and geometric isomers
112	(N-[3-methyl-1-(2-phenylethyl)-4-piperidyl]- N-phenylpropanamide);
113	(MMM) 3-methylthiofentanyl
114	(N-[(3-methyl-1-(2-thienyl)ethyl-4-piperidinyl]-N-phenylpropanamide);
115	(NNN) 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide also
116	known as U-47700; and
117	(OOO) 4-cyano CUMYL-BUTINACA.
118	(ii) Unless specifically excepted or unless listed in another schedule, any of the
119	following opium derivatives, their salts, isomers, and salts of isomers when the existence of the
120	salts, isomers, and salts of isomers is possible within the specific chemical designation:

121	$(\Lambda)$ A setembine.
	<ul><li>(A) Acetorphine;</li><li>(B) A setuldibudge of deine;</li></ul>
122	(B) Acetyldihydrocodeine;
123	(C) Benzylmorphine;
124	<ul><li>(D) Codeine methylbromide;</li><li>(E) Codeine N. Codeine</li></ul>
125	(E) Codeine-N-Oxide;
126	(F) Cyprenorphine;
127	(G) Desomorphine;
128	(H) Dihydromorphine;
129	(I) Drotebanol;
130	(J) Etorphine (except hydrochloride salt);
131	(K) Heroin;
132	(L) Hydromorphinol;
133	(M) Methyldesorphine;
134	(N) Methylhydromorphine;
135	(O) Morphine methylbromide;
136	(P) Morphine methylsulfonate;
137	(Q) Morphine-N-Oxide;
138	(R) Myrophine;
139	(S) Nicocodeine;
140	(T) Nicomorphine;
141	(U) Normorphine;
142	(V) Pholcodine; and
143	(W) Thebacon.
144	(iii) Unless specifically excepted or unless listed in another schedule, any material,
145	compound, mixture, or preparation which contains any quantity of the following hallucinogenic
146	substances, or which contains any of their salts, isomers, and salts of isomers when the
147	existence of the salts, isomers, and salts of isomers is possible within the specific chemical
148	designation; as used in this Subsection (2)(a)(iii) only, "isomer" includes the optical, position,
149	and geometric isomers:
150	(A) Alpha-ethyltryptamine, some trade or other names: etryptamine; Monase;
151	$\alpha$ -ethyl-1H-indole-3-ethanamine; 3-(2-aminobutyl) indole; $\alpha$ -ET; and AET;

152	(B) 4-bromo-2,5-dimethoxy-amphetamine, some trade or other names:
153	4-bromo-2,5-dimethoxy-α-methylphenethylamine; 4-bromo-2,5-DMA;
154	(C) 4-bromo-2,5-dimethoxyphenethylamine, some trade or other names:
155	2-(4-bromo-2,5-dimethoxyphenyl)-1-aminoethane; alpha-desmethyl DOB; 2C-B, Nexus;
156	(D) 2,5-dimethoxyamphetamine, some trade or other names:
157	2,5-dimethoxy-α-methylphenethylamine; 2,5-DMA;
158	(E) 2,5-dimethoxy-4-ethylamphetamine, some trade or other names: DOET;
159	(F) 4-methoxyamphetamine, some trade or other names:
160	4-methoxy-α-methylphenethylamine; paramethoxyamphetamine, PMA;
161	(G) 5-methoxy-3,4-methylenedioxyamphetamine;
162	(H) 4-methyl-2,5-dimethoxy-amphetamine, some trade and other names:
163	4-methyl-2,5-dimethoxy-α-methylphenethylamine; "DOM"; and "STP";
164	(I) 3,4-methylenedioxy amphetamine;
165	(J) 3,4-methylenedioxymethamphetamine (MDMA);
166	(K) 3,4-methylenedioxy-N-ethylamphetamine, also known as N-ethyl-
167	alpha-methyl-3,4(methylenedioxy)phenethylamine, N-ethyl MDA, MDE, MDEA;
168	(L) N-hydroxy-3,4-methylenedioxyamphetamine, also known as
169	N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine, and N-hydroxy MDA;
170	(M) 3,4,5-trimethoxy amphetamine;
171	(N) Bufotenine, some trade and other names:
172	3-(β-Dimethylaminoethyl)-5-hydroxyindole; 3-(2-dimethylaminoethyl)-5-indolol; N,
173	N-dimethylserotonin; 5-hydroxy-N,N-dimethyltryptamine; mappine;
174	(O) Diethyltryptamine, some trade and other names: N,N-Diethyltryptamine; DET;
175	(P) Dimethyltryptamine, some trade or other names: DMT;
176	(Q) Ibogaine, some trade and other names:
177	7-Ethyl-6,6β,7,8,9,10,12,13-octahydro-2-methoxy-6,9-methano-5H-pyrido [1', 2':1,2] azepino
178	[5,4-b] indole; Tabernanthe iboga;
179	(R) Lysergic acid diethylamide;
180	(S) Marijuana;
181	(T) Mescaline;
182	(U) Parahexyl, some trade or other names:

184(V) Peyote, meaning all parts of the plant presently classified botanically as185Lophophora williamsii Lemaire, whether growing or not, the seeds thereof, any extract from186any part of such plant, and every compound, manufacture, salts, derivative, mixture, or187preparation of such plant, its seeds or extracts (Interprets 21 USC 812(c), Schedule I(c) (12));188(W) N-ethyl-3-piperidyl benzilate;199(Y) Psilocybin;191(Z) Psilocyn;192(AA) Tetrahydrocannabinols, naturally contained in a plant of the genus Cannabis193(cannabis plant), except for marijuana as defined in Subsection 58-37-2(1)(aa)(i)(E), as well as194synthetic equivalents of the substances contained in the cannabis plant, or in the resinous195extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with196similar chemical structure and pharmacological activity to those substances contained in the197plant, such as the following: $\Delta 1$ cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 3$ , 4 cis or trans198tetrahydrocannabinol, and their optical isomers $\Delta 3$ , 4 cis or trans199tetrahydrocannabinol, and their optical isomers (CC) Pyrrolidine analog of phencyclidine, some trade or other names:1061-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;107(DD) Thiophene analog of phencyclidine, some trade or other names:1081-(1-(2-thienyl)-cyclohexyl)-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and109(EE) 1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, any material <th>183</th> <th>3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran; Synhexyl;</th>	183	3-Hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran; Synhexyl;
186any part of such plant, and every compound, manufacture, salts, derivative, mixture, or187preparation of such plant, its seeds or extracts (Interprets 21 USC 812(c), Schedule I(c) (12));188(W) N-ethyl-3-piperidyl benzilate;190(Y) Psilocybin;191(Z) Psilocyp;192(AA) Tetrahydrocannabinols, naturally contained in a plant of the genus Cannabis193(cannabis plant), except for marijuana as defined in Subsection 58-37-2(1)(aa)(i)(E), as well as194synthetic equivalents of the substances contained in the cannabis plant, or in the resinous195extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with196similar chemical structure and pharmacological activity to those substances contained in the197plant, such as the following: A1 cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 6$ 198cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 3.4$ cis or trans199tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not201intermationally standardized, compounds of these structures, regardless of numerical202(BB) Ethylamine analog of phencyclidine, some trade or other names:203N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;204(CC) Pyrrolidine analog of phencyclidine, some trade or other names:205(CC) Pyrrolidine, PCPy, PHP;207(DD) Thiophene analog of phencyclidine, some trade or other names:2081-[1-(2-thienyl)-cyclohexyl]-pyrrolidine, some other names: TCPy.209(EE) 1-[1-(2-	184	(V) Peyote, meaning all parts of the plant presently classified botanically as
187preparation of such plant, its seeds or extracts (Interprets 21 USC 812(c), Schedule I(c) (12));188(W) N-ethyl-3-piperidyl benzilate;189(X) N-methyl-3-piperidyl benzilate;190(Y) Psilocybin;191(Z) Psilocyn;192(AA) Tetrahydrocannabinols, naturally contained in a plant of the genus Cannabis193(cannabis plant), except for marijuana as defined in Subsection 58-37-2(1)(aa)(i)(E), as well as194synthetic equivalents of the substances contained in the cannabis plant, or in the resinous195extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with196similar chemical structure and pharmacological activity to those substances contained in the197plant, such as the following: A1 cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 6$ 198cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 3.4$ cis or trans199tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not100internationally standardized, compounds of these structures, regardless of numerical101designation of atomic positions covered;102(CC) Pytrolidine analog of phencyclidine, some trade or other names:103N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;104(DD) Thiophene analog of phencyclidine, some trade or other names:1051-(1-phenylcyclohexyl)-pytrolidine, PCPy, PHP;106(DD) Thiophene analog of phencyclidine, some trade or other names:1081-[1-(2-thienyl)-cyclohexyl]-pytrolidine, some other names: TCPy. <td>185</td> <td>Lophophora williamsii Lemaire, whether growing or not, the seeds thereof, any extract from</td>	185	Lophophora williamsii Lemaire, whether growing or not, the seeds thereof, any extract from
188(W) N-ethyl-3-piperidyl benzilate;189(X) N-methyl-3-piperidyl benzilate;190(Y) Psilocybin;191(Z) Psilocyn;192(AA) Tetrahydrocannabinols, naturally contained in a plant of the genus Cannabis193(cannabis plant), except for marijuana as defined in Subsection $58-37-2(1)(aa)(i)(E)$ , as well as194synthetic equivalents of the substances contained in the cannabis plant, or in the resinous195extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with196similar chemical structure and pharmacological activity to those substances contained in the197plant, such as the following: $\Delta 1$ cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 3, 4$ cis or trans198tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not100intermationally standardized, compounds of these structures, regardless of numerical201designation of atomic positions covered;202(BB) Ethylamine analog of phencyclidine, some trade or other names:203N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,204N-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;205(CC) Pyrrolidine analog of phencyclidine, some trade or other names:2061-(1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and209(EE) 1-[1-(2-thienyl)-cyclohexyl]pyrrolidine, some other names:201compound, mixture, or preparation which contains any quantity of the following substances	186	any part of such plant, and every compound, manufacture, salts, derivative, mixture, or
<ul> <li>(X) N-methyl-3-piperidyl benzilate;</li> <li>(Y) Psilocybin;</li> <li>(Z) Psilocypi;</li> <li>(AA) Tetrahydrocannabinols, naturally contained in a plant of the genus Cannabis</li> <li>(cannabis plant), except for marijuana as defined in Subsection 58-37-2(1)(aa)(i)(E), as well as</li> <li>synthetic equivalents of the substances contained in the cannabis plant, or in the resinous</li> <li>extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with</li> <li>similar chemical structure and pharmacological activity to those substances contained in the</li> <li>plant, such as the following: Δ1 cis or trans tetrahydrocannabinol, and their optical isomers Δ6</li> <li>cis or trans tetrahydrocannabinol, and their optical isomers Δ3,4 cis or trans</li> <li>tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not</li> <li>internationally standardized, compounds of these structures, regardless of numerical</li> <li>designation of atomic positions covered;</li> <li>(BB) Ethylamine analog of phencyclidine, some trade or other names:</li> <li>N-ethyl-1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)-cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	187	preparation of such plant, its seeds or extracts (Interprets 21 USC 812(c), Schedule I(c) (12));
<ul> <li>(Y) Psilocybin;</li> <li>(Z) Psilocyn;</li> <li>(AA) Tetrahydrocannabinols, naturally contained in a plant of the genus Cannabis</li> <li>(cannabis plant), except for marijuana as defined in Subsection 58-37-2(1)(aa)(i)(E), as well as</li> <li>synthetic equivalents of the substances contained in the cannabis plant, or in the resinous</li> <li>extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with</li> <li>similar chemical structure and pharmacological activity to those substances contained in the</li> <li>plant, such as the following: Δ1 cis or trans tetrahydrocannabinol, and their optical isomers Δ6</li> <li>cis or trans tetrahydrocannabinol, and their optical isomers Δ3,4 cis or trans</li> <li>tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not</li> <li>internationally standardized, compounds of these structures, regardless of numerical</li> <li>designation of atomic positions covered;</li> <li>(BB) Ethylamine analog of phencyclidine, some trade or other names:</li> <li>N-ethyl-1-phenylcyclohexyl)ethylamine, (1-phenylcyclohexyl)ethylamine,</li> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	188	(W) N-ethyl-3-piperidyl benzilate;
<ul> <li>(Z) Psilocyn;</li> <li>(AA) Tetrahydrocannabinols, naturally contained in a plant of the genus Cannabis</li> <li>(cannabis plant), except for marijuana as defined in Subsection 58-37-2(1)(aa)(i)(E), as well as</li> <li>synthetic equivalents of the substances contained in the cannabis plant, or in the resinous</li> <li>extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with</li> <li>similar chemical structure and pharmacological activity to those substances contained in the</li> <li>plant, such as the following: Δ1 cis or trans tetrahydrocannabinol, and their optical isomers Δ6</li> <li>cis or trans tetrahydrocannabinol, and their optical isomers Δ3,4 cis or trans</li> <li>tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not</li> <li>internationally standardized, compounds of these structures, regardless of numerical</li> <li>designation of atomic positions covered;</li> <li>(BB) Ethylamine analog of phencyclidine, some trade or other names:</li> <li>N-ethyl-1-phenylcyclohexyl)ethylamine, (1-phenylcyclohexyl)ethylamine,</li> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	189	(X) N-methyl-3-piperidyl benzilate;
192(AA) Tetrahydrocannabinols, naturally contained in a plant of the genus Cannabis193(cannabis plant), except for marijuana as defined in Subsection 58-37-2(1)(aa)(i)(E), as well as194synthetic equivalents of the substances contained in the cannabis plant, or in the resinous195extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with196similar chemical structure and pharmacological activity to those substances contained in the197plant, such as the following: $\Delta 1$ cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 6$ 198cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 3, 4$ cis or trans199tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not201internationally standardized, compounds of these structures, regardless of numerical202(BB) Ethylamine analog of phencyclidine, some trade or other names:203N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,204N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;205(CC) Pyrrolidine analog of phencyclidine, some trade or other names:2061-[1-(2-thienyl)-pyrrolidine, PCPy, PHP;207(DD) Thiophene analog of phencyclidine, some trade or other names:2081-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and209(EE) 1-[1-(2-thienyl)-cyclohexyl]pyrrolidine, some other names: TCPy.201(iv) Unless specifically excepted or unless listed in another schedule, any material201compound, mixture, or preparation which contains any quantity	190	(Y) Psilocybin;
193(cannabis plant), except for marijuana as defined in Subsection 58-37-2(1)(aa)(i)(E), as well as194synthetic equivalents of the substances contained in the cannabis plant, or in the resinous195extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with196similar chemical structure and pharmacological activity to those substances contained in the197plant, such as the following: $\Delta 1$ cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 5$ 198cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 3, 4$ cis or trans199tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not200internationally standardized, compounds of these structures, regardless of numerical201designation of atomic positions covered;202(BB) Ethylamine analog of phencyclidine, some trade or other names:203N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,204N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;205(CC) Pyrrolidine analog of phencyclidine, some trade or other names:2061-[1-(2-thienyl)-cyclohexyl]-pyrrolidine, PCPy, PHP;207(DD) Thiophene analog of phencyclidine, some other names:2081-[1-(2-thienyl)-cyclohexyl]pyrrolidine, some other names:209(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names:201(iv) Unless specifically excepted or unless listed in another schedule, any material202(iv) Unless specifically excepted or unless listed in another schedule, any material	191	(Z) Psilocyn;
194synthetic equivalents of the substances contained in the cannabis plant, or in the resinous195extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with196similar chemical structure and pharmacological activity to those substances contained in the197plant, such as the following: $\Delta 1$ cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 6$ 198cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 3, 4$ cis or trans199tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not200internationally standardized, compounds of these structures, regardless of numerical201designation of atomic positions covered;202(BB) Ethylamine analog of phencyclidine, some trade or other names:203N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,204N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;205(CC) Pyrrolidine analog of phencyclidine, some trade or other names:2061-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;207(DD) Thiophene analog of phencyclidine, some trade or other names:2081-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and209(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names:201(iv) Unless specifically excepted or unless listed in another schedule, any material201compound, mixture, or preparation which contains any quantity of the following substances	192	(AA) Tetrahydrocannabinols, naturally contained in a plant of the genus Cannabis
195extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with196similar chemical structure and pharmacological activity to those substances contained in the197plant, such as the following: $\Delta 1$ cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 6$ 198cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 3, 4$ cis or trans199tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not200internationally standardized, compounds of these structures, regardless of numerical201designation of atomic positions covered;202(BB) Ethylamine analog of phencyclidine, some trade or other names:203N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,204N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;205(CC) Pyrrolidine analog of phencyclidine, some trade or other names:2061-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;207(DD) Thiophene analog of phencyclidine, some trade or other names:2081-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and209(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names:2081-[1-(2-thienyl)-cyclohexyl]pyrrolidine, some other names:209(iv) Unless specifically excepted or unless listed in another schedule, any material201compound, mixture, or preparation which contains any quantity of the following substances	193	(cannabis plant), except for marijuana as defined in Subsection 58-37-2(1)(aa)(i)(E), as well as
<ul> <li>similar chemical structure and pharmacological activity to those substances contained in the</li> <li>plant, such as the following: Δ1 cis or trans tetrahydrocannabinol, and their optical isomers Δ6</li> <li>cis or trans tetrahydrocannabinol, and their optical isomers Δ3,4 cis or trans</li> <li>tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not</li> <li>internationally standardized, compounds of these structures, regardless of numerical</li> <li>designation of atomic positions covered;</li> <li>(BB) Ethylamine analog of phencyclidine, some trade or other names:</li> <li>N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,</li> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	194	synthetic equivalents of the substances contained in the cannabis plant, or in the resinous
<ul> <li>plant, such as the following: Δ1 cis or trans tetrahydrocannabinol, and their optical isomers Δ6</li> <li>cis or trans tetrahydrocannabinol, and their optical isomers Δ3,4 cis or trans</li> <li>tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not</li> <li>internationally standardized, compounds of these structures, regardless of numerical</li> <li>designation of atomic positions covered;</li> <li>(BB) Ethylamine analog of phencyclidine, some trade or other names:</li> <li>N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,</li> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	195	extractives of Cannabis, sp. and/or synthetic substances, derivatives, and their isomers with
<ul> <li>cis or trans tetrahydrocannabinol, and their optical isomers Δ3,4 cis or trans</li> <li>tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not</li> <li>internationally standardized, compounds of these structures, regardless of numerical</li> <li>designation of atomic positions covered;</li> <li>(BB) Ethylamine analog of phencyclidine, some trade or other names:</li> <li>N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,</li> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	196	similar chemical structure and pharmacological activity to those substances contained in the
<ul> <li>tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not</li> <li>internationally standardized, compounds of these structures, regardless of numerical</li> <li>designation of atomic positions covered;</li> <li>(BB) Ethylamine analog of phencyclidine, some trade or other names:</li> <li>N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,</li> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	197	plant, such as the following: $\Delta 1$ cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 6$
<ul> <li>internationally standardized, compounds of these structures, regardless of numerical</li> <li>designation of atomic positions covered;</li> <li>(BB) Ethylamine analog of phencyclidine, some trade or other names:</li> <li>N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,</li> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	198	cis or trans tetrahydrocannabinol, and their optical isomers $\Delta 3,4$ cis or trans
<ul> <li>designation of atomic positions covered;</li> <li>(BB) Ethylamine analog of phencyclidine, some trade or other names:</li> <li>N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,</li> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	199	tetrahydrocannabinol, and its optical isomers, and since nomenclature of these substances is not
<ul> <li>(BB) Ethylamine analog of phencyclidine, some trade or other names:</li> <li>N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,</li> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	200	internationally standardized, compounds of these structures, regardless of numerical
<ul> <li>N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,</li> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	201	designation of atomic positions covered;
<ul> <li>N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;</li> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	202	(BB) Ethylamine analog of phencyclidine, some trade or other names:
<ul> <li>(CC) Pyrrolidine analog of phencyclidine, some trade or other names:</li> <li>1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	203	N-ethyl-1-phenylcyclohexylamine, (1-phenylcyclohexyl)ethylamine,
<ul> <li>206 1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;</li> <li>207 (DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>208 1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>209 (EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>210 (iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>211 compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	204	N-(1-phenylcyclohexyl)ethylamine, cyclohexamine, PCE;
<ul> <li>(DD) Thiophene analog of phencyclidine, some trade or other names:</li> <li>1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	205	(CC) Pyrrolidine analog of phencyclidine, some trade or other names:
<ul> <li>208 1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and</li> <li>209 (EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>210 (iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>211 compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	206	1-(1-phenylcyclohexyl)-pyrrolidine, PCPy, PHP;
<ul> <li>(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.</li> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	207	(DD) Thiophene analog of phencyclidine, some trade or other names:
<ul> <li>(iv) Unless specifically excepted or unless listed in another schedule, any material</li> <li>compound, mixture, or preparation which contains any quantity of the following substances</li> </ul>	208	1-[1-(2-thienyl)-cyclohexyl]-piperidine, 2-thienylanalog of phencyclidine, TPCP, TCP; and
211 compound, mixture, or preparation which contains any quantity of the following substances	209	(EE) 1-[1-(2-thienyl)cyclohexyl]pyrrolidine, some other names: TCPy.
	210	(iv) Unless specifically excepted or unless listed in another schedule, any material
212 having a depressant effect on the central nervous system including its salts isomers and salts	211	compound, mixture, or preparation which contains any quantity of the following substances
212 naving a depressant effect on the central hervous system, including its saits, isolifets, and saits	212	having a depressant effect on the central nervous system, including its salts, isomers, and salts
213 of isomers when the existence of the salts, isomers, and salts of isomers is possible within the	213	of isomers when the existence of the salts, isomers, and salts of isomers is possible within the

214	specific chemical designation:
215	(A) Mecloqualone; and
216	(B) Methaqualone.
217	(v) Any material, compound, mixture, or preparation containing any quantity of the
218	following substances having a stimulant effect on the central nervous system, including their
219	salts, isomers, and salts of isomers:
220	(A) Aminorex, some other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; or
221	4,5-dihydro-5-phenyl-2-oxazolamine;
222	(B) Cathinone, some trade or other names: 2-amino-1-phenyl-1-propanone,
223	alpha-aminopropiophenone, 2-aminopropiophenone, and norephedrone;
224	(C) Fenethylline;
225	(D) Methcathinone, some other names: 2-(methylamino)-propiophenone;
226	alpha-(methylamino)propiophenone; 2-(methylamino)-1-phenylpropan-1-one;
227	alpha-N-methylaminopropiophenone; monomethylpropion; ephedrone; N-methylcathinone;
228	methylcathinone; AL-464; AL-422; AL-463 and UR1432, its salts, optical isomers, and salts of
229	optical isomers;
230	(E) ( $\pm$ )cis-4-methylaminorex (( $\pm$ )cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);
231	(F) N-ethylamphetamine; and
232	(G) N,N-dimethylamphetamine, also known as
233	N,N-alpha-trimethyl-benzeneethanamine; N,N-alpha-trimethylphenethylamine.
234	(vi) Any material, compound, mixture, or preparation which contains any quantity of
235	the following substances, including their optical isomers, salts, and salts of isomers, subject to
236	temporary emergency scheduling:
237	(A) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl); and
238	(B) N-[1- (2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl).
239	(vii) Unless specifically excepted or unless listed in another schedule, any material,
240	compound, mixture, or preparation which contains any quantity of gamma hydroxy butyrate
241	(gamma hydrobutyric acid), including its salts, isomers, and salts of isomers.
242	(b) Schedule II:
243	(i) Unless specifically excepted or unless listed in another schedule, any of the
244	following substances whether produced directly or indirectly by extraction from substances of

- 245 vegetable origin, or independently by means of chemical synthesis, or by a combination of
- 246 extraction and chemical synthesis:
- 247 (A) Opium and opiate, and any salt, compound, derivative, or preparation of opium or
- 248 opiate, excluding apomorphine, dextrorphan, nalbuphine, nalmefene, naloxone, and naltrexone,
- 249 and their respective salts, but including:
- 250 (I) Raw opium;
  251 (II) Opium extracts;
  252 (III) Opium fluid;
  253 (IV) Powdered opium;
- 254 (V) Granulated opium;
- 255 (VI) Tincture of opium;
- 256 (VII) Codeine;
- 257 (VIII) Ethylmorphine;
- 258 (IX) Etorphine hydrochloride;
- 259 (X) Hydrocodone;
- 260 (XI) Hydromorphone;
- 261 (XII) Metopon;
- 262 (XIII) Morphine;
- 263 (XIV) Oxycodone;
- 264 (XV) Oxymorphone; and
- 265 (XVI) Thebaine;
- (B) Any salt, compound, derivative, or preparation which is chemically equivalent or
  identical with any of the substances referred to in Subsection (2)(b)(i)(A), except that these
  substances may not include the isoquinoline alkaloids of opium;
- 269 (C) Opium poppy and poppy straw;
- 270 (D) Coca leaves and any salt, compound, derivative, or preparation of coca leaves, and 271 any salt, compound, derivative, or preparation which is chemically equivalent or identical with 272 any of these substances, and includes cocaine and ecgonine, their salts, isomers, derivatives.
- and salts of isomers and derivatives, whether derived from the coca plant or synthetically
- 274 produced, except the substances may not include decocainized coca leaves or extraction of coca
- 275 leaves, which extractions do not contain cocaine or ecgonine; and

276	(E) Concentrate of poppy straw, which means the crude extract of poppy straw in either
277	liquid, solid, or powder form which contains the phenanthrene alkaloids of the opium poppy.
278	(ii) Unless specifically excepted or unless listed in another schedule, any of the
279	following opiates, including their isomers, esters, ethers, salts, and salts of isomers, esters, and
280	ethers, when the existence of the isomers, esters, ethers, and salts is possible within the specific
281	chemical designation, except dextrorphan and levopropoxyphene:
282	(A) Alfentanil;
283	(B) Alphaprodine;
284	(C) Anileridine;
285	(D) Bezitramide;
286	(E) Bulk dextropropoxyphene (nondosage forms);
287	(F) Carfentanil;
288	(G) Dihydrocodeine;
289	(H) Diphenoxylate;
290	(I) Fentanyl;
291	(J) Isomethadone;
292	(K) Levo-alphacetylmethadol, some other names: levo-alpha-acetylmethadol,
293	levomethadyl acetate, or LAAM;
294	(L) Levomethorphan;
295	(M) Levorphanol;
296	(N) Metazocine;
297	(O) Methadone;
298	(P) Methadone-Intermediate, 4-cyano-2-dimethylamino-4, 4-diphenyl butane;
299	(Q) Moramide-Intermediate, 2-methyl-3-morpholino-1, 1-diphenylpropane-carboxylic
300	acid;
301	(R) Pethidine (meperidine);
302	(S) Pethidine-Intermediate-A, 4-cyano-1-methyl-4-phenylpiperidine;
303	(T) Pethidine-Intermediate-B, ethyl-4-phenylpiperidine-4-carboxylate;
304	(U) Pethidine-Intermediate-C, 1-methyl-4-phenylpiperidine-4-carboxylic acid;
305	(V) Phenazocine;
306	(W) Piminodine;

307	(X) Racemethorphan;
308	(Y) Racemorphan;
308 309	
	(Z) Remifentanil; and
310	(AA) Sufentanil.
311	(iii) Unless specifically excepted or unless listed in another schedule, any material,
312	compound, mixture, or preparation which contains any quantity of the following substances
313	having a stimulant effect on the central nervous system:
314	(A) Amphetamine, its salts, optical isomers, and salts of its optical isomers;
315	(B) Methamphetamine, its salts, isomers, and salts of its isomers;
316	(C) Phenmetrazine and its salts; and
317	(D) Methylphenidate.
318	(iv) Unless specifically excepted or unless listed in another schedule, any material,
319	compound, mixture, or preparation which contains any quantity of the following substances
320	having a depressant effect on the central nervous system, including its salts, isomers, and salts
321	of isomers when the existence of the salts, isomers, and salts of isomers is possible within the
322	specific chemical designation:
323	(A) Amobarbital;
324	(B) Glutethimide;
325	(C) Pentobarbital;
326	(D) Phencyclidine;
327	(E) Phencyclidine immediate precursors: 1-phenylcyclohexylamine and
328	1-piperidinocyclohexanecarbonitrile (PCC); and
329	(F) Secobarbital.
330	(v) (A) Unless specifically excepted or unless listed in another schedule, any material,
331	compound, mixture, or preparation which contains any quantity of Phenylacetone.
332	(B) Some of these substances may be known by trade or other names:
333	phenyl-2-propanone; P2P; benzyl methyl ketone; and methyl benzyl ketone.
334	(vi) Nabilone, another name for nabilone:
335	(±)-trans-3-(1,1-dimethylheptyl)-6,6a,7,8,10,10a-hexahydro-1-hydroxy-6,
336	6-dimethyl-9H-dibenzo[b,d]pyran-9-one.
337	(vii) A drug product or preparation that contains any component of marijuana,

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including tetrahydrocannabinol, and is approved by the United States Food and Drug
Administration and scheduled by the Drug Enforcement Administration in Schedule II of the
federal Controlled Substances Act, Title II, P.L. 91-513.

341 (c) Schedule III:

(i) Unless specifically excepted or unless listed in another schedule, any material,
compound, mixture, or preparation which contains any quantity of the following substances
having a stimulant effect on the central nervous system, including its salts, isomers whether
optical, position, or geometric, and salts of the isomers when the existence of the salts, isomers,
and salts of isomers is possible within the specific chemical designation:

(A) Those compounds, mixtures, or preparations in dosage unit form containing any
stimulant substances listed in Schedule II, which compounds, mixtures, or preparations were
listed on August 25, 1971, as excepted compounds under Section 1308.32 of Title 21 of the
Code of Federal Regulations, and any other drug of the quantitive composition shown in that
list for those drugs or which is the same except that it contains a lesser quantity of controlled
substances;

- 353 (B) Benzphetamine;
- 354 (C) Chlorphentermine;
- 355 (D) Clortermine; and

#### 356 (E) Phendimetrazine.

357 (ii) Unless specifically excepted or unless listed in another schedule, any material,

compound, mixture, or preparation which contains any quantity of the following substanceshaving a depressant effect on the central nervous system:

360 (A) Any compound, mixture, or preparation containing amobarbital, secobarbital,
361 pentobarbital, or any salt of any of them, and one or more other active medicinal ingredients
362 which are not listed in any schedule;

363 (B) Any suppository dosage form containing amobarbital, secobarbital, or
364 pentobarbital, or any salt of any of these drugs which is approved by the Food and Drug
365 Administration for marketing only as a suppository;

366 (C) Any substance which contains any quantity of a derivative of barbituric acid or any367 salt of any of them;

368 (D) Chlorhexadol;

369	(E) Buprenorphine;
370	(F) Any drug product containing gamma hydroxybutyric acid, including its salts,
371	isomers, and salts of isomers, for which an application is approved under the federal Food,
372	Drug, and Cosmetic Act, Section 505;
373	(G) Ketamine, its salts, isomers, and salts of isomers, some other names for ketamine:
374	± -2-(2-chlorophenyl)-2-(methylamino)-cyclohexanone;
375	(H) Lysergic acid;
376	(I) Lysergic acid amide;
377	(J) Methyprylon;
378	(K) Sulfondiethylmethane;
379	(L) Sulfonethylmethane;
380	(M) Sulfonmethane; and
381	(N) Tiletamine and zolazepam or any of their salts, some trade or other names for a
382	tiletamine-zolazepam combination product: Telazol, some trade or other names for tiletamine:
383	2-(ethylamino)-2-(2-thienyl)-cyclohexanone, some trade or other names for zolazepam:
384	4-(2-fluorophenyl)-6,8-dihydro-1,3,8-trimethylpyrazolo-[3,4-e] [1,4]-diazepin-7(1H)-one,
385	flupyrazapon.
386	(iii) Dronabinol (synthetic) in sesame oil and encapsulated in a soft gelatin capsule in a
387	U.S. Food and Drug Administration approved drug product, some other names for dronabinol:
388	(6aR-trans)-6a,7,8,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1-ol, or
389	(-)-delta-9-(trans)-tetrahydrocannabinol.
390	(iv) Nalorphine.
391	(v) Unless specifically excepted or unless listed in another schedule, any material,
392	compound, mixture, or preparation containing limited quantities of any of the following
393	narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid:
394	(A) Not more than 1.8 grams of codeine per 100 milliliters or not more than 90
395	milligrams per dosage unit, with an equal or greater quantity of an isoquinoline alkaloid of
396	opium;
397	(B) Not more than 1.8 grams of codeine per 100 milliliters or not more than 90
398	milligrams per dosage unit, with one or more active non-narcotic ingredients in recognized
399	therapeutic amounts;

400	(C) Not more than 300 milligrams of dihydrocodeinone per 100 milliliters or not more
401	than 15 milligrams per dosage unit, with a fourfold or greater quantity of an isoquinoline
402	alkaloid of opium;
403	(D) Not more than 300 milligrams of dihydrocodeinone per 100 milliliters or not more
404	than 15 milligrams per dosage unit, with one or more active, non-narcotic ingredients in
405	recognized therapeutic amounts;
406	(E) Not more than 1.8 grams of dihydrocodeine per 100 milliliters or not more than 90
407	milligrams per dosage unit, with one or more active non-narcotic ingredients in recognized
408	therapeutic amounts;
409	(F) Not more than 300 milligrams of ethylmorphine per 100 milliliters or not more
410	than 15 milligrams per dosage unit, with one or more active, non-narcotic ingredients in
411	recognized therapeutic amounts;
412	(G) Not more than 500 milligrams of opium per 100 milliliters or per 100 grams, or not
413	more than 25 milligrams per dosage unit, with one or more active, non-narcotic ingredients in
414	recognized therapeutic amounts; and
415	(H) Not more than 50 milligrams of morphine per 100 milliliters or per 100 grams with
416	one or more active, non-narcotic ingredients in recognized therapeutic amounts.
417	(vi) Unless specifically excepted or unless listed in another schedule, anabolic steroids
418	including any of the following or any isomer, ester, salt, or derivative of the following that
419	promotes muscle growth:
420	(A) Boldenone;
421	(B) Chlorotestosterone (4-chlortestosterone);
422	(C) Clostebol;
423	(D) Dehydrochlormethyltestosterone;
424	(E) Dihydrotestosterone (4-dihydrotestosterone);
425	(F) Drostanolone;
426	(G) Ethylestrenol;
427	(H) Fluoxymesterone;
428	(I) Formebulone (formebolone);
429	(J) Mesterolone;
430	(K) Methandienone;

431	(L) Methandranone;
432	(M) Methandriol;
433	(N) Methandrostenolone;
434	(O) Methenolone;
435	(P) Methyltestosterone;
436	(Q) Mibolerone;
437	(R) Nandrolone;
438	(S) Norethandrolone;
439	(T) Oxandrolone;
440	(U) Oxymesterone;
441	(V) Oxymetholone;
442	(W) Stanolone;
443	(X) Stanozolol;
444	(Y) Testolactone;
445	(Z) Testosterone; and
446	(AA) Trenbolone.
447	(vii) Anabolic steroids expressly intended for administration through implants to cattle
448	or other nonhuman species, and approved by the Secretary of Health and Human Services for
449	use, may not be classified as a controlled substance.
450	(viii) A drug product or preparation that contains any component of marijuana,
451	including tetrahydrocannabinol, and is approved by the United States Food and Drug
452	Administration and scheduled by the Drug Enforcement Administration in Schedule III of the
453	federal Controlled Substances Act, Title II, P.L. 91-513.
454	(ix) Nabiximols.
455	(d) Schedule IV:
456	(i) Unless specifically excepted or unless listed in another schedule, any material,
457	compound, mixture, or preparation containing not more than 1 milligram of difenoxin and not
458	less than 25 micrograms of atropine sulfate per dosage unit, or any salts of any of them.
459	(ii) Unless specifically excepted or unless listed in another schedule, any material,
460	compound, mixture, or preparation which contains any quantity of the following substances,
461	including its salts, isomers, and salts of isomers when the existence of the salts, isomers, and

- 462 salts of isomers is possible within the specific chemical designation:
- 463 (A) Alprazolam;464 (B) Barbital;
- 465 (C) Bromazepam;
- 466 (D) Butorphanol;
- 467 (E) Camazepam;
- 468 (F) Carisoprodol;
- 469 (G) Chloral betaine;
- 470 (H) Chloral hydrate;
- 471 (I) Chlordiazepoxide;
- 472 (J) Clobazam;
- 473 (K) Clonazepam;
- 474 (L) Clorazepate;
- 475 (M) Clotiazepam;
- 476 (N) Cloxazolam;
- 477 (O) Delorazepam;
- 478 (P) Diazepam;
- 479 (Q) Dichloralphenazone;
- 480 (R) Estazolam;
- 481 (S) Ethchlorvynol;
- 482 (T) Ethinamate;
- 483 (U) Ethyl loflazepate;
- 484 (V) Fludiazepam;
- 485 (W) Flunitrazepam;
- 486 (X) Flurazepam;
- 487 (Y) Halazepam;
- 488 (Z) Haloxazolam;
- 489 (AA) Ketazolam;
- 490 (BB) Loprazolam;
- 491 (CC) Lorazepam;
- 492 (DD) Lormetazepam;

493	(EE) Mebutamate;
494	(FF) Medazepam;
495	(GG) Meprobamate;
496	(HH) Methohexital;
497	(II) Methylphenobarbital (mephobarbital);
498	(JJ) Midazolam;
499	(KK) Nimetazepam;
500	(LL) Nitrazepam;
501	(MM) Nordiazepam;
502	(NN) Oxazepam;
503	(OO) Oxazolam;
504	(PP) Paraldehyde;
505	(QQ) Pentazocine;
506	(RR) Petrichloral;
507	(SS) Phenobarbital;
508	(TT) Pinazepam;
509	(UU) Prazepam;
510	(VV) Quazepam;
511	(WW) Temazepam;
512	(XX) Tetrazepam;
513	(YY) Tramadol;
514	(ZZ) Triazolam;
515	(AAA) Zaleplon; and
516	(BBB) Zolpidem.
517	(iii) Any material, compound, mixture, or preparation of fenfluramine which contains
518	any quantity of the following substances, including its salts, isomers whether optical, position,
519	or geometric, and salts of the isomers when the existence of the salts, isomers, and salts of
520	isomers is possible.
521	(iv) Unless specifically excepted or unless listed in another schedule, any material.

(iv) Unless specifically excepted or unless listed in another schedule, any material,
compound, mixture, or preparation which contains any quantity of the following substances
having a stimulant effect on the central nervous system, including its salts, isomers whether

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524 optical, position, or geometric isomers, and salts of the isomers when the existence of the salts,

525 isomers, and salts of isomers is possible within the specific chemical designation:

- 526 (A) Cathine ((+)-norpseudoephedrine);
- 527 (B) Diethylpropion;
- 528 (C) Fencamfamine;
- 529 (D) Fenproprex;
- 530 (E) Mazindol;
- 531 (F) Mefenorex;
- 532 (G) Modafinil;
- 533 (H) Pemoline, including organometallic complexes and chelates thereof;
- 534 (I) Phentermine;
- 535 (J) Pipradrol;
- 536 (K) Sibutramine; and
- 537 (L) SPA ((-)-1-dimethylamino-1,2-diphenylethane).
- (v) Unless specifically excepted or unless listed in another schedule, any material,
  compound, mixture, or preparation which contains any quantity of dextropropoxyphene
  (alpha-(+)-4-dimethylamino-1, 2-diphenyl-3-methyl-2-propionoxybutane), including its salts.
- (vi) A drug product or preparation that contains any component of marijuana and is
  approved by the United States Food and Drug Administration and scheduled by the Drug
  Enforcement Administration in Schedule IV of the federal Controlled Substances Act, Title II,
  P.L. 91-513.
- 545 (e) Schedule V:

(i) Any compound, mixture, or preparation containing any of the following limited
quantities of narcotic drugs, or their salts calculated as the free anhydrous base or alkaloid,
which includes one or more non-narcotic active medicinal ingredients in sufficient proportion
to confer upon the compound, mixture, or preparation valuable medicinal qualities other than
those possessed by the narcotic drug alone:

551

(A) not more than 200 milligrams of codeine per 100 milliliters or per 100 grams;

- (B) not more than 100 milligrams of dihydrocodeine per 100 milliliters or per 100grams;
- (C) not more than 100 milligrams of ethylmorphine per 100 milliliters or per 100

555 grams;

- 556 (D) not more than 2.5 milligrams of diphenoxylate and not less than 25 micrograms of 557 atropine sulfate per dosage unit;
- (E) not more than 100 milligrams of opium per 100 milliliters or per 100 grams;
- 559 (F) not more than 0.5 milligram of difenoxin and not less than 25 micrograms of 560 atropine sulfate per dosage unit; and
- (G) unless specifically exempted or excluded or unless listed in another schedule, any
   material, compound, mixture, or preparation which contains Pyrovalerone having a stimulant
   effect on the central nervous system, including its salts, isomers, and salts of isomers.
- 564 (ii) A drug product or preparation that contains any component of marijuana, including
- 565 cannabidiol, and is approved by the United States Food and Drug Administration and
- 566 scheduled by the Drug Enforcement Administration in Schedule V of the federal Controlled
- 567 Substances Act, Title II, P.L. 91-513.
- 568 <u>(iii) Gabapentin.</u>
- 569 Section 2. Effective date.
- 570 <u>This bill takes effect on May 1, 2024.</u>