99TH GENERAL ASSEMBLY

State of Illinois

2015 and 2016

HB3588

by Rep. Michael J. Zalewski

SYNOPSIS AS INTRODUCED:

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Creates the Synthetic Drug Identification Pilot Program Act. Provides that at least once every 90 days, and in consultation with the Division of Forensic Services of the Department of State Police, the Department of Human Services shall submit a report to the Governor, Attorney General, and General Assembly outlining whether the Division of Forensic Services of the Department of State Police has identified any new chemical formulas that are used to make synthetic cannabinoids or cathinones (synthetic drugs) that are not currently illegal under State law. Provides that if the Department of Human Services' report to the Governor, Attorney General, and General Assembly confirms the Department of Human Services has identified new chemical formulas that are used to make synthetic drugs, the Department Human Services shall as soon as practicable propose an emergency rule to add any new chemical formulas to the current list of chemical formulas that are listed in the Illinois Controlled Substances Act as Schedule I controlled substances, and adopt the proposed rule as quickly as allowed for under the Illinois Administrative Procedure Act. Provides that if the Department of Human Services adopts the emergency rule, the new rule shall be recognized as law under the Illinois Controlled Substances Act. Provides that the Department of State Police, by rule, shall create a pilot program that uses technologies and protocols to instantly identify synthetic cannabinoids and cathinones (synthetic drugs), as well as other designer drugs. Provides that the pilot program shall focus on using technology capable of presumptive identification of illicit drugs in the field. Repeals the Act on July 30, 2018. Amends the Illinois Controlled Substances Act to make conforming changes. Effective immediately.

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FISCAL NOTE ACT MAY APPLY

A BILL FOR

1

AN ACT concerning criminal law.

2 Be it enacted by the People of the State of Illinois, 3 represented in the General Assembly:

Section 1. Short title. This Act may be cited as the
Synthetic Drug Identification Pilot Program Act.

6 Section 5. Legislative intent. The General Assembly 7 recognizes the recent growth of synthetic drugs, such as 8 Spice/K2 and Bath Salts, and the dangers caused by these 9 substances. The concern is exemplified by a Substance Abuse and 10 Mental Health Services Administration report which summarizes the frequency and trends of abuse for these substances. The 11 12 General Assembly further recognizes that better methods and 13 strategies that appropriately respond to new synthetic drugs as 14 soon as they are made known to the State is of particular importance. The General Assembly further recognizes that law 15 16 enforcement is in need of presumptive testing tools capable of 17 quickly identifying substances as illegal synthetic drugs defined under State law. It is the intent of this legislation 18 19 to create a process by which synthetic drugs can be quickly 20 outlawed under State law and instantly identified by police in 21 the field.

22

Section 10. Department of Human Services; emergency rule.

(a) At least once every 90 days, and in consultation with 1 2 the Division of Forensic Services of the Department of State 3 Police, the Department of Human Services shall submit a report to the Governor, Attorney General, and General Assembly 4 5 outlining whether the Division of Forensic Services of the Department of State Police has identified any new chemical 6 formulas that are used to make synthetic cannabinoids or 7 8 cathinones (synthetic drugs) that are not currently illegal 9 under State law. To identify new chemical formulas, the 10 Department of Human Services shall, among other activities, 11 routinely communicate with the Division of Forensic Services of 12 the Department of State Police, the United States Drug 13 Enforcement Administration, the United States Office of National Drug Control Policy, and the Scientific Working Group 14 15 for the Analysis of Seized Drugs (SWDRUG), and agencies of 16 other states involved in drug schedule classifications.

17 (b) If the Department of Human Services' report to the Governor, Attorney General, and General Assembly confirms the 18 Department of Human Services has identified new chemical 19 20 formulas that are used to make synthetic drugs, the Department of Human Services shall as soon as practicable propose an 21 22 emergency rule to add any new chemical formulas to the current 23 list of chemical formulas that are listed in Section 204 of the 24 Illinois Controlled Substances Act as Schedule I controlled 25 substances, and adopt the proposed rule as quickly as allowed for under the Illinois Administrative Procedure Act. If the 26

Department of Human Services adopts the emergency rule under
 this subsection, the rule shall have the force of law under
 Section 204 of the Illinois Controlled Substances Act.

4 (c) Any emergency rule adopted under this Section shall be 5 inoperative in 12 months from the date that the emergency rule 6 becomes effective, or when the General Assembly takes action to 7 ratify, change, or reject the emergency rule enacted by the 8 Department of Human Services.

9 (d) Nothing in this Section shall interfere with the 10 exemptions provided for under State law to any person or entity 11 that possesses a chemical formula defined as a scheduled drug 12 for a lawful purpose.

Section 15. Law enforcement field testing to instantly identify synthetic drugs.

(a) The Department of State Police, by rule, shall create a
pilot program that uses technologies and protocols to instantly
identify synthetic cannabinoids and cathinones (synthetic
drugs), as well as other designer drugs. The pilot program
shall focus on using technology capable of presumptive
identification of illicit drugs in the field.

(b) A minimum of 5 separate State Police districts in the State shall be equipped to operate the pilot program. The Department of State Police may choose to have one or more of the pilot sites located within a municipal or county police department.

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(c) Before July 30, 2018, the Department of State Police 1 2 shall submit a report to the General Assembly outlining the 3 findings of the pilot program, and make any recommendation on whether the technologies and protocols selected for the pilot 4 5 program can be used to help prevent the growth of synthetic illicit drugs 6 drugs and other throughout the State. 7 Specifically, the evaluation shall include, but is not limited 8 to the following:

9 (1) Review of technical capabilities and accuracy 10 rates of technologies and protocols selected for the pilot 11 program.

12 (2) Describe the impact to State and local crime 13 laboratory backlogs if the technologies and protocols 14 selected could eliminate the need to send synthetic drugs, 15 or other illicit drugs to the crime laboratory for 16 presumptive testing, including the potential cost savings 17 to State and local government.

18 (3) Describe the status of court acceptance of the 19 technologies and protocols selected for the pilot program 20 for the presumptive identification of synthetic drugs and 21 other illicit drugs.

22

Section 20. Repeal. This Act is repealed on July 30, 2018.

23 Section 105. The Illinois Controlled Substances Act is 24 amended by changing Section 204 as follows:

1	(720 ILCS 570/204) (from Ch. 56 1/2, par. 1204)
2	Sec. 204. (a) The controlled substances listed in this
3	Section are included in Schedule I.
4	(b) Unless specifically excepted or unless listed in
5	another schedule, any of the following opiates, including their
6	isomers, esters, ethers, salts, and salts of isomers, esters,
7	and ethers, whenever the existence of such isomers, esters,
8	ethers and salts is possible within the specific chemical
9	designation:
10	(1) Acetylmethadol;
11	(1.1) Acetyl-alpha-methylfentanyl
12	(N-[1-(1-methyl-2-phenethyl)-
13	4-piperidinyl]-N-phenylacetamide);
14	(2) Allylprodine;
15	(3) Alphacetylmethadol, except
16	levo-alphacetylmethadol (also known as levo-alpha-
17	acetylmethadol, levomethadyl acetate, or LAAM);
18	(4) Alphameprodine;
19	(5) Alphamethadol;
20	(6) Alpha-methylfentanyl
21	(N-(1-alpha-methyl-beta-phenyl) ethyl-4-piperidyl)
22	propionanilide; 1-(1-methyl-2-phenylethyl)-4-(N-
23	propanilido) piperidine;
24	(6.1) Alpha-methylthiofentanyl
25	(N-[1-methyl-2-(2-thienyl)ethyl-

1	4-piperidinyl]-N-phenylpropanamide);
2	(7) 1-methyl-4-phenyl-4-propionoxypiperidine (MPPP);
3	(7.1) PEPAP
4	(1-(2-phenethyl)-4-phenyl-4-acetoxypiperidine);
5	(8) Benzethidine;
6	(9) Betacetylmethadol;
7	(9.1) Beta-hydroxyfentanyl
8	(N-[1-(2-hydroxy-2-phenethyl)-
9	4-piperidinyl]-N-phenylpropanamide);
10	(10) Betameprodine;
11	(11) Betamethadol;
12	(12) Betaprodine;
13	(13) Clonitazene;
14	(14) Dextromoramide;
15	(15) Diampromide;
16	(16) Diethylthiambutene;
17	(17) Difenoxin;
18	(18) Dimenoxadol;
19	(19) Dimepheptanol;
20	(20) Dimethylthiambutene;
21	(21) Dioxaphetylbutyrate;
22	(22) Dipipanone;
23	(23) Ethylmethylthiambutene;
24	(24) Etonitazene;
25	(25) Etoxeridine;
26	(26) Furethidine;

1	(27) Hydroxpethidine;
2	(28) Ketobemidone;
3	(29) Levomoramide;
4	(30) Levophenacylmorphan;
5	(31) 3-Methylfentanyl
6	(N-[3-methyl-1-(2-phenylethyl)-
7	4-piperidyl]-N-phenylpropanamide);
8	(31.1) 3-Methylthiofentanyl
9	(N-[(3-methyl-1-(2-thienyl)ethyl-
10	4-piperidinyl]-N-phenylpropanamide);
11	(32) Morpheridine;
12	(33) Noracymethadol;
13	(34) Norlevorphanol;
14	(35) Normethadone;
15	(36) Norpipanone;
16	(36.1) Para-fluorofentanyl
17	(N-(4-fluorophenyl)-N-[1-(2-phenethyl)-
18	4-piperidinyl]propanamide);
19	(37) Phenadoxone;
20	(38) Phenampromide;
21	(39) Phenomorphan;
22	(40) Phenoperidine;
23	(41) Piritramide;
24	(42) Proheptazine;
25	(43) Properidine;
26	(44) Propiram;

1	(45) Racemoramide;
2	(45.1) Thiofentanyl
3	(N-phenyl-N-[1-(2-thienyl)ethyl-
4	4-piperidinyl]-propanamide);
5	(46) Tilidine;
6	(47) Trimeperidine;
7	(48) Beta-hydroxy-3-methylfentanyl (other name:
8	N-[1-(2-hydroxy-2-phenethyl)-3-methyl-4-piperidinyl]-
9	N-phenylpropanamide).
10	(c) Unless specifically excepted or unless listed in
11	another schedule, any of the following opium derivatives, its
12	salts, isomers and salts of isomers, whenever the existence of
13	such salts, isomers and salts of isomers is possible within the
14	specific chemical designation:
15	(1) Acetorphine;
16	(2) Acetyldihydrocodeine;
17	<pre>(3) Benzylmorphine;</pre>
18	(4) Codeine methylbromide;
19	(5) Codeine-N-Oxide;
20	(6) Cyprenorphine;
21	(7) Desomorphine;
22	(8) Diacetyldihydromorphine (Dihydroheroin);
23	(9) Dihydromorphine;
24	(10) Drotebanol;
25	(11) Etorphine (except hydrochloride salt);
26	(12) Heroin;

1	(13) Hydromorphinol;
2	(14) Methyldesorphine;
3	(15) Methyldihydromorphine;
4	(16) Morphine methylbromide;
5	(17) Morphine methylsulfonate;
6	(18) Morphine-N-Oxide;
7	(19) Myrophine;
8	(20) Nicocodeine;
9	(21) Nicomorphine;
10	(22) Normorphine;
11	(23) Pholcodine;
12	(24) Thebacon.
13	(d) Unless specifically excepted or unless listed in
14	another schedule, any material, compound, mixture, or
15	preparation which contains any quantity of the following
16	hallucinogenic substances, or which contains any of its salts,
17	isomers and salts of isomers, whenever the existence of such
18	salts, isomers, and salts of isomers is possible within the
19	specific chemical designation (for the purposes of this
20	paragraph only, the term "isomer" includes the optical,
21	position and geometric isomers):

22

(1) 3,4-methylenedioxyamphetamine

23 (alpha-methyl, 3, 4-methylenedioxyphenethylamine,

methylenedioxyamphetamine, MDA); 24

(1.1) Alpha-ethyltryptamine 25

26 (some trade or other names: etryptamine;

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1	MONASE; alpha-ethyl-1H-indole-3-ethanamine;
2	3-(2-aminobutyl)indole; a-ET; and AET);
3	(2) 3,4-methylenedioxymethamphetamine (MDMA);
4	(2.1) 3,4-methylenedioxy-N-ethylamphetamine
5	(also known as: N-ethyl-alpha-methyl-
6	3,4(methylenedioxy) Phenethylamine, N-ethyl MDA, MDE,
7	and MDEA);
8	(2.2) N-Benzylpiperazine (BZP);
9	(3) 3-methoxy-4,5-methylenedioxyamphetamine, (MMDA);
10	(4) 3,4,5-trimethoxyamphetamine (TMA);
11	(5) (Blank);
12	(6) Diethyltryptamine (DET);
13	(7) Dimethyltryptamine (DMT);
14	(7.1) 5-Methoxy-diallyltryptamine;
15	(8) 4-methyl-2,5-dimethoxyamphetamine (DOM, STP);
16	(9) Ibogaine (some trade and other names:
17	7-ethyl-6,6,beta,7,8,9,10,12,13-octahydro-2-methoxy-
18	6,9-methano-5H-pyrido [1',2':1,2] azepino [5,4-b]
19	<pre>indole; Tabernanthe iboga);</pre>
20	(10) Lysergic acid diethylamide;
21	(10.1) Salvinorin A;
22	(10.5) Salvia divinorum (meaning all parts of the plant
23	presently classified botanically as Salvia divinorum,
24	whether growing or not, the seeds thereof, any extract from
25	any part of that plant, and every compound, manufacture,
26	salts, isomers, and salts of isomers whenever the existence

of such salts, isomers, and salts of isomers is possible 1 2 within the specific chemical designation, derivative, 3 mixture, or preparation of that plant, its seeds or extracts); 4 5 (11) 3,4,5-trimethoxyphenethylamine (Mescaline); (12) Peyote (meaning all parts of the plant presently 6 7 classified botanically as Lophophora williamsii Lemaire, 8 whether growing or not, the seeds thereof, any extract from 9 any part of that plant, and every compound, manufacture, 10 salts, derivative, mixture, or preparation of that plant, 11 its seeds or extracts); 12 (13) N-ethyl-3-piperidyl benzilate (JB 318); 13 (14) N-methyl-3-piperidyl benzilate; (14.1) N-hydroxy-3, 4-methylenedioxyamphetamine 14 15 (also known as N-hydroxy-alpha-methyl-16 3,4 (methylenedioxy) phenethylamine and N-hydroxy MDA); 17 (15) Parahexyl; some trade or other names: 3-hexyl-1-hydroxy-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-18 dibenzo (b,d) pyran; Synhexyl; 19 20 (16) Psilocybin; 21 (17) Psilocyn; 22 (18) Alpha-methyltryptamine (AMT); 23 (19) 2,5-dimethoxyamphetamine (2,5-dimethoxy-alpha-methylphenethylamine; 2,5-DMA); 24 (20) 4-bromo-2,5-dimethoxyamphetamine 25 26 (4-bromo-2,5-dimethoxy-alpha-methylphenethylamine;

4-bromo-2, 5-DMA); 1 2 (20.1) 4-Bromo-2,5 dimethoxyphenethylamine. Some trade or other names: 2-(4-bromo-3 4 2,5-dimethoxyphenyl)-1-aminoethane; 5 alpha-desmethyl DOB, 2CB, Nexus; (21) 4-methoxyamphetamine 6 7 (4-methoxy-alpha-methylphenethylamine; 8 paramethoxyamphetamine; PMA); 9 (22) (Blank); 10 (23) Ethylamine analog of phencyclidine. 11 Some trade or other names: 12 N-ethyl-1-phenylcyclohexylamine, 13 (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine, cyclohexamine, PCE; 14 15 (24) Pyrrolidine analog of phencyclidine. Some trade 16 or other names: 1-(1-phenylcyclohexyl) pyrrolidine, PCPy, 17 PHP; (25) 5-methoxy-3, 4-methylenedioxy-amphetamine; 18 (26) 2,5-dimethoxy-4-ethylamphetamine 19 20 (another name: DOET); (27) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine 21 22 (another name: TCPy); 23 (28) (Blank); (29) Thiophene analog of phencyclidine (some trade 24 25 or other names: 1-[1-(2-thienyl)-cyclohexyl]-piperidine; 26 2-thienyl analog of phencyclidine; TPCP; TCP);

1	(30) Bufotenine (some trade or other names:
2	3-(Beta-Dimethylaminoethyl)-5-hydroxyindole;
3	3-(2-dimethylaminoethyl)-5-indolol;
4	5-hydroxy-N,N-dimethyltryptamine;
5	N,N-dimethylserotonin; mappine);
6	(31) 1-Pentyl-3-(1-naphthoyl)indole
7	Some trade or other names: JWH-018;
8	(32) 1-Butyl-3-(1-naphthoyl)indole
9	Some trade or other names: JWH-073;
10	(33) 1-[(5-fluoropentyl)-1H-indol-3-yl]-
11	(2-iodophenyl)methanone
12	Some trade or other names: AM-694;
13	(34) 2-[(1R,3S)-3-hydroxycyclohexyl]-5-
14	(2-methyloctan-2-yl)phenol
15	Some trade or other names: CP 47,497
16	and its C6, C8 and C9 homologs;
17	(34.5) 2-[(1R,3S)-3-hydroxycyclohexyl]-5-
18	(2-methyloctan-2-yl) phenol), where side chain n=5;
19	and homologues where side chain $n=4$, 6, or 7; Some
20	trade or other names: CP 47,497;
21	(35) (6aR,10aR)-9-(hydroxymethyl)-6,6-dimethyl-3-
22	(2-methyloctan-2-yl)-6a,7,
23	10,10a-tetrahydrobenzo[c] chromen-1-ol
24	Some trade or other names: HU-210;
25	(35.5) (6aS,10aS)-9-(hydroxymethyl)-6,6-
26	dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-

tetrahydrobenzo[c] chromen-1-ol, its isomers, 1 2 salts, and salts of isomers; Some trade or other names: HU-210, Dexanabinol; 3 (36) Dexanabinol, (6aS,10aS)-9-(hydroxymethyl)-4 5 6,6-dimethyl-3-(2-methyloctan-2-yl)-6a, 7, 10, 10a-tetrahydrobenzo[c] chromen-1-ol 6 Some trade or other names: HU-211; 7 (37) (2-methyl-1-propyl-1H-indol-8 3-yl)-1-naphthalenyl-methanone 9 Some trade or other names: JWH-015; 10 11 (38) 4-methoxynaphthalen-1-yl-12 (1-pentylindol-3-yl)methanone 13 Some trade or other names: JWH-081; (39) 1-Pentyl-3-(4-methyl-1-naphthoyl) indole 14 15 Some trade or other names: JWH-122; (40) 2-(2-methylphenyl)-1-(1-pentyl-16 17 1H-indol-3-yl)-ethanone Some trade or other names: JWH-251; 18 (41) 1-(2-cyclohexylethyl)-3-19 20 (2-methoxyphenylacetyl) indole Some trade or other names: RCS-8, BTW-8 and SR-18; 21 22 (42) Any compound structurally derived from 23 3-(1-naphthoyl)indole or 1H-indol-3-yl-(1-naphthyl) methane by substitution at the 24 25 nitrogen atom of the indole ring by alkyl, haloalkyl, 26 alkenyl, cycloalkylmethyl, cycloalkylethyl or

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- 1 2

3

2-(4-morpholinyl)ethyl whether or not further substituted in the indole ring to any extent, whether or not substituted in the naphthyl ring to any extent;

- 4 (43) Any compound structurally derived from
 5 3-(1-naphthoyl)pyrrole by substitution at the nitrogen
 6 atom of the pyrrole ring by alkyl, haloalkyl, alkenyl,
 7 cycloalkylmethyl, cycloalkylethyl or
 8 2-(4-morpholinyl)ethyl, whether or not further
 9 substituted in the pyrrole ring to any extent, whether
 10 or not substituted in the naphthyl ring to any extent;
- 11 (44) Any compound structurally derived from
 12 1-(1-naphthylmethyl)indene by substitution
 13 at the 3-position of the indene ring by alkyl, haloalkyl,
 14 alkenyl, cycloalkylmethyl, cycloalkylethyl or
 15 2-(4-morpholinyl)ethyl whether or not further
 16 substituted in the indene ring to any extent, whether
 17 or not substituted in the naphthyl ring to any extent;

(45) Any compound structurally derived from 18 3-phenylacetylindole by substitution at the 19 20 nitrogen atom of the indole ring with alkyl, haloalkyl, 21 alkenyl, cycloalkylmethyl, cycloalkylethyl or 22 2-(4-morpholinyl)ethyl, whether or not further 23 substituted in the indole ring to any extent, whether or not substituted in the phenyl ring to any extent; 24 25 (46) Any compound structurally derived from

26 2-(3-hydroxycyclohexyl)phenol by substitution

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1	at the 5-position of the phenolic ring by alkyl,
2	haloalkyl, alkenyl, cycloalkylmethyl, cycloalkylethyl
3	or 2-(4-morpholinyl)ethyl, whether or not substituted
4	in the cyclohexyl ring to any extent;
5	(47) 3,4-Methylenedioxymethcathinone
6	Some trade or other names: Methylone;
7	(48) 3,4-Methyenedioxypyrovalerone
8	Some trade or other names: MDPV;
9	(49) 4-Methylmethcathinone
10	Some trade or other names: Mephedrone;
11	(50) 4-methoxymethcathinone;
12	(51) 4-Fluoromethcathinone;
13	(52) 3-Fluoromethcathinone;
14	(53) 2,5-Dimethoxy-4-(n)-propylthio-
15	phenethylamine;
16	(54) 5-Methoxy-N,N-diisopropyltryptamine;
17	(55) Pentedrone;
18	(56) 4-iodo-2,5-dimethoxy-N-((2-methoxy
19	phenyl)methyl)-benzeneethanamine
20	(trade or other name: 25I-NBOMe);
21	(57) 4-chloro-2,5-dimethoxy-N-[(2-methoxyphenyl)
22	methyl]-benzeneethanamine (trade or other name:
23	25C-NBOMe);
24	(58) 4-bromo-2,5-dimethoxy-N-[(2-methoxyphenyl)
25	methyl]-benzeneethanamine (trade or other name:
26	25B-NBOMe).

Unless specifically excepted or unless listed in 1 (e) 2 schedule, any material, compound, mixture, another or preparation which contains any quantity of the following 3 substances having a depressant effect on the central nervous 4 5 system, including its salts, isomers, and salts of isomers 6 whenever the existence of such salts, isomers, and salts of isomers is possible within the specific chemical designation: 7

8

mecloqualone;

9

(2) methaqualone; and

10

(3) gamma hydroxybutyric acid.

(f) 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, and salts of isomers:

16

17

(1) Fenethylline;

(2) N-ethylamphetamine;

18 (3) Aminorex (some other names:

19 2-amino-5-phenyl-2-oxazoline; aminoxaphen;

20 4-5-dihydro-5-phenyl-2-oxazolamine) and its

21 salts, optical isomers, and salts of optical isomers;

(4) Methcathinone (some other names:
23 2-methylamino-1-phenylpropan-1-one;
24 Ephedrone; 2-(methylamino)-propiophenone;
25 alpha-(methylamino)propiophenone; N-methylcathinone;
26 methycathinone; Monomethylpropion; UR 1431) and its

1	salts, optical isomers, and salts of optical isomers;
2	(5) Cathinone (some trade or other names:
3	2-aminopropiophenone; alpha-aminopropiophenone;
4	2-amino-1-phenyl-propanone; norephedrone);
5	(6) N,N-dimethylamphetamine (also known as:
6	N,N-alpha-trimethyl-benzeneethanamine;
7	N,N-alpha-trimethylphenethylamine);
8	(7) (+ or -) cis-4-methylaminorex ((+ or -) cis-
9	4,5-dihydro-4-methyl-4-5-phenyl-2-oxazolamine);
10	(8) 3,4-Methylenedioxypyrovalerone (MDPV).
11	(g) Temporary listing of substances subject to emergency
12	scheduling. Any material, compound, mixture, or preparation
13	that contains any quantity of the following substances:
14	(1) N-[1-benzyl-4-piperidyl]-N-phenylpropanamide
15	(benzylfentanyl), its optical isomers, isomers, salts,
16	and salts of isomers;
17	(2) N-[1(2-thienyl)
18	methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl),
19	its optical isomers, salts, and salts of isomers.
20	(h) If the Department adopts an emergency rule under
21	Section 10 of the Synthetic Drug Identification Pilot Program
22	Act adding a synthetic drug to the list of Schedule I
23	controlled substances that drug shall be automatically added as
24	a Schedule I controlled substance upon the effective date of
25	the emergency rule. This subsection (h) is inoperative on July
26	30, 2018.

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1 (Source: P.A. 97-192, eff. 7-22-11; 97-193, eff. 1-1-12; 97-194, eff. 7-22-11; 97-334, eff. 1-1-12; 97-813, eff.
3 7-13-12; 97-872, eff. 7-31-12; 98-987, eff. 1-1-15.)
4 Section 999. Effective date. This Act takes effect upon

5 becoming law.

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4 720 ILCS 570/204 from Ch. 56 1/2, par. 1204