

**Decree No**

**.../2023 of the Minister of Interior of [ date ]**

**amending Ministry of Interior Decree No 78/2022 of 28 December 2022 on controlled substances**

Based on the authorisation granted under Section 32(5a) of Act XCV of 2005 on medicinal products for human use and on the amendment of other regulations related to medicinal products, and acting within my duties as defined in Section 66(1)(26) of Government Decree No 182/2022 of 24 May 2022 on the scope of duties and powers of members of the Government, I hereby order the following:

**Section 1**

In Decree No 78/2022 of the Ministry of Interior of 28 December 2022 (hereinafter: Decree) the following Section 4 shall be inserted:

‘Section 4 The requirement for the prior notification of this draft regulation, as stipulated in Articles 5–7 of Directive (EU) 2015/1535 of the European Parliament and of the Council of 9 September 2015 laying down a procedure for the provision of information in the field of technical regulations and of rules on Information Society services, has been met.’

**Section 2**

Annex 1 to the Decree shall be replaced by *Annex 1* hereto.

**Section 3**

Annex 2 to the Decree shall be replaced by *Annex 2* hereto.

**Section 4**

Annex 3 to the Decree is amended in accordance with *Annex 3* hereto.

**Section 5**

This Decree shall enter into force on the third day following its publication.

**Section 6**

The requirement for the prior notification of this draft Decree, as stipulated in Articles 5–7 of Directive (EU) 2015/1535 of the European Parliament and of the Council of 9 September 2015 laying down a procedure for the provision of information in the field of technical regulations and of rules on Information Society services, has been met.

## **LISTS OF NARCOTIC DRUGS**

*Column A* of the tables in the lists of narcotic drugs in this Annex contains the International Non-proprietary Name (hereinafter: INN) of the compound or substance, as the official name. The English name is in bold, the Hungarian name is in brackets below that. If the current list published by the International Narcotic Control Board (hereinafter: INCB) contains no INN for the compound or substance, the first other name published in this current list will be taken as the official name. *Column B* of the tables in this Annex shows the other names published by INCB, separated by a comma. *Column C* of Lists K1 and K2 contains the chemical name and the description in italics, whereas the molecular formula for the chemical name is indicated in *column D*, and the chemical identifier InChIKey in *column E*. For the purposes of listing, the data in columns A to C shall prevail, and columns D and E ensure the computer-based searchability of the compounds.

### **1. List 1 of narcotic drugs (List K1)**

1.1. The following substances and compounds are classified as narcotics:

|   | A  | B   | C  | D   | E                            |
|---|--|---|--|---|------------------------------|
| 1 | <b>Official name</b><br>(Hungarian name)               | Other name or abbreviation, spelling commonly used abroad | Chemical name / <i>Description</i>   | Molecular formula                                 | InChIKey chemical identifier |
| 2 | 3-methylfentanyl*<br>(3-metilfentanil)                 |   | <i>N</i> -(3-methyl-1-phenethyl-4-piperidyl)propionanilide                     | C <sub>23</sub> H <sub>30</sub> N <sub>2</sub> O  | MLQRZXNZHAOCHQ-UHFFFAOYSA-N  |
| 3 | 3-methylthiofentanyl*<br>(3-metiltiofentanil)          |   | <i>N</i> -[3-methyl-1-[2-(2-thienyl)ethyl]-4-piperidyl]propionanilide          | C <sub>21</sub> H <sub>28</sub> N <sub>2</sub> OS | SRARDYUHGVMEQI-UHFFFAOYSA-N  |
| 4 | 4-fluoroisobutyrfentanyl<br>(4-fluoroizobutirfentanil) | 4-FIBF, pFIBF   | <i>N</i> -(4-fluorophenyl)- <i>N</i> -(1-phenethylpiperidin-4-yl)isobutyramide | C <sub>23</sub> H <sub>29</sub> FN <sub>2</sub> O | OZDOSQNUJIXEOR-UHFFFAOYSA-N  |
| 5 | Acetorphine*   |   | 3- <i>O</i> -acetyltetrahydro-7 $\alpha$ -(1-hydroxy-1-                        | C <sub>27</sub> H <sub>35</sub> NO <sub>5</sub>   | LFYBMMHFJIAKFE-              |

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|----|--|--|--|---|-----------------------------|
|    | (acetorfin)  |  | methylbutyl)-6,14- <i>endo</i> -ethenooripavine (derivative of thebaine)   |   | PMEKXCSPSA-N                |
| 6  | Acetyl-alpha-methylfentanyl* (acetil-alfa-metilfentanil) |  | <i>N</i> -[1-( $\alpha$ -methylphenethyl)-4-piperidyl]acetanilide  | C <sub>22</sub> H <sub>28</sub> N <sub>2</sub> O              | OKTLVZBUKMRPLL-UHFFFAOYSA-N |
| 7  | Acetylfentanyl* (acetilfentanil)                         | Desmethyl fentanyl (dezmetil fentanil) | <i>N</i> -phenyl- <i>N</i> -[1-(2-phenylethyl)-4-piperidiny]acetamide  | C <sub>21</sub> H <sub>26</sub> N <sub>2</sub> O              | FYIUUQUPOKIKNI-UHFFFAOYSA-N |
| 8  | Acetylmethadol (acetilmetadol)                           |  | 3-acetoxy-6-dimethylamino-4,4-diphenylheptane  | C <sub>23</sub> H <sub>31</sub> NO <sub>2</sub>               | XBMIVRRWGCYBTQ-UHFFFAOYSA-N |
| 9  | Acryloylfentanyl (akrilfentanil)                         | Acrylfentanyl                          | <i>N</i> -phenyl- <i>N</i> -[1-(2-phenylethyl)piperidin-4-yl]prop-2-enamide  | C <sub>22</sub> H <sub>26</sub> N <sub>2</sub> O              | RFQNLMWUIJJEQF-UHFFFAOYSA-N |
| 10 | Alfentanil (alfentanil)                                  |  | <i>N</i> -[1-[2-(4-ethyl-4,5-dihydro-5-oxo-1 <i>H</i> -tetrazol-1-yl)ethyl]-4-(methoxymethyl)-4-piperidiny]- <i>N</i> -phenylpropanamide | C <sub>21</sub> H <sub>32</sub> N <sub>6</sub> O <sub>3</sub> | IDBPHNDTYPBSNI-UHFFFAOYSA-N |
| 11 | Alphacetylmethadol (alfacetilmetadol)                    |  | $\alpha$ -3-acetoxy-6-dimethylamino-4,4-diphenylheptane  | C <sub>23</sub> H <sub>31</sub> NO <sub>2</sub>               | XBMIVRRWGCYBTQ-XMSQKQJNSA-N |
| 12 | Alphameprodine (alfameprodin)                            |  | $\alpha$ -3-ethyl-1-methyl-4-phenyl-4-propionoxypiperidine   | C <sub>17</sub> H <sub>25</sub> NO <sub>2</sub>               | ODEGQXRCQDVXSJ-WMLDXEAASA-N |
| 13 | Alphamethadol (alfametadol)                              |  | $\alpha$ -6-dimethylamino-4,4-diphenyl-3-heptanol  | C <sub>21</sub> H <sub>29</sub> NO                            | QIRAYNIFEOXSPW-YLJYHZDGSAN  |
| 14 | Alpha-methylfentanyl* (alfa-metilfentanil)               |  | <i>N</i> -[1-( $\alpha$ -methylphenethyl)-4-piperidyl]propionanilide   | C <sub>23</sub> H <sub>30</sub> N <sub>2</sub> O              | NGTVDHYUFBKWID-UHFFFAOYSA-N |
| 15 | Alpha-methylthiofentanyl* (alfa-metiltiofentanil)        |  | <i>N</i> -[1-[1-methyl-2-(2-thienyl)ethyl]-4-piperidyl]propionanilide  | C <sub>21</sub> H <sub>28</sub> N <sub>2</sub> OS             | YPOXDUYRRSUFFG-UHFFFAOYSA-N |
| 16 | Alfaprodine (alfaprodin)                                 |  | $\alpha$ -1,3-dimethyl-4-phenyl-4-propionoxypiperidine   | C <sub>16</sub> H <sub>23</sub> NO <sub>2</sub>               | UVAZQQHAVMNMHE-CJNGLKHVSA-N |
| 17 | Allylprodine (allilprodin)                               |  | 3-allyl-1-methyl-4-phenyl-4-propionoxypiperidine   | C <sub>18</sub> H <sub>25</sub> NO <sub>2</sub>               | KGYFOSCXVAXULR-UHFFFAOYSA-N |

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|----|--|--|--|---|-----------------------------|
| 18 | Anileridine<br>(anileridin)  |  | 1- <i>p</i> -aminophenethyl-4-phenylpiperidine-4-carboxylic acid ethyl ester             | C <sub>22</sub> H <sub>28</sub> N <sub>2</sub> O <sub>2</sub> | LKYQLAWMNBFNJT-UHFFFAOYSA-N |
| 19 | Benzethidine<br>(benzetidin)   |  | 1-(2-benzyloxyethyl)-4-phenylpiperidine-4-carboxylic acid ethyl ester                    | C <sub>23</sub> H <sub>29</sub> NO <sub>3</sub>               | UVTBZAWTRVBTMK-UHFFFAOYSA-N |
| 20 | Benzylmorphine<br>(benzilmorfin)   |  | 3-benzylmorphine   | C <sub>24</sub> H <sub>25</sub> NO <sub>3</sub>               | RDJGWRFTDZZXSM-RNWLQCGYSA-N |
| 21 | Betacetylmethadol<br>(bétacetilmetadol)  |  | $\beta$ -3-acetoxy-6-dimethylamino-4,4-diphenylheptane                                   | C <sub>23</sub> H <sub>31</sub> NO <sub>2</sub>               | XBMIVRRWGCYBTQ-GCJKJVERSA-N |
| 22 | Beta-hydroxy-3-methylfentanyl*<br>(béta-hidroxi-3-metilfentanil)                             |  | <i>N</i> -[1-( $\beta$ -hydroxyphenethyl)-3-methyl-4-piperidyl]propionanilide            | C <sub>23</sub> H <sub>30</sub> N <sub>2</sub> O <sub>2</sub> | FRPRNNRJTCONEC-UHFFFAOYSA-N |
| 23 | Beta-hydroxyfentanyl*<br>(béta-hidroxifentanil)  |  | <i>N</i> -[1-( $\beta$ -hydroxyphenethyl)-4-piperidyl]propionanilide                     | C <sub>22</sub> H <sub>28</sub> N <sub>2</sub> O <sub>2</sub> | JEFVHLMGRUJLET-UHFFFAOYSA-N |
| 24 | Betameprodine<br>(bétameprodin)  |  | $\beta$ -3-ethyl-1-methyl-4-phenyl-4-propionoxypiperidine                                | C <sub>17</sub> H <sub>25</sub> NO <sub>2</sub>               | ODEGQXRCQDVXSJ-RHSMWYFYSA-N |
| 25 | Betamethadol<br>(bétametadol)  |  | $\beta$ -6-dimethylamino-4,4-diphenyl-3-heptanol   | C <sub>21</sub> H <sub>29</sub> NO                            | QIRAYNIFEOXSPW-XLIONFOSSA-N |
| 26 | Betaprodine<br>(bétaprodin)  |  | $\beta$ -1,3-dimethyl-4-phenyl-4-propionoxypiperidine                                    | C <sub>16</sub> H <sub>23</sub> NO <sub>2</sub>               | UVAZQQHAVMNMHE-CZUORRHYSAN  |
| 27 | Bezitramide<br>(bezitramid)  |  | 1-(3-cyano-3,3-diphenylpropyl)-4-(2-oxo-3-propionyl-1-benzimidazolyl)-piperidine         | C <sub>31</sub> H <sub>32</sub> N <sub>4</sub> O <sub>2</sub> | FLKWNFFCSSJANB-UHFFFAOYSA-N |
| 28 | Brorphine<br>(brorfin)   |  | 1-{1-[1-(4-bromophenyl)ethyl]-piperidin-4-yl}-1,3-dihydro-2 <i>H</i> -benzimidazol-2-one | C <sub>20</sub> H <sub>22</sub> BrN <sub>3</sub> O            | CNOFBGYRMCBVLO-UHFFFAOYSA-N |
| 29 | Butyrfentanyl<br>(butirfentanil)   |  | <i>N</i> -phenyl- <i>N</i> -[1-(2-phenylethyl)-4-piperidinyl]butanamide                  | C <sub>23</sub> H <sub>30</sub> N <sub>2</sub> O              | QQOMYEQLWQJRKK-UHFFFAOYSA-N |
| 30 | Cannabis*, cannabis resin* and extracts and tinctures of cannabis<br>(Kannabisz*, kannabisz- |  |  | -   | -                           |

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|----|--|-----|---|---|----------------------------------|
|    | gyanta*, -extraktum és -<br>tinktúra)                  |     |   |   |                                  |
| 31 | Carfentanil*<br>(karfentanil)                          |     | methyl 1-(2-phenylethyl)-4-<br>[phenyl(propanoyl)amino]piperidine-4-<br>carboxylate   | C <sub>24</sub> H <sub>30</sub> N <sub>2</sub> O <sub>3</sub>   | YDSDEBIZUNNPOB-<br>UHFFFAOYSA-N  |
| 32 | Clonitazene<br>(klonitazén)                            |     | 2-( <i>p</i> -chlorobenzyl)-1-diethylaminoethyl-5-<br>nitrobenzimidazole  | C <sub>20</sub> H <sub>23</sub> ClN <sub>4</sub> O <sub>2</sub> | GPZLDQAEBHTMPG-<br>UHFFFAOYSA-N  |
| 33 | Coca leaf<br>(koka levél)                              |     |   | -   | -                                |
| 34 | Cocaine<br>(kokain)                                    |     | <i>methyl ester of benzoylecgonine</i>  | C <sub>17</sub> H <sub>21</sub> NO <sub>4</sub>                 | ZPUCINDJBIVPJ-<br>LJISPDSOSA-N   |
| 35 | Codoxime<br>(kodoxim)                                  |     | dihydrocodeinone-6-carboxymethyloxime   | C <sub>20</sub> H <sub>24</sub> N <sub>2</sub> O <sub>5</sub>   | WKJYCUVUZIIMJA-<br>HJWMAQEXSA-N  |
| 36 | Concentrate of poppy straw<br>(mákszalma koncentrátum) | CPS | <i>Material intended for commercial distribution,<br/>which was obtained from processing to enrich<br/>the alkaloid content of poppy straw.</i> | -   | -                                |
| 37 | Crotonylfentanyl<br>(krotonilfentanil)                 |     | ( <i>E</i> )- <i>N</i> -(1-phenethylpiperidin-4-yl)- <i>N</i> -phenylbut-<br>2-enamide  | C <sub>23</sub> H <sub>28</sub> N <sub>2</sub> O                | VDYXGPCGBKLRDA-<br>XNWCZRBMISA-N |
| 38 | Cyclopropylfentanyl<br>(ciklopropilfentanil)           |     | <i>N</i> -phenyl- <i>N</i> -[1-(2-phenylethyl)piperidin-4-<br>yl]cyclopropanecarboxamide  | C <sub>23</sub> H <sub>28</sub> N <sub>2</sub> O                | OIQSKDSKROTEMN-<br>UHFFFAOYSA-N  |
| 39 | Desomorphine*<br>(dezomorfin)                          |     | dihydrodeoxymorphine  | C <sub>17</sub> H <sub>21</sub> NO <sub>2</sub>                 | LNNWVNGFPYWNQE-<br>GMIGKAJZSA-N  |
| 40 | Dextromoramide<br>(dextromoramid)                      |     | (+)-4-[2-methyl-4-oxo-3,3-diphenyl-4-(1-<br>pyrrolidiny)butyl]-morpholine   | C <sub>25</sub> H <sub>32</sub> N <sub>2</sub> O <sub>2</sub>   | INUNXTSAACVKJS-<br>OAQYLSRUSA-N  |
| 41 | Diampromide<br>(diampromid)                            |     | <i>N</i> -[2-(methylphenethylamino)-<br>propyl]propionanilide   | C <sub>21</sub> H <sub>28</sub> N <sub>2</sub> O                | RXTHKWVSOIHJS-<br>UHFFFAOYSA-N   |
| 42 | Diethylthiambutene<br>(dietiltiambutén)                |     | 3-diethylamino-1,1-di-(2'-thienyl)-1-butene   | C <sub>16</sub> H <sub>21</sub> NS <sub>2</sub>                 | CBYWMRHUUVRIAF-<br>UHFFFAOYSA-N  |
| 43 | Difenoxin<br>(difenoxin)                               |     | 1-(3-cyano-3,3-diphenylpropyl)-4-<br>phenylisonipectic acid   | C <sub>28</sub> H <sub>28</sub> N <sub>2</sub> O <sub>2</sub>   | UFIVBRCCIRTJTN-<br>UHFFFAOYSA-N  |

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|----|--|--|---|---|-----------------------------|
| 44 | Dihydroetorphine<br>(dihidroetorfin)           |  | 7,8-dihydro-7 $\alpha$ -[1-( <i>R</i> )-hydroxy-1-methylbutyl]-6,14- <i>endo</i> -ethanotetrahydrooripavine | C <sub>25</sub> H <sub>35</sub> NO <sub>4</sub>               | BRTSNYPDACNMIP-FAWZKKEFSA-N |
| 45 | Dihydromorphine<br>(dihidromorfin)             |  | <i>morphine derivative</i>  | C <sub>17</sub> H <sub>21</sub> NO <sub>3</sub>               | IJVCMSMFSCRME-KBQPJGBKSA-N  |
| 46 | Dimenoxadol<br>(dimenoxadol)                   |  | 2-dimethylaminoethyl-1-ethoxy-1,1-diphenylacetate   | C <sub>20</sub> H <sub>25</sub> NO <sub>3</sub>               | RHUWRJWFHUKVED-UHFFFAOYSA-N |
| 47 | Dimepheptanol<br>(dimefeptanol)                |  | 6-dimethylamino-4,4-diphenyl-3-heptanol   | C <sub>21</sub> H <sub>29</sub> NO                            | QIRAYNIFEOXSPW-UHFFFAOYSA-N |
| 48 | Dimethylthiambutene<br>(dimetiltiambutén)      |  | 3-dimethylamino-1,1-di-(2'-thienyl)-1-butene  | C <sub>14</sub> H <sub>17</sub> NS <sub>2</sub>               | CANBGVXYBPOLRR-UHFFFAOYSA-N |
| 49 | Dioxaphetyl butyrate<br>(dioxafetil-butirát)   |  | ethyl-4-morpholino-2,2-diphenylbutyrate   | C <sub>22</sub> H <sub>27</sub> NO <sub>3</sub>               | LQGIXNQCOXNCRP-UHFFFAOYSA-N |
| 50 | Diphenoxylate<br>(difenoxilát)                 |  | 1-(3-cyano-3,3-diphenylpropyl)-4-phenylpiperidine-4-carboxylic acid ethyl ester                             | C <sub>30</sub> H <sub>32</sub> N <sub>2</sub> O <sub>2</sub> | HYPPXZBJBPSRLK-UHFFFAOYSA-N |
| 51 | Dipipanone<br>(dipipanon)                      |  | 4,4-diphenyl-6-piperidine-3-heptanone   | C <sub>24</sub> H <sub>31</sub> NO                            | SVDHSZFEQYXRDC-UHFFFAOYSA-N |
| 52 | Drotebanol<br>(drotebanol)                     |  | 3,4-dimethoxy-17-methylmorphinan-6 $\beta$ ,14-diol   | C <sub>19</sub> H <sub>27</sub> NO <sub>4</sub>               | LCAHPIFLPICNRW-SVYNMNNPSA-N |
| 53 | Ecgonine<br>(ekgonin)                          |  | <i>esters and derivatives thereof, which can be converted into ecgonine and cocaine</i>                     | C <sub>9</sub> H <sub>15</sub> NO <sub>3</sub>                | PHMBVCPLDPDESM-UHFFFAOYSA-N |
| 54 | Ethylmethylthiambutene<br>(etilmetiltiambutén) |  | 3-ethylmethylamino-1,1-di-(2'-thienyl)-1-butene   | C <sub>15</sub> H <sub>19</sub> NS <sub>2</sub>               | MORSAEFGQPDBKM-UHFFFAOYSA-N |
| 55 | Etonitazene<br>(etonitazén)                    |  | 1-diethylaminoethyl-2- <i>p</i> -ethoxybenzyl-5-nitrobenzimidazole  | C <sub>22</sub> H <sub>28</sub> N <sub>4</sub> O <sub>3</sub> | PXDBZSCGSQSKST-UHFFFAOYSA-N |
| 56 | Etorphine*<br>(etorfin)                        |  | tetrahydro-7 $\alpha$ -(1-hydroxy-1-methylbutyl)-6,14- <i>endo</i> -ethenoripavine                          | C <sub>25</sub> H <sub>33</sub> NO <sub>4</sub>               | CAHCBJPUTCKATP-UHFFFAOYSA-N |

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|----|---|--|--|---|------------------------------|
| 57 | Etoxidine<br>(etoxidin)                       |  | 1-[2-(2-hydroxyethoxy)-ethyl]-4-phenylpiperidine-4-carboxylic acid ethyl ester                                 | C <sub>18</sub> H <sub>27</sub> NO <sub>4</sub>               | KJTKYGFQPQRRA-UHFFFAOYSA-N   |
| 58 | Fentanyl<br>(fentanil)                        |  | 1-phenethyl-4- <i>N</i> -propionylanilinopiperidine  | C <sub>22</sub> H <sub>28</sub> N <sub>2</sub> O              | PJMPHNIQZUBGLI-UHFFFAOYSA-N  |
| 59 | Furanylfentanil<br>(furanilfentanil)          |  | <i>N</i> -phenyl- <i>N</i> -[1-(2-phenylethyl)piperidin-4-yl]furan-2-carboxamide                               | C <sub>24</sub> H <sub>26</sub> N <sub>2</sub> O <sub>2</sub> | FZJVHWISUGFFQV-UHFFFAOYSA-N  |
| 60 | Furethidine<br>(furetidin)                    |  | 1-(2-tetrahydrofurfuryloxyethyl)-4-phenylpiperidine-4-carboxylic acid ethyl ester                              | C <sub>21</sub> H <sub>31</sub> NO <sub>4</sub>               | NNCOZXNZFLUYGG-UHFFFAOYSA-N  |
| 61 | Heroin*<br>(heroin)                           |  | diacetylmorphine   | C <sub>21</sub> H <sub>23</sub> NO <sub>5</sub>               | GVGLGOZIDCSQPN-PVHGPHFFSA-N  |
| 62 | Hydrocodone<br>(hidrokodon)                   |  | dihydrocodeinone   | C <sub>18</sub> H <sub>21</sub> NO <sub>3</sub>               | LLPOLZWFYMWKNCMCMFDCUSA-N    |
| 63 | Hydromorfinol<br>(hidromorfinol)              |  | 14-hydroxydihydromorphine  | C <sub>17</sub> H <sub>21</sub> NO <sub>4</sub>               | AABLHGPVOULICIBRJGLHKUSA-N   |
| 64 | Hydromorphone<br>(hidromorfon)                |  | dihydromorphinone  | C <sub>17</sub> H <sub>19</sub> NO <sub>3</sub>               | WVLOADHCBXTIJKYNHQPCIGSA-N   |
| 65 | Hydroxypethidine<br>(hidroxipetidín)          |  | 4- <i>m</i> -hydroxyphenyl-1-methylpiperidine-4-carboxylic acid ethyl ester                                    | C <sub>15</sub> H <sub>21</sub> NO <sub>3</sub>               | WTJBNMUWRKPFERS-UHFFFAOYSA-N |
| 66 | Isomethadone<br>(izometadon)                  |  | 6-dimethylamino-5-methyl-4,4-diphenyl-3-hexanone   | C <sub>21</sub> H <sub>27</sub> NO                            | IFKPLJWIEQBPGG-UHFFFAOYSA-N  |
| 67 | Isotonitazene<br>(izotonitazén)               |  | <i>N,N</i> -diethyl-2-[2-(4-isopropoxybenzyl)-5-nitro-1 <i>H</i> -benzo[ <i>d</i> ]imidazol-1-yl]ethan-1-amine | C <sub>23</sub> H <sub>30</sub> N <sub>4</sub> O <sub>3</sub> | OIOQREYBGDAYGT-UHFFFAOYSA-N  |
| 68 | Ketobemidone*<br>(ketobemidon)                |  | 4- <i>m</i> -hydroxyphenyl-1-methyl-4-propionylpiperidine  | C <sub>15</sub> H <sub>21</sub> NO <sub>2</sub>               | ALFGKMXHOUSVAD-UHFFFAOYSA-N  |
| 69 | Levomethorphan <sup>1</sup><br>(levometorfán) |  | (-)-3-methoxy- <i>N</i> -methylmorphinan   | C <sub>18</sub> H <sub>25</sub> NO                            | MKXZASYAUGDDCJ-CGTJXYLNSA-N  |

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|----|---|--|---|---|---------------------------------|
| 70 | Levomoramide<br>(levomoramid)                   |  | (-)-4-[2-methyl-4-oxo-3,3-diphenyl-4-(1-pyrrolidinyl)-butyl]morpholine                                      | C <sub>25</sub> H <sub>32</sub> N <sub>2</sub> O <sub>2</sub> | INUNXTSAACVKJS-NR<br>FANRHFSA-N |
| 71 | Levophenacylmorphan<br>(levofenacilmorfán)      |  | (-)-3-hydroxy- <i>N</i> -phenacylmorphinan  | C <sub>24</sub> H <sub>27</sub> NO <sub>2</sub>               | RCYBMSQOSGJZLO-BG<br>WNEDDSSA-N |
| 72 | Levorphanol <sup>1</sup><br>(levorfanol)        |  | (-)-3-hydroxy- <i>N</i> -methyilmorphinan   | C <sub>17</sub> H <sub>23</sub> NO                            | JAQUASYNZVUNQP-US<br>XIJHARSA-N |
| 73 | Metazocine<br>(metazocin)                       |  | 2'-hydroxy-2,5,9-trimethyl-6,7-benzomorphan   | C <sub>15</sub> H <sub>21</sub> NO                            | YGSVZRIZCHZUHB-VR<br>FXOJNMSA-N |
| 74 | Methadone<br>(metadon)                          |  | 6-dimethylamino-4,4-diphenyl-3-heptanone  | C <sub>21</sub> H <sub>27</sub> NO                            | USSIQXCVUWKGNF-UH<br>FFFAOYSA-N |
| 75 | Methadone intermediate<br>(metadon intermediér) |  | 4-cyano-2-dimethylamino-4,4-diphenylbutane  | C <sub>19</sub> H <sub>22</sub> N <sub>2</sub>                | GJJQIGFCGLPOQK-UH<br>FFFAOYSA-N |
| 76 | Methoxyacetylfentanyl<br>(metoxiacetilfentanil) |  | 2-methoxy- <i>N</i> -phenyl- <i>N</i> -[1-(2-phenylethyl)piperidin-4-yl]acetamide                           | C <sub>22</sub> H <sub>28</sub> N <sub>2</sub> O <sub>2</sub> | SADNVKRDSWWFTK-UH<br>FFFAOYSA-N |
| 77 | Methyldesorphine<br>(metildezorfin)             |  | 6-methyl- $\Delta$ 6-deoxymorphine (derivative of morphine)   | C <sub>18</sub> H <sub>21</sub> NO <sub>2</sub>               | CUFWYVOFDYVCPM-GG<br>NLRSJOSA-N |
| 78 | Methyldihydromorphine<br>(metildihidromorfin)   |  | 6-methyldihydromorphine   | C <sub>18</sub> H <sub>23</sub> NO <sub>3</sub>               | NBKVWIJQJMEQLE-NG<br>TWOADLSA-N |
| 79 | Metonitazene<br>(metonitazén)                   |  | <i>N,N</i> -diethyl-2-[2-(4-methoxybenzyl)-5-nitro-1 <i>H</i> -benzo[ <i>d</i> ]imidazol-1-yl]ethan-1-amine | C <sub>21</sub> H <sub>26</sub> N <sub>4</sub> O <sub>3</sub> | HNGZTLMRQTVPBH-UH<br>FFFAOYSA-N |
| 80 | Metopon<br>(metopon)                            |  | 5-methyldihydromorphinone   | C <sub>18</sub> H <sub>21</sub> NO <sub>3</sub>               | NPZXCTIHHUUEEJ-CM<br>KMFDCUSA-N |
| 81 | Moramide intermediate<br>(moramid intermediér)  |  | 2-methyl-3-morpholino-1,1-diphenylpropane carboxylic acid   | C <sub>21</sub> H <sub>25</sub> NO <sub>3</sub>               | AWLNVHVUYACOMZ-UH<br>FFFAOYSA-N |
| 82 | Morpheridine<br>(morferidin)                    |  | 1-(2-morpholinoethyl)-4-phenylpiperidine-4-carboxylic acid ethyl ester                                      | C <sub>20</sub> H <sub>30</sub> N <sub>2</sub> O <sub>3</sub> | JDEDMCKQPKGSAX-UH<br>FFFAOYSA-N |
| 83 | Morphine<br>(morfin)                            |  | <i>the main alkaloid of opium and opium poppy</i>   | C <sub>17</sub> H <sub>19</sub> NO <sub>3</sub>               | BQJCRHHNABKAKU-KB<br>QPJGBKSA-N |
| 84 | Morphine methobromide                           |  | <i>morphine methobromide and other pentavalent</i>  | C <sub>18</sub> H <sub>22</sub> BrNO <sub>3</sub>             | KQUQZJSQMSHWHP-                 |



## DRAFT

|    |   |  |   |  |                                  |
|----|---|--|---|--|----------------------------------|
|    | (morfin metobromid)                         |  | <i>nitrogen-containing morphine derivatives including in particular the morphine-N-oxide derivatives, one of which is codeine-N-oxide</i> |  | SCLAZZCHSA-N                     |
| 85 | Morphine-N-oxide<br>(morfin-N-oxid)         |  | <i>morphine derivative</i>  | C <sub>17</sub> H <sub>19</sub> NO <sub>4</sub>                | AMAPEXTUMXQULJ-<br>APQDOHRLSA-N  |
| 86 | MPPP*                                       |  | 1-methyl-4-phenyl-4-piperidinol propionate<br>(ester)   | C <sub>15</sub> H <sub>21</sub> NO <sub>2</sub>                | BCQMRZRAWHNSBF-<br>UHFFFAOYSA-N  |
| 87 | Myrophine<br>(mirofin)                      |  | myristylbenzylmorphine  | C <sub>38</sub> H <sub>51</sub> NO <sub>4</sub>                | GODGZZGKTZQSAL-<br>VXFFQEMOSA-N  |
| 88 | Nicomorphine<br>(nikomorfin)                |  | 3,6-dinicotinylmorphine   | C <sub>29</sub> H <sub>25</sub> N <sub>3</sub> O <sub>5</sub>  | HNDXBGYRMHRUFN-<br>CIVUWBIHSA-N  |
| 89 | Noracymethadol<br>(noracimetadol)           |  | (±)-α-3-acetoxy-6-methylamino-4,4-<br>diphenylheptane   | C <sub>22</sub> H <sub>29</sub> NO <sub>2</sub>                | VWCUGCYZZGRKEE-<br>UHFFFAOYSA-N  |
| 90 | Norlevorphanol<br>(norlevorfanol)           |  | (-)-3-hydroxymorphinan  | C <sub>16</sub> H <sub>21</sub> NO                             | IYNWSQDZXMGGGI-<br>NUEKZKHPSA-N  |
| 91 | Normethadone<br>(normetadon)                |  | 6-dimethylamino-4,4-diphenyl-3-hexanone   | C <sub>20</sub> H <sub>25</sub> NO                             | WCJFBSYALHQBSK-<br>UHFFFAOYSA-N  |
| 92 | Normorphine<br>(normorfin)                  |  | demethylmorphine  | C <sub>16</sub> H <sub>17</sub> NO <sub>3</sub>                | ONBWJWYUHXVEJS-<br>ZTYRTETDSA-N  |
| 93 | Norpipanone<br>(norpipanon)                 |  | 4,4-diphenyl-6-piperidino-3-hexanone  | C <sub>23</sub> H <sub>29</sub> NO                             | WCDSHELZWCOTMI-<br>UHFFFAOYSA-N  |
| 94 | Ocfentanyl<br>(okfentanil)                  |  | <i>N</i> -(2-fluorophenyl)-2-methoxy- <i>N</i> -[1-(2-<br>phenylethyl)piperidin-4-yl]acetamide  | C <sub>22</sub> H <sub>27</sub> FN <sub>2</sub> O <sub>2</sub> | NYISTOZKVCMLVEL-<br>UHFFFAOYSA-N |
| 95 | Opium<br>(ópium)                            |  |   | -  | -                                |
| 96 | Oripavine<br>(oripavin)                     |  | <i>O</i> 3-demethylthebaine vagy 6,7,8,14-<br>tetrahydro-4,5- <i>alpha</i> -epoxy-6-methoxy-17-<br>methylmorphinan-3-ol                   | C <sub>18</sub> H <sub>19</sub> NO <sub>3</sub>                | ZKLXUUYLEHCAMF-<br>UWFMWQGSAN    |
| 97 | Orthofluorofentanyl<br>(ortofluorofentanil) |  | <i>N</i> -(2-fluorophenyl)- <i>N</i> -[1-(2-<br>phenylethyl)piperidin-4-yl]propanamide  | C <sub>22</sub> H <sub>27</sub> FN <sub>2</sub> O              | BKUWDIVZCJNXRA-<br>UHFFFAOYSA-N  |

## DRAFT

|     |  |  |   |   |                                  |
|-----|--|--|---|---|----------------------------------|
| 98  | Oxycodone<br>(oxikodon)                                  |  | 14-hydroxydihydrocodeinone  | C <sub>18</sub> H <sub>21</sub> NO <sub>4</sub>   | BRUQQQPBMZOVGD-<br>XFKAJCMBSA-N  |
| 99  | Oxymorphone<br>(oximorfon)                               |  | 14-hydroxydihydromorphinone   | C <sub>17</sub> H <sub>19</sub> NO <sub>4</sub>   | UQCNKQCJZOFTQ-<br>ISWURRPUA-N    |
| 100 | Parafluorobutyrylfentanyl<br>(parafluorobutirilfentanil) |  | <i>N</i> -(4-fluorophenyl)- <i>N</i> -[1-(2-phenylethyl)piperidin-4-yl]butanamide | C <sub>23</sub> H <sub>29</sub> FN <sub>2</sub> O | QZFMICYUBPSLOBP-<br>UHFFFAOYSA-N |
| 101 | Para-fluorofentanyl*<br>(para-fluorofentanil)            |  | 4'-fluoro- <i>N</i> -(1-phenethyl-4-piperidyl)propionanilide                      | C <sub>22</sub> H <sub>27</sub> FN <sub>2</sub> O | KXUBAVLIJFTASZ-<br>UHFFFAOYSA-N  |
| 102 | PEPAP*   |  | 1-phenethyl-4-phenyl-4-piperidinol acetate<br>(ester)                             | C <sub>21</sub> H <sub>25</sub> NO <sub>2</sub>   | BVURVTVDNWSNFN-<br>UHFFFAOYSA-N  |
| 103 | Pethidine<br>(petidin)                                   |  | 1-methyl-4-phenylpiperidine-4-carboxylic acid<br>ethyl ester                      | C <sub>15</sub> H <sub>21</sub> NO <sub>2</sub>   | XADCESSVHJOZHK-<br>UHFFFAOYSA-N  |
| 104 | Pethidine intermediate A<br>(petidin A intermediar)      |  | 4-cyano-1-methyl-4-phenylpiperidine   | C <sub>13</sub> H <sub>16</sub> N <sub>2</sub>    | ZLFQTZYFXYOGLS-<br>UHFFFAOYSA-N  |
| 105 | Pethidine intermediate B<br>(petidin B intermediar)      |  | 4-phenylpiperidine-4-carboxylic acid ethyl ester                                  | C <sub>14</sub> H <sub>19</sub> NO <sub>2</sub>   | QKHMFBKXTNQCTM-<br>UHFFFAOYSA-N  |
| 106 | Pethidine intermediate C<br>(petidin C intermediar)      |  | 1-methyl-4-phenylpiperidine-4-carboxylic acid                                     | C <sub>13</sub> H <sub>17</sub> NO <sub>2</sub>   | KHUPPYUUMRDAAX-<br>UHFFFAOYSA-N  |
| 107 | Phenadoxone<br>(fenadoxon)                               |  | 6-morpholino-4,4-diphenyl-3-heptanone   | C <sub>23</sub> H <sub>29</sub> NO <sub>2</sub>   | LOXCOAXRHYDLOW-<br>UHFFFAOYSA-N  |
| 108 | Phenampromide<br>(fenampromid)                           |  | <i>N</i> -(1-methyl-2-piperidinoethyl)propionanilide                              | C <sub>17</sub> H <sub>26</sub> N <sub>2</sub> O  | DHTRHEVNFFZCNU-<br>UHFFFAOYSA-N  |
| 109 | Phenazocine<br>(fenazocin)                               |  | 2'-hydroxy-5,9-dimethyl-2-phenethyl-6,7-<br>benzomorphan                          | C <sub>22</sub> H <sub>27</sub> NO                | ZQHYKVKNPWDQSL-<br>UHFFFAOYSA-N  |
| 110 | Phenomorphane<br>(fenomorfan)                            |  | 3-hydroxy- <i>N</i> -phenethylmorphinan   | C <sub>24</sub> H <sub>29</sub> NO                | CFBQYWXPZVQQTN-<br>QPTUXGOLSA-N  |
| 111 | Phenoperidine<br>(fenoperidin)                           |  | 1-(3-hydroxy-3-phenylpropyl)-4-<br>phenylpiperidine-4-carboxylic acid ethyl ester | C <sub>23</sub> H <sub>29</sub> NO <sub>3</sub>   | IPOPQVVNCFQFRK-<br>UHFFFAOYSA-N  |

## DRAFT

|     |  |       |   |   |                              |
|-----|--|-------|---|---|------------------------------|
| 112 | Piminodine<br>(piminodin)                                |       | 4-phenyl-1-(3-phenylaminopropyl)-piperidine-4-carboxylic acid ethyl ester                                   | C <sub>23</sub> H <sub>30</sub> N <sub>2</sub> O <sub>2</sub>   | PXXKIYPSXYFATG-UHFFFAOYSA-N  |
| 113 | Piritramide<br>(piritramid)                              |       | 1-(3-cyano-3,3-diphenylpropyl)-4-(1-piperidino)piperidine-4-carboxylic acid amide                           | C <sub>27</sub> H <sub>34</sub> N <sub>4</sub> O                | IHEHEFLXQFOQJO-UHFFFAOYSA-N  |
| 114 | Proheptazine<br>(proheptazin)                            |       | 1,3-dimethyl-4-phenyl-4-propionoxyazacycloheptane   | C <sub>17</sub> H <sub>25</sub> NO <sub>2</sub>                 | ZXWAUWBYASJEOE-UHFFFAOYSA-N  |
| 115 | Properidine<br>(properidin)                              |       | 1-methyl-4-phenylpiperidine-4-carboxylic acid isopropyl ester   | C <sub>16</sub> H <sub>23</sub> NO <sub>2</sub>                 | XJKQCILVUHXXVIQ-UHFFFAOYSA-N |
| 116 | Racemethorphan<br>(racemetorfán)                         |       | (±)-3-methoxy- <i>N</i> -methylmorphinan  | C <sub>18</sub> H <sub>25</sub> NO                              | MKXZASYAUGDDCJ-CGTJXYLNSA-N  |
| 117 | Racemoramide<br>(racemoramid)                            |       | (±)-4-[2-methyl-4-oxo-3,3-diphenyl-4-(1-pyrrolidinyl)-butyl]-morpholine                                     | C <sub>25</sub> H <sub>32</sub> N <sub>2</sub> O <sub>2</sub>   | INUNXTSAACVKJS-UHFFFAOYSA-N  |
| 118 | Racemorphan<br>(racemorfán)                              |       | (±)-3-hydroxy- <i>N</i> -methylmorphinan  | C <sub>17</sub> H <sub>23</sub> NO                              | JAQUASYNZVUNQP-USXIJHARSA-N  |
| 119 | Remifentanil<br>(remifentanil)                           |       | 1-(2-methoxy-carbonyl-ethyl)-4-(phenyl- <i>N</i> -propionylamino)-piperidine-4-carboxylic acid methyl ester | C <sub>20</sub> H <sub>28</sub> N <sub>2</sub> O <sub>5</sub>   | ZTVQQQVZCWLTFD-UHFFFAOYSA-N  |
| 120 | Sufentanil<br>(szufentanil)                              |       | <i>N</i> -[4-(methoxymethyl)-1-[2-(2-thienyl)-ethyl]-4-piperidyl]propionanilide                             | C <sub>22</sub> H <sub>30</sub> N <sub>2</sub> O <sub>2</sub> S | GGCSSNBKKAUURC-UHFFFAOYSA-N  |
| 121 | Tetrahydrofuranylfentanyl<br>(tetrahidrofuranilfentanil) | THF-F | <i>N</i> -phenyl- <i>N</i> -[1-(2-phenylethyl)piperidin-4-yl]tetrahydrofuran-2-carboxamide                  | C <sub>24</sub> H <sub>30</sub> N <sub>2</sub> O <sub>2</sub>   | OHJNHKUFKAANI-UHFFFAOYSA-N   |
| 122 | Thebacon<br>(tebakon)                                    |       | acetyldihydrocodeinone  | C <sub>20</sub> H <sub>23</sub> NO <sub>4</sub>                 | RRJQTGHQFYTZOW-ILWKUFEGSA-N  |
| 123 | Thebaine<br>(tebain)                                     |       | <i>opium alkaloid, found in Papaver bracteatum Lindl.</i>   | C <sub>19</sub> H <sub>21</sub> NO <sub>3</sub>                 | FQXXSQDCDRQNQE-VMDGZTHMSA-N  |
| 124 | Thiofentanyl*<br>(tiofentanil)                           |       | <i>N</i> -[1-[2-(2-thienyl)ethyl]-4-piperidyl]propionanilide  | C <sub>20</sub> H <sub>26</sub> N <sub>2</sub> OS               | YMRFZDHYDKZXP-UHFFFAOYSA-N   |
| 125 | Tilidine<br>(tilidin)                                    |       | (±)-ethyl- <i>trans</i> -2-(dimethylamino)-1-phenyl-3-cyclohexene-1-carboxylate                             | C <sub>17</sub> H <sub>23</sub> NO <sub>2</sub>                 | WDEFBBTXULIOBB-UHFFFAOYSA-N  |

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|     |                                     |  |   |  |                             |
|-----|-------------------------------------|--|---|--|-----------------------------|
| 126 | Trimeperidine<br>(trimeperidin)     |  | 1,2,5-trimethyl-4-phenyl-4-propionoxypiperidine                                   | C <sub>17</sub> H <sub>25</sub> NO <sub>2</sub>                  | UVITTYOJFDLOGI-UHFFFAOYSA-N |
| 127 | U-47700                             |  | 3,4-dichloro- <i>N</i> -[2-(dimethylamino)cyclohexyl]- <i>N</i> -methyl-benzamide | C <sub>16</sub> H <sub>22</sub> Cl <sub>2</sub> N <sub>2</sub> O | JGPNMZWFVRQNGU-UHFFFAOYSA-N |
| 128 | Valerylentanyl<br>(valerilfentanil) |  | <i>N</i> -(1-phenethylpiperidin-4-yl)- <i>N</i> -phenylpentanamide                | C <sub>24</sub> H <sub>32</sub> N <sub>2</sub> O                 | VCCPXHWAJYWQMR-UHFFFAOYSA-N |

1.2. In addition, isomers of the above substances, provided that they correspond to the chemical name indicated, exist on the basis of their chemical structure and are not expressly covered by exceptional provisions, and their esters and ethers, in so far as such esters and ethers exist and do not appear on any other list, and their salts, including those of esters, ethers, isomers, where such salts exist.

This List K1 is identical in content to the updated Schedule I of the Single Convention on Narcotic Drugs signed in New York on 30 March 1961 (hereinafter: Narcotics Convention).

1.3. In the case of substances marked with the symbol <sup>1</sup>, the dextromethorphan (dextrometorfán) [(+)-3-methoxy-*N*-methylmorphinan] and dextrorphan (dextrorfán) [(+)-3-hydroxy-*N*-methylmorphinan] isomers are not subject to international control.

1.4. Substances marked with \* are subject to increased international control and are also included in Schedule IV of the Narcotics Convention.

## 2. List 2 of narcotic drugs (List K2)

2.1. The following substances and compounds are classified as narcotics:

|   | <b>A</b>                                 | <b>B</b>  | <b>C</b>                           | <b>D</b>          | <b>E</b>                     |
|---|--|---|------------------------------------|-------------------|------------------------------|
| 1 | <b>Official name</b><br>(Hungarian name) | Other name or abbreviation, spelling commonly used abroad | Chemical name / <i>Description</i> | Molecular formula | InChIKey chemical identifier |

## DRAFT

|    |   |  |  |   |                                 |
|----|---|--|--|---|---------------------------------|
| 2  | Acetyldihydrocodeine<br>(acetildihidrokodein) |  | <i>codeine derivative</i>  | C <sub>20</sub> H <sub>25</sub> NO <sub>4</sub>               | LGGDXXJAGWBUSL-<br>BKRJIHRRSA-N |
| 3  | Codeine<br>(kodein)                           |  | 3-methylmorphine   | C <sub>18</sub> H <sub>21</sub> NO <sub>3</sub>               | OROGSEYTTFOCAN-<br>DNJOTXNNSA-N |
| 4  | Dextropropoxyphene<br>(dextropropoxifén)      |  | $\alpha$ -(+)-4-dimethylamino-1,2-diphenyl-3-methyl-2-<br>butanol propionate | C <sub>22</sub> H <sub>29</sub> NO <sub>2</sub>               | XLMALTXPSGQGBX-<br>GCJKJVERSA-N |
| 5  | Dihydrocodeine<br>(dihidrokodein)             |  | <i>morphine derivative</i>   | C <sub>18</sub> H <sub>23</sub> NO <sub>3</sub>               | RBOXVHNMFORY-<br>UHFFFAOYSA-N   |
| 6  | Ethylmorphine<br>(etilmorfin)                 |  | 3-ethylmorphine  | C <sub>19</sub> H <sub>23</sub> NO <sub>3</sub>               | OGDVEMNWJVYAJL-<br>LEPYJNQMSA-N |
| 7  | Nicocodeine<br>(nikokodin)                    |  | 6-nicotinylcodeine   | C <sub>24</sub> H <sub>24</sub> N <sub>2</sub> O <sub>4</sub> | RYBGRHAWFUVMS-<br>MJFIPZRTSA-N  |
| 8  | Nicodicodine<br>(nikodikodin)                 |  | 6-nicotinyldihydrocodeine  | C <sub>24</sub> H <sub>26</sub> N <sub>2</sub> O <sub>4</sub> | GTGRMWCOZHEYRL-<br>MJFIPZRTSA-N |
| 9  | Norcodeine<br>(norkodein)                     |  | <i>N</i> -demethylcodeine  | C <sub>17</sub> H <sub>19</sub> NO <sub>3</sub>               | HKOIXWVRNLGFOR-<br>KOFBORESSA-N |
| 10 | Pholcodine<br>(folkodin)                      |  | morpholinylethylmorphine   | C <sub>23</sub> H <sub>30</sub> N <sub>2</sub> O <sub>4</sub> | GPFJKDEDBRFOS-<br>FKQDBXSBSA-N  |
| 11 | Propiram<br>(propirám)                        |  | <i>N</i> -(1-methyl-2-piperidinoethyl)- <i>N</i> -2-<br>pyridylpropionamide  | C <sub>16</sub> H <sub>25</sub> N <sub>3</sub> O              | ZBAFFZBKCMWUHM-<br>UHFFFAOYSA-N |

2.2. In addition, the isomers of the above substances, if they correspond to the chemical name indicated, exist on the basis of their chemical structure and are not expressly covered by exceptional provisions, and their salts, including those of their isomers, in so far as such salts may exist.

2.3. List K2 is identical in content to the updated Schedule II of the Narcotics Convention. Preparations of the active substances listed here may be included in List K3.

### 3. List of exceptions containing narcotic drugs (List K3)

3.1. Narcotic drug-containing preparations as defined below shall be excluded in accordance with the provisions laid down in this regulation.

3.2. Single or multiple component medicinal products containing the following active substances:

3.2.1. acetyldihydrocodeine (acetildihidrokodein),

3.2.2. codeine (kodein),

3.2.3. **dihydrocodeine** (dihidrokodein),

3.2.4. ethylmorphine (etilmorfin),

3.2.5. **nicocodeine** (nikokodin),

3.2.6. **nicodicodine** (nikodikodin),

3.2.7. **norcodeine** (norkodein),

3.2.8. **pholcodine** (folkodin),

where these preparations contain one or more components and the amount of the narcotic drug does not exceed 100 mg per dosage unit and the concentration of the drug in undivided preparations (e.g. solution) is not higher than 2.5 %, except for injectable preparations.

3.3. Complex medicinal products containing up to 100 mg of **propiram** (propiram) and methylcellulose in a quantity at least identical to propiram.

3.4. Medicinal products containing **dextropropoxyphene** (dextropropoxifén) which are administered orally and contain no more than 135 mg dextropropoxyphene (calculated as base) per dosage unit, or undivided medicinal products (e.g. solution) in which the dextropropoxyphene concentration, calculated as base, does not exceed 2.5 %, provided that these preparations do not contain any other active substances covered by the Convention on Psychotropic Substances signed in Vienna on 21 February 1971 (hereinafter: the Psychotropic Convention).

3.5. Those complex **cocaine** (kokain) medicinal products which contain up to 0.1 % of cocaine (calculated as base) and those complex opium (ópium) or morphine (morfin) products which contain up to 0.2 % of morphine (calculated as anhydrous morphine base) and which contain one or more components in a way that disables the re-extraction of narcotics by an easily implementable method or in such a quantity that would pose a public health risk.

3.6. Those medicinal products containing **difenoxin** (difenoxin) in which the quantity of difenoxin does not exceed 0.5 mg per dosage unit and contain atropin-sulphate in a quantity identical to at least 5 % of the difenoxin content.

3.7. Those medicinal products containing diphenoxylate (difenoxilát) in which the diphenoxylate content (calculated as base) per dosage unit does not exceed 2.5 mg and which contain atropin-sulphate in a quantity identical to at least 1 % of the diphenoxylate content.

3.8. Medicinal products with the following composition:

3.8.1. 10 g Pulvis opii

3.8.2. 10 g *Ipecacuanhae radix et rhizoma*

3.8.3. 80 g *Saccharosum* (or other powder vehicle not containing active substance).

3.9. Medicinal products with identical product characteristics to any of those included in the present Schedule, and mixtures thereof with any substance that does not contain narcotic drugs. '

**LISTS OF PSYCHOTROPIC SUBSTANCES**

*Column A* of the table in the list of psychotropic substances in this Annex contains the International Non-Proprietary Name (hereinafter: INN) of the compound or substance, as the official name. The English name is written in bold, the Hungarian name is in brackets below that. If the current list published by the International Narcotic Control Board (hereinafter: INCB) contains no INN for the compound or substance, the first other name published in this current list will be taken as the official name. *Column B* of the tables in this Annex shows the other names published by INCB, any other names if appropriate, abbreviations, and spelling commonly used abroad, separated by a comma. In *Column C* of the lists the chemical name is shown together with the description in italics, whereas the molecular formula for the chemical name is indicated in *column D* and the chemical identifier InChIKey in *column E*. For the purposes of listing, the data in columns A to C shall prevail, and columns D and E ensure the computer-based searchability of the compounds.

**1. List 1 of psychotropic substances (List P1)**

1.1. The following substances and compounds are classified as psychotropic substances:

|   | <b>A</b>                                 | <b>B</b>  | <b>C</b>   | <b>D</b>                           | <b>E</b>                        |
|---|--|---|--|------------------------------------|---------------------------------|
| 1 | <b>Official name</b><br>(Hungarian name) | Other name or<br>abbreviation, spelling<br>commonly used abroad | Chemical name / <i>Description</i>                           | Molecular<br>formula               | InChIKey chemical<br>identifier |
| 2 | 1-naphyrone***<br>(1-nafiron)            |   | 1-(naphthalen-1-yl)-2-(pyrrolidin-1-yl)pentan-1-one          | C <sub>19</sub> H <sub>23</sub> NO | PDYIKONOBSEMS-<br>UHFFFAOYSA-N  |
| 3 | 1-PEA***                                 |   | 1-amino-1-phenylethane; <i>1-phenethylamine</i> <sup>o</sup> | C <sub>8</sub> H <sub>11</sub> N   | RQEUFKEYXDPUK-<br>UHFFFAOYSA-N  |



## DRAFT

|    |   |                 |   |   |                              |
|----|---|-----------------|---|---|------------------------------|
| 4  | 25I-NBOMe   | 2C-I-NBOMe      | 2-(4-iodo-2,5-dimethoxyphenyl)- <i>N</i> -(2-methoxybenzyl)ethanamine   | C <sub>18</sub> H <sub>22</sub> INO <sub>3</sub>                | ZFUOLNAKPBFDIJ-UHFFFAOYSA-N  |
| 5  | 2-aminoindane***<br>(2-aminoindán)                              |                 | 2,3-dihydro-1 <i>H</i> -inden-2-amine   | C <sub>9</sub> H <sub>11</sub> N                                | LMHHFZAXSANGGM-UHFFFAOYSA-N  |
| 6  | 2C-I <sup>EU</sup>  |                 | 2,5-dimethoxy-4-iodophenethylamine <sup>EU2</sup> ; 4-Iod-2,5-dimethoxyphenethylazan <sup>oo</sup>                                | C <sub>10</sub> H <sub>14</sub> INO <sub>2</sub>                | PQHQB RJAAZQXHL-UHFFFAOYSA-N |
| 7  | 2C-T-2 <sup>EU</sup>  |                 | 2,5-dimethoxy-4-ethylthiophenethylamine <sup>EU2</sup> ; 4-Ethylsulfanyl-2,5-dimethoxy-phenethylazan <sup>oo</sup>                | C <sub>12</sub> H <sub>19</sub> NO <sub>2</sub> S               | HCWQGDLBIKOJPM-UHFFFAOYSA-N  |
| 8  | 2C-T-7 <sup>EU</sup>  |                 | 2,5-dimetoxi-4-( <i>n</i> )-propil-tio-fenetil-amin <sup>EU2</sup> ; 2,5-Dimethoxy-4-(propylsulfanyl)-phenethylazan <sup>oo</sup> | C <sub>13</sub> H <sub>21</sub> NO <sub>2</sub> S               | OLEVEPDJOF PJTF-UHFFFAOYSA-N |
| 9  | 2-DPMP***<br>(dezoxipipradrol)                                  | Desoxypipradrol | 2-(diphenylmethyl)piperidine  | C <sub>18</sub> H <sub>21</sub> N                               | RWTNXJXZVGHMGI-UHFFFAOYSA-N  |
| 10 | 2-naphyrone***<br>(2-nafiron)                                   |                 | 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one   | C <sub>19</sub> H <sub>23</sub> NO                              | DTNUPBSOODGRKW-UHFFFAOYSA-N  |
| 11 | 3,4-dichloromethylphenidate***                                  | 3,4-CTMP        | methyl-2-(3,4-dichlorophenyl)-2-(piperidin-2-yl)acetate   | C <sub>14</sub> H <sub>17</sub> Cl <sub>2</sub> NO <sub>2</sub> | JUKMAYKVHWKRKY-UHFFFAOYSA-N  |
| 12 | 3,4-methylenedioxyprovalerone**<br>(3,4-metiléndioxiprovaleron) | MDPV            | ( <i>RS</i> )-1-(benzo[ <i>d</i> ][1,3]dioxol-5-yl)-2-(pyrrolidin-1-yl)pentan-1-one   | C <sub>16</sub> H <sub>21</sub> NO <sub>3</sub>                 | SYHGEUNFJIGTRX-UHFFFAOYSA-N  |

## DRAFT

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|----|--|---|--|--|-----------------------------|
| 13 | 3F-phenmetrazine***                      | meta-fluoro-phenmetrazine, meta-F-phenmetrazine | 2-(3-fluorophenyl)-3-methylmorpholine  | C <sub>11</sub> H <sub>14</sub> FNO                            | VHYVKJAQSJCYCK-UHFFFAOYSA-N |
| 14 | 3-MeO-PCE***                             |   | <i>N</i> -ethyl-1-(3-methoxyphenyl)cyclohexanamine   | C <sub>15</sub> H <sub>23</sub> NO                             | OFGOOZLOGUNDFS-UHFFFAOYSA-N |
| 15 | 3-MeO-PCP**                              | 3-Methoxyphencyclidine                          | 1-[1-(3-methoxyphenyl)cyclohexyl]-piperidine   | C <sub>18</sub> H <sub>27</sub> NO                             | BQQSZHHKGPOXLN-UHFFFAOYSA-N |
| 16 | 4-MA <sup>EU</sup><br>(4-metilamfetamin) | 4-methylamphetamine                             | 1-(4-methylphenyl)propan-2-amine   | C <sub>10</sub> H <sub>15</sub> N                              | ZDHZDWSHLNBTEB-UHFFFAOYSA-N |
| 17 | 4-benzylpiperidine***                    |   | 4-(phenylmethyl)piperidine   | C <sub>12</sub> H <sub>17</sub> N                              | ABGXADJDTPFFSZ-UHFFFAOYSA-N |
| 18 | 4F-MDMB-BICA <sup>EU</sup>               |   | methyl 2-{{1-(4-fluorobutyl)-1 <i>H</i> -indole-3-carbonyl}amino}-3,3-dimethylbutanoate            | C <sub>20</sub> H <sub>27</sub> FN <sub>2</sub> O <sub>3</sub> | QIKHYQCWUGFBB-UHFFFAOYSA-N  |
| 19 | 4-MEC**                                  | 4-methylethcathinone                            | 2-(ethylamino)-1-(4-methylphenyl)propan-1-one  | C <sub>12</sub> H <sub>17</sub> NO                             | ZOXZWYWOECCBSH-UHFFFAOYSA-N |
| 20 | 4-MeO-PCP***                             |   | 1-[1-(4-methoxyphenyl)cyclohexyl]piperidine  | C <sub>18</sub> H <sub>27</sub> NO                             | MUZGGFNYPVLGUF-UHFFFAOYSA-N |
| 21 | 4-methylaminorex<br>(4-metilaminorex)    |   | (±)- <i>cis</i> -2-amino-4-methyl-5-phenyl-2-oxazoline   | C <sub>10</sub> H <sub>12</sub> N <sub>2</sub> O               | LJQBMYDFWFGESC-CBAPKCEASA-N |
| 22 | 4-MTA                                    |   | α-methyl-4-methylthiophenethylamine  | C <sub>10</sub> H <sub>15</sub> NS                             | OLEWMKVPSUCNLG-UHFFFAOYSA-N |
| 23 | 5F-AB-PINACA***                          |   | <i>N</i> -(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1 <i>H</i> -indazole-3-carboxamide | C <sub>18</sub> H <sub>25</sub> FN <sub>4</sub> O <sub>2</sub> | WCBYXIBEPFZUBG-UHFFFAOYSA-N |

## DRAFT

|    |                            |   |   |  |                             |
|----|----------------------------|---|---|--|-----------------------------|
| 24 | 5-IAI***                   |   | 5-iodo-2,3-dihydro-1 <i>H</i> -inden-2-amine  | C <sub>9</sub> H <sub>10</sub> IN                                | BIHPYCDDPGNWQO-UHFFFAOYSA-N |
| 25 | 5-IT <sup>EU</sup>         |   | 5-(2-aminopropyl)indole   | C <sub>11</sub> H <sub>14</sub> N <sub>2</sub>                   | AULGMISRJWGTBA-UHFFFAOYSA-N |
| 26 | 5-MeO-AMT***               |   | 1-(5-methoxy-1 <i>H</i> -indol-3-yl)propan-2-amine  | C <sub>12</sub> H <sub>16</sub> N <sub>2</sub> O                 | OGNJZVNNKBZFRM-UHFFFAOYSA-N |
| 27 | A-836,339***               |   | <i>N</i> -[3-(2-methoxyethyl)-4,5-dimethyl-1,3-thiazol-2-ylidene]-2,2,3,3-tetramethylcyclopropane-carboxamide | C <sub>16</sub> H <sub>26</sub> N <sub>2</sub> O <sub>2</sub> S  | JKGIMVBQKSRTGX-UHFFFAOYSA-N |
| 28 | AH-7921 <sup>EU</sup>      |   | 3,4-dichloro- <i>N</i> -{[1-(dimethylamino)cyclohexyl]methyl}benzamide  | C <sub>16</sub> H <sub>22</sub> Cl <sub>2</sub> N <sub>2</sub> O | JMZROFPPEXCTST-UHFFFAOYSA-N |
| 29 | AL***                      |   | 4-allyloxy-3,5-dimethoxyphenethylamine °  | C <sub>13</sub> H <sub>19</sub> NO <sub>3</sub>                  | JNUAYHHGCXYBHX-UHFFFAOYSA-N |
| 30 | alpha-PBT***<br>(alfa-PBT) | α-PBT, α-Pyrrolidinobutiothiophenone      | 2-(pyrrolidin-1-yl)-1-(thiophen-2-yl)butan-1-one  | C <sub>12</sub> H <sub>17</sub> NOS                              | NGVNNJYFJYTCCO-UHFFFAOYSA-N |
| 31 | alpha-PVP**                | α-pyrrolidinovalerophenone                | 1-phenyl-2-(1-pyrrolidinyl)-1-pentanone   | C <sub>15</sub> H <sub>21</sub> NO                               | YDIIDRWHPFMLGR-UHFFFAOYSA-N |
| 32 | alpha-PVT***               | α-PVT, alpha-pyrrolidinopentiothiophenone | 2-(pyrrolidin-1-yl)-1-(thiophen-2-yl)pentan-1-one   | C <sub>13</sub> H <sub>19</sub> NOS                              | OOSRPGUQJAKBLV-UHFFFAOYSA-N |
| 33 | AM-1248 azepane isomer***  |   | adamant-1-yl[1-(1-methylazepan-3-yl)-1 <i>H</i> -indol-3-yl] methanone  | C <sub>26</sub> H <sub>34</sub> N <sub>2</sub> O                 | HSCSEKGAOWTVDH-UHFFFAOYSA-N |

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|    |                                   |                    |   |   |                             |
|----|-----------------------------------|--------------------|---|---|-----------------------------|
| 34 | AM-2201**                         | JWH-2201           | 1-[(5-fluoropentyl)-1 <i>H</i> -indol-3-yl]-(naphthalen-1-yl)methanone                            | C <sub>24</sub> H <sub>22</sub> FNO               | ALQFAGFPQCBPED-UHFFFAOYSA-N |
| 35 | AMT***<br>(alfa-metiltriptamin)   |                    | 1-(1 <i>H</i> -indol-3-yl)propan-2-amine  | C <sub>11</sub> H <sub>14</sub> N <sub>2</sub>    | QSQQQURBVYWZKJ-UHFFFAOYSA-N |
| 36 | BDB***                            |                    | 1-(1,3-benzodioxol-5-yl)-2-butanamine; 1-(1,3-benzodioxol-5-yl)butan-2-ylazan <sup>oo</sup>       | C <sub>11</sub> H <sub>15</sub> NO <sub>2</sub>   | VHMRXGAIDDCGDU-UHFFFAOYSA-N |
| 37 | Brolamfetamine<br>(brolamfetamin) | DOB                | (±)-4-bromo-2,5-dimethoxy- $\alpha$ -methylphenethylamine   | C <sub>11</sub> H <sub>16</sub> BrNO <sub>2</sub> | FXMWUTGUCAKGQL-UHFFFAOYSA-N |
| 38 | BZP <sup>EU3</sup>                | N-Benzylpiperazine | 1-benzylpiperazine <sup>EU3</sup> ;<br>1-benzyl-1,4-diaza-cyclohexane; <i>N</i> -benzylpiperazine | C <sub>11</sub> H <sub>16</sub> N <sub>2</sub>    | IQXXEPZFOOTTBA-UHFFFAOYSA-N |
| 39 | $\beta$ -Me-PEA***                |                    | 2-phenylpropan-1-amine  | C <sub>9</sub> H <sub>13</sub> N                  | AXORVIZLPOGIRG-UHFFFAOYSA-N |
| 40 | Camfetamine***<br>(kamfetamin)    |                    | <i>N</i> -methyl-3-phenylbicyclo[2.2.1]heptan-2-amine   | C <sub>14</sub> H <sub>19</sub> N                 | CTVMYAZECFXZLN-UHFFFAOYSA-N |
| 41 | Cathinone<br>(katinon)            |                    | (-)-( <i>S</i> )-2-amino-propiofenone   | C <sub>9</sub> H <sub>11</sub> NO                 | PUAQLLVFLMYYJJ-ZETCQYMHSA-N |
| 42 | CP 47,497 C8-homológ***           |                    | 2-(3-hydroxycyclohexyl)-5-(2-methylnonan-2-yl)phenol  | C <sub>22</sub> H <sub>36</sub> O <sub>2</sub>    | HNMJDLVMIUDJNH-UHFFFAOYSA-N |
| 43 | CRA 13***                         |                    | naphthalen-1-yl [4-(pentyloxy)naphthalen-1-yl]methanone   | C <sub>26</sub> H <sub>24</sub> O <sub>2</sub>    | RSUMDJRTAFBISX-UHFFFAOYSA-N |
| 44 | D2PM***<br>(difenilprolinol)      | Diphenylprolinol   | diphenyl(pyrrolidin-2-yl)methanol   | C <sub>17</sub> H <sub>19</sub> NO                | OGCGXUGBDJGFFY-UHFFFAOYSA-N |
| 45 | DBZP***                           |                    | 1,4-dibenzylpiperazine  | C <sub>18</sub> H <sub>22</sub> N <sub>2</sub>    | YPUGLZQRXQQCSX-UHFFFAOYSA-N |
| 46 | Desoxy-D2PM***<br>(dezoxi-D2PM)   |                    | 2-(diphenylmethyl)pyrrolidine   | C <sub>17</sub> H <sub>19</sub> N                 | OXOBKZZXZVFOBB-UHFFFAOYSA-N |
| 47 | DET                               |                    | 3-[2-(diethyl-amino)ethyl]indole  | C <sub>14</sub> H <sub>20</sub> N <sub>2</sub>    | LSSUMOWDTKZHHT              |

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|    |                                  |  |   |   | -UHFFFAOYSA-N                   |
| 48 | Dimethocaine****<br>(dimetokain) |  | 3-(diethylamino)-2,2-dimethylpropyl-4-aminobenzoate   | C <sub>16</sub> H <sub>26</sub> N <sub>2</sub> O <sub>2</sub> | OWQIUQKMPDHQ<br>Q-UHFFFAOYSA-N  |
| 49 | Diphenidine**                    | 1,2-diphenylethylpiperidine,<br>DPD, 1,2-DEP, DIPH | 1-(1,2-diphenylethyl)piperidine   | C <sub>19</sub> H <sub>23</sub> N                             | JQWJJYHVHNXJH-<br>UHFFFAOYSA-N  |
| 50 | DMA                              |  | (±)-2,5-dimethoxy-α-methylphenethylamine  | C <sub>11</sub> H <sub>17</sub> NO <sub>2</sub>               | LATVFDIBMDBSY-<br>UHFFFAOYSA-N  |
| 51 | DMAA***                          |  | 4-methylhexan-2-amine   | C <sub>7</sub> H <sub>17</sub> N                              | YAHRDLICUYEDAU-<br>UHFFFAOYSA-N |
| 52 | DMHP                             |  | 3-(1,2-dimethylheptyl)-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo[ <i>b,d</i> ]pyran-1-ol | C <sub>25</sub> H <sub>38</sub> O <sub>2</sub>                | QBEFIFWEOSUTKV-<br>UHFFFAOYSA-N |
| 53 | DMT                              |  | 3-[2-(dimethylamino)ethyl]indole  | C <sub>12</sub> H <sub>16</sub> N <sub>2</sub>                | DMULVCHRPCFFGV-<br>UHFFFAOYSA-N |
| 54 | DOC                              |  | 4-chloro-2,5-dimethoxyamfetamine; 2,5-dimethoxy-4-chloroamphetamine°                          | C <sub>11</sub> H <sub>16</sub> ClNO <sub>2</sub>             | ACRITBNCBMTINK-<br>UHFFFAOYSA-N |
| 55 | DOET                             |  | (±)-4-ethyl-2,5-dimethoxy-α-methylphenethylamine  | C <sub>13</sub> H <sub>21</sub> NO <sub>2</sub>               | HXJKWPGVENNMCC-<br>UHFFFAOYSA-N |
| 56 | EG-018****                       |  | (naphthalen-1-yl)(9-pentyl-9H-carbazol-3-yl)-methanone  | C <sub>28</sub> H <sub>25</sub> NO                            | FJMMDJDPNLZYLA-<br>UHFFFAOYSA-N |
| 57 | Etaqualone****<br>(etakvalon)    |  | 3-(2-ethylphenyl)-2-methylquinazolin-4(3H)-one  | C <sub>17</sub> H <sub>16</sub> N <sub>2</sub> O              | UVTJKLLUVOTSOB-<br>UHFFFAOYSA-N |
| 58 | Eticyclidine<br>(eticiklidin)    | PCE  | N-ethyl-1-phenylcyclohexylamine   | C <sub>14</sub> H <sub>21</sub> N                             | IFYLVUHLOOCYBG-<br>UHFFFAOYSA-N |
| 59 | Etryptamine<br>(etriptamin)      |  | 3-(2-aminobutyl)indole  | C <sub>12</sub> H <sub>16</sub> N <sub>2</sub>                | ZXUMUPVQYAFTLF-<br>UHFFFAOYSA-N |
| 60 | FLEA***                          |  | N-hydroxy-N-methyl-3,4-   | C <sub>11</sub> H <sub>15</sub> NO <sub>3</sub>               | ORADFQZOLNHWRQ                  |

## DRAFT

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|----|-----------------------------------|---------------------|--|--|----------------------------------|
|    |                                   |                     | methylenedioxyamphetamine;<br><i>N</i> -hydroxy-3,4-<br>methylenedioxymethamphetamine°;<br><i>N</i> -[1-(1,3-benzodioxol-5-yl)propan-2-yl]- <i>N</i> -<br>methylhydroxylamine°°  |  | -UHFFFAOYSA-N                    |
| 61 | GBL***<br>(gamma-butirolakton)    | gamma-butyrolactone | dihydrofuran-2(3 <i>H</i> )-one  | C <sub>4</sub> H <sub>6</sub> O <sub>2</sub>     | YEJRWHA VMIAJKC-<br>UHFFFAOYSA-N |
| 62 | Homoamphetamine***                |                     | 3-amino-1-phenyl-butane  | C <sub>10</sub> H <sub>15</sub> N                | WECUIGDEWBNQJJ-<br>UHFFFAOYSA-N  |
| 63 | HU-210***                         |                     | 9-(hydroxymethyl)-6,6-dimethyl-3-(2-<br>methyloctan-2-yl)-6a,7,10,10a-<br>tetrahydrobenzo[ <i>c</i> ]chromen-1-ol  | C <sub>25</sub> H <sub>38</sub> O <sub>3</sub>   | SSQJFGMEZBFMNV-<br>UHFFFAOYSA-N  |
| 64 | Ibogain***                        |                     | (6 <i>R</i> ,6a <i>S</i> ,7 <i>S</i> ,9 <i>R</i> )-7-ethyl-2-methoxy-<br>6,6a,7,8,9,10,12,13-octahydro-5 <i>H</i> -6,9-<br>methanopyrido[10,20:1,2]azepino[4,5- <i>b</i> ]indole | C <sub>20</sub> H <sub>26</sub> N <sub>2</sub> O | HSIBGVUMFOSJPD-<br>CFDPKNGZSA-N  |
| 65 | JWH-018**                         | AM-678              | naphthalen-1-yl (1-pentyl-1 <i>H</i> -indol-3-<br>yl)methanone   | C <sub>24</sub> H <sub>23</sub> NO               | JDNLPKCAXICMBW-<br>UHFFFAOYSA-N  |
| 66 | JWH-073***                        |                     | (1-butyl-1 <i>H</i> -indol-3-yl)(naphthalen-1-<br>yl)methanone   | C <sub>23</sub> H <sub>21</sub> NO               | VCHHHSMPMLNVGS-<br>UHFFFAOYSA-N  |
| 67 | JWH-081***                        |                     | (4-methoxynaphthalen-1-yl)(1-pentyl-1 <i>H</i> -indol-<br>3-yl)methanone   | C <sub>25</sub> H <sub>25</sub> NO <sub>2</sub>  | UBMPKJKGUQDHRM-<br>UHFFFAOYSA-N  |
| 68 | JWH-122***                        |                     | (4-methylnaphthalen-1-yl)(1-pentyl-1 <i>H</i> -indol-3-<br>yl)methanone  | C <sub>25</sub> H <sub>25</sub> NO               | HUKJQMKQFWYIHS-<br>UHFFFAOYSA-N  |
| 69 | JWH-210***                        |                     | (4-ethylnaphthalen-1-yl)(1-pentyl-1 <i>H</i> -indol-3-<br>yl)methanone   | C <sub>26</sub> H <sub>27</sub> NO               | LACIUQLUNACUKC-<br>UHFFFAOYSA-N  |
| 70 | (+)-Lysergide<br>( (+)-lizergid ) | LSD, LSD-25         | 9,10-didehydro- <i>N,N</i> -diethyl-6-methylergoline-<br>8β-carboxamide  | C <sub>20</sub> H <sub>25</sub> N <sub>3</sub> O | VAYOSLLFUXYJDT-<br>RDTXWAMCSA-N  |

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|----|--|--|---|---|---------------------------------|
| 71 | MAL***                                 |  | 3,5-dimethoxy-4-metallyloxyphenethylamine°;<br>3,5-dimethoxy-4-(2-methylallyloxy)-<br>phenethylazan°°   | C <sub>14</sub> H <sub>21</sub> NO <sub>3</sub>               | FOXJFBFFGULACD-<br>UHFFFAOYSA-N |
| 72 | MBDB***                                |  | 2-(methylamino)-1-(3,4-methylenedioxyphenyl)-<br>butane or N-methyl-1-(1,3-benzodioxol-5-yl)-2-<br>butanamine°;<br>[1-(1,3-Benzodioxol-5-yl)butan-2-yl]<br>(methyl)azan°° | C <sub>12</sub> H <sub>17</sub> NO <sub>2</sub>               | USWVWJSAJAEHQ-<br>UHFFFAOYSA-N  |
| 73 | MBZP***                                |  | 1-benzyl-4-methylpiperazine   | C <sub>12</sub> H <sub>18</sub> N <sub>2</sub>                | MLJOKPBESJWYGL-<br>UHFFFAOYSA-N |
| 74 | mCPP***<br>(meta-klorofenilpiperazin)  |  | meta-chlorophenylpiperazine°;<br>[1-(3-chlorophenyl)piperazine]°°   | C <sub>10</sub> H <sub>13</sub> ClN <sub>2</sub>              | VHFVKMTVMIZMIK-<br>UHFFFAOYSA-N |
| 75 | MDAI***                                |  | 6,7-dihydro-5H-indeno[5,6-d][1,3]dioxol-6-<br>amine   | C <sub>10</sub> H <sub>11</sub> NO <sub>2</sub>               | FQDRMHHCWZAXJM-<br>UHFFFAOYSA-N |
| 76 | MDE<br>(N-etil-MDA)                    | N-ethyl MDA, MDEA                                      | (±)-N-ethyl-α-methyl-3,4-<br>(methylenedioxy)phenethylamine   | C <sub>12</sub> H <sub>17</sub> NO <sub>2</sub>               | PVXVWWANJIWJOO-<br>UHFFFAOYSA-N |
| 77 | MDMA                                   |  | (±)-N,α-dimethyl-3,4-<br>(methylenedioxy)phenethylamine   | C <sub>11</sub> H <sub>15</sub> NO <sub>2</sub>               | SHXWCVYOXRDMC<br>X-UHFFFAOYSA-N |
| 78 | MDMB-CHMICA**                          |  | Methyl 2-{{1-(cyclohexylmethyl)indole-3-<br>carbonyl}amino}-3,3-dimethylbutanoate   | C <sub>23</sub> H <sub>32</sub> N <sub>2</sub> O <sub>3</sub> | SRJKCVHWIDFUBO-<br>UHFFFAOYSA-N |
| 79 | Mephedrone <sup>EU</sup><br>(mefedron) | 4-methylmethcathinone,<br>4-methylephedrone, 4-<br>MMC | (RS)-2-methylamino-1-(4-<br>methylphenyl)propan-1-one   | C <sub>11</sub> H <sub>15</sub> NO                            | YELGFTGWJGBAQU-<br>UHFFFAOYSA-N |
| 80 | Mescaline<br>(meszkalin)               |  | 3,4,5-trimethoxyphenethylamine  | C <sub>11</sub> H <sub>17</sub> NO <sub>3</sub>               | RHCSKNNOAZULRK-<br>UHFFFAOYSA-N |
| 81 | Methcathinone<br>(metkatinon)          |  | 2-(methylamino)-1-phenylpropan-1-one  | C <sub>10</sub> H <sub>13</sub> NO                            | LPLLVINFLBSFRP-<br>UHFFFAOYSA-N |

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|----|---|-------------------------------------|---|---|------------------------------|
| 82 | Methoxetamine**<br>(metoxetamin)  | MXE                                 | ( <i>RS</i> )-2-(3-methoxyphenyl)-2-(ethylamino)-cyclohexanone  | C <sub>15</sub> H <sub>21</sub> NO <sub>2</sub>   | LPKTWLVEGBNOOX-UHFFFAOYSA-N  |
| 83 | Methylone**<br>(Metilon)  | beta-keto-MDMA                      | ( <i>RS</i> )-2-methylamino-1-(3,4-methylenedioxyphenyl)propan-1-one  | C <sub>11</sub> H <sub>13</sub> NO <sub>3</sub>   | VKEQBMCQRQDSRET-UHFFFAOYSA-N |
| 84 | Methiopropamine**<br>(metiltienil-propamin)   | MPA,<br>Methylthietnylpropamine     | 1-(thiophen-2-yl)-2-methylaminopropane  | C <sub>8</sub> H <sub>13</sub> NS                 | HPHUWHKFQXTZPS-UHFFFAOYSA-N  |
| 85 | MMDA  |                                     | 5-methoxy- $\alpha$ -methyl-3,4-(methylenedioxy)phenylethylamine  | C <sub>11</sub> H <sub>15</sub> NO <sub>3</sub>   | YQYUWUKDEVZFDB-UHFFFAOYSA-N  |
| 86 | MT-45 <sup>EU</sup>   |                                     | 1-cyclohexyl-4-(1,2-diphenylethyl)piperazine  | C <sub>24</sub> H <sub>32</sub> N <sub>2</sub>    | IGBRRSIHEGCUEN-UHFFFAOYSA-N  |
| 87 | MTP***  | methcathinone thiophen analog       | 2-(methylamino)-1-(thiophen-2-yl)propan-1-one   | C <sub>8</sub> H <sub>11</sub> NOS                | DOZQPYDMJMLVKX-UHFFFAOYSA-N  |
| 88 | MTTA***   | MTA, Mephtetramine                  | 2-[(methylamino)methyl]-3,4-dihydronaphthalen-1(2 <i>H</i> )-one  | C <sub>12</sub> H <sub>15</sub> NO                | FTRWLSZFILOOD-UHFFFAOYSA-N   |
| 89 | MXP***  | methoxyphenidine, 2-MeO-diphenidine | 1-[1-(2-methoxyphenyl)-2-phenylethyl]piperidine   | C <sub>20</sub> H <sub>25</sub> NO                | QXXCUXIRBHSITD-UHFFFAOYSA-N  |
| 90 | N-(2-methoxyethyl)-N-(1-methylethyl)-2-(1-pentyl-1 <i>H</i> -indol-3-yl)-4-thiazol-methanamine*** |                                     | <i>N</i> -(2-methoxyethyl)- <i>N</i> -(1-methylethyl)-2-(1-pentyl-1 <i>H</i> -indol-3-yl)-4-thiazol-methanamine | C <sub>23</sub> H <sub>33</sub> N <sub>3</sub> OS | PSAKYYVEVVAWJL-UHFFFAOYSA-N  |
| 91 | N,N-diethyl-2-(1-pentyl-1 <i>H</i> -indol-3-yl)-4-thiazol-methanamine***                          |                                     | <i>N,N</i> -diethyl-2-(1-pentyl-1 <i>H</i> -indol-3-yl)-4-thiazol-methanamine                                   | C <sub>21</sub> H <sub>29</sub> N <sub>3</sub> S  | PCNLLVFKBKMRDB-UHFFFAOYSA-N  |
| 92 | N-ethylnorketamine***   |                                     | 2-(2-chlorophenyl)-2-(ethylamino)cyclohexanone  | C <sub>14</sub> H <sub>18</sub> ClNO              | ITBBBZIIFJJMDU-UHFFFAOYSA-N  |
| 93 | N-hydroxy MDA<br>(N-hidroxi-MDA)  |                                     | ( $\pm$ )- <i>N</i> -[ $\alpha$ -methyl-3,4-(methylenedioxy)phenethyl] hydroxylamine                            | C <sub>10</sub> H <sub>13</sub> NO <sub>3</sub>   | FNDCTJYFKOQGTL-UHFFFAOYSA-N  |



## DRAFT

|     |                                 |  |  |   |                             |
|-----|---------------------------------|--|--|---|-----------------------------|
| 94  | Nitracaine***                   |  | 3-(diethylamino)-2,2-dimethylpropyl 4-nitrobenzoate                            | C <sub>16</sub> H <sub>24</sub> N <sub>2</sub> O <sub>4</sub> | SPTIETJWCCCJSE-UHFFFAOYSA-N |
| 95  | N-Me-1-PEA***                   |  | 1-(N-methylamino)-1-phenylethane;<br>N-methyl-1-phenethylamine                 | C <sub>9</sub> H <sub>13</sub> N                              | RCSHZGQHHEHPZ-UHFFFAOYSA-N  |
| 96  | ODT***<br>(O-desmetiltramadol)  | O-desmethyltramadol  | 3-{2-[(dimethylamino)methyl]-1-hydroxycyclohexyl}phenol                        | C <sub>15</sub> H <sub>23</sub> NO <sub>2</sub>               | UWJUQVWARXYRCG-UHFFFAOYSA-N |
| 97  | Parahexyl<br>(parahexil)        |  | 3-hexyl-7,8,9,10-tetrahydro-6,6,9-trimethyl-6H-dibenzo[ <i>b,d</i> ]pyran-1-ol | C <sub>22</sub> H <sub>32</sub> O <sub>2</sub>                | OORFXDSWECAQLI-UHFFFAOYSA-N |
| 98  | para-Methyl-4-methylaminorex ** | 4,4'-DMAR  | 4-methyl-5-(4-methylphenyl)-4,5-dihydro-1,3-oxazol-2-amine                     | C <sub>11</sub> H <sub>14</sub> N <sub>2</sub> O              | NPILLHMQNMXXTL-UHFFFAOYSA-N |
| 99  | Pentedron**                     | $\beta$ -ethyl-methcathinone<br>( $\beta$ -etil-metkatinon),<br>pentedrone | ( $\pm$ )-2-(methylamino)-1-phenylpentan-1-one                                 | C <sub>12</sub> H <sub>17</sub> NO                            | WLIWIUNEJRETFX-UHFFFAOYSA-N |
| 100 | pFBT***<br>(fluortropakokain)   | 4-fluorotropacocaine   | 8-methyl-8-azabicyclo[3.2.1]oct-3-yl 4-fluorobenzoate                          | C <sub>15</sub> H <sub>18</sub> FNO <sub>2</sub>              | YXDfSLsXLYAAPF-UHFFFAOYSA-N |
| 101 | pFPP***                         |  | 1-(4-fluorophenyl)piperazine   | C <sub>10</sub> H <sub>13</sub> FN <sub>2</sub>               | AVJKDKWRVSSJPK-UHFFFAOYSA-N |
| 102 | Phenazepam**<br>(fenazepam)     |  | 7-bromo-5-(2-chlorophenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one              | C <sub>15</sub> H <sub>10</sub> BrClN <sub>2</sub> O          | CGMJQJQSWIRRL-UHFFFAOYSA-N  |
| 103 | PMA                             |  | 1-(4-methoxyphenyl)propan-2-amine  | C <sub>10</sub> H <sub>15</sub> NO                            | NEGYEDYHPMHGK-UHFFFAOYSA-N  |

## DRAFT

|     |  |  |  |   |                                 |
|-----|--|--|--|---|---------------------------------|
| 104 | PMMA <sup>EU</sup><br>(parametoxi-<br>metilamfetamin)  | para-<br>methoxymethylampheta-<br>mine | paramethoxymethylamphetamine <sup>EU1</sup> ;<br><i>N</i> -methyl-1-(4-methoxyphenyl)-2-<br>aminopropane or <i>p</i> -methoxy- <i>N</i> , $\alpha$ -<br>dimethylphenethylamine or 1-(4-<br>methoxyphenyl)- <i>N</i> -methylpropan-2-amine                          | C <sub>11</sub> H <sub>17</sub> NO                              | UGFMBZYKVQSQFX-<br>UHFFFAOYSA-N |
| 105 | Psilocine<br>(pszilocin)   | Psilotsin                              | 3-[2-(dimethylamino)ethyl]indol-4-ol   | C <sub>12</sub> H <sub>16</sub> N <sub>2</sub> O                | SPCIYGNTAMCTRO-<br>UHFFFAOYSA-N |
| 106 | Psilocybine<br>(pszilocibin)   |  | 3-[2-(dimethylamino)ethyl]indol-4-yl<br>dihydrogen phosphate   | C <sub>12</sub> H <sub>17</sub> N <sub>2</sub> O <sub>4</sub> P | QVDSEJDULKHCG-<br>UHFFFAOYSA-N  |
| 107 | Rolicyclidine<br>(rolciklidin)   | PHP, PCPY                              | 1-(1-phenylcyclohexyl)pyrrolidine  | C <sub>16</sub> H <sub>23</sub> N                               | FYOWWXMGDATDQ<br>Y-UHFFFAOYSA-N |
| 108 | Salvinorin A****<br>(Szalvinorin A)  |  | (2 <i>S</i> ,4 <i>aR</i> ,6 <i>aR</i> ,7 <i>R</i> ,9 <i>S</i> ,10 <i>aS</i> ,10 <i>bR</i> )-9-(acetyloxy)-2-<br>(3-furanyl)dodecahydro-6 <i>a</i> ,10 <i>b</i> -dimethyl-4,10-<br>dioxo-2 <i>H</i> -naphtho[2,1- <i>c</i> ]pyran-7-carboxylic<br>acid methyl ester | C <sub>23</sub> H <sub>28</sub> O <sub>8</sub>                  | OBSYBRPAKCASQB-<br>AGQYDFLVSA-N |
| 109 | STP  | DOM                                    | 2,5-dimethoxy- $\alpha$ ,4-dimethylphenethylamine  | C <sub>12</sub> H <sub>19</sub> NO <sub>2</sub>                 | NTJQREUGJKIARY-<br>UHFFFAOYSA-N |
| 110 | Tenamfetamine<br>(tenamfetamin)  | MDA                                    | $\alpha$ -methyl-3,4-(methylenedioxy)phenethylamine  | C <sub>10</sub> H <sub>13</sub> NO <sub>2</sub>                 | NGBBVGZWCFBGO<br>-UHFFFAOYSA-N  |
| 111 | Tenocyclidine<br>(tenociklidin)  | TCP                                    | 1-[1-(2-thienyl)cyclohexyl]piperidine  | C <sub>15</sub> H <sub>23</sub> NS                              | JUZZEWSCNBCFRL-<br>UHFFFAOYSA-N |
| 112 | Tetrahydrocannabinol, the<br>following isomers and their<br>stereochemical variants<br>(a következő tetrahydro-<br>kannabinol izomerek és<br>szterokémiai variánsaik)<br>[THC] | <i>delta</i> -6a(10a)-THC              | 7,8,9,10-tetrahydro-6,6,9-trimethyl-3-pentyl-<br>6 <i>H</i> -dibenzo[ <i>b,d</i> ]pyran-1-ol   | C <sub>21</sub> H <sub>30</sub> O <sub>2</sub>                  | NEBZNJDFIPBXCS-<br>UHFFFAOYSA-N |
|     |  | <i>delta</i> -6a(7)-THC                | 8,9,10,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-<br>6 <i>H</i> -dibenzo-[ <i>b,d</i> ]pyran-1-ol  | C <sub>21</sub> H <sub>30</sub> O <sub>2</sub>                  | UQOUHXDCXBITSF-<br>UHFFFAOYSA-N |
|     |  | <i>delta</i> -7-THC                    | 6a,9,10,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-<br>6 <i>H</i> -dibenzo[ <i>b,d</i> ]pyran-1-ol  | C <sub>21</sub> H <sub>30</sub> O <sub>2</sub>                  | WWYMYGIVLCKTBL<br>-UHFFFAOYSA-N |
|     |  | <i>delta</i> -8-THC                    | 6a,7,10,10a-tetrahydro-6,6,9-trimethyl-3-pentyl-   | C <sub>21</sub> H <sub>30</sub> O <sub>2</sub>                  | HCAWPGARWVBULJ                  |

|     |  |                         |   |   |                                 |
|-----|--|-------------------------|---|---|---------------------------------|
| 113 | TFMPP***   |                         | 6 <i>H</i> -dibenzo- <i>[b,d]</i> pyran-1-ol  |   | -UHFFFAOYSA-N                   |
|     |  | <i>delta</i> -10-THC    | 6a,7,8,9-tetrahydro-6,6,9-trimethyl-3-pentyl-6 <i>H</i> -dibenzo <i>[b,d]</i> pyran-1-ol                | C <sub>21</sub> H <sub>30</sub> O <sub>2</sub>                | YLTWYAXWDLZZCU<br>-UHFFFAOYSA-N |
|     |  | <i>delta</i> -9(11)-THC | 6a,7,8,9,10,10a-hexahydro-6,6-dimethyl-9-methylene-3-pentyl-6 <i>H</i> -dibenzo <i>[b,d]</i> pyran-1-ol | C <sub>21</sub> H <sub>30</sub> O <sub>2</sub>                | AOYYFUGUUIRBML-<br>UHFFFAOYSA-N |
|     |  |                         | 1-[3-(trifluoromethyl)phenyl]piperazine   | C <sub>11</sub> H <sub>13</sub> F <sub>3</sub> N <sub>2</sub> | KKIMDKMETPPURN-<br>UHFFFAOYSA-N |
| 114 | TMA<br>(trimetoxiamfetamin)                        |                         | (±)-3,4,5-trimethoxy- $\alpha$ -methylphenethylamine  | C <sub>12</sub> H <sub>19</sub> NO <sub>3</sub>               | WGTASENVNYJZBK-<br>UHFFFAOYSA-N |
| 115 | TMA-2 <sup>EU</sup><br>(2,4,5- trimetoxiamfetamin) |                         | 2,4,5-trimethoxyamphetamine <sup>EU2</sup> ;<br>1-(2,4,5-trimethoxyphenyl)propan-2-ylazan <sup>oo</sup> | C <sub>12</sub> H <sub>19</sub> NO <sub>3</sub>               | TVSIMAWGATVNGK<br>-UHFFFAOYSA-N |
| 116 | 3-methylmethcathinone, 3-MMC                       |                         | 2-(methylamino)-1-(3-methylphenyl)propan-1-one  | C <sub>11</sub> H <sub>15</sub> NO                            | QDNXSIYWHYGMCD<br>-UHFFFAOYSA-N |
| 117 | 3-chloromethcathinone, 3-CMC                       |                         | 1-(3-chlorophenyl)-2-(methylamino)propan-1-one  | C <sub>10</sub> H <sub>12</sub> ClNO                          | VOEFELLSAAJCHJ-<br>UHFFFAOYSA-N |

1.2. In addition, the stereoisomers of the above substances, if they correspond to the chemical name indicated, exist on the basis of their chemical structure and are not expressly covered by exceptional provisions, and their salts, including those of their stereoisomers, where such salts exist.

1.3. The following substances are included in List P1:

1.3.1. substances included in the updated Schedule I of the Psychotropic Convention,

1.3.2. substances listed in Council Framework Decision 2004/757/JHA of 25 October 2004 laying down minimum provisions on the constituent elements of criminal acts and penalties in the field of illicit drug trafficking—they are marked with the symbol <sup>EU</sup>,

1.3.3. substances whose national control is stricter, on the basis of national and international law enforcement experience, than the requirement of the Psychotropic Convention—they are marked with \*\*,

1.3.4. substances whose national control is stricter based on national and international law enforcement experience and which are not included in the Schedules of the Narcotics Convention and the Psychotropic Convention—they are marked with \*\*\*.

1.4. In *column C* of this list, the former Hungarian official names are marked with the symbol ° and in italics, while the IUPAC name used by the German BtMG (Betäubungsmittelgesetz, 22 December 2003) is marked with ° ° and in italics, if available.

## 2. List 2 of psychotropic substances (List P2)

2.1. The following substances and compounds are classified as psychotropic substances:

|   | <b>A</b>                                 | <b>B</b>  | <b>C</b>   | <b>D</b>   | <b>E</b>                        |
|---|--|---|--|--|---------------------------------|
| 1 | <b>Official name</b><br>(Hungarian name) | Other name or<br>abbreviation, spelling<br>commonly used abroad | Chemical name / <i>Description</i>   | Molecular formula  | InChIKey chemical<br>identifier |
| 2 | 2C-B                                     |   | 4-bromo-2,5-dimethoxyphenethylamine  | C <sub>10</sub> H <sub>14</sub> BrNO <sub>2</sub>              | YMHOBZXQZVXHBM-<br>UHFFFAOYSA-N |
| 3 | 4-CMC                                    | 4-<br>chloromethcathinone,<br>clephedrone                       | 1-(4-chlorophenyl)-2-(methylamino)-1-<br>propanone   | C <sub>10</sub> H <sub>12</sub> ClNO                           | UEJBEOXRNGPEI-<br>UHFFFAOYSA-N  |
| 4 | 4-FA                                     | 4-fluoroamphetamine   | 1-(4-fluorophenyl)propan-2-amine   | C <sub>9</sub> H <sub>12</sub> FN                              | DGXWNDGLEOIEGT-<br>UHFFFAOYSA-N |
| 5 | 4F-MDMB-BINACA                           |   | Methyl 2-(1-(4-fluorobutyl)-1 <i>H</i> -indazole-3-<br>carboxamido)-3,3-dimethylbutanoate    | C <sub>19</sub> H <sub>26</sub> FN <sub>3</sub> O <sub>3</sub> | GZGKSDAMWRWYOZ-<br>UHFFFAOYSA-N |
| 6 | 5F-AKB-48                                | 5F-APINACA  | <i>N</i> -(adamantan-1-yl)-1-(5-fluoropentyl)-1 <i>H</i> -<br>indazole-3-carboxamide         | C <sub>23</sub> H <sub>30</sub> FN <sub>3</sub> O              | UCMFSGVIEPXYIV-<br>UHFFFAOYSA-N |
| 7 | 5F-AMB                                   | 5F-AMB-PINACA,<br>5F-MMB-PINACA                                 | Methyl 2-([1-(5-fluoropentyl)-1 <i>H</i> -indazol-3-<br>yl]carbonyl)amino)-3-methylbutanoate | C <sub>19</sub> H <sub>26</sub> FN <sub>3</sub> O <sub>3</sub> | SAFXSUZMRLTBMM-<br>UHFFFAOYSA-N |
| 8 | 5F-MDMB-PICA                             | 5F-MDMB-2201  | Methyl 2-(1-(5-fluoropentyl)-1 <i>H</i> -indole-3-   | C <sub>21</sub> H <sub>29</sub> FN <sub>2</sub> O <sub>3</sub> | CHSUEEBESACQDV-                 |

## DRAFT

|    |                               |                |  |  |                              |
|----|-------------------------------|----------------|--|--|------------------------------|
|    |                               |                | carboxamido)-3,3-dimethylbutanoate   |  | UHFFFAOYSA-N                 |
| 9  | 5F-ADB                        | 5F-MDMB-PINACA | Methyl 2-{{1-(fluoropentyl)-1 <i>H</i> -indazole-3-carbonyl}amino}-3,3-dimethylbutanoate                                 | C <sub>20</sub> H <sub>28</sub> FN <sub>3</sub> O <sub>3</sub> | PWEKNGSNNAKWBL-UHFFFAOYSA -N |
| 10 | 5F-PB-22                      |                | Quinolin-8-yl 1-(5-fluoropentyl)-1 <i>H</i> -indole-3-carboxylate  | C <sub>23</sub> H <sub>21</sub> FN <sub>2</sub> O <sub>2</sub> | MBOCMBFDYVSGLJ-UHFFFAOYSA-N  |
| 11 | AB-CHMINACA                   |                | <i>N</i> -[1-amino-3-methyl-1-oxobutan-2-yl]-1-(cyclohexylmethyl)-1 <i>H</i> -indazole-3-carboxamide                     | C <sub>20</sub> H <sub>28</sub> N <sub>4</sub> O <sub>2</sub>  | KJNZIEGLNLCWTQ-UHFFFAOYSA -N |
| 12 | AB-FUBINACA                   |                | <i>N</i> -[1-amino-3-methyl-1-oxobutan-2-yl]-1-[(4-fluorophenyl)methyl]indazole-3-carboxamide                            | C <sub>20</sub> H <sub>21</sub> FN <sub>4</sub> O <sub>2</sub> | AKOOIMKXADOPDA-UHFFFAOYSA-N  |
| 13 | AB-PINACA                     |                | <i>N</i> -[1-amino-3-methyl-1-oxobutan-2-yl]-1-pentyl-1 <i>H</i> -indazole-3-carboxamide                                 | C <sub>18</sub> H <sub>26</sub> N <sub>4</sub> O <sub>2</sub>  | GIMHPAQOAAZSHS-UHFFFAOYSA -N |
| 14 | ADB-CHMINACA                  | MAB-CHMINACA   | <i>N</i> -[1-amino-3,3-dimethyl-1-oxobutan-2-yl]-1-(cyclohexylmethyl)-1 <i>H</i> -indazole-3-carboxamide                 | C <sub>21</sub> H <sub>30</sub> N <sub>4</sub> O <sub>2</sub>  | ZWCCSIUBHCZKOY-UHFFFAOYSA -N |
| 15 | ADB-FUBINACA                  |                | <i>N</i> -[1-amino-3,3-dimethyl-1-oxobutan-2-yl]-1-[(4-fluorophenyl)methyl]-1 <i>H</i> -indazole-3-carboxamide           | C <sub>21</sub> H <sub>23</sub> FN <sub>4</sub> O <sub>2</sub> | ZSSGCSINPVBLQD-UHFFFAOYSA -N |
| 16 | alpha-PHP                     |                | ( <i>RS</i> )-1-phenyl-2-(pyrrolidine-1-yl)hexan-1-one   | C <sub>16</sub> H <sub>23</sub> NO                             | KYIJLDDXQWBNGX-UHFFFAOYSA-N  |
| 17 | Amfetamine (amfetamin)        | amphetamine    | (±)- $\alpha$ -methylphenethylamine  | C <sub>9</sub> H <sub>13</sub> N                               | KWTSXDURSIMDCE-UHFFFAOYSA-N  |
| 18 | Amineptine (amineptin)        |                | 7-[(10,11-dihydro-5 <i>H</i> -dibenzo[ <i>a,d</i> ]cyclohepten-5-yl)amino]heptanoic acid                                 | C <sub>22</sub> H <sub>27</sub> NO <sub>2</sub>                | ONNOFKFOZAJDHT-UHFFFAOYSA-N  |
| 19 | Buprenorphine** (buprenorfin) |                | 21-cyclopropyl-7- $\alpha$ -[( <i>S</i> )-1-hydroxy-1,2,2-trimethylpropyl]-6,14-endo-ethano-6,7,8,14-tetrahydrooripavine | C <sub>29</sub> H <sub>41</sub> NO <sub>4</sub>                | RMRJXGBAOAMLHD-IHFGGWKQSA-N  |
| 20 | CUMYL-4CN-BINACA              |                | 1-(4-cyanobutyl)- <i>N</i> -(2-phenylpropan-2-yl)-1 <i>H</i> -indazole-3-carboxamide                                     | C <sub>22</sub> H <sub>24</sub> N <sub>4</sub> O               | JGTSOWOPISVAHG-UHFFFAOYSA-N  |
| 21 | CUMYL-                        |                | 5-pentyl-2-(2-phenylpropan-2-yl)-2,5-dihydro-  | C <sub>25</sub> H <sub>28</sub> N <sub>2</sub> O               | AWHWTKXMUJLSRM-              |

## DRAFT

|    |   |  |  |  |                                  |
|----|---|--|--|--|----------------------------------|
|    | PEGACLONE   |  | 1 <i>H</i> -pyrido[4,3- <i>b</i> ]indol-1-one  |  | UHFFFAOYSA-N                     |
| 22 | Dexamfetamine<br>(dexamfetamin)   | dexamphetamine   | (+)- $\alpha$ -methylphenethylamine  | C <sub>9</sub> H <sub>13</sub> N                               | KWTSXDURSIMDCE-<br>QMMMGPBSA-N   |
| 23 | Dronabinol <sup>2</sup><br>(delta-9-<br>tetrahidrokannabinol<br>[delta-9-THC] és<br>sztereokémiai varánsai) | delta-9-<br>tetrahydrocannabinol<br>and its stereochemical<br>variants | (6 <i>aR</i> , 10 <i>aR</i> )-6 <i>a</i> ,7,8,10 <i>a</i> -tetrahydro-6,6,9-<br>trimethyl-3-pentyl-6 <i>H</i> -dibenzo[ <i>b,d</i> ]pyran-1-ol | C <sub>21</sub> H <sub>30</sub> O <sub>2</sub>                 | CYQFCXCEBYINGO-<br>IAGOWNOFSA-N  |
| 24 | Ethylone<br>(etilon)  |  | ( <i>RS</i> )-1-(1,3-benzodioxol-5-yl)-2-<br>(ethylamino)propan-1-one  | C <sub>12</sub> H <sub>15</sub> NO <sub>3</sub>                | MJEMIOXXNCZZFK-<br>UHFFFAOYSA-N  |
| 25 | Ethylphenidate<br>(etilfenidát)   | „Nopaine”, „Fake<br>cocaine”   | Ethyl 2-phenyl-2-(piperidin-2-yl)acetate   | C <sub>15</sub> H <sub>21</sub> NO <sub>2</sub>                | AIVSIRYZIBXTMM-<br>UHFFFAOYSA-N  |
| 26 | Eutylone (eutilon)  |  | 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)butan-<br>1-one   | C <sub>13</sub> H <sub>17</sub> NO <sub>3</sub>                | YERSNXHEOIYEGX-<br>UHFFFAOYSA-N  |
| 27 | Fenetylline (fenetillin)  |  | 7-{2-[( $\alpha$ -<br>methylphenethyl)amino]ethyl}theophylline   | C <sub>18</sub> H <sub>23</sub> N <sub>5</sub> O <sub>2</sub>  | NMCHYWGKBADVMK-<br>UHFFFAOYSA-N  |
| 28 | FUB-AMB   | MMB-FUBINACA,<br>AMB-FUBINACA  | Methyl 2-({1-[(4-fluorophenyl)methyl]-1 <i>H</i> -<br>indazole-3-carbonyl}amino)-3-<br>methylbutanoate   | C <sub>21</sub> H <sub>22</sub> FN <sub>3</sub> O <sub>3</sub> | FRFFLYJQPCIIQB-<br>UHFFFAOYSA -N |
| 29 | gamma-Hydroxybutiric<br>acid (gamma-hidroxi-<br>vajsav)   | GHB  | $\gamma$ -hydroxybutyric acid  | C <sub>4</sub> H <sub>8</sub> O <sub>3</sub>                   | SJZRECIVHVDYJC-<br>UHFFFAOYSA-N  |
| 30 | Ketamine***<br>(ketamin)  |  | 2-(2-chlorophenyl)-2-(methylamino)-<br>cyclohexanone   | C <sub>13</sub> H <sub>16</sub> ClNO                           | YQEZLKZALYSWHR-<br>UHFFFAOYSA-N  |
| 31 | Levamfetamine<br>(levamfetamin)   | levamphetamine   | (-)-( <i>R</i> )- $\alpha$ -methylphenethylamine   | C <sub>9</sub> H <sub>13</sub> N                               | KWTSXDURSIMDCE-<br>MRVPVSSYSA-N  |
| 32 | Levomethamphetamine<br>(levometamfetamin)   | levomethamphetamine  | (-)- <i>N</i> , $\alpha$ -dimethylphenethylamine   | C <sub>10</sub> H <sub>15</sub> N                              | MYWUZJCMWCOHBA-<br>SECBINFHSA-N  |
| 33 | MDMB-4en-PINACA   |  | Methyl 3,3-dimethyl-2-[1-(pent-4-en-1-yl)-1 <i>H</i> -<br>indazole-3-carboxamido]butanoate   | C <sub>20</sub> H <sub>27</sub> N <sub>3</sub> O <sub>3</sub>  | LWOCBHBFWNGPGM-<br>UHFFFAOYSA-N  |
| 34 | Mecloqualone<br>(meklokvalon)   |  | 3-( <i>o</i> -chlorophenyl)-2-methyl-4-(3 <i>H</i> )-<br>quinazolinone   | C <sub>15</sub> H <sub>11</sub> ClN <sub>2</sub> O             | SFITWQDBYUMAPS-<br>UHFFFAOYSA-N  |

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|    |  |                            |  |   |                                 |
|----|--|----------------------------|--|---|---------------------------------|
| 35 | Metamfetamine<br>(metamfetamin)                        | metamphetamine             | (+)-( <i>S</i> )- <i>N</i> , $\alpha$ -dimethylphenethylamine  | C <sub>10</sub> H <sub>15</sub> N                             | MYWUZJCMWCOHBA-<br>VIFPVBQESA-N |
| 36 | Metamfetamine<br>racemate<br>(metamfetamin<br>racemát) | metamphetamine<br>racemate | ( $\pm$ )- <i>N</i> , $\alpha$ -dimethylphenethylamine   | C <sub>10</sub> H <sub>15</sub> N                             | MYWUZJCMWCOHBA-<br>UHFFFAOYSA-N |
| 37 | Methaqualone<br>(metakvalon)                           |                            | 2-methyl-3- <i>o</i> -tolyl-4-(3 <i>H</i> )-quinazolinone  | C <sub>16</sub> H <sub>14</sub> N <sub>2</sub> O              | JEYCTXHKTXCGPB-<br>UHFFFAOYSA-N |
| 38 | Methylphenidate<br>(metilfenidát)                      |                            | Methyl $\alpha$ -phenyl-2-piperidine acetate   | C <sub>14</sub> H <sub>19</sub> NO <sub>2</sub>               | DUGOZIWVEXMGBE-<br>UHFFFAOYSA-N |
| 39 | N-ethylhexedrone                                       |                            | 2-(ethylamino)-1-phenyl-1-hexanone   | C <sub>14</sub> H <sub>21</sub> NO                            | CWNKMHIETKEBCA-<br>UHFFFAOYSA-N |
| 40 | N-ethyl norpentylone                                   | ephylone                   | 1-(2 <i>H</i> -1,3-benzodioxol-5-yl)-2-(ethylamino)pentan-1-one  | C <sub>14</sub> H <sub>19</sub> NO <sub>3</sub>               | VERDHJIMZYXGIW-<br>UHFFFAOYSA-N |
| 41 | Pentazocine**<br>(pentazocin)                          |                            | (2 <i>R</i> ,*6 <i>R</i> ,*11 <i>R</i> *)-1,2,3,4,5,6-hexahydro-6,11-dimethyl-3-(3-methyl-2-butenyl)-2,6-methano-3-benzazocin-8-ol | C <sub>19</sub> H <sub>27</sub> NO                            | VOKSWYLNZZRQPF-<br>GDIGMMSISA-N |
| 42 | Phencyclidine<br>(fenciklidin)                         | PCP                        | 1-(1-phenylcyclohexyl)piperidine   | C <sub>17</sub> H <sub>25</sub> N                             | JTJMJGYZQZDUJJ-<br>UHFFFAOYSA-N |
| 43 | Phenmetrazine<br>(fenmetrazin)                         |                            | 3-methyl-2-phenylmorpholine  | C <sub>11</sub> H <sub>15</sub> NO                            | OOBHFESNSZDWIU-<br>UHFFFAOYSA-N |
| 44 | Poppy straw**<br>(mákszalma)                           |                            |  | -   | -                               |
| 45 | Secobarbital<br>(szekobarbitál)                        |                            | 5-allyl-5-(1-methylbutyl)barbituric acid   | C <sub>12</sub> H <sub>18</sub> N <sub>2</sub> O <sub>3</sub> | KQPKPCNLIDLUMF-<br>UHFFFAOYSA-N |
| 46 | Tapentadol ***<br>(tapentadol)                         |                            | 3-[(2 <i>R</i> ,3 <i>R</i> )-1-(dimethylamino)-2-methylpentan-3-yl]- phenol  | C <sub>14</sub> H <sub>23</sub> NO                            | KWTWDQCKEHXFFR-<br>SMDDNHRTSA-N |
| 47 | UR-144   |                            | (1-pentyl-1 <i>H</i> -indol-3-yl)(2,2,3,3-tetramethylcyclopropyl)methanone   | C <sub>21</sub> H <sub>29</sub> NO                            | NBMMIBNZVQFQEO-<br>UHFFFAOYSA-N |
| 48 | XLR-11   |                            | [1-(5-fluoropentyl)-1 <i>H</i> -indol-3-yl](2,2,3,3-tetramethylcyclopropyl)methanone   | C <sub>21</sub> H <sub>28</sub> FNO                           | PXLDPUUMIHVLEC-<br>UHFFFAOYSA-N |
| 49 | Zipeprol   |                            | $\alpha$ -( $\alpha$ -methoxybenzyl)-4-( $\beta$ -methoxyphenethyl)-   | C <sub>23</sub> H <sub>32</sub> N <sub>2</sub> O <sub>3</sub> | VSTNNAYSCJQCQI-                 |

|            |  |                     |  |              |
|------------|--|---------------------|--|--------------|
| (zipeprol) |  | 1-piperazineethanol |  | UHFFFAOYSA-N |
|------------|--|---------------------|--|--------------|

2.2. In addition, the stereoisomers of the above substances, if they correspond to the chemical name indicated, exist on the basis of their chemical structure and are not expressly covered by exceptional provisions, and their salts, where such salts exist.

2.3. The following substances are included in List P2:

2.3.1. substances included in the updated Schedule II of the Psychotropic Convention,

2.3.2. substances whose national control is stricter, on the basis of national and international law enforcement experience, than the requirement of the Psychotropic Convention—they are marked with \*\*,

2.3.3. substances whose national control is stricter on the basis of national and international law enforcement experience and which are not included in the Schedules of the Narcotics Convention and the Psychotropic Convention—they are marked with \*\*\*.

2.4. For a substance marked with the symbol <sup>2</sup>, dronabinol refers only to (-)-trans-delta-9-tetrahydrocannabinol stereoisomer.

### 3. List 3 of psychotropic substances (List P3)

3.1. The following substances and compounds are classified as psychotropic substances:

|   | A  | B   | C  | D   | E                               |
|---|--|---|--|---|---------------------------------|
| 1 | <b>Official name</b><br>(Hungarian name) | Other name or<br>abbreviation, spelling<br>commonly used abroad | Chemical name / <i>Description</i>         | Molecular formula   | InChIKey chemical<br>identifier |
| 2 | Amobarbital<br>(amobarbitál)             |   | 5-ethyl-5-isopentylbarbituric acid         | C <sub>11</sub> H <sub>18</sub> N <sub>2</sub> O <sub>3</sub> | VIROVYVQCGLCII-<br>UHFFFAOYSA-N |
| 3 | Butalbital<br>(butalbitál)               |   | 5-allyl-5-isobutylbarbituric acid          | C <sub>11</sub> H <sub>16</sub> N <sub>2</sub> O <sub>3</sub> | UZVHFVZFNXBMQJ-<br>UHFFFAOYSA-N |
| 4 | Cathine<br>(katin)                       | (+)-<br>norpseudoephedrine                                      | (+)-(S)-α-[(S)-1-aminoethyl]benzyl alcohol | C <sub>9</sub> H <sub>13</sub> NO                             | DLNKOYKMWOXYQA-<br>IONNQARKSA-N |



|   |                                  |  |  |  |                                 |
|---|----------------------------------|--|--|--|---------------------------------|
| 5 | Cyclobarbitál<br>(ciklobarbitál) |  | 5-(1-cyclohexen-1-yl)-5-ethylbarbituric acid   | C <sub>12</sub> H <sub>16</sub> N <sub>2</sub> O <sub>3</sub>  | WTYGAUXICFETTC-<br>UHFFFAOYSA-N |
| 6 | Flunitrazepam<br>(flunitrazepám) |  | 5-( <i>o</i> -fluorophenyl)-1,3-dihydro-1-methyl-7-nitro-2 <i>H</i> -1,4-benzodiazepin-2-one | C <sub>16</sub> H <sub>12</sub> FN <sub>3</sub> O <sub>3</sub> | PPTYJKAXVCCBDU-<br>UHFFFAOYSA-N |
| 7 | Glutethimid<br>(glutetimid)      |  | 2-ethyl-2-phenylglutarimide  | C <sub>13</sub> H <sub>15</sub> NO <sub>2</sub>                | JMBQKKAJIKAWKF-<br>UHFFFAOYSA-N |
| 8 | Pentobarbitál<br>(pentobarbitál) |  | 5-ethyl-5-(1-methylbutyl)barbituric acid   | C <sub>11</sub> H <sub>18</sub> N <sub>2</sub> O <sub>3</sub>  | WEXRUCMBJFQVBZ-<br>UHFFFAOYSA-N |

3.2. In addition, the stereoisomers of the above substances, if they correspond to the chemical name indicated, exist on the basis of their chemical structure and are not expressly covered by exceptional provisions, and their salts, where such salts exist.

3.3. This P3 list is based on the updated Schedule III of the Psychotropic Convention.

#### 4. List 4 of psychotropic substances (List P4)

4.1. The following substances and compounds are classified as psychotropic substances:

|   | <b>A</b>                                 | <b>B</b>  | <b>C</b>   | <b>D</b>  | <b>E</b>                        |
|---|--|---|--|---|---------------------------------|
| 1 | <b>Official name</b><br>(Hungarian name) | Other name or<br>abbreviation, spelling<br>commonly used abroad | Chemical name / <i>Description</i>   | Molecular formula   | InChIKey chemical<br>identifier |
| 2 | Allobarbitál<br>(allobarbitál)           |   | 5,5-diallylbarbituric acid   | C <sub>10</sub> H <sub>12</sub> N <sub>2</sub> O <sub>3</sub> | FDQGNLOWMMVRQL-<br>UHFFFAOYSA-N |
| 3 | Alprazolám<br>(alprazolám)               |   | 8-chloro-1-methyl-6-phenyl-4 <i>H</i> -s-triazolo[4,3- <i>a</i> ][1,4]benzodiazepine | C <sub>17</sub> H <sub>13</sub> ClN <sub>4</sub>              | VREFGVBLTWBCJP-<br>UHFFFAOYSA-N |
| 4 | Amfepramón<br>(amfepramón)               | diethylpropion  | 2-(diethylamino)propiofenone   | C <sub>13</sub> H <sub>19</sub> NO                            | XXEPPPIWZFIQJ-<br>UHFFFAOYSA-N  |
| 5 | Aminorex                                 |   | 2-amino-5-phenyl-2-oxazoline   | C <sub>9</sub> H <sub>10</sub> N <sub>2</sub> O               | SYAKTDIEAPMBAL-                 |

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|    |                                      |               |   |   |                                  |
|----|--------------------------------------|---------------|---|---|----------------------------------|
|    | (aminorex)                           |               |   |   | UHFFFAOYSA-N                     |
| 6  | Barbital<br>(barbital)               |               | 5,5-diethylbarbituric acid  | C <sub>8</sub> H <sub>12</sub> N <sub>2</sub> O <sub>3</sub>                  | FTOAOBMCPZCFFF-<br>UHFFFAOYSA-N  |
| 7  | Benzfetamine<br>(benzfetamin)        | benzphetamine | <i>N</i> -benzyl- <i>N</i> , $\alpha$ -dimethylphenethylamine   | C <sub>17</sub> H <sub>21</sub> N   | YXKTVDFXDRQTKV-<br>UHFFFAOYSA-N  |
| 8  | Bromazepam<br>(bromazepám)           |               | 7-bromo-1,3-dihydro-5-(2-pyridyl)-2 <i>H</i> -1,4-benzodiazepin-2-one   | C <sub>14</sub> H <sub>10</sub> BrN <sub>3</sub> O                            | VMIYHDSEFNYSJL-<br>UHFFFAOYSA-N  |
| 9  | Brotizolam<br>(brotizolám)           |               | 2-bromo-4-( <i>o</i> -chlorophenyl)-9-methyl-6 <i>H</i> -thieno[3,2- <i>f</i> ]- <i>s</i> -triazolo[4,3- <i>a</i> ][1,4]diazepine | C <sub>15</sub> H <sub>10</sub> BrClN <sub>4</sub> S                          | UMSGKTJDUHERQW-<br>UHFFFAOYSA-N  |
| 10 | Butobarbital<br>(butobarbitál)       |               | 5-butyl-5-ethylbarbituric acid  | C <sub>10</sub> H <sub>16</sub> N <sub>2</sub> O <sub>3</sub>                 | STDBAQMTJLUMFW-<br>UHFFFAOYSA-N  |
| 11 | Camazepam<br>(kamazepám)             |               | 7-chloro-1,3-dihydro-3-hydroxy-1-methyl-5-phenyl-2 <i>H</i> -1,4-benzodiazepin-2-one dimethylcarbamate (ester)                    | C <sub>19</sub> H <sub>18</sub> ClN <sub>3</sub> O <sub>3</sub>               | PXBVEXGRHZFEOf-<br>UHFFFAOYSA-N  |
| 12 | Chlordiazepoxide<br>(klórdiazepoxid) |               | 7-chloro-2-(methylamino)-5-phenyl-3 <i>H</i> -1,4-benzodiazepine-4-oxide  | C <sub>16</sub> H <sub>14</sub> ClN <sub>3</sub> O                            | ANTSCNMPPGJYLG-<br>UHFFFAOYSA-N  |
| 13 | Clobazam<br>(klobazám)               |               | 7-chloro-1-methyl-5-phenyl-1 <i>H</i> -1,5-benzodiazepine-2,4(3 <i>H</i> ,5 <i>H</i> )-dione                                      | C <sub>16</sub> H <sub>13</sub> ClN <sub>2</sub> O <sub>2</sub>               | CXOXHMZGEKVPMT-<br>UHFFFAOYSA-N  |
| 14 | Clonazepam<br>(klonazepám)           |               | 5-( <i>o</i> -chlorophenyl)-1,3-dihydro-7-nitro-2 <i>H</i> -1,4-benzodiazepin-2-one   | C <sub>15</sub> H <sub>10</sub> ClN <sub>3</sub> O <sub>3</sub>               | DGBIGWXXNGSACT-<br>UHFFFAOYSA-N  |
| 15 | Clonazolam<br>(klonazolám)           |               | 6-(2-chlorophenyl)-1-methyl-8-nitro-4 <i>H</i> -benzo[ <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine                    | C <sub>17</sub> H <sub>12</sub> ClN <sub>5</sub> O <sub>2</sub>               | XJRGLCAWBRZUFC-<br>UHFFFAOYSA-N  |
| 16 | Clorazepate<br>(klorazepát)          |               | 7-chloro-2,3-dihydro-2-oxo-5-phenyl-1 <i>H</i> -1,4-benzodiazepine-3-carboxylic acid  | C <sub>16</sub> H <sub>11</sub> ClN <sub>2</sub> O <sub>3</sub>               | XDDJGVMJFWAHJX-<br>UHFFFAOYSA-N  |
| 17 | Clotiazepam<br>(klotiazepám)         |               | 5-( <i>o</i> -chlorophenyl)-7-ethyl-1,3-dihydro-1-methyl-2 <i>H</i> -thieno[2,3- <i>e</i> ]-1,4-diazepin-2-one                    | C <sub>16</sub> H <sub>15</sub> ClN <sub>2</sub> OS                           | CHBRHODLKOZEPZ-<br>UHFFFAOYSA-N  |
| 18 | Cloxazolam<br>(kloxazolám)           |               | 10-chloro-11b-( <i>o</i> -chlorophenyl)-2,3,7,11b-tetrahydro-oxazolo-[3,2- <i>d</i> ][1,4]benzodiazepin-6(5 <i>H</i> )-one        | C <sub>17</sub> H <sub>14</sub> Cl <sub>2</sub> N <sub>2</sub> O <sub>2</sub> | ZIXNZOBDFKSQTC-<br>UHFFFAOYSA-N  |
| 19 | Delorazepam<br>(delorazepám)         |               | 7-chloro-5-( <i>o</i> -chlorophenyl)-1,3-dihydro-2 <i>H</i> -1,4-benzodiazepin-2-one  | C <sub>15</sub> H <sub>10</sub> Cl <sub>2</sub> N <sub>2</sub> O              | CHIFCDOIIPRCHCF-<br>UHFFFAOYSA-N |
| 20 | Diazepam                             |               | 7-chloro-1,3-dihydro-1-methyl-5-phenyl-2 <i>H</i> -   | C <sub>16</sub> H <sub>13</sub> ClN <sub>2</sub> O                            | AAOVKJBEBIDNHE-                  |

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|----|--|-------------------------------|---|---|---------------------------------|
|    | (diazepám)                             |                               | 1,4-benzodiazepin-2-one   |   | UHFFFAOYSA-N                    |
| 21 | Diclazepam<br>(diklazepám)             | 2-chlorodiazepam,<br>Ro5-3448 | 7-Chloro-5-(2-chlorophenyl)-1-methyl-1,3-dihydro-2 <i>H</i> -benzo[ <i>e</i> ][1,4]diazepin-2-one                   | C <sub>16</sub> H <sub>12</sub> Cl <sub>2</sub> N <sub>2</sub> O  | VPAYQWRBBOGGPY-<br>UHFFFAOYSA-N |
| 22 | Estazolam<br>(esztazolám)              |                               | 8-chloro-6-phenyl-4 <i>H</i> -s-triazolo[4,3- <i>a</i> ][1,4]benzodiazepine   | C <sub>16</sub> H <sub>11</sub> ClN <sub>4</sub>                  | CDCHDCWJMGXXRH-<br>UHFFFAOYSA-N |
| 23 | Ethchlorvynol<br>(etklórvinol)         |                               | 1-chloro-3-ethyl-1-penten-4-yn-3-ol   | C <sub>7</sub> H <sub>9</sub> ClO                                 | ZEHYJZXQEQOSON-<br>UHFFFAOYSA-N |
| 24 | Ethinamate<br>(etinamát)               |                               | 1-ethynylcyclohexanolcarbamate  | C <sub>9</sub> H <sub>13</sub> NO <sub>2</sub>                    | GXRZIMHKGDIBEW-<br>UHFFFAOYSA-N |
| 25 | Ethyl loflazepate<br>(etil-loflazepát) |                               | ethyl 7-chloro-5-( <i>o</i> -fluorophenyl)-2,3-dihydro-2-oxo-1 <i>H</i> -1,4-benzodiazepine-3-carboxylate           | C <sub>18</sub> H <sub>14</sub> ClFN <sub>2</sub> O <sub>3</sub>  | CUCHJCMWNFEYOM-<br>UHFFFAOYSA-N |
| 26 | Etilamfetamine<br>(etilamfetamin)      | N-ethylamphetamine            | <i>N</i> -ethyl- $\alpha$ -methylphenethylamine   | C <sub>11</sub> H <sub>17</sub> N                                 | YAGBSNMZQKEFCO-<br>UHFFFAOYSA-N |
| 27 | Etizolam<br>(etizolám)                 |                               | 4-(2-chlorophenyl)-2-ethyl-9-methyl-6 <i>H</i> -thieno[3,2- <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine | C <sub>17</sub> H <sub>15</sub> ClN <sub>4</sub> S                | VMZUTJCNQWMAGF-<br>UHFFFAOYSA-N |
| 28 | Fencamfamin<br>(fenkamfamin)           |                               | <i>N</i> -ethyl-3-phenyl-2-norbornanamine   | C <sub>15</sub> H <sub>21</sub> N                                 | IKFBPFGUINLYQI-<br>UHFFFAOYSA-N |
| 29 | Fenproporex<br>(fenproporex)           |                               | ( $\pm$ )-3-[( $\alpha$ -methylphenylethyl)amino]propionitrile  | C <sub>12</sub> H <sub>16</sub> N <sub>2</sub>                    | IQUFSXIQAFFIMR-<br>UHFFFAOYSA-N |
| 30 | Flualprazolam<br>(flualprazolám)       |                               | 8-chloro-6-(2-fluorophenyl)-1-methyl-4 <i>H</i> -benzo[ <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine     | C <sub>17</sub> H <sub>12</sub> ClFN <sub>4</sub>                 | MPZVLJCMGPYWQQ-<br>UHFFFAOYSA-N |
| 31 | Flubromazolam<br>(flubromazolám)       |                               | 8-bromo-6-(2-fluorophenyl)-1-methyl-4 <i>H</i> -benzo[ <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine      | C <sub>17</sub> H <sub>12</sub> BrFN <sub>4</sub>                 | VXGSZBZQCBNUIP-<br>UHFFFAOYSA-N |
| 32 | Fludiazepam<br>(fludiazepám)           |                               | 7-chloro-5-( <i>o</i> -fluorophenyl)-1,3-dihydro-1-methyl-2 <i>H</i> -1,4-benzodiazepin-2-one                       | C <sub>16</sub> H <sub>12</sub> ClFN <sub>2</sub> O               | ROYOYTLGDLIGBX-<br>UHFFFAOYSA-N |
| 33 | Flurazepam<br>(flurazepám)             |                               | 7-chloro-1-[2-(diethylamino)ethyl]-5-( <i>o</i> -fluorophenyl)-1,3-dihydro-2 <i>H</i> -1,4-benzodiazepin-2-one      | C <sub>21</sub> H <sub>23</sub> ClFN <sub>3</sub> O               | SAADBVGJQAEFS-<br>UHFFFAOYSA-N  |
| 34 | Halazepam<br>(halazepám)               |                               | 7-chloro-1,3-dihydro-5-phenyl-1-(2,2,2-trifluoroethyl)-2 <i>H</i> -1,4-benzodiazepin-2-one                          | C <sub>17</sub> H <sub>12</sub> ClF <sub>3</sub> N <sub>2</sub> O | WYCLKVQLVUQKNZ-<br>UHFFFAOYSA-N |
| 35 | Haloxazolam                            |                               | 10-bromo-11 <i>b</i> -( <i>o</i> -fluorophenyl)-2,3,7,11 <i>b</i> -   | C <sub>17</sub> H <sub>14</sub> BrFN <sub>2</sub> O <sub>2</sub>  | XDKCGKQHVBBOHC-                 |

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|    |   |     |   |   |                             |
|----|---|-----|---|---|-----------------------------|
|    | (haloxazolám)                           |     | tetrahydrooxazolo[3,2- <i>d</i> ][1,4]benzodiazepin-6(5 <i>H</i> )-one  |   | UHFFFAOYSA-N                |
| 36 | Ketazolam (ketazolám)                   |     | 11-chloro-8,12 <i>b</i> -dihydro-2,8-dimethyl-12 <i>b</i> -phenyl-4 <i>H</i> -[1,3]oxazino[3,2- <i>d</i> ][1,4]benzodiazepin-4,7(6 <i>H</i> )-dione | C <sub>20</sub> H <sub>17</sub> ClN <sub>2</sub> O <sub>3</sub>               | PWAJCNITSBZRBL-UHFFFAOYSA-N |
| 37 | Lefetamine (lefetamin)                  | SPA | (-)- <i>N,N</i> -dimethyl-1,2-diphenylethylamine  | C <sub>16</sub> H <sub>19</sub> N   | YEJZJVJJPVZXGX-MRXNPFEDSA-N |
| 38 | Loprazolam (loprazolám)                 |     | 6-( <i>o</i> -chlorophenyl)-2,4-dihydro-2-[(4-methyl-1-piperazinyl)methylene]-8-nitro-1 <i>H</i> -imidazo[1,2- <i>a</i> ][1,4]benzodiazepin-1-one   | C <sub>23</sub> H <sub>21</sub> ClN <sub>6</sub> O <sub>3</sub>               | UTEFBSAVJNEPTR-RGEXLXHISA-N |
| 39 | Lorazepam (lorazepám)                   |     | 7-chloro-5-( <i>o</i> -chlorophenyl)-1,3-dihydro-3-hydroxy-2 <i>H</i> -1,4-benzodiazepin-2-one  | C <sub>15</sub> H <sub>10</sub> Cl <sub>2</sub> N <sub>2</sub> O <sub>2</sub> | DIWRORZWFLOCLC-UHFFFAOYSA-N |
| 40 | Lormetazepam (lormetazepám)             |     | 7-chloro-5-( <i>o</i> -chlorophenyl)-1,3-dihydro-3-hydroxy-1-methyl-2 <i>H</i> -1,4-benzodiazepin-2-one   | C <sub>16</sub> H <sub>12</sub> Cl <sub>2</sub> N <sub>2</sub> O <sub>2</sub> | FJKWGRGCXUCUIG-UHFFFAOYSA-N |
| 41 | Mazindol (mazindol)                     |     | 5-( <i>p</i> -chlorophenyl)-2,5-dihydro-3 <i>H</i> -imidazo[2,1- <i>a</i> ]isoindol-5-ol  | C <sub>16</sub> H <sub>13</sub> ClN <sub>2</sub> O                            | ZPXSCAKFGYXMGA-UHFFFAOYSA-N |
| 42 | Medazepam (medazepám)                   |     | 7-chloro-2,3-dihydro-1-methyl-5-phenyl-1 <i>H</i> -1,4-benzodiazepine   | C <sub>16</sub> H <sub>15</sub> ClN <sub>2</sub>                              | YLCXGBZIZBEVPZ-UHFFFAOYSA-N |
| 43 | Mefenorex (mefenorex)                   |     | <i>N</i> -(3-chloropropyl)- $\alpha$ -methylphenethylamine  | C <sub>12</sub> H <sub>18</sub> ClN   | XXVROGAVTTXONC-UHFFFAOYSA-N |
| 44 | Meprobamate (meprobamát)                |     | 2-methyl-2-propyl-1,3-propanedioldicarbamate  | C <sub>9</sub> H <sub>18</sub> N <sub>2</sub> O <sub>4</sub>                  | NPPQSCRMBWNHMW-UHFFFAOYSA-N |
| 45 | Mesocarb (mezokarb)                     |     | 3-( $\alpha$ -methylphenethyl)- <i>N</i> -(phenylcarbamoyl)sydnone imine  | C <sub>18</sub> H <sub>18</sub> N <sub>4</sub> O <sub>2</sub>                 | OWFUPROYPKGHMH-UHFFFAOYSA-N |
| 46 | Methylphenobarbital (metilfenobarbitál) |     | 5-ethyl-1-methyl-5-phenylbarbituric acid  | C <sub>13</sub> H <sub>14</sub> N <sub>2</sub> O <sub>3</sub>                 | ALARQZQTBTVLJV-UHFFFAOYSA-N |
| 47 | Methyprylon (metiprilon)                |     | 3,3-diethyl-5-methyl-2,4-piperidine-dione   | C <sub>10</sub> H <sub>17</sub> NO <sub>2</sub>                               | SIDLZWOQUZRBRU-UHFFFAOYSA-N |
| 48 | Midazolam (midazolám)                   |     | 8-chloro-6-( <i>o</i> -fluorophenyl)-1-methyl-4 <i>H</i> -imidazo[1,5- <i>a</i> ][1,4]benzodiazepine  | C <sub>18</sub> H <sub>13</sub> ClFN <sub>3</sub>                             | DDLIGBOFAVUZHB-UHFFFAOYSA-N |
| 49 | Nimetazepam                             |     | 1,3-dihydro-1-methyl-7-nitro-5-phenyl-2 <i>H</i> -  | C <sub>16</sub> H <sub>13</sub> N <sub>3</sub> O <sub>3</sub>                 | GWUSZQUVEVMBPI-             |

## DRAFT

|    |                                    |  |  |   |                             |
|----|------------------------------------|--|--|---|-----------------------------|
|    | (nimetazepám)                      |  | 1,4-benzodiazepin-2-one  |   | UHFFFAOYSA-N                |
| 50 | Nitrazepam (nitrazepám)            |  | 1,3-dihydro-7-nitro-5-phenyl-2 <i>H</i> -1,4-benzodiazepin-2-one   | C <sub>15</sub> H <sub>11</sub> N <sub>3</sub> O <sub>3</sub>   | KJONHKAYOJNZEC-UHFFFAOYSA-N |
| 51 | Nordazepam (nordazepám)            |  | 7-chloro-1,3-dihydro-5-phenyl-2 <i>H</i> -1,4-benzodiazepin-2-one  | C <sub>15</sub> H <sub>11</sub> ClN <sub>2</sub> O              | AKPLHCDWDRPJGD-UHFFFAOYSA-N |
| 52 | Oxazepam (oxazepám)                |  | 7-chloro-1,3-dihydro-3-hydroxy-5-phenyl-2 <i>H</i> -1,4-benzodiazepin-2-one  | C <sub>15</sub> H <sub>11</sub> ClN <sub>2</sub> O <sub>2</sub> | ADIMAYPTOBDMTL-UHFFFAOYSA-N |
| 53 | Oxazolam (oxazolám)                |  | 10-chloro-2,3,7,11 <i>b</i> -tetrahydro-2-methyl-11 <i>b</i> -phenyloxazolo[3,2- <i>d</i> ][1,4]benzodiazepin-6(5 <i>H</i> )-one | C <sub>18</sub> H <sub>17</sub> ClN <sub>2</sub> O <sub>2</sub> | VCCZBYPHZRWKFY-UHFFFAOYSA-N |
| 54 | Pemoline (pemolin)                 |  | 2-amino-5-phenyl-2-oxazolin-4-one  | C <sub>9</sub> H <sub>8</sub> N <sub>2</sub> O <sub>2</sub>     | NRNCYVBFPDDJNE-UHFFFAOYSA-N |
| 55 | Phendimetrazine (fendimetrazin)    |  | (+)-(2 <i>S</i> ,3 <i>S</i> )-3,4-dimethyl-2-phenylmorpholine  | C <sub>12</sub> H <sub>17</sub> NO                              | MFOCDFTXLCYLKU-CMPLNLGQSA-N |
| 56 | Phenobarbital (fenobarbitál)       |  | 5-ethyl-5-phenylbarbituric acid  | C <sub>12</sub> H <sub>12</sub> N <sub>2</sub> O <sub>3</sub>   | DDBREPKUVSBGFI-UHFFFAOYSA-N |
| 57 | Phentermine (fentermin)            |  | $\alpha,\alpha$ -dimethylphenethylamine  | C <sub>10</sub> H <sub>15</sub> N                               | DHHVAGZRUROJKS-UHFFFAOYSA-N |
| 58 | Pinazepam (pinazepám)              |  | 7-chloro-1,3-dihydro-5-phenyl-1-(2-propynyl)-2 <i>H</i> -1,4-benzodiazepin-2-one   | C <sub>18</sub> H <sub>13</sub> ClN <sub>2</sub> O              | MFZOSKPPVCIFMT-UHFFFAOYSA-N |
| 59 | Pipradrol (pipradrol)              |  | 1,1-diphenyl-1-(2-piperidyl)methanol   | C <sub>18</sub> H <sub>21</sub> NO                              | XSWHNYGMWWVAIE-UHFFFAOYSA-N |
| 60 | Prazepam (prazepám)                |  | 7-chloro-1-(cyclopropylmethyl)-1,3-dihydro-5-phenyl-2 <i>H</i> -1,4-benzodiazepin-2-one  | C <sub>19</sub> H <sub>17</sub> ClN <sub>2</sub> O              | MWQCHHACWWAQLJ-UHFFFAOYSA-N |
| 61 | Pyrovalerone (pirovaleron)         |  | 4'-methyl-2-(1-pyrrolidinyl)valerophenone  | C <sub>16</sub> H <sub>23</sub> NO                              | SWUVZKWCBOGPTH-UHFFFAOYSA-N |
| 62 | Secbutabarbital (szekbutabarbitál) |  | 5-sec-butyl-5-ethylbarbituric acid   | C <sub>10</sub> H <sub>16</sub> N <sub>2</sub> O <sub>3</sub>   | ZRIHAIZYIMGOAB-UHFFFAOYSA-N |
| 63 | Temazepam (temazepám)              |  | 7-chloro-1,3-dihydro-3-hydroxy-1-methyl-5-phenyl-2 <i>H</i> -1,4-benzodiazepin-2-one   | C <sub>16</sub> H <sub>13</sub> ClN <sub>2</sub> O <sub>2</sub> | SEQDDYPDSLOBDC-UHFFFAOYSA-N |
| 64 | Tetrazepam (tetrazepám)            |  | 7-chloro-5-(1-cyclohexen-1-yl)-1,3-dihydro-1-methyl-2 <i>H</i> -1,4-benzodiazepin-2-one  | C <sub>16</sub> H <sub>17</sub> ClN <sub>2</sub> O              | IQWYAQCHYZHJOS-UHFFFAOYSA-N |

## DRAFT

|    |                         |  |   |  |                             |
|----|-------------------------|--|---|--|-----------------------------|
| 65 | Triazolam (triazolám)   |  | 8-chloro-6-( <i>o</i> -chlorophenyl)-1-methyl-4 <i>H</i> - <i>s</i> -triazolo[4,3- <i>a</i> ][1,4] benzodiazepine | C <sub>17</sub> H <sub>12</sub> Cl <sub>2</sub> N <sub>4</sub> | JOFWLTCLBGQGBO-UHFFFAOYSA-N |
| 66 | Vinylbital (vinilbitál) |  | 5-(1-methylbutyl)-5-vinylbarbituric acid  | C <sub>11</sub> H <sub>16</sub> N <sub>2</sub> O <sub>3</sub>  | KGKJZEKQJQQOTD-UHFFFAOYSA-N |
| 67 | Zolpidem (zolpidem)     |  | <i>N,N</i> ,6-trimethyl-2- <i>p</i> -tolylimidazo[1,2- <i>a</i> ]pyridine-3-acetamide                             | C <sub>19</sub> H <sub>21</sub> N <sub>3</sub> O               | ZAFYATHCZYHLPB-UHFFFAOYSA-N |

4.2. In addition, the stereoisomers of the above substances, if they correspond to the chemical name indicated, exist on the basis of their chemical structure and are not expressly covered by exceptional provisions, and their salts, where such salts exist.

4.3. This P4 list is based on the updated Schedule IV of the Psychotropic Convention.

1. Point 6.1 of Annex 3 to the Decree is replaced by the following point:

‘6.1. the following compounds:

|    | A   | B   | C   | D                            |
|----|---|---|---|------------------------------|
| 1  | Official name (or other name, abbreviation and spelling commonly used abroad) | Chemical name                                   | Molecular formula   | InChIKey chemical identifier |
| 2  | 2C-B-BZP  | 1-(4-bromo-2,5-dimethoxybenzyl)-4-piperazine    | C <sub>13</sub> H <sub>19</sub> BrN <sub>2</sub> O <sub>2</sub> | OHXVYXBOJDDYJS-UHFFFAOYSA-N  |
| 3  | CPCPP, Gelbes   | 1-(3-chlorophenyl)-4-(3-chloropropyl)piperazine | C <sub>13</sub> H <sub>18</sub> Cl <sub>2</sub> N <sub>2</sub>  | NDQKGEFMUGSRNS-UHFFFAOYSA-N  |
| 4  | Dichlorophenyl-piperazine (diklórfenil-piperazin)                             | 1-(2,3-dichlorophenyl)piperazine                | C <sub>10</sub> H <sub>12</sub> Cl <sub>2</sub> N <sub>2</sub>  | UDQMXYJSNNCRAS-UHFFFAOYSA-N  |
| 5  | Harmin (harmine)  | 7-methoxy-1-methyl-9H-β-carboline               | C <sub>13</sub> H <sub>12</sub> N <sub>2</sub> O                | BXNJHAXVSOCGBA-UHFFFAOYSA-N  |
| 6  | 5-MeO-triptamin (5-MeO-tryptamine)  | 2-(5-methoxy-1H-indol-3-yl)ethanamine           | C <sub>11</sub> H <sub>14</sub> N <sub>2</sub> O                | JTEJPPKMYBDEMY-UHFFFAOYSA-N  |
| 7  | N-benzyl-1-PEA (N-benzil-1-PEA)   | N-benzyl-1-phenylethanamine                     | C <sub>15</sub> H <sub>17</sub> N                               | ZYZHMSJNPCYUTB-UHFFFAOYSA-N  |
| 8  | pCPP  | 1-(4-chlorophenyl)piperazine                    | C <sub>10</sub> H <sub>13</sub> ClN <sub>2</sub>                | UNEIHNMKASENIG-UHFFFAOYSA-N  |
| 9  | 1-phenyl-propyl-amine (fenilpropilamin, 1-phenyl-1-propanamine)               | 1-phenylpropan-1-amine                          | C <sub>9</sub> H <sub>13</sub> N                                | AQFLVLHRZFLDDV-UHFFFAOYSA-N  |
| 10 | pMeOPP  | 1-(4-methoxyphenyl)piperazine                   | C <sub>11</sub> H <sub>16</sub> N <sub>2</sub> O                | MRDGZSKYFPGAKP-              |

## DRAFT

|    |   |   |   |                              |
|----|---|---|---|------------------------------|
|    |   |   |   | UHFFFAOYSA-N                 |
| 11 | URB-754   | 6-methyl-2-[(4-methylphenyl)amino]-4H-3,1-benzoxazin-4-one                                | C <sub>16</sub> H <sub>14</sub> N <sub>2</sub> O <sub>2</sub> | GFWNGVKCDGYFKG-UHFFFAOYSA-N  |
| 12 | RH-34   | 3-[2-(2-methoxybenzylamino)ethyl]-1H-quinazoline-2,4-dione                                | C <sub>18</sub> H <sub>19</sub> N <sub>3</sub> O <sub>3</sub> | NUAJBITWGGTZCM-UHFFFAOYSA-N  |
| 13 | MEOP; Methoxypiperamide; metoxipiperamid; MEXP; 1(4methoxybenzoyl)-4-methylpiperazine | (4-methoxyphenyl)(4-methylpiperazin-1-yl)methanone  | C <sub>13</sub> H <sub>18</sub> N <sub>2</sub> O <sub>2</sub> | DWPVVZZGGGCRM-UHFFFAOYSA-N   |
| 14 | bk-2C-B   | 2-amino-1-(4-bromo-2,5-dimethoxyphenyl)ethanone   | C <sub>10</sub> H <sub>12</sub> BrNO <sub>3</sub>             | HFYJGAIIOBIDRPX-UHFFFAOYSA-N |
| 15 | 25I-NBF   | 2-(4-iodo-2,5-dimethoxyphenyl)-N-[(2-fluorophenyl)methyl]ethanamine                       | C <sub>17</sub> H <sub>19</sub> FINO <sub>2</sub>             | LPBKNBHMWRBPHT-UHFFFAOYSA-N  |
| 16 | Mexedrone; 4-MMC-oMe; 'MEX'   | 3-methoxy-2-(methylamino)-1-(4-methylphenyl)propan-1-one                                  | C <sub>12</sub> H <sub>17</sub> NO <sub>2</sub>               | JHGDCSPZKQLBOP-UHFFFAOYSA-N  |
| 17 | TH-PVP  | 2-(pyrrolidin-1-yl)-1-(5,6,7,8-tetrahydronaphthalen-2-yl)pentan-1-one                     | C <sub>19</sub> H <sub>27</sub> NO                            | MMIKQWIZKBYLKZ-UHFFFAOYSA-N  |
| 18 | izopropilfenidát (isopropylphenidate)   | propan-2-yl phenyl(piperidin-2-yl)acetate   | C <sub>16</sub> H <sub>23</sub> NO <sub>2</sub>               | AZVPADMEIMLODT-UHFFFAOYSA-N  |
| 19 | metamnetamin (methylnaphetamine; N-methyl-PAL-287; MNT; MNA)                          | N-methyl-1-(naphthalen-2-yl)propan-2-amine  | C <sub>14</sub> H <sub>17</sub> N                             | BWWWOLYZMKACSB-UHFFFAOYSA-N  |
| 20 | dezklorketamin (deschloroketamine; 2-(phenyl)-2-(methylamino)-cyclohexanone)          | 2-(methylamino)-2-phenylcyclohexanone   | C <sub>13</sub> H <sub>17</sub> NO                            | ZAGBSZSITDFFAF-UHFFFAOYSA-N  |
| 21 | 1P-LSD (1-propionyl-d-lysergic acid diethylamide)                                     | N,N-diethyl-7-methyl-4-propanoyl-6,6a,8,9-tetrahydroindolo[4,3-fg]quinoline-9-carboxamide | C <sub>23</sub> H <sub>29</sub> N <sub>3</sub> O <sub>2</sub> | JSMQOVGXBIDBIE-UHFFFAOYSA-N  |
| 22 | Ethyl-naphthidate; HDEP-28  | ethyl 2-(naphthalen-2-yl)-2-(piperidin-2-yl)acetate                                       | C <sub>19</sub> H <sub>23</sub> NO <sub>2</sub>               | OTQVTBPHZRARTL-UHFFFAOYSA-N  |
| 23 | 4-methylmethylphenidate; 4Me-   | methyl 2-(piperidin-2-yl)-2-(p-tolyl)acetate  | C <sub>15</sub> H <sub>21</sub> NO <sub>2</sub>               | WJZNCJIOIACDBR-              |



## DRAFT

|    |  |   |   |                              |
|----|--|---|---|------------------------------|
|    | TMP  |   |   | UHFFFAOYSA-N                 |
| 24 | 4-fluoromethylphenidate; 4F-TMP; 4F-MPH; 4-FMPH  | methyl 2-(4-fluorophenyl)-2-(piperidin-2-yl)acetate   | C <sub>14</sub> H <sub>18</sub> FNO <sub>2</sub>                | XISBAJBPDVRSFG-UHFFFAOYSA-N  |
| 25 | CUMYL-5F-P7AICA; CUMYL-5F-PAICA; SGT-263   | 1-(5-fluoropentyl)- <i>N</i> -(2-phenylpropan-2-yl)-1 <i>H</i> -pyrrolo[2,3- <i>b</i> ]pyridine-3-carboxamide                                 | C <sub>22</sub> H <sub>26</sub> FN <sub>3</sub> O               | MXJYOUUMYJGNQEY-UHFFFAOYSA-N |
| 26 | (Iso)butyryl-F-fentanyl <i>N</i> -benzyl analogue                                      | 2-methyl- <i>N</i> -(1-benzylpiperidin-4-yl)- <i>N</i> -(4-fluorophenyl)propanamide   | C <sub>22</sub> H <sub>27</sub> FN <sub>2</sub> O               | XNQGKYHSTDKIKG-UHFFFAOYSA-N  |
| 27 | Despropionyl-2-fluoro fentanyl; despropionyl- <i>o</i> -fluoro fentanyl                | <i>N</i> -(2-fluorophenyl)-1-(2-phenylethyl)piperidin-4-amine   | C <sub>19</sub> H <sub>23</sub> FN <sub>2</sub>                 | WUNLGTOLOUTCPE-UHFFFAOYSA-N  |
| 28 | Flubromazepam  | 7-bromo-5-(2-fluorophenyl)-1,3-dihydro-2 <i>H</i> -1,4-benzodiazepin-2-one  | C <sub>15</sub> H <sub>10</sub> BrFN <sub>2</sub> O             | ZRKDDZBVSZLOFS-UHFFFAOYSA-N  |
| 29 | 5-fluoropentyl-3-pyridinoylindole  | [1-(5-fluoropentyl)-1 <i>H</i> -indol-3-yl](pyridin-3-yl)methanone  | C <sub>19</sub> H <sub>19</sub> FN <sub>2</sub> O               | CNMQLCYLJPWHEW-UHFFFAOYSA-N  |
| 30 | Phenibut   | 4-amino-3-phenylbutanoic acid   | C <sub>10</sub> H <sub>13</sub> NO <sub>2</sub>                 | DAFOCGYVTAOKAJ-UHFFFAOYSA-N  |
| 31 | 2-fluorodeschloroketamine; 2-Fl-2'-Oxo-PCM; fluoroketamine; 2-FDCK; 2F-DK; 2-FDK; 2-FK | 2-(2-fluorophenyl)-2-(methylamino)cyclohexanone   | C <sub>13</sub> H <sub>16</sub> FNO                             | PHFAGYYTDLITB-UHFFFAOYSA-N   |
| 32 | Dichloropane; RTI-111; RTI-4229-111; O-401   | methyl 3-(3,4-dichlorophenyl)-8-methyl-8-azabicyclo[3.2.1]octane-2-carboxylate  | C <sub>16</sub> H <sub>19</sub> Cl <sub>2</sub> NO <sub>2</sub> | AMIHUYQKNJHXPT-UHFFFAOYSA-N  |
| 33 | AMAPN  | 2-(methylamino)-1-(naphthalen-1-yl)propan-1-one   | C <sub>14</sub> H <sub>15</sub> NO                              | QIACKSHQBOUATI-UHFFFAOYSA-N  |
| 34 | 5Cl-bk-MPA; 5Cl-bk-methylthienylpropamine; 5Cl-bk-methiopropamine                      | 1-(5-chlorothiophen-2-yl)-2-(methylamino)propan-1-one   | C <sub>8</sub> H <sub>10</sub> ClNOS                            | VTSPXQPERGVFBV-UHFFFAOYSA-N  |
| 35 | 5F-Cumyl-PeGaClone; 5F-SGT-151   | 5-(5-fluoropentyl)-2-(2-phenylpropan-2-yl)-2,5-dihydro-1 <i>H</i> -pyrido[4,3- <i>b</i> ]indol-1-one  | C <sub>25</sub> H <sub>27</sub> FN <sub>2</sub> O               | SMRRORRDOWXERZ-UHFFFAOYSA-N  |
| 36 | ALD-52; 1-acetyl-LSD   | (6 <i>aR</i> ,9 <i>R</i> )-4-acetyl- <i>N,N</i> -diethyl-7-methyl-6,6 <i>a</i> ,8,9-tetrahydroindolo[4,3- <i>f,g</i> ]quinoline-9-carboxamide | C <sub>22</sub> H <sub>27</sub> N <sub>3</sub> O <sub>2</sub>   | FJOWXGYLIWJFCH-OXQOHEQNSA-N  |
| 37 | ETH-LAD; 6-ethyl-6-nor-lysergic acid diethylamide                                      | (6 <i>aR</i> ,9 <i>R</i> )- <i>N,N</i> -diethyl-7-ethyl-4,6,6 <i>a</i> ,7,8,9-hexahydroindolo-[4,3- <i>fg</i> ]quinoline-9-                   | C <sub>21</sub> H <sub>27</sub> N <sub>3</sub> O                | MYNOUXJLOHVSMQ-DNVCBOLYSA-N  |

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|    |   |  |   |                              |
|----|---|--|---|------------------------------|
|    |   | carboxamide  |   |                              |
| 38 | diphenyl-PBP                              | 1-(1,3-diphenylpropan-2-yl)pyrrolidine   | C <sub>19</sub> H <sub>23</sub> N                 | VQECHRQHFMUVRS-UHFFFAOYSA-N  |
| 39 | benzoil-fentanil (benzoylfentanyl)        | N-phenyl-N-[1-(2-phenylethyl)-4-piperidyl]benzamide  | C <sub>26</sub> H <sub>28</sub> N <sub>2</sub> O  | BJPDWVPQDSVQKD-UHFFFAOYSA-N  |
| 40 | Cumyl-CH-MegaClone                        | 5-cyclohexylmethyl-2-(2-phenylpropan-2-yl)-2,5-dihydro-1 <i>H</i> -pyrido[4,3- <i>b</i> ]indol-1-one | C <sub>27</sub> H <sub>30</sub> N <sub>2</sub> O  | CGHCGYCTOLWAPL-UHFFFAOYSA-N  |
| 41 | 1-Aminoindan                              | 2,3-dihydro-1 <i>H</i> -inden-1-amine  | C <sub>9</sub> H <sub>11</sub> N                  | XJEVHMGJSYVQBQ-UHFFFAOYSA-N  |
| 42 | N-methyl-2AI                              | <i>N</i> -methyl-2,3-dihydro-1 <i>H</i> -inden-2-amine   | C <sub>10</sub> H <sub>13</sub> N                 | SXWZQUCTTOBHJT-UHFFFAOYSA-N  |
| 43 | M-ALPHA                                   | 1-(1,3-benzodioxol-5-yl)- <i>N</i> -methylpropan-1-amine   | C <sub>11</sub> H <sub>15</sub> NO <sub>2</sub>   | NLINVDHEDVEOMJ-UHFFFAOYSA-N  |
| 44 | Thiopropamine                             | 1-(thiophen-2-yl)propan-2-amine  | C <sub>7</sub> H <sub>11</sub> NS                 | NYVQQTTOGYLBBDQ-UHFFFAOYSA-N |
| 45 | 2-APB                                     | 1-(1-benzofuran-2-yl)propan-2-amine  | C <sub>11</sub> H <sub>13</sub> NO                | QGLBWEFCBFEAPH-UHFFFAOYSA-N  |
| 46 | 2-MAPB                                    | 1-(1-benzofuran-2-yl)- <i>N</i> -methylpropan-2-amine  | C <sub>12</sub> H <sub>15</sub> NO                | ANJIDHKQUCZNQY-UHFFFAOYSA-N  |
| 47 | 2-EAPB                                    | 1-(1-benzofuran-2-yl)- <i>N</i> -ethylpropan-2-amine   | C <sub>13</sub> H <sub>17</sub> NO                | SGGKRTSTBXBERJ-UHFFFAOYSA-N  |
| 48 | 5-MeO-DIBF                                | <i>N</i> -[2-(5-methoxy-1-benzofuran-3-yl)ethyl]- <i>N</i> -(propan-2-yl)propan-2-amine              | C <sub>17</sub> H <sub>25</sub> NO <sub>2</sub>   | NBFMSQBTYHYVKP-UHFFFAOYSA-N  |
| 49 | 2-MABB                                    | 1-(1-benzofuran-2-yl)- <i>N</i> -methylbutan-2-amine   | C <sub>13</sub> H <sub>17</sub> NO                | YBPPNDUCEMDPJZ-UHFFFAOYSA-N  |
| 50 | 1-(4-Bromo-2,5-dimethoxyphenyl)ethanamine | 1-(4-bromo-2,5-dimethoxyphenyl)ethanamine  | C <sub>10</sub> H <sub>14</sub> BrNO <sub>2</sub> | JKOMOTQHGDWZAQ-UHFFFAOYSA-N  |
| 51 | BOH-PHP                                   | 1-phenyl-2-(pyrrolidin-1-yl)hexan-1-ol   | C <sub>16</sub> H <sub>25</sub> NO                | GJHNPZVIGWPXBH-UHFFFAOYSA-N  |
| 52 | 2-MeO-Ketamine                            | 2-(2-methoxyphenyl)-2-(methylamino)cyclohexanone   | C <sub>14</sub> H <sub>19</sub> NO <sub>2</sub>   | OYAUVHORXFUVAJ-UHFFFAOYSA-N  |
| 53 | Methoxetamine brominated                  | 2-(2-bromo-5-methoxyphenyl)-2-   | C <sub>15</sub> H <sub>20</sub> BrNO <sub>2</sub> | YNEMYNGTFPWRHN-              |

## DRAFT

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|    | derivative                                 | (ethylamino)cyclohexanone   |   | UHFFFAOYSA-N                |
| 54 | 3MeO-PCMo                                  | 4-[1-(3-methoxyphenyl)cyclohexyl]morpholine   | C <sub>17</sub> H <sub>25</sub> NO <sub>2</sub>                 | BOGOEDFWPOXWQE-UHFFFAOYSA-N |
| 55 | Benocyclidine                              | 1-[1-(benzothiophen-2-yl)cyclohexyl]piperidine  | C <sub>19</sub> H <sub>25</sub> NS                              | RGSVXQJPSWZXOP-UHFFFAOYSA-N |
| 56 | Tiletamine                                 | 2-(ethylamino)-2-(2-thienyl)cyclohexanone   | C <sub>12</sub> H <sub>17</sub> NOS                             | QAXBVGVDCAVLV-UHFFFAOYSA-N  |
| 57 | deschloro-N-ethyl-ketamine (O-PCE)         | 2-(ethylamino)-2-phenyl-cyclohexanone   | C <sub>14</sub> H <sub>19</sub> NO                              | IDLBAANXISGEI-UHFFFAOYSA-N  |
| 58 | 3-MeO-PCMMo                                | 4-{{1-(3-methoxyphenyl)cyclohexyl}methyl}morpholine   | C <sub>18</sub> H <sub>27</sub> NO <sub>2</sub>                 | KVDDTOKOCUZIFC-UHFFFAOYSA-N |
| 59 | 3-HO-PCE                                   | 3-[1-(ethylamino)cyclohexyl]phenol  | C <sub>14</sub> H <sub>21</sub> NO                              | MIKNPNLBFHVMKK-UHFFFAOYSA-N |
| 60 | 3-HO-PCP                                   | 3-(1-piperidin-1-ylcyclohexyl)phenol  | C <sub>17</sub> H <sub>25</sub> NO                              | AMSXTZUCNOKUEN-UHFFFAOYSA-N |
| 61 | Methoxpropamine                            | 2-(3-methoxyphenyl)-2-(propylamino)cyclohexan-1-one   | C <sub>16</sub> H <sub>23</sub> NO <sub>2</sub>                 | AAVOSBAXDRASAH-UHFFFAOYSA-N |
| 62 | Pyrazolam                                  | 8-bromo-1-methyl-6-(pyridin-2-yl)-4 <i>H</i> -[1,2,4]triazolo[4,3- <i>a</i> ][1,4]benzodiazepine                          | C <sub>16</sub> H <sub>12</sub> BrN <sub>5</sub>                | BGRWSFIQPVEML-UHFFFAOYSA-N  |
| 63 | Alprazolam triazolobenzophenone derivative | {2-[3-(aminomethyl)-5-methyl-4 <i>H</i> -1,2,4-triazol-4-yl]-5-chlorophenyl}phenyl-methanone                              | C <sub>17</sub> H <sub>15</sub> ClN <sub>4</sub> O              | WWADXOXMCNJJKR-UHFFFAOYSA-N |
| 64 | Meclonazepam                               | 5-(2-chlorophenyl)-3-methyl-7-nitro-1,3-dihydro-1,4-benzodiazepin-2-one   | C <sub>16</sub> H <sub>12</sub> ClN <sub>3</sub> O <sub>3</sub> | LMUVYJCAFVGNSY-UHFFFAOYSA-N |
| 65 | Deschloroetizolam                          | 2-ethyl-9-methyl-4-phenyl-6 <i>H</i> -thieno[3,2- <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine                 | C <sub>17</sub> H <sub>16</sub> N <sub>4</sub> S                | JIOBORXCOGMHSV-UHFFFAOYSA-N |
| 66 | Nifoxipam                                  | 5-(2-fluorophenyl)-3-hydroxy-7-nitro-1,3-dihydro-2 <i>H</i> -1,4-benzodiazepin-2-one                                      | C <sub>15</sub> H <sub>10</sub> FN <sub>3</sub> O <sub>4</sub>  | UHFIFTRHLBAWGY-UHFFFAOYSA-N |
| 67 | Adinazolam                                 | 1-(8-chloro-6-phenyl-4 <i>H</i> -[1,2,4]triazolo[4,3- <i>a</i> ][1,4]benzodiazepin-1-yl)- <i>N,N</i> -dimethylmethanamine | C <sub>19</sub> H <sub>18</sub> ClN <sub>5</sub>                | GJSLOMWRLALDCT-UHFFFAOYSA-N |
| 68 | Metizolam                                  | 4-(2-chlorophenyl)-2-ethyl-6 <i>H</i> -thieno[3,2- <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine                | C <sub>16</sub> H <sub>13</sub> ClN <sub>4</sub> S              | NQSSWDKQLVBUQN-UHFFFAOYSA-N |

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|----|---------------------|--|---|-----------------------------|
| 69 | Nitrazolam          | 1-methyl-8-nitro-6-phenyl-4 <i>H</i> -1,2,4]triazolo[4,3- <i>a</i> ][1,4]benzodiazepine                              | C <sub>17</sub> H <sub>13</sub> N <sub>5</sub> O <sub>2</sub>     | OYRPNABWTHDOFK-UHFFFAOYSA-N |
| 70 | Cloniprazepam       | 5-(2-chlorophenyl)-1-(cyclopropylmethyl)-7-nitro-1,3-dihydro-2 <i>H</i> -1,4-benzodiazepin-2-one                     | C <sub>19</sub> H <sub>16</sub> ClN <sub>3</sub> O <sub>3</sub>   | CCSYKGYLSFXNTA-UHFFFAOYSA-N |
| 71 | Cinazepam           | 4-{{[7-bromo-5-(2-chlorophenyl)-2-oxo-1,3-dihydro-1,4- benzodiazepin-3-yl]oxy}-4-oxo-butanoic acid                   | C <sub>19</sub> H <sub>14</sub> BrClN <sub>2</sub> O <sub>5</sub> | NQTRBZXDWMDXQA-UHFFFAOYSA-N |
| 72 | 3-hydroxyphenazepam | 7-bromo-5-(2-chlorophenyl)-3-hydroxy-1,3-dihydro-1,4- benzodiazepin-2-one  | C <sub>15</sub> H <sub>10</sub> BrClN <sub>2</sub> O <sub>2</sub> | KRJKJUWAZOWXNV-UHFFFAOYSA-N |
| 73 | Fonazepam           | 5-(2-fluorophenyl)-1,3-dihydro-7-nitro-2 <i>H</i> -1,4-benzodiazepin-2-one   | C <sub>15</sub> H <sub>10</sub> FN <sub>3</sub> O <sub>3</sub>    | KNGIGRDYBQPXKQ-UHFFFAOYSA-N |
| 74 | 4-chlorodiazepam    | 7-chloro-5-(4-chlorophenyl)-1-methyl-3 <i>H</i> -1,4-benzodiazepin-2-one   | C <sub>16</sub> H <sub>12</sub> Cl <sub>2</sub> N <sub>2</sub> O  | PUMYFTJOWAJIKF-UHFFFAOYSA-N |
| 75 | Flunitrazolam       | 6-(2-fluorophenyl)-1-methyl-8-nitro-4 <i>H</i> -[1,2,4]triazolo[4,3- <i>a</i> ][1,4]benzodiazepine                   | C <sub>17</sub> H <sub>12</sub> FN <sub>5</sub> O <sub>2</sub>    | RDLAGIOILLWVTM-UHFFFAOYSA-N |
| 76 | Bromazolam          | 8-bromo-1-methyl-6-phenyl-4 <i>H</i> -[1,2,4]triazolo[4,3- <i>a</i> ][1,4]benzodiazepine                             | C <sub>17</sub> H <sub>13</sub> BrN <sub>4</sub>                  | KCEIOBKDDQAYCM-UHFFFAOYSA-N |
| 77 | Norfludiazepam      | 7-chloro-5-(2-fluorophenyl)-1,3-dihydro-1,4-benzodiazepin-2-one  | C <sub>15</sub> H <sub>10</sub> ClFN <sub>2</sub> O               | UVCOILFBWYKHHB-UHFFFAOYSA-N |
| 78 | Ro 07-4065          | 7-chloro-5-(2,6-difluorophenyl)-1-methyl-3 <i>H</i> -1,4-benzodiazepin-2-one   | C <sub>16</sub> H <sub>11</sub> ClF <sub>2</sub> N <sub>2</sub> O | DUNFPASORLTEGN-UHFFFAOYSA-N |
| 79 | Thionordazepam      | 7-chloro-5-phenyl-1,3-dihydro-1,4-benzodiazepine-2-thione  | C <sub>15</sub> H <sub>11</sub> ClN <sub>2</sub> S                | ULILTJWAJZIROM-UHFFFAOYSA-N |
| 80 | Methyl clonazepam   | 5-(2-chlorophenyl)-1-methyl-7-nitro-3 <i>H</i> -1,4-benzodiazepin-2-one  | C <sub>16</sub> H <sub>12</sub> ClN <sub>3</sub> O <sub>3</sub>   | AZVBJDUDXZLTM-UHFFFAOYSA-N  |
| 81 | Fluclotizolam       | 2-chloro-4-(2-fluorophenyl)-9-methyl-6 <i>H</i> -thieno[3,2- <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine | C <sub>15</sub> H <sub>10</sub> ClFN <sub>4</sub> S               | ZDYRCUZZLRLMHG-UHFFFAOYSA-N |
| 82 | Clobromazolam       | 8-bromo-6-(2-chlorophenyl)-1-methyl-4 <i>H</i> -[1,2,4]triazolo[4,3- <i>a</i> ][1,4]benzodiazepine                   | C <sub>17</sub> H <sub>12</sub> BrClN <sub>4</sub>                | BUTCFAZTKZDYCN-UHFFFAOYSA-N |
| 83 | Bentazepam          | 5-phenyl-1,3,6,7,8,9-hexahydro-2 <i>H</i> -[1]benzothieno[2,3- <i>e</i> ][1,4]diazepin-2-one                         | C <sub>17</sub> H <sub>16</sub> N <sub>2</sub> OS                 | AIZFEOPQVZBNGH-UHFFFAOYSA-N |
| 84 | CP 47,497           | 5-(1,1-dimethylheptyl)-2-[(1 <i>R</i> ,3 <i>S</i> )-3-   | C <sub>21</sub> H <sub>34</sub> O <sub>2</sub>                    | ZWWRREXSUJTKNN-             |

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|----|---|--|--|-----------------------------|
|    |   | hydroxycyclohexyl]phenol   |  | AEFFLSMTSA-N                |
| 85 | AM-1220 Azepane Isomer  | [1-(hexahydro-1-methyl-1 <i>H</i> -azepin-3-yl)-1 <i>H</i> -indol-3-yl]-1-naphthalenyl-methanone                                       | C <sub>26</sub> H <sub>26</sub> N <sub>2</sub> O                             | ZDCZZWAEXISRJF-UHFFFAOYSA-N |
| 86 | Org 27569   | 5-chloro-3-ethyl- <i>N</i> -[2-[4-(1-piperidinyl)phenyl]ethyl]-1 <i>H</i> -indole-2-carboxamide  | C <sub>24</sub> H <sub>28</sub> ClN <sub>3</sub> O                           | AHFZDNYNXFMRFQ-UHFFFAOYSA-N |
| 87 | Org 27759   | <i>N</i> -[2-[4-(dimethylamino)phenyl]ethyl]-3-ethyl-5-fluoro-1 <i>H</i> -indole-2-carboxamide   | C <sub>21</sub> H <sub>24</sub> FN <sub>3</sub> O                            | MUYUEZAKMLKZSO-UHFFFAOYSA-N |
| 88 | Org 29647   | <i>N</i> -(1-benzylpyrrolidin-3-yl)-5-chloro-3-ethyl-1 <i>H</i> -indole-2-carboxamide  | C <sub>22</sub> H <sub>24</sub> ClN <sub>3</sub> O                           | MYJJVFCPCDPOSM-UHFFFAOYSA-N |
| 89 | HU-331  | 3-hydroxy-2-(6-isopropenyl-3-methyl-cyclohex-2-en-1-yl)-5-pentyl-1,4-benzoquinone  | C <sub>21</sub> H <sub>28</sub> O <sub>3</sub>                               | WDXXEUARVHTWQF-UHFFFAOYSA-N |
| 90 | WIN 55212-2   | [(3 <i>R</i> )-2,3-dihydro-5-methyl-3-(4-morpholinylmethyl)pyrrolo[1,2,3- <i>de</i> ]-1,4-benzoxazin-6-yl]-1-naphthalen-1-yl-methanone | C <sub>27</sub> H <sub>26</sub> N <sub>2</sub> O <sub>3</sub>                | HQVHOQAKMCMIIM-HXUWFJFHSA-N |
| 91 | AB-005 azepane isomer   | [1-(1-methylazepan-2-yl)indol-3-yl]-(2,2,3,3-tetramethylcyclopropyl)methanone  | C <sub>23</sub> H <sub>32</sub> N <sub>2</sub> O                             | VBJSVQSAWSBQA-UHFFFAOYSA-N  |
| 92 | 4-HTMPIPO   | 4-hydroxy-3,3,4-trimethyl-1-(1-pentyl-1 <i>H</i> -indol-3-yl)pentan-1-one  | C <sub>21</sub> H <sub>31</sub> NO <sub>2</sub>                              | GWHGUYKAGQLPTQ-UHFFFAOYSA-N |
| 93 | URB-597   | 3'-carbamoylbiphenyl-3-yl cyclohexylcarbamate  | C <sub>20</sub> H <sub>22</sub> N <sub>2</sub> O <sub>3</sub>                | ROFVXGGUISEHAM-UHFFFAOYSA-N |
| 94 | JTE-907   | <i>N</i> -(1,3-benzodioxol-5-ylmethyl)-2-hydroxy-7-methoxy-8-pentoxo-quinoline-3-carboxamide   | C <sub>24</sub> H <sub>26</sub> N <sub>2</sub> O <sub>6</sub>                | GRAJFFFJYFVOC-UHFFFAOYSA-N  |
| 95 | LY2183240   | 5-([1,1'-biphenyl]-4-ylmethyl)- <i>N,N</i> -dimethyl-1 <i>H</i> -tetrazole-1-carboxamide   | C <sub>17</sub> H <sub>17</sub> N <sub>5</sub> O                             | GZNIYOXWFCDBBJ-UHFFFAOYSA-N |
| 96 | 1-(Cyclohexylmethyl)-2-[(4-ethoxyphenyl)methyl]- <i>N,N</i> -diethyl-1 <i>H</i> -benzimidazol-5-carboxamide | 1-(cyclohexylmethyl)-2-[(4-ethoxyphenyl)methyl]- <i>N,N</i> -diethyl-benzimidazole-5-carboxamide                                       | C <sub>28</sub> H <sub>37</sub> N <sub>3</sub> O <sub>2</sub>                | WRVZBXHTUOPQJS-UHFFFAOYSA-N |
| 97 | AM-2201 benzimidazole analogue (FUBIMINA)   | [1-(5-fluoropentyl)-1 <i>H</i> -benzimidazol-2-yl]-(naphthalen-1-yl)methanone  | C <sub>23</sub> H <sub>21</sub> FN <sub>2</sub> O                            | KUESSZMROAFKQJ-UHFFFAOYSA-N |
| 98 | 5F-AB-FUPPYCA (5F-5,3-AB-   | 2-[[1-(5-fluoropentyl)-5-(4-fluorophenyl)-1 <i>H</i> -   | C <sub>20</sub> H <sub>26</sub> F <sub>2</sub> N <sub>4</sub> O <sub>2</sub> | GSXRDTDYPSATDE-             |

## DRAFT

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|-----|-----------------------------------|---|--|-----------------------------|
|     | PFUPPYCA)                         | pyrazol-3-yl]formamido}-3-methylbutanamide  |  | UHFFFAOYSA-N                |
| 99  | 5F-PY-PICA                        | [1-(5-fluoropentyl)-1 <i>H</i> -indol-3-yl](pyrrolidin-1-yl)methanone   | C <sub>18</sub> H <sub>23</sub> FN <sub>2</sub> O                            | AJOAHRJLOXOZKX-UHFFFAOYSA-N |
| 100 | 5F-PY-PINACA                      | [1-(5-fluoropentyl)-1 <i>H</i> -indazol-3-yl](pyrrolidin-1-yl)methanone   | C <sub>17</sub> H <sub>22</sub> FN <sub>3</sub> O                            | GSCLIRQNUBFUJA-UHFFFAOYSA-N |
| 101 | AB-CHMFUPPYCA (3,5-AB-CHMFUPPYCA) | <i>N</i> -(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-3-(4-fluorophenyl)-1 <i>H</i> -pyrazole-5-carboxamide | C <sub>22</sub> H <sub>29</sub> FN <sub>4</sub> O <sub>2</sub>               | NDYOOVJIZQTGHY-UHFFFAOYSA-N |
| 102 | MDMB-CHMCZCA                      | 9-(cyclohexylmethyl)- <i>N</i> -(1-methoxycarbonyl-2,2-dimethyl-propyl)carbazole-3-carboximidic acid                    | C <sub>27</sub> H <sub>34</sub> N <sub>2</sub> O <sub>3</sub>                | FAWVRKNYDPKTDZ-UHFFFAOYSA-N |
| 103 | 5F-PCN                            | 1-(5-fluoropentyl)- <i>N</i> -(naphthalen-1-yl)-1 <i>H</i> -pyrrolo[3,2- <i>c</i> ]pyridine-3-carboxamide               | C <sub>23</sub> H <sub>22</sub> FN <sub>3</sub> O                            | BRRZRRZUERBDQL-UHFFFAOYSA-N |
| 104 | EG-2201                           | [9-(5-fluoropentyl)-9 <i>H</i> -carbazol-3-yl](naphthalen-1-yl)methanone  | C <sub>28</sub> H <sub>24</sub> FNO  | LYDDINAZVHIBGP-UHFFFAOYSA-N |
| 105 | MDA 19 (BZO-HEXOXIZID)            | <i>N</i> -[( <i>Z</i> )-(1-hexyl-2-oxo-indolin-3-ylidene)amino]benzamide  | C <sub>21</sub> H <sub>23</sub> N <sub>3</sub> O <sub>2</sub>                | ZGQHMZCITJHYOW-QOCHGBHMSA-N |
| 106 | MO-CHMINACA                       | 1-methoxy-3,3-dimethyl-1-oxobutan-2-yl 1-(cyclohexylmethyl)-1 <i>H</i> -indazole-3-carboxylate                          | C <sub>22</sub> H <sub>30</sub> N <sub>2</sub> O <sub>4</sub>                | SUEOBRAXHJBVGY-UHFFFAOYSA-N |
| 107 | MDMB-PCZCA                        | methyl 3,3-dimethyl-2-[(9-pentylcarbazole-3-carbonyl)amino] butanoate   | C <sub>25</sub> H <sub>32</sub> N <sub>2</sub> O <sub>3</sub>                | GILRPTXNJMVJBM-UHFFFAOYSA-N |
| 108 | 5F-3,5-AB-PFUPPYCA                | <i>N</i> -(1-carbamoyl-2-methyl-propyl)-1-(5-fluoropentyl)-3-(4-fluorophenyl)pyrazole-5-carboxamide                     | C <sub>20</sub> H <sub>26</sub> F <sub>2</sub> N <sub>4</sub> O <sub>2</sub> | JPKXVUNTSWGYKJ-UHFFFAOYSA-N |
| 109 | CUMYL-4CN-B7AICA                  | 1-(4-cyanobutyl)- <i>N</i> -(1-methyl-1-phenylethyl)pyrrolo[2,3- <i>b</i> ]pyridine-3-carboxamide                       | C <sub>22</sub> H <sub>24</sub> N <sub>4</sub> O                             | DYPZVGXELGJCHD-UHFFFAOYSA-N |
| 110 | 5F-MDMB-P4AICA                    | methyl 2-{{1-(5-fluoropentyl)pyrrolo[3,2- <i>b</i> ]pyridine-3-carbonyl}amino}-3,3-dimethylbutanoate                    | C <sub>20</sub> H <sub>28</sub> FN <sub>3</sub> O <sub>3</sub>               | YUNKAZHULJISDA-UHFFFAOYSA-N |
| 111 | 5F-MDMB-P7AICA                    | methyl 2-{{1-(5-fluoropentyl)-1 <i>H</i> -pyrrolo[2,3- <i>b</i> ]pyridin-3-yl]formamido}-3,3-dimethylbutanoate          | C <sub>20</sub> H <sub>28</sub> FN <sub>3</sub> O <sub>3</sub>               | LIRBKFHBIDESHO-UHFFFAOYSA-N |

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|-----|--|---|--|-------------------------------|
| 112 | 5F-AB-P7AICA   | <i>N</i> -(1-amino-3-methyl-1-oxobutan-2-yl)-1-(5-fluoropentyl)-1 <i>H</i> -pyrrolo[2,3- <i>b</i> ]pyridine-3-carboxamide | C <sub>18</sub> H <sub>25</sub> FN <sub>4</sub> O <sub>2</sub>                 | PEAJNPAIVYHNDQ-UHFFFAOYSA-N   |
| 113 | 2F-QMPSB   | quinolin-8-yl 3-[(4,4-difluoropiperidin-1-yl)sulfonyl]-4-methylbenzoate   | C <sub>22</sub> H <sub>20</sub> F <sub>2</sub> N <sub>2</sub> O <sub>4</sub> S | JOSWCKYCXJMLNM-UHFFFAOYSA-N   |
| 114 | 5F-A-P7AICA  | <i>N</i> -(adamantan-1-yl)-1-(5-fluoropentyl)-1 <i>H</i> -pyrrolo[2,3- <i>b</i> ]pyridine-3-carboxamide                   | C <sub>23</sub> H <sub>30</sub> FN <sub>3</sub> O                              | JYMCNFJLBFSGH-UHFFFAOYSA-N    |
| 115 | 3',4'-Methylenedioxy-alpha-methylPPP (MDMPP)                                     | 1-(1,3-benzodioxol-5-yl)-2-methyl-2-(pyrrolidin-1-yl)propan-1-one   | C <sub>15</sub> H <sub>19</sub> NO <sub>3</sub>                                | QEVPOSURBZYOOK-UHFFFAOYSA-N   |
| 116 | 1-(2,3-dihydro-1 <i>H</i> -inden-5-yl)-2-phenyl-2-(pyrrolidinyl-1-yl)ethan-1-one | 1-(2,3-dihydro-1 <i>H</i> -inden-5-yl)-2-phenyl-2-(pyrrolidin-1-yl)ethanone   | C <sub>21</sub> H <sub>23</sub> NO   | JFDZZTSFBTYDF-UHFFFAOYSA-N    |
| 117 | alpha-PPP-MeO  | 3-methoxy-1-phenyl-2-(pyrrolidin-1-yl)propan-1-one  | C <sub>14</sub> H <sub>19</sub> NO <sub>2</sub>                                | TVBNCCGCIPGJQO-UHFFFAOYSA-N   |
| 118 | 3,4-Dichloro- <i>N,N</i> -cyclohexylmethylmethcathinone                          | 2-[cyclohexyl(methyl)amino]-1-(3,4-dichlorophenyl)propan-1-one  | C <sub>16</sub> H <sub>21</sub> Cl <sub>2</sub> NO                             | YOJYFWVGILSSIW-UHFFFAOYSA-N   |
| 119 | TH-PBP   | 2-(pyrrolidin-1-yl)-1-(5,6,7,8-tetrahydronaphthalen-2-yl)butan-1-one  | C <sub>18</sub> H <sub>25</sub> NO   | DNKYSKGGKPOHTKY-UHFFFAOYSA-N  |
| 120 | alpha-pyrrolidinocyclohexylphenone (alpha-PCYP)                                  | 2-cyclohexyl-1-phenyl-2-(pyrrolidin-1-yl)ethan-1-one  | C <sub>18</sub> H <sub>25</sub> NO   | FKEHRWJWTWDTDB-UHFFFAOYSA-N   |
| 121 | 2-Me-DMT   | <i>N,N</i> -dimethyl-2-(2-methyl-1 <i>H</i> -indol-3-yl)ethanamine  | C <sub>13</sub> H <sub>18</sub> N <sub>2</sub>                                 | NDGCOWDSLVLNGLGE-UHFFFAOYSA-N |
| 122 | AL-LAD   | (8β)-9,10-didehydro- <i>N,N</i> -diethyl-6-(2-propenyl)-ergoline-8-carboxamide  | C <sub>22</sub> H <sub>27</sub> N <sub>3</sub> O                               | JCQLEPDZFXGHHQ-OXQOHEQNSA-N   |
| 123 | McPT   | <i>N</i> -[2-(1 <i>H</i> -indol-3-yl)ethyl]- <i>N</i> -methylcyclopropanamine   | C <sub>14</sub> H <sub>18</sub> N <sub>2</sub>                                 | LVOBNSVSQLXACL-UHFFFAOYSA-N   |
| 124 | alpha-TMT  | 1-(1 <i>H</i> -indol-3-yl)- <i>N,N</i> -dimethylpropan-2-amine  | C <sub>13</sub> H <sub>18</sub> N <sub>2</sub>                                 | XQFCCTPWINMCQJ-UHFFFAOYSA-N   |
| 125 | 1P-ETH-LAD   | <i>N,N</i> ,7-triethyl-4-propanoyl-6,6a,8,9-tetrahydroindolo[4,3- <i>fg</i> ]quinoline-9-carboxamide                      | C <sub>24</sub> H <sub>31</sub> N <sub>3</sub> O <sub>2</sub>                  | MLOFCBXSOAYCIF-UHFFFAOYSA-N   |
| 126 | 5-MeO-pyr-T  | 5-methoxy-3-(2-pyrrolidin-1-ylethyl)-1 <i>H</i> -indole   | C <sub>15</sub> H <sub>20</sub> N <sub>2</sub> O                               | KAASYKNZNPWPQG-               |

## DRAFT

|     |                                 |   |  |                              |
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|     |                                 |   |  | UHFFFAOYSA-N                 |
| 127 | Lysergic acid methyl ester      | methyl 7-methyl-6,6a,8,9-tetrahydro-4 <i>H</i> -indolo[4,3- <i>fg</i> ]quinoline-9-carboxylate          | C <sub>17</sub> H <sub>18</sub> N <sub>2</sub> O <sub>2</sub>    | RNHDWLRHUIJZABX-UHFFFAOYSA-N |
| 128 | 4-HO-McPT                       | 3-{2-[cyclopropyl(methyl)amino]ethyl}-1 <i>H</i> -indol-4-ol  | C <sub>14</sub> H <sub>18</sub> N <sub>2</sub> O                 | GFVJBFIXZYL VPO-UHFFFAOYSA-N |
| 129 | 4-PrO-DMT                       | 3-[2-(dimethylamino)ethyl]-1 <i>H</i> -indol-4-yl propanoate  | C <sub>15</sub> H <sub>20</sub> N <sub>2</sub> O <sub>2</sub>    | KUOGXPDQORRHED-UHFFFAOYSA-N  |
| 130 | Butorphanol                     | 17-(cyclobutylmethyl)morphinan-3,14-diol  | C <sub>21</sub> H <sub>29</sub> NO <sub>2</sub>                  | IFKLAQQSCNHLHL-QHAWAJNXSA-N  |
| 131 | U-49900                         | 3,4-dichloro- <i>N</i> -[2-(diethylamino)cyclohexyl]- <i>N</i> -methylbenzamide                         | C <sub>18</sub> H <sub>26</sub> Cl <sub>2</sub> N <sub>2</sub> O | AXACJBKFKCCIOR-UHFFFAOYSA-N  |
| 132 | U-51754                         | 2-(3,4-dichlorophenyl)- <i>N</i> -[2-(dimethylamino)cyclohexyl]- <i>N</i> -methyl-acetamide             | C <sub>17</sub> H <sub>24</sub> Cl <sub>2</sub> N <sub>2</sub> O | ISJUYFBACBKWBV-UHFFFAOYSA-N  |
| 133 | Benzodioxole-fentanyl           | <i>N</i> -phenyl- <i>N</i> -[1-(2-phenylethyl)-4-piperidyl]-1,3-benzodioxole-5-carboxamide              | C <sub>27</sub> H <sub>28</sub> N <sub>2</sub> O <sub>3</sub>    | ZFAAZMIOHJNKGD-UHFFFAOYSA-N  |
| 134 | 3-phenylpropanoylfentanyl       | <i>N</i> ,3-diphenyl- <i>N</i> -[1-(2-phenylethyl)-4-piperidyl]propanamide                              | C <sub>28</sub> H <sub>32</sub> N <sub>2</sub> O                 | DIRAGWDYMRIDIO-UHFFFAOYSA-N  |
| 135 | Tetramethylcyclopropanefentanyl | 2,2,3,3-tetramethyl- <i>N</i> -phenyl- <i>N</i> -[1-(2-phenylethyl)-4-piperidyl]cyclopropanecarboxamide | C <sub>27</sub> H <sub>36</sub> N <sub>2</sub> O                 | BYCDHAVFKDTVAM-UHFFFAOYSA-N  |
| 136 | U-48800                         | 2-(2,4-dichlorophenyl)- <i>N</i> -[2-(dimethylamino)cyclohexyl]- <i>N</i> -methyl-acetamide             | C <sub>17</sub> H <sub>24</sub> Cl <sub>2</sub> N <sub>2</sub> O | FKUWIGXXBMULOI-UHFFFAOYSA-N  |
| 137 | Thiophenefentanyl               | <i>N</i> -phenyl- <i>N</i> -[1-(2-phenylethyl)-4-piperidyl]thiophene-2-carboxamide                      | C <sub>24</sub> H <sub>26</sub> N <sub>2</sub> OS                | CCHPKGYUIHSQIE-UHFFFAOYSA-N  |
| 138 | Benzylfentanyl                  | <i>N</i> -(1-benzyl-4-piperidyl)- <i>N</i> -phenyl-propanamide  | C <sub>21</sub> H <sub>26</sub> N <sub>2</sub> O                 | POQDXIFVWVZVML-UHFFFAOYSA-N  |
| 139 | Bromadoline                     | 4-bromo- <i>N</i> -[2-(dimethylamino)cyclohexyl]benzamide   | C <sub>15</sub> H <sub>21</sub> BrN <sub>2</sub> O               | UFDJFJYMMIZKLG-UHFFFAOYSA-N  |
| 140 | Acetylbenzylfentanyl            | <i>N</i> -(1-benzyl-4-piperidyl)- <i>N</i> -phenyl-acetamide  | C <sub>20</sub> H <sub>24</sub> N <sub>2</sub> O                 | UKGXYSOSRSCGJB-UHFFFAOYSA-N  |
| 141 | Benzoylbenzylfentanyl           | <i>N</i> -(1-benzyl-4-piperidyl)- <i>N</i> -phenyl-benzamide  | C <sub>25</sub> H <sub>26</sub> N <sub>2</sub> O                 | TVPYSIMEYHFHLN-UHFFFAOYSA-N  |



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|-----|------------------------------------|---|--|-----------------------------|
| 142 | 4-hydroxybutyrfentanyl; 4-HO-BF    | <i>N</i> -(4-hydroxyphenyl)- <i>N</i> -[1-(2-phenylethyl)piperidin-4-yl]butanamide          | C <sub>23</sub> H <sub>30</sub> N <sub>2</sub> O <sub>2</sub>                | HMPNQEXOJZKKKE-UHFFFAOYSA-N |
| 143 | isopropyl-U-47700                  | 3,4-dichloro- <i>N</i> -[2-(dimethylamino)cyclohexyl]- <i>N</i> -(propan-2-yl)benzamide     | C <sub>18</sub> H <sub>26</sub> Cl <sub>2</sub> N <sub>2</sub> O             | LGYSYWASFBSHJ-UHFFFAOYSA-N  |
| 144 | U-50488                            | 2-(3,4-dichlorophenyl)- <i>N</i> -methyl- <i>N</i> -(2-pyrrolidin-1-yl)cyclohexyl)acetamide | C <sub>19</sub> H <sub>26</sub> Cl <sub>2</sub> N <sub>2</sub> O             | VQLPLYSROCPWFF-UHFFFAOYSA-N |
| 145 | 3,4-methylenedioxy-U-47700         | <i>N</i> -[2-(dimethylamino)cyclohexyl]- <i>N</i> -methyl-1,3-benzodioxole-5-carboxamide    | C <sub>17</sub> H <sub>24</sub> N <sub>2</sub> O <sub>3</sub>                | UUAVKYBZWVMWSM-UHFFFAOYSA-N |
| 146 | 4-fluoro-cyclopropylbenzylfentanyl | <i>N</i> -(1-benzyl-4-piperidyl)- <i>N</i> -(4-fluorophenyl)cyclopropanecarboxamide         | C <sub>22</sub> H <sub>25</sub> FN <sub>2</sub> O                            | DGFCHXNCEJHFDB-UHFFFAOYSA-N |
| 147 | Furanylbenzylfentanyl              | <i>N</i> -(1-benzyl-4-piperidyl)- <i>N</i> -phenyl-furan-2-carboxamide                      | C <sub>23</sub> H <sub>24</sub> N <sub>2</sub> O <sub>2</sub>                | GDPJXHFCUEJBT-UHFFFAOYSA-N  |
| 148 | 2-methylacetylfentanyl             | <i>N</i> -(2-methylphenyl)- <i>N</i> -[1-(2-phenylethyl)-4-piperidinyl]-acetamide           | C <sub>22</sub> H <sub>28</sub> N <sub>2</sub> O                             | GRDWUDZBHWLHSH-UHFFFAOYSA-N |
| 149 | <i>N</i> -methyl U-47931E          | 4-bromo- <i>N</i> -[2-(dimethylamino)cyclohexyl]- <i>N</i> -methylbenzamide                 | C <sub>16</sub> H <sub>23</sub> BrN <sub>2</sub> O                           | XQCGUPNNNXRBDG-UHFFFAOYSA-N |
| 150 | Piperidylthiambutene               | 1-[4,4-di(thiophen-2-yl)but-3-en-2-yl]piperidine  | C <sub>17</sub> H <sub>21</sub> NS <sub>2</sub>                              | FQRWJLVMJCKSME-UHFFFAOYSA-N |
| 151 | 2-methyl-AP-237                    | 1-[2-methyl-4-(3-phenylprop-2-en-1-yl)piperazin-1-yl]butan-1-one                            | C <sub>18</sub> H <sub>26</sub> N <sub>2</sub> O                             | CRSFXYZFNAFVFC-UHFFFAOYSA-N |
| 152 | AP-237                             | 1-[4-(3-phenylprop-2-en-1-yl)piperazin-1-yl]butan-1-one                                     | C <sub>17</sub> H <sub>24</sub> N <sub>2</sub> O                             | ZQBMUHABRSEAIK-UHFFFAOYSA-N |
| 153 | Furanyl UF-17                      | <i>N</i> -[2-(dimethylamino)cyclohexyl]- <i>N</i> -phenyl-furan-2-carboxamide               | C <sub>19</sub> H <sub>24</sub> N <sub>2</sub> O <sub>2</sub>                | OSFBKUBOZUWERW-UHFFFAOYSA-N |
| 154 | 2F-viminol                         | 2-[di(butan-2-yl)amino]-1-[1-(2-fluorobenzyl)-1 <i>H</i> -pyrrol-2-yl]ethan-1-ol            | C <sub>21</sub> H <sub>31</sub> FN <sub>2</sub> O                            | FXQMNNHPOYWWKI-UHFFFAOYSA-N |
| 155 | Iso-ethcathinone                   | 1-(ethylamino)-1-phenylpropan-2-one   | C <sub>11</sub> H <sub>15</sub> NO   | GHVCNLDNRVGRJP-UHFFFAOYSA-N |
| 156 | 5-HTP                              | 2-amino-3-(5-hydroxy-1 <i>H</i> -indol-3-yl)propanoic acid                                  | C <sub>11</sub> H <sub>12</sub> N <sub>2</sub> O <sub>3</sub>                | LDCYZAJDBXYCGN-UHFFFAOYSA-N |
| 157 | Ostarine                           | 3-(4-cyanophenoxy)- <i>N</i> -[4-cyano-3-(trifluoromethyl)phenyl]-2-hydroxy-2-              | C <sub>19</sub> H <sub>14</sub> F <sub>3</sub> N <sub>3</sub> O <sub>3</sub> | JNGVJMBLXIUVRD-UHFFFAOYSA-N |

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|-----|---|--|---|-----------------------------|
|     |   | methylpropanamide  |   |                             |
| 158 | 4-Fluoroephedrine                         | 4-fluoro- $\alpha$ -[1-(methylamino)ethyl]-benzenemethanol   | C <sub>10</sub> H <sub>14</sub> FNO                               | SPEQHEOLWDGWML-UHFFFAOYSA-N |
| 159 | 4-methylphendimetrazine                   | 3,4-dimethyl-2-(4-methylphenyl)morpholine  | C <sub>13</sub> H <sub>19</sub> NO                                | UJDFQROPBYDNBJ-UHFFFAOYSA-N |
| 160 | Mebroqualone                              | 3-(2-bromophenyl)-2-methylquinazolin-4(3 <i>H</i> )-one  | C <sub>15</sub> H <sub>11</sub> BrN <sub>2</sub> O                | NBUSAPJNASSKBP-UHFFFAOYSA-N |
| 161 | W-15                                      | 4-chloro- <i>N</i> -[(2 <i>E</i> )-1-(2-phenylethyl)piperidin-2-ylidene]benzenesulfonamide   | C <sub>19</sub> H <sub>21</sub> ClN <sub>2</sub> O <sub>2</sub> S | VJHXSSVOCOBVMI-XUTLUUPISA-N |
| 162 | Sibutramine                               | 1-[1-(4-chlorophenyl)cyclobutyl]- <i>N,N</i> ,3-trimethyl-butan-1-amine  | C <sub>17</sub> H <sub>26</sub> ClN                               | UNAANXDKBXWMLN-UHFFFAOYSA-N |
| 163 | Embutramide                               | <i>N</i> -[2-ethyl-2-(3-methoxyphenyl)butyl]-4-hydroxybutanamide   | C <sub>17</sub> H <sub>27</sub> NO <sub>3</sub>                   | LMBMDLOSPKIWAP-UHFFFAOYSA-N |
| 164 | Lysergic acid 2,4-dimethylazetidide (LSZ) | [(6 <i>aS</i> ,9 <i>S</i> )-7-methyl-6,6 <i>a</i> ,8,9-tetrahydro-4 <i>H</i> -indolo[4,3- <i>f,g</i> ]quinoline-9-yl]-[(2 <i>S</i> ,4 <i>S</i> )-2,4-dimethylazetidid-1-yl]methanone | C <sub>21</sub> H <sub>25</sub> N <sub>3</sub> O                  | DUKNIHFTDAXJON-UHFFFAOYSA-N |
| 165 | Noopept                                   | ethyl ({[(2 <i>S</i> )-1-(phenylacetyl)pyrrolidin-2-yl]carbonyl}amino)acetate  | C <sub>17</sub> H <sub>22</sub> N <sub>2</sub> O <sub>4</sub>     | PJNSMUBMSNAEEN-AWEZNQCLSA-N |
| 166 | Mesembrine                                | 3 <i>a</i> -(3,4-dimethoxyphenyl)-1-methyl-2,3,4,5,7,7 <i>a</i> -hexahydroindol-6-one  | C <sub>17</sub> H <sub>23</sub> NO <sub>3</sub>                   | DAHIQPJTGIHDGO-UHFFFAOYSA-N |
| 167 | Orphenadrine                              | <i>N,N</i> -dimethyl-2-[(2-methylphenyl)(phenyl)methoxy]ethanamine   | C <sub>18</sub> H <sub>23</sub> NO                                | QVYRGXJSLMXQH-UHFFFAOYSA-N  |
| 168 | Ephedrine (NEDPA)                         | <i>N</i> -ethyl-1,2-diphenylethanamine   | C <sub>16</sub> H <sub>19</sub> N                                 | IGFZMQXEKIZPDR-UHFFFAOYSA-N |
| 169 | NPDPA                                     | <i>N</i> -(1,2-diphenylethyl)propan-2-amine  | C <sub>17</sub> H <sub>21</sub> N                                 | FBRJTEBLJRHAQX-UHFFFAOYSA-N |
| 170 | W-18                                      | 4-chloro- <i>N</i> -{(2 <i>E</i> )-1-[2-(4-nitrophenyl)ethyl]piperidin-2-ylidene}benzenesulfonamide  | C <sub>19</sub> H <sub>20</sub> ClN <sub>3</sub> O <sub>4</sub> S | BKRSVROQVRTSND-XUTLUUPISA-N |
| 171 | Adrafinil                                 | 2-[(diphenylmethyl)sulfinyl]- <i>N</i> -hydroxyacetamide   | C <sub>15</sub> H <sub>15</sub> NO <sub>3</sub> S                 | CGNMLOKEMNBUI-UHFFFAOYSA-N  |
| 172 | Afloqualone                               | 6-amino-2-(fluoromethyl)-3-(2-   | C <sub>16</sub> H <sub>14</sub> FN <sub>3</sub> O                 | VDOSWXIDETXFET-             |

|     |                                      |   |  |                              |
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|     |                                      | methylphenyl)quinazolin-4(3 <i>H</i> )-one  |  | UHFFFAOYSA-N                 |
| 173 | Modafienz                            | 2-[[bis(4-fluorophenyl)methyl]sulfinyl]- <i>N</i> -methylacetamide                                  | C <sub>16</sub> H <sub>15</sub> F <sub>2</sub> NO <sub>2</sub> S | MQZWTCIUUSDFCQ-UHFFFAOYSA-N  |
| 174 | Methylmethaqualone                   | 3-(2,4-dimethylphenyl)-2-methylquinazolin-4(3 <i>H</i> )-one  | C <sub>17</sub> H <sub>16</sub> N <sub>2</sub> O                 | MPMDMUROZIYIIM-UHFFFAOYSA-N  |
| 175 | Flibanserin                          | 1,3-dihydro-1-(2-{4-[3-(trifluoromethyl)phenyl]-1-piperazinyl}ethyl)-2 <i>H</i> -benzimidazol-2-one | C <sub>20</sub> H <sub>21</sub> F <sub>3</sub> N <sub>4</sub> O  | PPRRDFIXUUSXRA-UHFFFAOYSA-N  |
| 176 | 4-methylpentan-2-amine (DMBA)        | 4-methylpentan-2-amine  | C <sub>6</sub> H <sub>15</sub> N                                 | UNBMPKNTYKDYCG-UHFFFAOYSA-N  |
| 177 | Modafinil sulphone                   | 2-[(diphenylmethyl)sulfonyl]acetamide   | C <sub>15</sub> H <sub>15</sub> NO <sub>3</sub> S                | ZESNOWZYHYRSRY-UHFFFAOYSA-N  |
| 178 | <i>N</i> -methyl aminorex derivative | 3-methyl-5-phenyl-oxazolidin-2-imine  | C <sub>10</sub> H <sub>12</sub> N <sub>2</sub> O                 | PINRUEQFGKWBTO-UHFFFAOYSA-N  |
| 179 | 3,4-DMAR                             | 3,4-dimethyl-5-phenyl-1,3-oxazolidin-2-imine  | C <sub>11</sub> H <sub>14</sub> N <sub>2</sub> O                 | AINPTMZQNYUGDW-UHFFFAOYSA-N  |
| 180 | 4-MPH                                | 3-methyl-2-(4-methylphenyl)morpholine   | C <sub>12</sub> H <sub>17</sub> NO                               | NWNCIXFIIDVRKE-UHFFFAOYSA-N  |
| 181 | Phenetrazine                         | 3-ethyl-2-phenylmorpholine  | C <sub>12</sub> H <sub>17</sub> NO                               | DOMAVIHZFHQQHF-UHFFFAOYSA-N  |
| 182 | Epirocaine                           | 2-methyl-2-(propylamino)propyl benzoate   | C <sub>14</sub> H <sub>21</sub> NO <sub>2</sub>                  | VXJABHHJLXLNMP-UHFFFAOYSA-N  |
| 183 | Modafinil                            | 2-[(diphenylmethyl)sulfinyl]acetamide   | C <sub>15</sub> H <sub>15</sub> NO <sub>2</sub> S                | YFGHCGITMMYXAQ-UHFFFAOYSA-N  |
| 184 | Iso-phenmetrazine                    | 5-methyl-2-phenylmorpholine   | C <sub>11</sub> H <sub>15</sub> NO                               | LQHGEIOIBMBXJGV-UHFFFAOYSA-N |
| 185 | Phenmetetrazine                      | 4-ethyl-3-methyl-2-phenylmorpholine   | C <sub>13</sub> H <sub>19</sub> NO                               | ZOMTUQAEMGCZPK-UHFFFAOYSA-N  |
| 186 | 3F-Phenetrazine                      | 3-ethyl-2-(3-fluorophenyl)morpholine  | C <sub>12</sub> H <sub>16</sub> FNO                              | GPTHUZORYIVVNJ-UHFFFAOYSA-N  |
| 187 | Bromantane                           | <i>N</i> -(4-bromophenyl)tricyclo[3.3.1.1 <sup>3,7</sup> ]decan-2-amine                             | C <sub>16</sub> H <sub>20</sub> BrN                              | LWJALJDRFBXHKX-UHFFFAOYSA-N  |
| 188 | Viloxazine                           | 2-[(2-ethoxyphenoxy)methyl]morpholine   | C <sub>13</sub> H <sub>19</sub> NO <sub>3</sub>                  | YWPHCCPCQOJSGZ-              |

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|     |  |  |  | UHFFFAOYSA-N                 |
| 189 | Fladrafinil                            | 2-[[bis(4-fluorophenyl)methyl]sulfinyl]- <i>N</i> -hydroxyacetamide  | C <sub>15</sub> H <sub>13</sub> F <sub>2</sub> NO <sub>3</sub> S | VKGUUSVYPXTWMA-UHFFFAOYSA-N  |
| 190 | PDM-35                                 | 3,5-dimethyl-2-phenylmorpholine  | C <sub>12</sub> H <sub>17</sub> NO                               | YKCSYIYQRSVLAK-UHFFFAOYSA-N  |
| 191 | 3,6-DMPM                               | 3,6-dimethyl-2-phenylmorpholine  | C <sub>12</sub> H <sub>17</sub> NO                               | FZEIVUHEODGHML-UHFFFAOYSA-N  |
| 192 | 3-methylphenmetrazine                  | 3-methyl-2-(3-methylphenyl)morpholine  | C <sub>12</sub> H <sub>17</sub> NO                               | QEDQZYNGDXULGO-UHFFFAOYSA-N  |
| 193 | G-130                                  | 5,5-dimethyl-2-phenyl-morpholine   | C <sub>12</sub> H <sub>17</sub> NO                               | KJUOROGOOZJYAI-UHFFFAOYSA-N  |
| 194 | Methylmorphenate                       | methyl 2-(morpholin-3-yl)-2-phenylacetate  | C <sub>13</sub> H <sub>17</sub> NO <sub>3</sub>                  | FTSNQYGFVFKHY-UHFFFAOYSA-N   |
| 195 | PRE-084                                | 2-(morpholin-4-yl)ethyl 1-phenylcyclohexane-1-carboxylate  | C <sub>19</sub> H <sub>27</sub> NO <sub>3</sub>                  | RQHKZUBCUZVZEF-UHFFFAOYSA-N  |
| 196 | NDTDI                                  | <i>N,N</i> -diethyl-3-{methyl-[(4 <i>R</i> )-1,3,4,5-tetrahydrobenzo[ <i>c,d</i> ]indol-4-yl]amino}propanamide | C <sub>19</sub> H <sub>27</sub> N <sub>3</sub> O                 | JECGWOMOCQPQHDH-UHFFFAOYSA-N |
| 197 | Ru-28306                               | <i>N,N</i> -dimethyl-1,3,4,5-tetrahydrobenzo[ <i>c,d</i> ]indol-4-amine  | C <sub>13</sub> H <sub>16</sub> N <sub>2</sub>                   | BQOANWOQEHVATQ-UHFFFAOYSA-N  |
| 198 | Octodrine                              | 6-methylheptan-2-amine   | C <sub>8</sub> H <sub>19</sub> N                                 | QNIVIMYXGGFTAK-UHFFFAOYSA-N  |
| 199 | 1,4-DMAA                               | 5-methylhexan-2-amine  | C <sub>7</sub> H <sub>17</sub> N                                 | IZCBXLKODYZSDJ-UHFFFAOYSA-N  |
| 200 | 2F-Phenmetrazine                       | 2-(2-fluorophenyl)-3-methylmorpholine  | C <sub>11</sub> H <sub>14</sub> FNO                              | QTYLEXQVLJYJHT-UHFFFAOYSA-N  |
| 201 | Troparil                               | methyl 8-methyl-3-phenyl-8-azabicyclo[3.2.1]octane-4-carboxylate   | C <sub>16</sub> H <sub>21</sub> NO <sub>2</sub>                  | OMBOXYLBHWNWHL-UHFFFAOYSA-N  |
| 202 | <i>N</i> -methyl-cyclazodone           | 2-[cyclopropyl(methyl)amino]-5-phenyl-4(5 <i>H</i> )-oxazolone   | C <sub>13</sub> H <sub>14</sub> N <sub>2</sub> O <sub>2</sub>    | FFWGGFGJVZVGOW-UHFFFAOYSA-N  |
| 203 | para-fluoro-4-methylaminorex;<br>4-FPO | 5-(4-fluorophenyl)-4-methyl-4,5-dihydro-1,3-oxazol-2-amine   | C <sub>10</sub> H <sub>11</sub> FN <sub>2</sub> O                | UYKYWISHPDEDQRQ-UHFFFAOYSA-N |

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|     |                         |   |  |                                  |
|-----|-------------------------|---|--|----------------------------------|
| 204 | Bisfluoromodafinil      | 2-{{bis(4-fluorophenyl)methyl}sulfinyl}acetamide  | C <sub>15</sub> H <sub>13</sub> F <sub>2</sub> NO <sub>2</sub> S | YEAQNUMCWMRYM<br>U-UHFFFAOYSA-N  |
| 205 | WIN 35428               | methyl 3-(4-fluorophenyl)-8-methyl-8-azabicyclo[3.2.1]octane-2-carboxylate  | C <sub>16</sub> H <sub>20</sub> FNO <sub>2</sub>                 | QUSLQENMLDRCTO-<br>UHFFFAOYSA-N  |
| 206 | alpha-methylephedrine   | 3-(methylamino)-2-phenylbutan-2-ol  | C <sub>11</sub> H <sub>17</sub> NO                               | SIBUOXVRODUTKZ-<br>UHFFFAOYSA-N  |
| 207 | N-methylephedrine       | 2-(dimethylamino)-1-phenylpropan-1-ol   | C <sub>11</sub> H <sub>17</sub> NO                               | FMCGSUUBYTWNDP-<br>UHFFFAOYSA-N  |
| 208 | 1B-LSD                  | 4-butyryl- <i>N,N</i> -diethyl-7-methyl-4,6,6a,7,8,9-hexahydroindolo[4,3- <i>fg</i> ]quinoline-9-carboxamide                | C <sub>24</sub> H <sub>31</sub> N <sub>3</sub> O <sub>2</sub>    | SVRFNPSJPIDUBC-<br>UHFFFAOYSA-N  |
| 209 | 2C-B aminorex           | 5-(4-bromo-2,5-dimethoxy-phenyl)-4,5-dihydrooxazol-2-amine  | C <sub>11</sub> H <sub>13</sub> BrN <sub>2</sub> O <sub>3</sub>  | XUTCHZHTIDHQOP-<br>UHFFFAOYSA-N  |
| 210 | Pagoclone               | 2-(7-chloro-1,8-naphthyridin-2-yl)-2,3-dihydro-3-(5-methyl-2-oxohexyl)-1 <i>H</i> -isoindol-1-one                           | C <sub>23</sub> H <sub>22</sub> ClN <sub>3</sub> O <sub>2</sub>  | HIUPRQPBWVEQJJ-<br>UHFFFAOYSA-N  |
| 211 | 4-fluorophenibut        | 4-amino-3-(4-fluorophenyl)butanoic acid   | C <sub>10</sub> H <sub>12</sub> FNO <sub>2</sub>                 | QWHXHLDNSXLAPX-<br>UHFFFAOYSA-N  |
| 212 | Pregabalin methyl ester | methyl 3-(aminomethyl)-5-methylhexanoate  | C <sub>9</sub> H <sub>19</sub> NO <sub>2</sub>                   | JDFPGFRTUSUWNI-<br>UHFFFAOYSA-N  |
| 213 | Xylazine                | <i>N</i> -(2,6-dimethylphenyl)-5,6-dihydro-4 <i>H</i> -1,3-thiazin-2-amine  | C <sub>12</sub> H <sub>16</sub> N <sub>2</sub> S                 | BPICBUSOMSTKRF-<br>UHFFFAOYSA-N  |
| 214 | SL-164                  | 5-chloro-3-(4-chloro-2-methylphenyl)-2-methyl-4(3 <i>H</i> )-quinazolinone  | C <sub>16</sub> H <sub>12</sub> Cl <sub>2</sub> N <sub>2</sub> O | KUIHLOHNUGOCTO-<br>UHFFFAOYSA-N  |
| 215 | 1cP-LSD                 | 4-(cyclopropanecarbonyl)- <i>N,N</i> -diethyl-7-methyl-4,6,6a,7,8,9-hexahydroindolo[4,3- <i>fg</i> ]quinoline-9-carboxamide | C <sub>24</sub> H <sub>29</sub> N <sub>3</sub> O <sub>2</sub>    | RAFUPYYDHPFASC-<br>DYESRHJHSA-N  |
| 216 | Nitromethaqualone       | 3-(2-methoxy-4-nitrophenyl)-2-methylquinazolin-4-one  | C <sub>16</sub> H <sub>13</sub> N <sub>3</sub> O <sub>4</sub>    | RZHHDMJWDYJXAW-<br>UHFFFAOYSA-N  |
| 217 | Nefiracetam             | <i>N</i> -(2,6-dimethylphenyl)-2-(2-oxopyrrolidin-1-yl)acetamide  | C <sub>14</sub> H <sub>18</sub> N <sub>2</sub> O <sub>2</sub>    | NGHTXZCKLWZPGK-<br>UHFFFAOYSA-N  |
| 218 | 4-Hydroxyamphetamine    | 4-(2-aminopropyl)phenol   | C <sub>9</sub> H <sub>13</sub> NO                                | GIKNHHRFLCDOEU -<br>UHFFFAOYSA-N |
| 219 | 25I-NBMD                | <i>N</i> -(1,3-benzodioxol-4-ylmethyl)-2-(4-iodo-2,5-   | C <sub>18</sub> H <sub>20</sub> INO <sub>4</sub>                 | NJNMIPDEUMTYNV -                 |

## DRAFT

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|-----|---------------------------------|--|--|------------------------------|
|     |                                 | dimethoxyphenyl)ethanamine   |  | UHFFFAOYSA-N                 |
| 220 | 25I-NB34MD                      | <i>N</i> -(1,3-benzodioxol-5-ylmethyl)-2-(4-iodo-2,5-dimethoxyphenyl)ethanamine                    | C <sub>18</sub> H <sub>20</sub> INO <sub>4</sub>                             | FWEBGKDUEZRMRQ -UHFFFAOYSA-N |
| 221 | 25C-NBF                         | 2-(4-chloro-2,5-dimethoxyphenyl)- <i>N</i> -(2-fluorobenzyl)ethanamine                             | C <sub>17</sub> H <sub>19</sub> ClFNO <sub>2</sub>                           | AHIUIEOLKNDLSC -UHFFFAOYSA-N |
| 222 | 25C-NBOH                        | 2-({[2-(4-chloro-2,5-dimethoxyphenyl)ethyl]amino}methyl)phenol                                     | C <sub>17</sub> H <sub>20</sub> ClNO <sub>3</sub>                            | VHWXICYQMMZCW -UHFFFAOYSA-N  |
| 223 | 25I-NBOH                        | 2-{{[2-(4-iodo-2,5-dimethoxyphenyl)ethylamino]methyl}phenol  | C <sub>17</sub> H <sub>20</sub> INO <sub>3</sub>                             | FEUZHYRXGQTBRO -UHFFFAOYSA-N |
| 224 | 25B-NBF                         | 2-(4-bromo-2,5-dimethoxyphenyl)- <i>N</i> -[(2-fluorophenyl)methyl]ethanamine                      | C <sub>17</sub> H <sub>19</sub> BrFNO <sub>2</sub>                           | ATMBBMXJNJRST -UHFFFAOYSA-N  |
| 225 | 25B-NBOH                        | 2-{{[2-(4-bromo-2,5-dimethoxyphenyl)ethylamino]methyl}phenol                                       | C <sub>17</sub> H <sub>20</sub> BrNO <sub>3</sub>                            | RSUNJYKZRKIBNB -UHFFFAOYSA-N |
| 226 | 1-phenethyl-4-hydroxypiperidine | 1-(2-phenylethyl)piperidin-4-ol  | C <sub>13</sub> H <sub>19</sub> NO   | KYGMSTKBHJVPJK -UHFFFAOYSA-N |
| 227 | 7-CDMeOPPAE                     | 7-(2-{{[2-(4-chloro-2,5-dimethoxyphenyl)-1-methyl-ethyl]amino}ethyl}-1,3-dimethyl-purine-2,6-dione | C <sub>20</sub> H <sub>26</sub> ClN <sub>5</sub> O <sub>4</sub>              | CWRNNANKXUBLFV -UHFFFAOYSA-N |
| 228 | 25E-NBOH                        | 2-({[2-(4-ethyl-2,5-dimethoxyphenyl)ethyl]amino}methyl)phenol                                      | C <sub>19</sub> H <sub>25</sub> NO <sub>3</sub>                              | SYBINTRPEZWFLZ -UHFFFAOYSA-N |
| 229 | BOD                             | 2-(2,5-dimethoxy-4-methylphenyl)-2-methoxyethan-1-amine  | C <sub>12</sub> H <sub>19</sub> NO <sub>3</sub>                              | VTEIFHQZWBABDE -UHFFFAOYSA-N |
| 230 | Vanoxerine                      | 1-{2-[bis(4-fluorophenyl)methoxy]ethyl}-4-(3-phenylpropyl)piperazine                               | C <sub>28</sub> H <sub>32</sub> F <sub>2</sub> N <sub>2</sub> O              | NAUWTFJOPJWYOT -UHFFFAOYSA-N |
| 231 | 1-(3-Methylbenzyl)piperazine    | 1-(3-methylbenzyl)piperazine   | C <sub>12</sub> H <sub>18</sub> N <sub>2</sub>                               | VTEOTZPEMDQENX -UHFFFAOYSA-N |
| 232 | NSI-189                         | (4-benzylpiperazin-1-yl)-[2-(isopentylamino)-3-pyridyl]methanone                                   | C <sub>22</sub> H <sub>30</sub> N <sub>4</sub> O                             | DYTOQURYRYYNOR -UHFFFAOYSA-N |
| 233 | DB-MDBP                         | 1-[(2,2-difluoro-1,3-benzodioxol-5-yl)methyl]piperazine  | C <sub>12</sub> H <sub>14</sub> F <sub>2</sub> N <sub>2</sub> O <sub>2</sub> | MPZINTHEMFDGTH -UHFFFAOYSA-N |
| 234 | 3,4-CFP                         | 1-(3-chloro-4-fluorophenyl)piperazine  | C <sub>10</sub> H <sub>12</sub> ClFN <sub>2</sub>                            | MKXFXPRJCBUTLX -UHFFFAOYSA-N |

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|-----|--|---|--|-----------------------------|
| 235 | pBPP                                       | 1-(4-bromophenyl)piperazine   | C <sub>10</sub> H <sub>13</sub> BrN <sub>2</sub>   | PJHPFAFEJNBIDC-UHFFFAOYSA-N |
| 236 | HDMP-28 (methylnaphthidate)                | methyl naphthalen-2-yl(piperidin-2-yl)acetate   | C <sub>18</sub> H <sub>21</sub> NO <sub>2</sub>    | DNRNSIJBSCBESJ-UHFFFAOYSA-N |
| 237 | Propylphenidate                            | propyl phenyl(piperidin-2-yl)acetate  | C <sub>16</sub> H <sub>23</sub> NO <sub>2</sub>    | PRMWWEANNQSWAR-UHFFFAOYSA-N |
| 238 | 4-fluoroethylphenidate                     | ethyl 2-(4-fluorophenyl)-2-(2-piperidyl)acetate   | C <sub>15</sub> H <sub>20</sub> FNO <sub>2</sub>   | RKXQYWFDJDYSEN-UHFFFAOYSA-N |
| 239 | methyl 2-phenyl-2-(pyrrolidin-1-yl)acetate | methyl 2-phenyl-2-(pyrrolidin-1-yl)acetate  | C <sub>13</sub> H <sub>17</sub> NO <sub>2</sub>    | NVTLFQBLMSATNO-UHFFFAOYSA-N |
| 240 | Glaucine                                   | 1,2,9,10-tetramethoxy-6-methyl-5,6,6a,7-tetrahydro-4 <i>H</i> -dibenzo[ <i>de,g</i> ]quinoline                              | C <sub>21</sub> H <sub>25</sub> NO <sub>4</sub>    | RUZIUYOSRDWYQF-UHFFFAOYSA-N |
| 241 | LSA  | (6 <i>aS</i> ,9 <i>S</i> )-7-methyl-6,6 <i>a</i> ,8,9-tetrahydro-4 <i>H</i> -indolo[4,3- <i>fg</i> ]quinoline-9-carboxamide | C <sub>16</sub> H <sub>17</sub> N <sub>3</sub> O   | GENAHGKEFJLNJB-HZMBPMFUSA-N |
| 242 | Arecoline                                  | methyl 1-methyl-1,2,5,6-tetrahydropyridine-3-carboxylate  | C <sub>8</sub> H <sub>13</sub> NO <sub>2</sub>     | HJJPJSXJAXAIPN-UHFFFAOYSA-N |
| 243 | 5-Br-DMT                                   | [2-(5-bromo-1 <i>H</i> -indol-3-yl)ethyl]dimethylamine  | C <sub>12</sub> H <sub>15</sub> BrN <sub>2</sub>   | ATEYZYQLBQUZJE-UHFFFAOYSA-N |
| 244 | 5-Cl-DMT                                   | [2-(5-chloro-1 <i>H</i> -indol-3-yl)ethyl]dimethylamine   | C <sub>12</sub> H <sub>15</sub> ClN <sub>2</sub>   | LXATUVRMTAHHDX-UHFFFAOYSA-N |
| 245 | Mephedrene                                 | <i>N</i> -methyl-1-(5-methyl-2-thiophenyl)propan-2-amine  | C <sub>9</sub> H <sub>15</sub> NS                  | HZICDJQMPFRKCUHFFFAOYSA-N   |
| 246 | PTI-3                                      | <i>N</i> -({2-[1-(5-fluoropentyl)-1 <i>H</i> -indol-3-yl]-1,3-thiazol-4-yl}methyl)-2-methoxy- <i>N</i> -methylethan-1-amine | C <sub>21</sub> H <sub>28</sub> FN <sub>3</sub> OS | LXQIIHJBHSFWQW-UHFFFAOYSA-N |
| 247 | Cumyl-Cb-MeGaClone                         | 5-(cyclobutylmethyl)-2-(1-methyl-1-phenylethyl)pyrido[4,3- <i>b</i> ]indol-1-one  | C <sub>25</sub> H <sub>26</sub> N <sub>2</sub> O   | VOCGZWPYRQJUMY-UHFFFAOYSA-N |
| 248 | BOH-2C-B                                   | 2-amino-1-(4-bromo-2,5-dimethoxyphenyl)ethanol  | C <sub>10</sub> H <sub>14</sub> BrNO <sub>3</sub>  | PCSKDXWCLQXURQUHFFFAOYSA-N  |
| 249 | <i>N</i> -methylbenzedrone                 | 2-[benzyl(methyl)amino]-1-(4-methylphenyl)propan-1-one  | C <sub>18</sub> H <sub>21</sub> NO                 | UIOKRZKDMRJWFG-UHFFFAOYSA-N |
| 250 | 3F-PCP                                     | 1-[1-(3-fluorophenyl)cyclohexyl]piperidine  | C <sub>17</sub> H <sub>24</sub> FN                 | PFPLGKFWWBXTNP-             |

## DRAFT

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|     |   |   |  | UHFFFAOYSA-N                |
| 251 | 3-Cl-PCP  | 1-[1-(3-chlorophenyl)cyclohexyl]piperidine  | C <sub>17</sub> H <sub>24</sub> ClN                            | HUHBTESMMFLCAN-UHFFFAOYSA-N |
| 252 | A-D2PV (alpha-pyrrolidino-2-phenylacetophenone, alpha-D2PV) | 1,2-diphenyl-2-(pyrrolidin-1-yl)ethan-1-one   | C <sub>18</sub> H <sub>19</sub> NO                             | GQCCTZGWWWUYLS-UHFFFAOYSA-N |
| 253 | 3-Me-PCP  | 1-[1-(3-methylphenyl)cyclohexyl]piperidine  | C <sub>18</sub> H <sub>27</sub> N                              | BMFKUCGCXMDGBK-UHFFFAOYSA-N |
| 254 | MXiPR (methoxisopropamine)                                  | 2-(isopropylamino)-2-(3-methoxyphenyl)-cyclohexanone  | C <sub>16</sub> H <sub>23</sub> NO <sub>2</sub>                | FTQIVDGNXPEKP-UHFFFAOYSA-N  |
| 255 | Cumyl-BC-HpMeGaClone-221                                    | 5-(bicyclo[2.2.1]hept-2-ylmethyl)-2-(2-phenylpropan-2-yl)-2,5-dihydro-1 <i>H</i> -pyrido[4,3- <i>b</i> ]indol-1-one | C <sub>28</sub> H <sub>30</sub> N <sub>2</sub> O               | VFBNNQOHBYEVLU-UHFFFAOYSA-N |
| 256 | 3-Methoxyphenmetrazine                                      | 2-(3-methoxyphenyl)-3-methylmorpholine  | C <sub>12</sub> H <sub>17</sub> NO <sub>2</sub>                | QKAKYFBKVKSLLT-UHFFFAOYSA-N |
| 257 | butonitazene  | 2-[2-(4-butoxybenzyl)-5-nitro-1 <i>H</i> -benzimidazol-1-yl]- <i>N,N</i> -diethylethan-1-amine                      | C <sub>24</sub> H <sub>32</sub> N <sub>4</sub> O <sub>3</sub>  | UZZPOLCDCVWLAZ-UHFFFAOYSA-N |
| 258 | etonitazepyne   | 2-(4-ethoxybenzyl)-5-nitro-1-[2-(pyrrolidin-1-yl)ethyl]-1 <i>H</i> -benzimidazole                                   | C <sub>22</sub> H <sub>26</sub> N <sub>4</sub> O <sub>3</sub>  | LQZWZCJCEPUKCI-UHFFFAOYSA-N |
| 259 | BDMT  | 2,2'-(1 <i>H</i> ,1' <i>H</i> -[2,2'-biindole]-3,3'-diyl)bis( <i>N,N</i> -dimethylethan-1-amine)                    | C <sub>24</sub> H <sub>30</sub> N <sub>4</sub>                 | LDSNARXIXVOSTN-UHFFFAOYSA-N |
| 260 | ABO-4en-PINACA  | <i>N</i> -(1-amino-1-oxobutan-2-yl)-1-(pent-4-en-1-yl)-1 <i>H</i> -indazole-3-carboxamide                           | C <sub>17</sub> H <sub>22</sub> N <sub>4</sub> O <sub>2</sub>  | SHELRBQJCJWWQC-UHFFFAOYSA-N |
| 261 | 4F-deprenyl   | <i>N</i> -[1-(4-fluorophenyl)propan-2-yl]- <i>N</i> -methylprop-2-yn-1-amine  | C <sub>13</sub> H <sub>16</sub> FN                             | MUDUXRHPVDVWHU-UHFFFAOYSA-N |
| 262 | fluonitazene  | <i>N,N</i> -diethyl-2-[2-(4-fluorobenzyl)-5-nitro-1 <i>H</i> -benzimidazol-1-yl]ethan-1-amine                       | C <sub>20</sub> H <sub>23</sub> FN <sub>4</sub> O <sub>2</sub> | ZTWHIDCAGRMKTC-UHFFFAOYSA-N |
| 263 | M-alpha-HCMA  | 3-(1,3-benzodioxol-5-yl)-2-hydroxy- <i>N</i> ,2-dimethyl-3-(methylamino)propanamide                                 | C <sub>13</sub> H <sub>18</sub> N <sub>2</sub> O <sub>4</sub>  | GGNDZIOJQYAARA-UHFFFAOYSA-N |
| 264 | AP-238  | 1-[2,6-dimethyl-4-(3-phenylprop-2-enyl)piperazin-1-yl]propan-1-one  | C <sub>18</sub> H <sub>26</sub> N <sub>2</sub> O               | JELNWDOXWGBBLO-UHFFFAOYSA-N |
| 265 | carbonyl-bromadol   | (4-bromophenyl)-[1-(dimethylamino)-4-hydroxy-   | C <sub>23</sub> H <sub>28</sub> BrNO <sub>2</sub>              | SQKVCASVOTZNCS-             |



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|-----|--|---|---|-----------------------------|
|     |  | 4-phenethylcyclohexyl]methanone   |   | UHFFFAOYSA-N                |
| 266 | O-AMKD   | 3-(4-acetyl-1-methylpiperidin-4-yl)phenyl acetate   | C <sub>16</sub> H <sub>21</sub> NO <sub>3</sub>                 | ZXPASXBVYVTLRG-UHFFFAOYSA-N |
| 267 | nortilidine  | ethyl 2-methylamino-1-phenylcyclohex-3-ene-1-carboxylate  | C <sub>16</sub> H <sub>21</sub> NO <sub>2</sub>                 | PDJZPNKVLDWEKI-UHFFFAOYSA-N |
| 268 | 4,4-dimethyl-1-phenyl-1-pyrrolidin-1-yl-pentan-3-one | 4,4-dimethyl-1-phenyl-1-pyrrolidin-1-yl-pentan-3-one  | C <sub>17</sub> H <sub>25</sub> NO                              | JZBYSJKUYKWHMX-UHFFFAOYSA-N |
| 269 | phenylpiracetam                                      | 2-(2-oxo-4-phenylpyrrolidin-1-yl)acetamide  | C <sub>12</sub> H <sub>14</sub> N <sub>2</sub> O <sub>2</sub>   | LYONXVJRBWWGQO-UHFFFAOYSA-N |
| 270 | deoxymethoxetamine                                   | 2-(ethylamino)-2-(3-methylphenyl)-cyclohexanone   | C <sub>15</sub> H <sub>21</sub> NO                              | WIMLPRYZJQNQLE-UHFFFAOYSA-N |
| 271 | CHM-MDMB-CHMINACA                                    | cyclohexylmethyl 2-[1-(cyclohexylmethyl)-1 <i>H</i> -indazole-3-carboxamido]-3,3-dimethylbutanoate      | C <sub>28</sub> H <sub>41</sub> N <sub>3</sub> O <sub>3</sub>   | BJNHBUMERXHSSM-UHFFFAOYSA-N |
| 272 | protonitazene  | <i>N,N</i> -diethyl-5-nitro-2-[(4-propoxyphenyl)methyl]-1 <i>H</i> -benzimidazole-1-ethanamine          | C <sub>23</sub> H <sub>30</sub> N <sub>4</sub> O <sub>3</sub>   | SJHUJFHOXYDSJY-UHFFFAOYSA-N |
| 273 | 4Br-MAR (para-bromo-4-methylaminorex)                | 5-(4-bromophenyl)-4-methyl-4,5-dihydro-1,3-oxazol-2-amine   | C <sub>10</sub> H <sub>11</sub> BrN <sub>2</sub> O              | TUHDNALAVIDYHT-UHFFFAOYSA-N |
| 274 | 4Cl-MAR (Para-chloro-4-methylaminorex)               | 5-(4-chlorophenyl)-4-methyl-4,5-dihydro-1,3-oxazol-2-amine  | C <sub>10</sub> H <sub>11</sub> ClN <sub>2</sub> O              | PEMJVPLFSLEVII-UHFFFAOYSA-N |
| 275 | 2C-T-21  | 2-[4-(2-fluoroethylsulfanyl)-2,5-dimethoxyphenyl]ethanamine   | C <sub>12</sub> H <sub>18</sub> FNO <sub>2</sub> S              | ZBUUUKBTOCTOPW-UHFFFAOYSA-N |
| 276 | 5-chloro-alpha MT (5-Chloro-alpha-methyltryptamine)  | 1-(5-chloro-1 <i>H</i> -indol-3-yl)propan-2-amine   | C <sub>11</sub> H <sub>13</sub> ClN <sub>2</sub>                | QMKOQSCXSYPIPB-UHFFFAOYSA-N |
| 277 | 3-chlorophenmetrazine; 3-CPM                         | 2-(3-chlorophenyl)-3-methylmorpholine   | C <sub>11</sub> H <sub>14</sub> ClNO                            | BOFUZZAQNVYZFF-UHFFFAOYSA-N |
| 278 | Hydroxetamine; HXE                                   | 2-(ethylamino)-2-(3-hydroxyphenyl)-cyclohexanone  | C <sub>14</sub> H <sub>19</sub> NO <sub>2</sub>                 | CQERUJSORROCGH-UHFFFAOYSA-N |
| 279 | Dipyanone  | 4,4-diphenyl-6-(pyrrolidin-1-yl)heptan-3-one  | C <sub>23</sub> H <sub>29</sub> NO                              | LJIUPFDRFKFNJE-UHFFFAOYSA-N |
| 280 | Bretazenil   | 1,1-dimethylethyl 8-bromo-11,12,13,13a-tetrahydro-9-oxo-9 <i>H</i> -imidazo[1,5- <i>a</i> ]pyrrolo[2,1- | C <sub>19</sub> H <sub>20</sub> BrN <sub>3</sub> O <sub>3</sub> | LWUDDYHYNNIQL-UHFFFAOYSA-N  |

|     |   |   |   |                             |
|-----|---|---|---|-----------------------------|
|     |   | c][1,4]benzodiazepine-1-carboxylate   |   |                             |
| 281 | Deschloroclotizolam                                       | 2-chloro-9-methyl-4-phenyl-6 <i>H</i> -thieno[3,2- <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine          | C <sub>15</sub> H <sub>11</sub> ClN <sub>4</sub> S              | DBAZIULAFZCKIM-UHFFFAOYSA-N |
| 282 | BZO-POXIZID (MDA-19 pentyl analogue)                      | <i>N</i> -[( <i>Z</i> )-(2-oxo-1-pentyl-indolin-3-ylidene)amino]benzamide   | C <sub>20</sub> H <sub>21</sub> N <sub>3</sub> O <sub>2</sub>   | PCHOEXVGYICASA-UZYVYHOESA-N |
| 283 | iso-3-CMC   | 1-(3-chlorophenyl)-1-(methylamino)propan-2-one  | C <sub>10</sub> H <sub>12</sub> ClNO                            | CRNFRDNGAHBKFB-UHFFFAOYSA-N |
| 284 | BZO-4en-POXIZID (MDA-19 4en-pentyl analogue)              | <i>N</i> -[( <i>Z</i> )-(2-oxo-1-pent-4-enyl-indolin-3-ylidene)amino]benzamide                                      | C <sub>20</sub> H <sub>19</sub> N <sub>3</sub> O <sub>2</sub>   | DIVZUDBOOCSMQO-UZYVYHOESA-N |
| 285 | ADB-5Br-INACA   | 5-bromo- <i>N</i> -(1-carbamoyl-2,2-dimethyl-propyl)-1 <i>H</i> -indazole-3-carboxamide                             | C <sub>14</sub> H <sub>17</sub> BrN <sub>4</sub> O <sub>2</sub> | AJGASUCDTSLMNP-UHFFFAOYSA-N |
| 286 | BZO-ChMOXIZID   | <i>N</i> -{(Z)-[1-(cyclohexylmethyl)-2-oxo-indolin-3-ylidene]amino}benzamide  | C <sub>22</sub> H <sub>23</sub> N <sub>3</sub> O <sub>2</sub>   | HTPDZRIIOLCPPS-ATJXCDBQSA-N |
| 287 | 3,5-ADB-4en-PFUPPYCA                                      | <i>N</i> -(1-carbamoyl-2,2-dimethyl-propyl)-3-(4-fluorophenyl)-1-pent-4-enyl-pyrazole-5-carboxamide                 | C <sub>21</sub> H <sub>27</sub> FN <sub>4</sub> O <sub>2</sub>  | JPCQBOGNKSMAIE-UHFFFAOYSA-N |
| 288 | desmethylnoramide   | 4-(4-morpholinyl)-2,2-diphenyl-1-(1-pyrrolidinyl)-1-butanone  | C <sub>24</sub> H <sub>30</sub> N <sub>2</sub> O <sub>2</sub>   | JRPANCYSRUEJDY-UHFFFAOYSA-N |
| 289 | ADB-FUBIATA (ADB-FUBIACA)                                 | 2-[(2-{1-[1-(4-fluorophenyl)methyl]indol-3-yl}acetyl)amino]-3,3-dimethyl-butanamide                                 | C <sub>23</sub> H <sub>26</sub> FN <sub>3</sub> O <sub>2</sub>  | KHAUCCNSUMBFOT-UHFFFAOYSA-N |
| 290 | 5F-BZO-POXIZID; 5F-MDA-19; MDA-19 5-fluoropentyl analogue | <i>N</i> -{(Z)-[1-(5-fluoropentyl)-2-oxo-indolin-3-ylidene]amino}benzamide  | C <sub>20</sub> H <sub>20</sub> FN <sub>3</sub> O <sub>2</sub>  | CJINBVDNZDUMJM-PYCFMQQDSA-N |
| 291 | 3-Me-PCPy   | 1-[1-(3-methylphenyl)cyclohexyl]pyrrolidine   | C <sub>17</sub> H <sub>25</sub> N                               | JZVMREFYFTZXGN-UHFFFAOYSA-N |
| 292 | flubrotizolam   | 2-bromo-4-(2-fluorophenyl)-9-methyl-6 <i>H</i> -thieno[3,2- <i>f</i> ][1,2,4]triazolo[4,3- <i>a</i> ][1,4]diazepine | C <sub>15</sub> H <sub>10</sub> BrFN <sub>4</sub> S             | VOZDBDBHBXLWCG-UHFFFAOYSA-N |
| 293 | 5,3-ADB-4en-PFUPPYCA                                      | <i>N</i> -(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-5-(4-fluorophenyl)-1-pent-4-enyl-pyrazole-3-carboxamide            | C <sub>21</sub> H <sub>27</sub> FN <sub>4</sub> O <sub>2</sub>  | WGRQGKCMZJBNQD-UHFFFAOYSA-N |
| 294 | 5,3-AB-CHMFUPPYCA   | <i>N</i> -(1-amino-3-methyl-1-oxobutan-2-yl)-1-(cyclohexylmethyl)-5-(4-fluorophenyl)-1 <i>H</i> -                   | C <sub>22</sub> H <sub>29</sub> FN <sub>4</sub> O <sub>2</sub>  | XMZXGVYPIIOHAQ-UHFFFAOYSA-N |

## DRAFT

|     |                                       |   |   |                              |
|-----|---------------------------------------|---|---|------------------------------|
|     |                                       | pyrazole-3-carboxamide  |   |                              |
| 295 | fenzolone                             | 2-(ethylamino)-5-phenyl-4(5 <i>H</i> )-oxazolone  | C <sub>11</sub> H <sub>12</sub> N <sub>2</sub> O <sub>2</sub>     | RXOIEVSUURELPG-UHFFFAOYSA-N  |
| 296 | CUMYL-TsINACA                         | <i>N</i> -(2-phenylpropan-2-yl)-1-tosyl-1 <i>H</i> -indazole-3-carboxamide                                    | C <sub>24</sub> H <sub>23</sub> N <sub>3</sub> O <sub>3</sub> S   | NZHIGHLHQWQLPPC-UHFFFAOYSA-N |
| 297 | 1V-LSD                                | <i>N,N</i> -diethyl-7-methyl-4-pentanoyl-4,6,6a,7,8,9-hexahydroindolo[4,3- <i>fg</i> ]quinoline-9-carboxamide | C <sub>25</sub> H <sub>33</sub> N <sub>3</sub> O <sub>2</sub>     | GIIBVGJWUZNECE-XMSQKQJNSA-N  |
| 298 | etonitazepipne                        | 2-(4-ethoxybenzyl)-5-nitro-1-[2-(piperidin-1-yl)ethyl]-1 <i>H</i> -benzo[ <i>d</i> ]imidazole                 | C <sub>23</sub> H <sub>28</sub> N <sub>4</sub> O <sub>3</sub>     | UMGXRAISFRUVKD-UHFFFAOYSA-N  |
| 299 | iso-(meta-methyl-propcathinone)       | 1-(3-methylphenyl)-1-(propylamino)propan-2-one  | C <sub>13</sub> H <sub>19</sub> NO                                | UPNWTTPGWJZSBN-UHFFFAOYSA-N  |
| 300 | <i>N</i> -benzyl-isopropylamine       | <i>N</i> -benzylpropan-2-amine  | C <sub>10</sub> H <sub>15</sub> N                                 | LYBKPDDZTNUNNM-UHFFFAOYSA-N  |
| 301 | ADB-IACA (ADB-IATA)                   | 2-[2-(1 <i>H</i> -indol-3-yl)acetamido]-3,3-dimethylbutanamide  | C <sub>16</sub> H <sub>21</sub> N <sub>3</sub> O <sub>2</sub>     | WUWCLBKDNAZHPN-UHFFFAOYSA-N  |
| 302 | Cumyl-INACA                           | <i>N</i> -(1-methyl-1-phenyl-ethyl)-1 <i>H</i> -indazole-3-carboxamide  | C <sub>17</sub> H <sub>17</sub> N <sub>3</sub> O                  | COOPWWXIRLDJCP-UHFFFAOYSA-N  |
| 303 | Cumyl-CHSINACA                        | 1-(cyclohexylsulfonyl)- <i>N</i> -(2-phenylpropan-2-yl)-1 <i>H</i> -indazole-3-carboxamide                    | C <sub>23</sub> H <sub>27</sub> N <sub>3</sub> O <sub>3</sub> S   | ZNRPDVCQZGICAF-UHFFFAOYSA-N  |
| 304 | Cumyl-CLCHSINACA (CUMYL-1Cl-CHSINACA) | 1-(1-chlorocyclohexyl)sulfonyl- <i>N</i> -(1-methyl-1-phenyl-ethyl)indazole-3-carboxamide                     | C <sub>23</sub> H <sub>26</sub> ClN <sub>3</sub> O <sub>3</sub> S | MNMREJDPBHYPPIR-UHFFFAOYSA-N |
| 305 | CH-FUBIACA                            | <i>N</i> -cyclohexyl-2-[1-(4-fluorobenzyl)-1 <i>H</i> -indol-3-yl]acetamide                                   | C <sub>23</sub> H <sub>25</sub> FN <sub>2</sub> O                 | GSTACBDFNQHWHP-UHFFFAOYSA-N  |
| 306 | CH-PIACA                              | <i>N</i> -cyclohexyl-2-(1-pentyl-1 <i>H</i> -indol-3-yl)acetamide   | C <sub>21</sub> H <sub>30</sub> N <sub>2</sub> O                  | SYOOLIGHZEOKJ-UHFFFAOYSA-N   |
| 307 | MDMB-5Br-INACA                        | methyl 2-(5-bromo-1 <i>H</i> -indazole-3-carboxamido)-3,3-dimethylbutanoate                                   | C <sub>15</sub> H <sub>18</sub> BrN <sub>3</sub> O <sub>3</sub>   | QGEVEXPJOKFMAN-UHFFFAOYSA-N  |
| 308 | ADB-D-5Br-INACA                       | <i>N</i> -(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-5-bromo-1-decyl-1 <i>H</i> -indazole-3-carboxamide           | C <sub>24</sub> H <sub>37</sub> BrN <sub>4</sub> O <sub>2</sub>   | YDOQIEUYWRSIGV-UHFFFAOYSA-N  |
| 309 | A-PONASA                              | <i>N</i> -(adamantan-1-yl)-4-(pentyloxy)naphthalene-1-sulfonamide   | C <sub>25</sub> H <sub>33</sub> NO <sub>3</sub> S                 | KDLJELWGBJUNBO-UHFFFAOYSA-N  |

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|-----|--|---|---|-----------------------------|
| 310 | fluetizolam                              | 2-ethyl-4-(2-fluorophenyl)-9-methyl-6H-thieno[3,2-f][1,2,4]triazolo[4,3-a][1,4]diazepine                              | C <sub>17</sub> H <sub>15</sub> FN <sub>4</sub> S                             | BCKPHENWWQCRCG-UHFFFAOYSA-N |
| 311 | ADB-4en-P-5Br-INACA                      | N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-5-bromo-1-(pent-4-en-1-yl)-1H-indazole-3-carboxamide                         | C <sub>19</sub> H <sub>25</sub> BrN <sub>4</sub> O <sub>2</sub>               | SKMMFULKOGGVOT-UHFFFAOYSA-N |
| 312 | Desalkylgidazepam                        | 7-bromo-5-phenyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one   | C <sub>15</sub> H <sub>11</sub> BrN <sub>2</sub> O                            | ATCCWKYKHCKDGT-UHFFFAOYSA-N |
| 313 | N-cyclohexyl methylone                   | 1-(1,3-benzodioxol-5-yl)-2-(cyclohexylamino)propan-1-one  | C <sub>16</sub> H <sub>21</sub> NO <sub>3</sub>                               | WZDQUORQDHVTKD-UHFFFAOYSA-N |
| 314 | N-ethyl zolpidem                         | N-ethyl-2-[6-methyl-2-(4-methylphenyl)imidazo[1,2-a]pyridin-3-yl]acetamide  | C <sub>19</sub> H <sub>21</sub> N <sub>3</sub> O                              | GQJSOTATHCNTSW-UHFFFAOYSA-N |
| 315 | 2-fluoro-deschloro-N-ethylketamine       | 2-(ethylamino)-2-(2-fluorophenyl)cyclohexanone  | C <sub>14</sub> H <sub>18</sub> FNO   | RTXKYSLDFKUESF-UHFFFAOYSA-N |
| 316 | A-FUBIACA                                | N-(1-adamantyl)-2-[1-(4-fluorophenyl)methyl]indol-3-yl]acetamide  | C <sub>27</sub> H <sub>29</sub> FN <sub>2</sub> O                             | UGPWISFILITTLX-UHFFFAOYSA-N |
| 317 | hexahydrocannabinol (HHC)                | 6a,7,8,9,10,10a-hexahydro-6,6,9-trimethyl-3-pentyl-6H-dibenzo[b,d]pyran-1-ol  | C <sub>21</sub> H <sub>32</sub> O <sub>2</sub>                                | XKRHRBJLCLXSGE-UHFFFAOYSA-N |
| 318 | 4en-PDMB-4en-PINACA                      | pent-4-en-1-yl 3,3-dimethyl-2-(1-(pent-4-en-1-yl)-1H-indazole-3-carboxamido)-butanoate                                | C <sub>24</sub> H <sub>33</sub> N <sub>3</sub> O <sub>3</sub>                 | CQZNZGBBRCLQTI-UHFFFAOYSA-N |
| 319 | Acetyl-hexahydrocannabinol (HHC acetate) | (6,6,9-trimethyl-3-pentyl-6a,7,8,9,10,10a-hexahydrobenzo[c]chromen-1-yl) acetate                                      | C <sub>23</sub> H <sub>34</sub> O <sub>3</sub>                                | ZAZIHGFBNRVMAI-UHFFFAOYSA-N |
| 320 | Hexahydrocannabiphorol (HHC-P)           | 3-Heptyl-6a,7,8,9,10,10a-hexahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran-1-ol  | C <sub>23</sub> H <sub>36</sub> O <sub>2</sub>                                | USZILQYXSONCHH-UHFFFAOYSA-N |
| 321 | Fluorexetamine (FXE)                     | 2-(ethylamino)-2-(3-fluorophenyl)cyclohexan-1-one   | C <sub>14</sub> H <sub>18</sub> FNO   | FCETYWCLCUZFI-UHFFFAOYSA-N  |
| 322 | Rilmazafone                              | 5-[[[(2-aminoacetyl)amino]methyl]-1-[4-chloro-2-(2-chlorobenzoyl)phenyl]-N,N-dimethyl-1H-1,2,4-triazole-3-carboxamide | C <sub>21</sub> H <sub>20</sub> Cl <sub>2</sub> N <sub>6</sub> O <sub>3</sub> | KYHFRCPLIGODFH-UHFFFAOYSA-N |
| 323 | 5-MeO-TMT                                | 2-(5-methoxy-2-methyl-1H-indol-3-yl)-N,N-dimethylethanamine   | C <sub>14</sub> H <sub>20</sub> N <sub>2</sub> O                              | ACEHBQPPDDGCGZ-UHFFFAOYSA-N |
| 324 | ADMB-INACA                               | N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1H-  | C <sub>14</sub> H <sub>18</sub> N <sub>4</sub> O <sub>2</sub>                 | UFECWXZSRFBSHC-             |

## DRAFT

|     |         |  |  |                                 |
|-----|---------|--|--|---------------------------------|
|     |         | indazole-3-carboxamide                             |  | UHFFFAOYSA-N                    |
| 325 | FUBIAT  | 1-[(4-fluorophenyl)methyl]-1H-indole-3-acetic acid | C <sub>17</sub> H <sub>14</sub> FNO <sub>2</sub> | MSECUWYFWKAGLD<br>-UHFFFAOYSA-N |
| 326 | 4F-MBZP | 1-[(4-fluorophenyl)methyl]-4-methylpiperazine      | C <sub>12</sub> H <sub>17</sub> FN <sub>2</sub>  | KFMDNJJTGINMCS-<br>UHFFFAOYSA-N |

”

2. In Annex 3 to the Decree, the following point 7 shall be inserted:

‘7. For the purposes of listing, the data in columns A and B of the table in point 6.1 shall prevail, and the data in columns C and D shall ensure the computer-based searchability of the compounds.’