

RESEARCH RESOURCES

DRUG SUPPLY PROGRAM CATALOG

25TH EDITION

MAY 2016

CHEMISTRY AND PHARMACEUTICS BRANCH

DIVISION OF THERAPEUTICS AND MEDICAL CONSEQUENCES

NATIONAL INSTITUTE ON DRUG ABUSE

NATIONAL INSTITUTES OF HEALTH

DEPARTMENT OF HEALTH AND HUMAN SERVICES

6001 EXECUTIVE BOULEVARD

ROCKVILLE, MARYLAND 20852



On the cover: CPK rendering of nalfurafine.

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A. INTRODUCTION

The National Institute on Drug Abuse (NIDA) Drug Supply Program (DSP) is administered by the Chemistry and Pharmaceutics Branch, Division of Therapeutics and Medical Consequences. As part of its mission to lead the nation in bringing the power of science to bear on drug abuse and addiction, NIDA provides researchers with chemicals and research probes that are either unavailable, difficult to obtain, or very expensive to purchase. In addition, this program also provides analytical services of experimental samples used in research.

Drugs and Research Compounds. The NIDA DSP provides a variety of research chemicals and controlled substances for research purposes to all investigators working in the area of drug abuse, addiction, and related disciplines. The availability of controlled substances is regulated by the United States Drug Enforcement Administration (DEA), Department of Justice under the Controlled Substances Act (CSA), and Psychotropic Convention. These substances include hallucinogens, stimulants, sedatives and hypnotics, narcotics, designer drugs, cannabinoids, marijuana, and other miscellaneous categories of drugs. The NIDA DSP maintains an inventory of such drugs and other chemical substances. In addition, continuous efforts are made to synthesize new compounds and add them to the inventory. The stability and purity of all such compounds are monitored and maintained.

Marijuana and Marijuana Cigarettes. Marijuana is grown, harvested, processed and analyzed for delta-9-THC and other cannabinoids, stored under controlled conditions to preserve its purity and stability, and distributed for approved research purposes. This substance is subject to control under Schedule I of the CSA (21 U.S.C. 801 et seq.), the most restrictive of the five federally-regulated classes of controlled substances. Persons who wish to conduct research using marijuana for purposes other than research, such as forensic analytical standards, or any other law enforcement purposes must first obtain a special DEA registration under the CSA and then submit a request to NIDA along with required documents for consideration. It should be noted that priority for supply is given to researchers conducting federally-funded studies.

Nicotine Research Cigarettes. The NIDA DSP also houses a variety of nicotine research cigarettes (NRC) with specific characteristics for nicotine content and yield, tar, menthol, and ventilation (NOT-DA-13-002). As with other research chemicals or controlled substances supplied by the DSP, NRCs can be provided to investigators working in the area of drug abuse, addiction, or related disciplines at no cost. At this time, NRCs are of limited supply and priority is given to federally-funded grantees including

those working on areas related to the tobacco regulatory science authority of the NIDA Drug Supply Program and the United States Food & Drug Administration Center for Tobacco Products (FDA/CTP). A separate administrative review by the FDA/CTP, in collaboration with NIDA, is required for those working on projects within the FDA/CTP priority areas.

In order to obtain controlled substances, other research chemicals, marijuana, or marijuana and nicotine research cigarettes, research investigators are required to submit their requests along with necessary documents to the NIDA DSP for consideration. Applicants (domestic and foreign) without an NIH grant should note that their research protocol will require additional scientific review. Furthermore, research investigators who are not funded by NIH and plan to use marijuana or marijuana cigarettes in human research should be aware that their research protocol is subject to additional review as required by the US Department of Health and Human Services, Public Health Services. If marijuana is to be used for purposes other than human research, the request is forwarded to NIDA's Office of the Director for review and recommendation.

Foreign applicants must provide necessary documentation to demonstrate that they are allowed to import controlled substances, research chemicals, or marijuana cigarettes into their respective countries.

Approved research investigators are provided with pure drugs, compounds, or marijuana cigarettes along with their respective analytical data sheets.

B. NIDA DRUG SUPPLY PROGRAM (DSP) ORDERING GUIDELINES

To obtain research chemicals and controlled substances from the NIDA DSP, research investigators should prepare and submit a Request Package consisting of the following items for consideration:

1. A cover letter including:
 - a. Name, phone number and e-mail address of Research Investigator (and consignee, if applicable),
 - Provide a current and complete address that would allow shipment by a suitable carrier such as Federal Express (FedEx) (*i.e.* street address, Building name or number, room number, city and state) - This address should coincide with the address on the DEA order form.
 - For radio-labeled drugs or chemical substances, indicate the address at which the shipment can be made . In case the shipping address is different from the researcher's address, a current copy of the radioactive material license must be submitted.
 - b. If applicable, provide the NIH grant number, project title, and name and contact information of NIH Program Officer. If a non-grantee, no information on program Officer and grant number is required.
 - c. Name and quantity of compounds or other materials being requested, The total radioactivity (preferred unit of measurement, mg per vial, etc.)
 - If multiple studies/tasks are planned, combine projected needs into a single request rather than placing several separate requests in a short time interval. The request should generally be limited to four items or drugs/compounds per request to avoid delay.
 - Avoid drug abbreviations, and include specifications such as (+), (-), (*d/l*), base, or salt. You may consult the NIDA DSP Catalog for such information.
2. The Research Investigator's *curriculum vitae* (if not a funded grantee)
3. A detailed research protocol clearly indicating:
 - a. The specific aims and goals of proposed study (preferably the study abstract)
 - b. The number of experiments and experimental subjects
 - c. The dosage or concentration of drugs

- Calculate required amount of drugs or chemical substances for your project and submit your request well in advance of your planned experiments or tasks, or 6-8 weeks prior to depletion of stock on hand for ongoing studies/tasks.
- d. Justification for quantity of compounds/drugs, or other substances in request requested.
- If the request is related to a previously submitted protocol, provide a reference to this protocol and a brief description of progress along with references to resulting publications, if any.
4. A statement or commitment that NIDA will be acknowledged in research publications resulting from the use of supplies received from the NIDA Drug Supply Program.
 5. A completed DEA Order Form-222 for Schedule I-II controlled substances.
 - a. DEA Form-222 is not necessary for drugs in Schedules III-V, but a valid registration for the appropriate schedule is required.
 - b. Under the third column of DEA Form-222 (Size of Package), list quantities as bulk weight. Radio-labeled compounds must be listed by weight and NOT by units of activity.
 6. A copy of current DEA registration, Form DEA-223 for controlled substances.
 - a. It is the Research Investigator's responsibility to keep his/her registration current and to verify the drug code for requested drug.
 - b. Investigators who request a Schedule I drug and/or Etorphine HCl or Diprenorphine must provide DEA documentation under which the requested drug is covered under their current DEA registration.
 - c. Avoid drug abbreviations, and include specifications such as (+), (-), (dl), base, or salt, as appropriate.
 7. A copy of a Nuclear Regulatory Commission license for radioactive compounds.
For more information, you may visit the Nuclear Regulatory Commission licensing website.
 8. Research projects involving human subjects (Clinical research)
 - a. Investigational New Drug (IND) number and a copy of approved IND letter from the FDA. For more information, please visit the FDA IND Application website.
 - b. A copy of your Institutional Review Board (IRB) approval letter

- c. A copy of your Data Safety and Monitoring Plan
 - d. Proof of registration with clinicaltrials.gov (NCT #)
 - e. A copy of the study consent form(s)
9. For basic (non-human) research projects.
A copy of the document demonstrating that the research is approved by the Animal Care & Use Committee and that adequate care in conducting animal research will be exercised (if applicable). For more information, please visit the Animal Care & Use Committee Animal Study Proposal website.
10. For ongoing research projects (This includes research previously supported by the NIH prior to request)
- a. Reference information pertaining the previous protocol / FDA approval to the previous protocol
 - b. A brief description of progress (500 words or less)
 - c. A list of relevant publications.
11. Billing Information for Shipping (if applicable):
- a. US investigators should provide their FedEx account number to bill shipping charges,
 - b. International requests for controlled compounds should provide their United Parcel Service (UPS) Supply Chain Solutions account number to bill shipping charges. For non-controlled compounds the investigator should provide their FedEx account number.
12. Foreign research investigators must submit an Import Permit (preferably in English) issued by an appropriate agency of their government for internationally-controlled drug substances.
- a. The name of the supplier on the import permit must be listed as follows:

Research Triangle Institute
Kenneth S. Rehder
E. Institute Dr., Hermann Bldg.
Room 106
Research Triangle Park, NC 27709
 - b. The import permit should be accompanied by a signed statement from the investigator to the effect that the drug will be used solely for the purpose of re-

search and will not be re-exported. The import permit must contain a clear address and individual name to which the requested drug is to be delivered (not a post box number). Shipment should be indicated on the import permit as Air Freight to this individual. Specify the appropriate port for customs clearance purposes and provide the name and address of the Customs Clearance Agent, if one is used.

- c. Please obtain the longest possible expiration date on import permits.
- d. Foreign investigators requesting drugs or other substances that are not very stable should be certain to make advance arrangements that would allow these drugs or other substances to be released to them expeditiously. This is particularly relevant when dealing with radioactive materials with high specific activity (>1.0 Ci/mmol). These compounds are shipped as solutions in ice and will not stay cold for more than 48 hours; they must be retrieved and put in cold storage immediately to prevent decomposition.
- e. Foreign investigators should also be aware that compounds are shipped with a listed value for insurance purposes. Since this may cause problems with customs officials, investigators should determine in advance what steps should be taken to avoid these complications. Often an official statement that the compounds are for research only and have no commercial value is sufficient.

Failure to comply with aforesaid guidelines may delay the processing of your request.

ADDITIONAL NOTES:

1. ETORPHINE AND DIPRENORPHINE - Request for either of these two compounds should be made on separate order forms when ordering additional compounds at the same time. Etorphine hydrochloride and diprenorphine (free base or hydrochloride) are Schedule II drugs, however etorphine free base is a Schedule I drug. Therefore, etorphine free base should be ordered using DEA Form-222 as is appropriate for all Schedule I drugs.
2. CARFENTANIL, ETORPHINE AND DIPRENORPHINE – The DEA registration of a research investigator requesting these compounds must show the proper registration for such compounds.
3. For more information on the Drug Enforcement Administration registration forms including Form 222 and Form 223, please see: <http://www.deadiversion.usdoj.gov/drugreg/index.html>.

4. For more information about the FDA Investigational New Drug Application, please see: <http://www.fda.gov/Drugs/DevelopmentApprovalProcess/HowDrugsareDevelopedandApproved/ApprovalApplications/InvestigationalNewDrugINDApplication/default.htm>.
5. For more information on the Nuclear Regulatory Commission License, please see: <http://www.nrc.gov/about-nrc/regulatory/licensing.html>. Foreign applicants are required to submit additional documentation, including an import permit from the country into which the drug(s) are being shipped.

CONTACT INFORMATION:

All requests or questions should be addressed to:

Richard Kline, Ph.D.

phone: (301) 827-5243

email: rkline@nida.nih.gov

Robert Walsh

Phone: (301) 443-9825

email: bob.walsh@nih.gov

Kevin Gormley

Drug Supply Specialist

phone: (301) 435-0264

email: kgormley@nida.nih.gov

Division of Therapeutics and Medical Consequences (DTMC)

Chemistry and Pharmaceutics Branch (CPB)

National Institute on Drug Abuse, NIH

6001 Executive Boulevard, Room 4119

Rockville, MD 20852

Fax: (301) 443-2599

Phone: (301) 827-5243

C. DRUG REQUEST CHECKLIST

1. A *cover letter* including the name and quantities of compounds or drugs being requested, grant number, and name, phone number and e-mail address of your program officer (for an NIH/NIDA grantee), your shipping address, e-mail address, phone and fax numbers.
2. A *recommendation letter* from your program officer in support of your request.
3. A *research protocol*, including justification for the requested quantity of compounds or drugs being requested.
4. A *DEA Form 222* (for controlled substances).
5. A copy of your current *DEA registration* (for controlled substances).
6. An *approved FDA letter* and *IND number* (for a clinical study).
7. A copy of an *NRC license* for radioactive compounds.
8. A *curriculum vitae* of the principal investigator, if applicable.
9. A *statement of commitment* that NIDA will be acknowledged in publications.

D. SAMPLE DEA ORDER FORM 222

(THIS IS A SAMPLE OF A CORRECTLY COMPLETED DEA 222 FORM.
PLEASE NOTE THAT ALL ENTRIES SHOULD BE IN THE APPROPRIATE COLUMNS.)

See Reverse of PURCHASER'S Copy for Instructions		No order form may be issued for Schedule I and II substances unless a completed application form has been received. (21 CFR 1305.04)		OMB APPROVAL No. 1117-0010																																																																																																							
TO: (Name of Supplier) NIDA		STREET ADDRESS 6001 Executive Blvd, Room 4282, MSC 9555																																																																																																									
CITY and STATE Bethesda, MD 20892-9555		DATE XX-XX-XXXX	TO BE FILLED IN BY SUPPLIER SUPPLIERS DEA REGISTRATION NO.																																																																																																								
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<hr/> <p style="text-align: center;">CONTROLLED SUBSTANCES REGISTRATION CERTIFICATE UNITED STATES DEPARTMENT OF JUSTICE DRUG ENFORCEMENT ADMINISTRATION WASHINGTON, D.C. 20537</p> <p>The Controlled Substances Act of 1970 reads in part as follows:</p> <p>Sec. 304. (2) A registration instrument "to manufacture, distribute, or dispense a controlled substance may be suspended or revoked by the Attorney General upon a finding that the registrant -</p> <ul style="list-style-type: none"> (1) has materially falsified any application filed pursuant to or required by this title or title III, (2) has been convicted of a felony under the title or title III of any other law of the United States, or of any State, relating to any substance defined in this title as a controlled substance; (3) has had his State license or registration suspended, revoked, or denied by competent State authority and is no longer authorized by State law to engage in the manufacturing, distribution, or dispensing of controlled substances. <table border="1"> <thead> <tr> <th>DEA REGISTRATION NUMBER</th> <th>THIS REGISTRATION EXPIRES</th> <th>FEES PAID</th> </tr> </thead> <tbody> <tr> <td>XXXXXXXXXXXX</td> <td>XX-XX-XX</td> <td>XXXXXXX</td> </tr> <tr> <td colspan="2">SCHEDULES</td> <td>BUSINESS ACTIVITY</td> </tr> <tr> <td>XXXX</td> <td colspan="2">XXXXXXXXXXXXXXXXXXXX</td> </tr> <tr> <td></td> <td colspan="2">XX-XX-XX</td> </tr> </tbody> </table> <div style="border: 1px solid black; padding: 10px; text-align: center;"> <p>SAMPLE ONLY!! A COPY MUST BE INCLUDED WITH EVERY ORDER!!</p> </div>									DEA REGISTRATION NUMBER	THIS REGISTRATION EXPIRES	FEES PAID	XXXXXXXXXXXX	XX-XX-XX	XXXXXXX	SCHEDULES		BUSINESS ACTIVITY	XXXX	XXXXXXXXXXXXXXXXXXXX			XX-XX-XX																																																																																					
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E. SUPPLY AND ANALYSIS OF STANDARD SOLUTIONS OF Δ⁹-THC

In discussions with investigators working in the area of quantitative analysis of Δ⁹-THC, the question on the availability of its standard has repeatedly arisen. Because of the difficulties associated with accurately weighing exact quantity of non-crystalline Δ⁹-THC, NIDA provided the following Δ⁹-THC and its standard in solution form upon special request to research investigators in past and it still continues to do so:

1. 1. Small quantity of two stock solutions of Δ⁹-THC, 5 mg/ml in ethanol, and 0.5 mg/ml in ethanol, and Δ⁹-THC 0.70 mg/ml and androst-4-ene-3,17-dione in ethanol. The latter material is an internal standard for GC analysis. This steroid can also be provided in solid form to researchers who want to perform their own GC analyses. Note that both of these stock solutions are analyzed for exact composition and analytical data are provided with the shipment
2. As a final check, NIDA can provide the analysis of your stock solution using standardized procedure. At least 1.0 mL of stock solution should be provided with approximate concentration. If the concentration is greater than 50 µg/mL Δ⁹-THC, a standard GLC or HPLC analysis is performed using a steroid internal standard. For concentrations lower than this, GC/MS technique is used.

For further information, contact:

Richard Kline, Ph.D.

Phone: (301) 827-5243

email:

rkline@nida.nih.gov

Robert Walsh

Phone: (301) 443-9825

email:

bob.walsh@nih.gov

Kevin Gormley (RTI)

Phone: (301) 435-0264

email:

kgormley@nida.nih.gov

Division of Therapeutics and Medical Consequences (DTMC)
Chemistry and Pharmaceutics Branch (CPB)
National Institute on Drug Abuse, NIH
6001 Executive Boulevard, Room 4119
Rockville, MD 20852

F. ALTERNATIVE SOURCES FOR PEPTIDES

Recognizing a need in the research community for peptides, NIDA has included many peptides in the list of compounds provided to researchers through the NIDA Drug Supply Program. There are, however, a number of commercial suppliers that have many compounds of interest which are currently unavailable from the NIDA program. The addresses and phone numbers of some of these suppliers are:

PolyPeptide Laboratories San Diego
9395 Cabot Drive
San Diego, CA 82126
(800) 338-4965

Pierce Biotechnology
P.O. Box 117
Rockford, IL 61105
(800) 874-3723

Bachem Americas, Inc.
3132 Kashiwa Street
Torrance, CA 90505
(888) 422-2436

Sigma Chemical Co.
P.O. Box 14508
St. Louis, MO 63178
(800) 325-3010

NIDA is always receptive to suggestions for new peptides that are of interest to the scientific community. These suggestions should be in writing accompanied by the structure of the peptide as well as any names by which it is known. These suggestions should be forwarded to:

Richard Kline, Ph.D.
Division of Therapeutics and Medical Consequences (DTMC)
Chemistry and Pharmaceutics Branch (CPB)
National Institute on Drug Abuse, NIH
6001 Executive Boulevard, Room 4119
Rockville, MD 20852
email: rkline@nida.nih.gov
Fax: (301) 443-2599
Phone: (301) 827-5243

G. INSTRUCTIONS FOR ANALYTICAL SERVICES

NIDA facilitates drug abuse research by providing analytical services to research investigators who do not have necessary analytical facilities in their own laboratories. Priority is given to those investigators who are funded by NIDA.

Experimental samples (such as tissue, plasma, urine, and saliva) are analyzed in a NIDA contract laboratory for determining the concentration of drugs of abuse and their metabolites or precursors. The result of analyses is sent to research investigator with a copy to NIDA program official.

Request for analytical services should be submitted to the NIDA Drug Supply Program official with the following information for consideration and approval:

1. Number and title of your funded NIDA research grant
2. Name and contact information of your program official
3. In case of no grant, a brief research protocol
4. Number and name of drug(s) or drug metabolite(s) for analysis
5. Nature and origin of matrix such as rat/mouse/or human serum, plasma, urine, tissue, or saliva
6. Minimum expected concentrations of each drug/metabolites, and
7. Any other information that could be useful for analysis
8. A commitment to acknowledge NIDA in research publications resulting from this service. Once a research paper is published, a reprint or reference must be provided to NIDA drug supply program official.

After request approval, researcher will be contacted to submit experimental samples with shipping instructions to the following NIDA contract laboratory:

David E. Moody, Ph.D.
Center for Human Toxicology
University of Utah
30 South 2000 East
(for regular mail) Room 105
(for courier) Room 3956
Salt Lake City, UT 84112
Phone: (801) 581-5117; Fax: (801) 581-5034
E-Mail:david.Moody@utah.edu

NOTE: It is recommended that the research investigator obtain authorization prior to conducting any experiments so they can consult with the analytical laboratory about sample preparation and appropriate analyses.

Request for analysis should be submitted to:

Richard Kline, Ph.D.

Division of Therapeutics and Medical Consequences (DTMC)

Chemistry and Pharmaceutics Branch (CPB)

National Institute on Drug Abuse, NIH

6001 Executive Boulevard, Room 4119

Rockville, MD 20852

email: rkline@nida.nih.gov

Fax: (301) 443-2599

Phone: (301) 827-5243

H. X-RAY DIFFRACTION ANALYSIS OF COMPOUNDS

The purpose of this service is to provide definitive three-dimensional structural coordinates of compounds in support of drug abuse research. X-Ray diffraction analysis is performed on two types of compounds. The first type consists of non-peptide drugs, their analogs, related probes or potential analytical therapeutics including substances related to cocaine, phencyclidine, cannabinoids, opiates, and others. The second type consists of opioids and related peptides (CCK related, FMRF related, and alpha MSH-related fragments, etc.).

X-ray diffraction results provide full characterization of a compound even when the empirical formula of the molecule is not known beforehand. It is only the analytical method that can clearly define the absolute configuration of a molecule. An X-ray diffraction experiment requires only one good single crystal ideally about 0.1 x 0.2 x 0.2 mm (~0.1 mg of starting material). Note that single crystals may contain solvent so the best results are often obtained when samples are shipped in their mother liquor. If the submitted sample does not contain suitable single crystals then a minimum of 10 mg of sample is required for crystallization experiments. For these non-crystalline compounds information about stability and solubility should be provided with the sample. In present day research environment growing good single crystal is often the slowest step in the process since the use of high speed computers and state of the art data collection systems have reduced the time needed to complete a single structure study from several weeks to a few days.

For all samples indicate: 1 – if absolute or relative conformation is required as part of the analysis; 2 – the chirality of any known centers; 3 – provide a diagram of the expected structure including a preferred numbering scheme if applicable.

To request X-ray diffraction analysis of compounds, research investigators should provide the following information for consideration:

REQUEST FOR X-RAY DIFFRACTION ANALYSIS

Name of Investigator: _____

Institution: _____

Mailing Address: _____

E-Mail Address: _____

Phone & Fax Numbers: _____

NIH/NIDA Grant Title: _____

NIH/NIDA Grant Number: _____

Name of Your Program Officer (NIH/NIDA): _____

Type of compound: _____

Empirical Formula: _____

Anticipated Structure/comments/suggestions:

Signature of Principal Investigator: _____

Provide a recommendation letter from your Program Officer in support of your request. If not a NIDA grantee, provide a research protocol of your study for review signifying the relevance of your study with the NIDA research programs. The priority will be given to research investigators who are funded by the NIDA. A commitment to recognize NIDA must be provided. Once a paper is published, a copy, reprint, or publication reference must be submitted to NIDA program official.

Note: An incomplete application may not be considered.

The signed request (or electronic version) for analysis should be submitted to the following for approval before submitting your sample to the Center for Crystallo-

graphic Studies, Laboratory for Structure of Matter, Naval Research Laboratory,
4555 Overlook Ave. S.W., Washington, D.C. 20375-5000:

Richard Kline, Ph.D.

Division of Therapeutics and Medical Consequences (DTMC)

Chemistry and Pharmaceutics Branch (CPB)

National Institute on Drug Abuse, NIH

6001 Executive Boulevard, Room 4119

Rockville, MD 20852

email: rkline@nida.nih.gov

Fax: (301) 443-2599

Phone: (301) 827-5243

I. Nicotine Research Cigarettes Drug Supply Program

The NIDA DSP now provides a variety of nicotine research cigarettes (NRC) to research investigators. For more information on the Notice of Availability, please see <http://grants.nih.gov/grants/guide/notice-files/NOT-DA-14-004.html>.

Nicotine Research Cigarette Drug Supply Program Application Process

To obtain NRCs from the National Institute of Drug Abuse Drug Supply Program, all research investigators will need to prepare a Request Package.

Stepwise Procedure (Basic Non-Human Research)

1. A Request Package for NRC is submitted to the NIDA Drug Supply Program.
2. Once all request package materials are received, the NIDA Drug Supply Program Director reviews request package for completeness.
3. Materials are reviewed by NIH and assigned to a Program Officer.
 - For non-Grantees, the Program Officer refers to an external Scientific Expert Committee for further review.
4. Once reviewed, the Program Officer provides a recommendation to the NIDA Drug Supply Director for approval.
5. Once approved, if there is adequate inventory of nicotine research cigarettes, the NIDA Drug Supply Director issues shipment authorization.

Stepwise Procedure (Clinical Research with Human subjects)

1. A Request Package for NRC is submitted to the NIDA Drug Supply Program.
2. Once all request package materials are received, the NIDA Drug Supply Program Director for consideration.
3. Materials are reviewed by NIH and assigned to a Program Officer
 - For non-Grantees, the Program Officer refers to FDA and the Scientific Expert Committee for further review.
4. Once reviewed, the Program Officer provides a recommendation to the NIDA Drug Supply Director for approval.
5. Once approved, and there is adequate inventory of nicotine research cigarettes, the NIDA Drug Supply Director issues shipment authorization.

FAQs for the Investigational Tobacco Product (ITP) application

For researchers, in order for the Nicotine Research Cigarette (NRC) order to be deemed complete by the NIDA Drug Supply Program (DSP) Director, an Investigational Tobacco Product (ITP) application needs to be completed and forward to the FDA for approval. Below are helpful notes which will further your application to approval status. **Please be consistent and as specific as possible throughout your ITP application.** If not, applications may be considered inappropriate and delayed until corrected.

- Include actual date of previously submitted ITPs for cross-reference.

Reference the actual Tobacco Product Master File (TPMF) code name for the requested nicotine research cigarette (NRC) (e.g., NRC 300, RN).

- As specifically as possible, provide the specific grant number, grant title (where appropriate), project and/or protocol number and title, and PI for the study.
- Please provide the study sponsor. Note the distinction between “sponsor” and “investigator”: sponsors are typically considered to be the awarded institution (for NIH grants) to which the funding will be designated. Investigators are the primary research contact who conducts the research at that particular institution.
- Please be sure that the title of SPECTRUM cigarettes should be in all caps and consistent throughout your ITP application.

All inquiries and requests for the Investigational Tobacco Product application should be forwarded to the:

**Center for Tobacco Products
Food and Drug Administration
Document Control Center, Room 020J
9200 Corporate Boulevard Rockville, MD 20850
(301) 796-0456**

Point of Contact. All other requests or questions should be addressed to:

**Richard Kline, Ph.D.
National Institute on Drug Abuse, NIH
6001 Executive Boulevard, Room 4119
Rockville, MD 20852
email: rkline@nida.nih.gov
Fax: (301) 443-2599
Phone: (301) 827-5243**

J. Ordering Guidelines for Nicotine Research Cigarettes (NRCs)

To obtain NRCs from the NIDA DSP, all research investigators will need to prepare a NRC Request Package. A Request Package should include the following items:

1. A cover letter including:
 - a. Name, phone number and e-mail address of Research Investigator (and consignee, if applicable),
 - Provide a current and complete address that would allow shipment by a suitable carrier such as Federal Express (FedEx) (i.e., street address, building name or number, room number, city and state),
 - b. If applicable, NIH grant number of project and name and contact information of project's NIDA/NIH Program Officer. If a non-NIH grantee, no information on Program Officer and grant number is required.
 - c. Type and quantity of NRCs being requested,
 - If multiple studies/tasks are planned, combine projected needs into a single order rather than placing several separate requests in a short time interval. The request should generally be limited to four items or drugs/compounds per order to avoid delay.
2. The Research Investigator's curriculum vitae (CV).
3. A detailed research protocol clearly indicating:
 - a. The specific aims and goals of proposed study (preferably the study abstract)
 - b. Number of experiments and experimental subjects
 - c. Number and type of NRCs (NOT-DA-14-004), including respective Tobacco Product Master File (TPMF) codes,
 - Calculate required amount of NRCs for your project and submit your request well in advance of your planned experiments or tasks, or 6-8 weeks prior to depletion of stock on hand for ongoing studies/tasks.
 - d. Justification for the quantity needed. If the request is related to a previously submitted protocol, provide a reference to this protocol and a brief statement of progress along with references to resulting publications.
 - e. A timeline indicating approximately when and what quantities of NRC shipments are needed over the duration of the protocol.

- f. Statement of commitment that NIDA will be acknowledged in research publications using the NIDA-supplied cigarettes.

4. For Clinical Research Projects Involving Human Subjects

- a. Investigational Tobacco Product (ITP) application information including: A copy of FDA letter (Advice/Information Request) in response to your ITP application. (NOTE: Suggested guidelines for submitting an Investigational Tobacco Product Application (ITPA) to FDA/CTP are given separately. For more information, please visit [here](#).)
- b. A copy of your Institutional Review Board (IRB) approval letter
- c. A copy of your Data Safety and Monitoring Plan
- d. Proof of registration with clinicaltrials.gov (NCT #)
- e. A copy of the study consent form(s)

5. For Basic (Non-human) Research Projects

- A copy of the document demonstrating that the research is approved by the Animal Care & Use Committee and that adequate care in conducting animal research will be exercised. For more information, please visit the Animal Care & Use Committee Animal Study Proposal website.

6. For Ongoing Research Projects (This includes research previously supported by the NIH prior to NRC request)

- a. Reference information pertaining the previous protocol / FDA approval to the previous protocol (if applicable).
- b. A brief statement of progress (500 words or less)
- c. A list of any relevant publications.

7. Billing Information for Shipping Costs (if applicable):

- a. US investigators should provide their FedEx account number to bill shipping charges,
- b. International requests should provide United Parcel Service (UPS) Supply Chain Solutions account number to bill shipping charges or FedEx account number as appropriated for the import of NRCs.

Contact Information:

All requests/questions should be sent to: NIDANRCSupply@mail.nih.gov.

Please specify materials requested in the subject line.

Address correspondence to:

Richard Kline, Ph.D.

Division of Therapeutics and Medical Consequences (DTMC)

Chemistry and Pharmaceutics Branch (CPB)

National Institute on Drug Abuse, NIH

6001 Executive Boulevard, Room 4119

Rockville, MD 20852

email: rkline@nida.nih.gov

Fax: (301) 443-2599

Phone: (301) 827-5243

K. Ordering Guidelines for Marijuana and Marijuana Cigarettes

To obtain marijuana and marijuana cigarettes from the NIDA DSP, all research investigators will need to prepare a Marijuana Cigarette Request Package. A Request Package should include the following items:

1. A cover letter including:

- a. Name, phone number and e-mail address of Research Investigator (and consignee, if applicable),
 - Provide a current and complete address that would allow shipment by a suitable carrier such as Federal Express(FedEx) (i.e. street address, building name or number, room number, city and state) - This address should coincide with the address on the DEA order form.
 - If requestor's address is outside the US, please provide all necessary documentation to verify whether particular compound requested is permitted in his/her country for importation.
- b. If applicable, NIH grant number of project and name and contact information of project's NIDA/NIH Program Officer (if applicable).
If a non-grantee, no information regarding the Program Officer and grant number is required.
- c. Name(s) and quantity of compounds or other substances being requested,
 - If multiple studies/tasks are planned, combine projected needs into a single order rather than placing several separate requests in a short time interval.

The request should generally be limited to no more than four items per order to avoid delay.

2. The Research Investigator's curriculum vitae (CV).

3. A detailed research protocol clearly indicating:

- a. The specific aims and goals of proposed study (preferably the study abstract),
- b. The number of experiments and experimental subjects,
- c. The strength or concentration of marijuana or marijuana cigarettes,
 - Calculate required amount of marijuana cigarettes for your project and submit your request well in advance of your planned experiments or tasks, or 6-8 weeks prior to depletion of stock on hand for ongoing studies/tasks.
- d. Justification for the quantity of marijuana or marijuana cigarettes requested.

If the request is related to a previously submitted protocol, provide a reference to this protocol and a brief statement of progress along with references to resulting publications.

4. Statement of commitment that NIDA will be acknowledged in research publications using the NIDA Drug Supply Program.

5. A completed [DEA Order Form-222](#).

- Under the third column of DEA Form-222 (Size of Package), list quantities as bulk weight.

6. A copy of current DEA registration, [Form DEA-223](#) for a controlled substance.

- It is the Research Investigator's responsibility to keep his/her registration current and verify the drug code for the requested materials.

7. For clinical research projects involving human subjects

1. Investigational New Drug (IND) number and a copy of approved IND letter from the FDA. For more information, please visit the [FDA IND Application website](#).
2. A copy of your Institutional Review Board (IRB) approval letter
3. A copy of your Data Safety and Monitoring Plan
4. Proof of registration with clinicaltrials.gov (NCT #)
5. A copy of the study consent form(s)

8. For basic (non-human) research projects

- A copy of the document demonstrating that the research is approved by the Animal Care & Use Committee and that adequate care in conducting animal research will be exercised (if applicable).

For more information, please visit the [Animal Care & Use Committee Animal Study Proposal website](#).

9. For ongoing research projects (This includes research previously supported by the NIH prior to request)

- a. Reference information pertaining to the previous protocol.
- b. A brief statement of progress (500 words or less)
- c. A list of any relevant publications.

10. Billing Information for Shipping Costs:

1. US investigators should provide a FedEx account number for shipping charges,
2. International requests for marijuana and marijuana cigarettes should provide their United Parcel Service (UPS) Supply Chain Solutions account number to bill shipping charges.

11. For Foreign Investigators, please submit the Import Permit (preferably in English) issued by an appropriate agency of your government. The name of the supplier on the import permit must be listed as follows:

Research Triangle Institute
Kenneth S. Rehder
E. Institute Dr., Hermann Bldg.
Room 106
Research Triangle Park, NC 27709

- The import permit should be accompanied by a signed statement from the investigator to the effect that the materials will be used solely for the purpose of research and will not be re-exported. The import permit must contain a clear address and individual name to which the requested material is to be delivered (not a post office box number). Shipment should be indicated on the import permit as Air Freight to this individual. Specify the appropriate port for customs clearance purposes and provide the name and address of the Customs Clearance Agent, if one is used.
- Please obtain the longest possible expiration date on import permits.
- Foreign investigators should also be aware that the material is shipped with a listed value (for insurance purposes). Since this can cause problems with customs officials, investigators should determine in advance what steps should be taken to avoid these complications. Often an official statement that the materials are for research only and have no commercial value is sufficient.

Failure to comply with these guidelines may delay the processing of your request.

Contact Information:

Richard Kline, Ph.D.

phone: (301) 827-5243

email: rkline@nida.nih.gov

Robert Walsh

Phone: (301) 443-9825

email: bob.walsh@nih.gov

Kevin Gormley

Drug Supply Specialist

phone: (301) 435-0264

email: kgormley@nida.nih.gov

Division of Therapeutics and Medical Consequences (DTMC)

Chemistry and Pharmaceutics Branch (CPB)

National Institute on Drug Abuse, NIH

6001 Executive Boulevard, Room 4119

Rockville, MD 20852

L. IMPORTANT ADDRESSES, TELEPHONE & FAX NUMBERS

NATIONAL INSTITUTE ON DRUG ABUSE

Richard Kline, Ph.D.	Robert Walsh	Kevin Gormley (RTI)
Phone: (301) 827-5243	Phone: (301) 443-9825	Phone: (301) 435-0264
email:	email:	email:
rkline@nida.nih.gov	bob.walsh@nih.gov	kgormley@nida.nih.gov

Division of Therapeutics and Medical Consequences (DTMC)
Chemistry and Pharmaceutics Branch (CPB)
National Institute on Drug Abuse, NIH
6001 Executive Boulevard, Room 4119
Rockville, MD 20852

DRUG ENFORCEMENT ADMINISTRATION (DEA)

Office of Diversion Control Online: <http://www.deadiversion.usdoj.gov/>
Drug Enforcement Administration
Office of Diversion Control
8701 Morrissette Drive
Springfield, Virginia 22152
Phone: (800) 882-9539

FOOD AND DRUG ADMINISTRATION (FDA)

U.S. Food & Drug Administration
Department of Health & Human Services
10903 New Hampshire Avenue
Silver Spring, Maryland 20993
Phone: (888) 463-6332
Website: <http://www.fda.gov>

M. AVAILABLE DRUGS, COMPOUNDS, & DOSAGE FORMS



Disclaimer: All compound descriptions in this catalog are provided for the purpose of general information only and are not intended to address specific issues or to be a complete or definitive source of such information. Nothing herein is to be considered as a warranty, expressed or otherwise. Users of any the materials provided by NIDA are obligated to understand the proper and safe research applications and to comply with the applicable federal, state, and local laws and regulations. The substances described in this catalog are suitable for basic research only and, unless otherwise indicated, are **not** for human use.

1 – New Compounds

(Compounds added to the NDSP catalog since the 24th edition)

Catalog number	Compound	Page
7023-001	AB-PINACA	37
7031-001	AB-CHMINACA	37
NOCD-142	(+)-PHT dihydrochloride	72
NICT-020	(-)–Nicotine ditartrate	74
9652-071	CNV-NLX	77
9739-001	Remifentanil hydrochloride	78
2270-012	Sodium pentobarbital	141
7535-002	(<i>R</i>)-MDPV hydrochloride	150
7535-003	(<i>S</i>)-MDPV hydrochloride	150
7545-001	α -PVP hydrochloride	150



2 – Cannabinoids and Cannabinoid-related

Cannabinoids: Allosteric Modulators

Catalog number : NOCD-133

CASRN : 868273-06-7

Name : ORG 27569

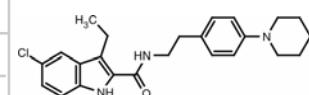
Mol. formula : C₂₄H₂₈ClNO₃

FW : 409.95

DEA schedule : 0

Notes : CB₁ receptor positive allosteric modulator.

References : Price et al (2005) Allosteric modulation of the Cannabinoid CB1 receptor. Mol.Pharmacol. 68 1484. PMID: 16113085.



Cannabinoids: Cannabichromene Class

Catalog number : 7360-007

CASRN : 20675-51-8

Name : Cannabichromene

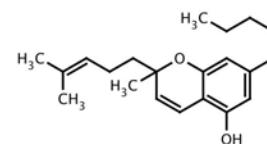
Mol. formula : C₂₁H₃₀O₂

FW : 314.46

DEA schedule : 1

Notes : Non-psychoactive constituent of cannabis.

References : Turner, CE; Elsohly, MA *J Clin Pharmacol* 1981, 21, 283S-291S.



Catalog number : 7360-008

CASRN : 20675-51-8 (parent cmpd)

Name : [1',2'-³H₂]Cannabichromene

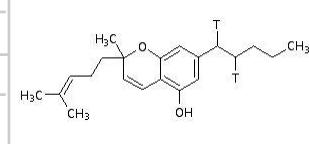
Mol. formula : C₂₁H₃₀O₂

FW : 314.46

DEA schedule : 1

Notes : Non-psychoactive constituent of cannabis (tritium-labeled).

References : Instrumental Data for Drug Analysis, 2nd Ed., 1996, Volume 1, p304.



Cannabinoids: Cannabicyclohexanol Class

Catalog number : 7927-001

CASRN : 114753-51-4

Name : (±)-CP-47497

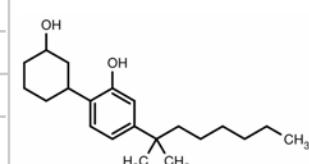
Mol. formula : C₂₁H₃₄O₂

FW : 318.50

DEA schedule : 1

Notes : Potent CB₁ receptor agonist.

References : Melvin, L. S., et al., *J Med Chem* 1984, 27(1), 67-71.



Catalog number : 7298-001

CASRN : 70434-92-3

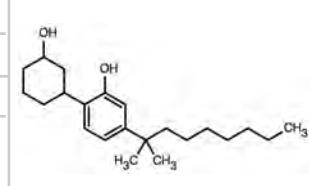
Name : Cannabicyclohexanol; (±)-CP-47,497 C8 homolog

Mol. formula : C₂₂H₃₆O₂

FW : 332.52

DEA schedule : 1

References : Compton DR, Johnson MR, Melvin LS, Martin BR, *J Pharm Exp Ther*, 1992, 260(1), 201-209.



2 – Cannabinoids and Cannabinoid-related

Catalog number : NOCD-091

CASRN : 83002-04-4

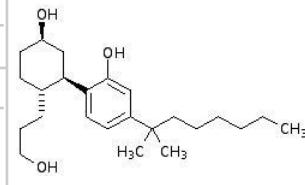
Name : (–)CP 55,940

Mol. formula : C₂₄H₄₀O₃

FW : 376.58 **DEA schedule :** 0

Notes : High-affinity cannabinoid CB1 and CB2 receptor agonist

References : Melvin, LS; et al. *J Med Chem* 1984, 27, 67–71.



Catalog number : NOCD-092

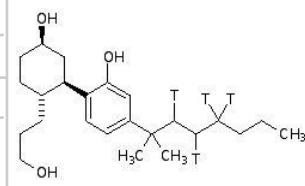
CASRN : 119095-48-6

Name : [2,3,4,4–³H₄] (–)CP 55,940

Mol. formula : C₂₄H₄₀O₃

FW : 376.58 **DEA schedule :** 0

Notes : High affinity CB1 and CB2 receptor radioligand (tritium-labeled).



Cannabinoids: Cannabidiol Class

Catalog number : 7360-021

CASRN : 13956-29-1

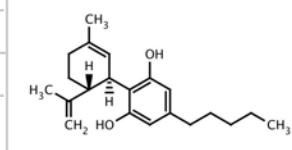
Name : (–)Cannabidiol

Mol. formula : C₂₁H₃₀O₂

FW : 314.46 **DEA schedule :** 1

Notes : CB₁ and CB₂ receptor antagonist

References : Petitet, F; et al. *Life Sci* 1998, 63, PL1–6.
Costa, B; et al. *Br J Pharmacol* 2004, 143, 247–50.
Thomas, A; et al. *Br J Pharmacol* 2007, 150, 613–23.



Catalog number : 7360-022

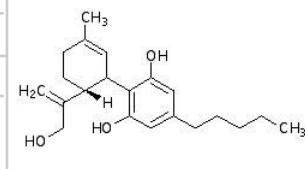
Name : 10-Hydroxycannabidiol

Mol. formula : C₂₁H₃₀O₃

FW : 330.46 **DEA schedule :** 1

Notes : Cannabidiol metabolite

References : Lander, N; et al. *J Chem Soc., Perkin Trans 1* 1976, 8–16.



Cannabinoids: Cannabigerol Class

Catalog number : 7360-010

CASRN : 25654-31-3

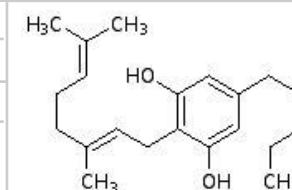
Name : Cannabigerol

Mol. formula : C₂₁H₃₂O₂

FW : 316.48 **DEA schedule :** 1

Notes : Analgesic; anti-inflammatory

References : Williamson, EM; Evans, FJ *Drugs* 2000, 60, 1303–14.



Cannabinoids: Cannabinol Class

Catalog number : 7360-013

CASRN : 21-35-7

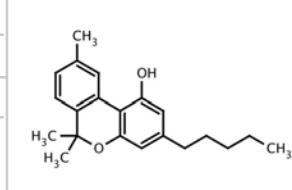
Name : Cannabinol; CBN

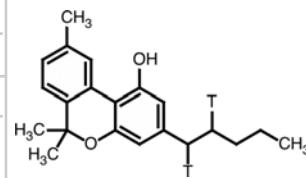
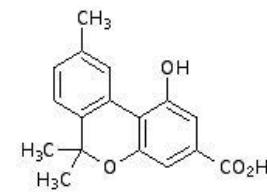
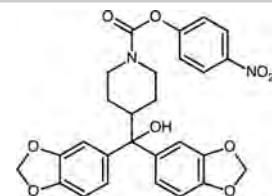
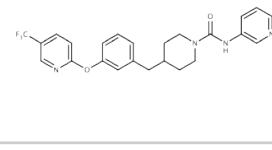
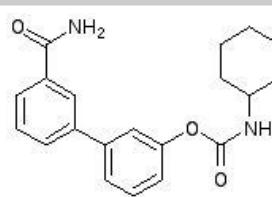
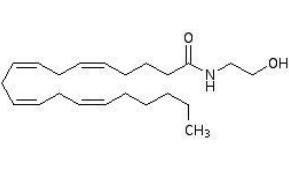
Mol. formula : C₂₁H₂₆O₂

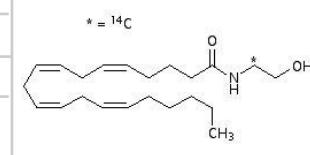
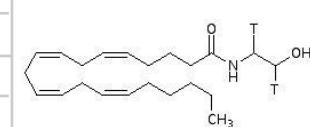
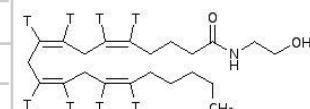
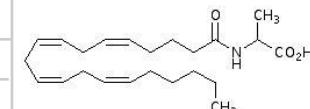
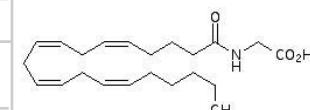
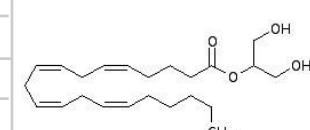
FW : 310.43 **DEA schedule :** 1

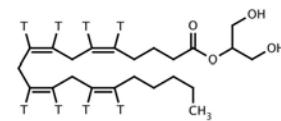
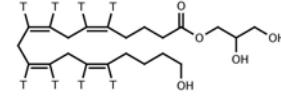
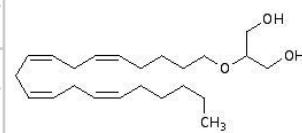
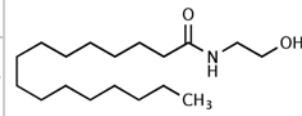
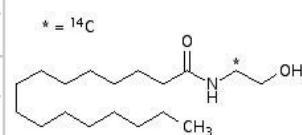
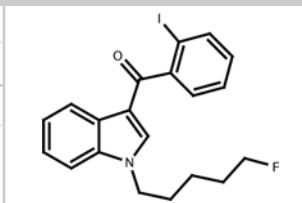
Notes : Inactive constituent of cannabis.

References : Mahadevan, A; et al. *J Med Chem* 2000, 43, 3778–85.

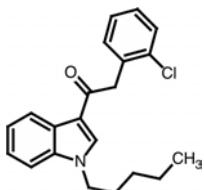
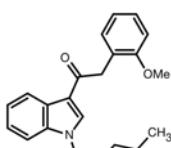
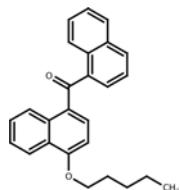
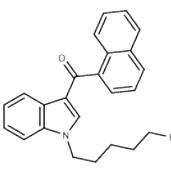
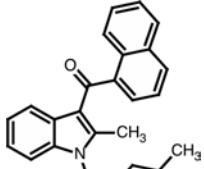
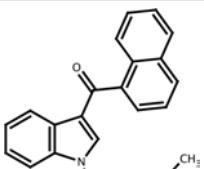


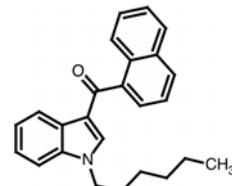
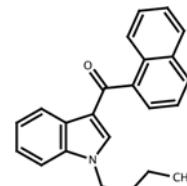
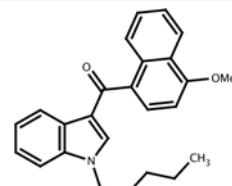
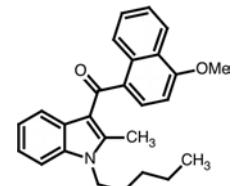
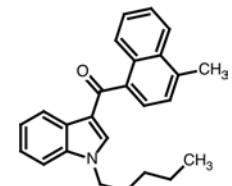
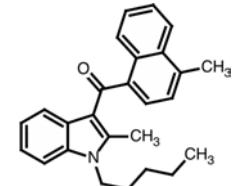
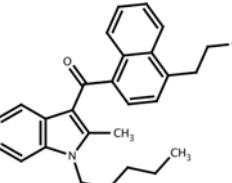
Catalog number : 7360-023**Name :** [1',2'-³H₂]Cannabinol**Mol. formula :** C₂₁H₂₆O₂**FW :** 310.43 **DEA schedule :** 1**Catalog number :** 7360-020**CASRN :** 60788-14-9**Name :** 1',2',3',4',5'-Pentanorcannabinol-3-carboxylic acid**Mol. formula :** C₁₇H₁₆O₄**FW :** 284.31 **DEA schedule :** 1**Cannabinoids: Enzyme Inhibitors****Catalog number :** NOCD-037**CASRN :** 1101854-58-3**Name :** JZL184**Mol. formula :** C₂₇H₂₄N₂O₉**FW :** 520.49 **DEA schedule :** 0**Notes :** Potent and selective inhibitor of monoacylglycerol lipase (MAGL).**References :** Long, JZ; et al. *Nat Chem Biol* 2009, 5, 37–44.**Catalog number :** NOCD-124**CASRN :** 1196109-52-0**Name :** PF-3845**Mol. formula :** C₂₄H₂₃F₃N₄O₂**FW :** 456.47 **DEA schedule :** 0**Notes :** Selective fatty acid amide hydrolase (FAAH) inhibitor.**References :** Ramesh D, Ross GR, Schlosburg JE, Owens RA, Abdullah RA, Kinsey SG, Long JZ, Nomura DK, Sim-Selley LJ, Cravatt BF, Akbarali HI, Lichtman AH, *J Pharmacol Exp Ther* 2011, 339(1), 173–85.**Catalog number :** NOCD-098**CASRN :** 546141-08-6**Name :** Cyclohexylcarbamic Acid 3'-carbamoylbiphenyl-3-yl ester; URB 597**Mol. formula :** C₂₀H₂₂N₂O₃**FW :** 338.41 **DEA schedule :** 0**Notes :** Fatty acid amide hydrolase (FAAH) inhibitor.**References :** Cravatt BF; et al. *Nature* 1996, 384, 83.
Kathuria S; et al. *Nat Med* 2003, 9, 76.**Cannabinoids: Fatty Acid Derivatives (Anandamides)****Catalog number :** NOCD-080**CASRN :** 94421-68-8**Name :** Arachidonylethanolamide; Anandamide**Mol. formula :** C₂₂H₃₇NO₂**FW :** 347.54 **DEA schedule :** 0**Notes :** Cannabinoid CB1 and CB2 receptor agonist.**References :** Devane, WA; et al. *Science* 1992, 258, 1946–9.

Catalog number : NOCD-007**Name :** Arachidonyl[1-¹⁴C]ethanolamide**Mol. formula :** C₂₃H₃₇NO₂**FW :** 361.56 **DEA schedule :** 0**Catalog number :** NOCD-008**Name :** Arachidonyl[1,2-³H]ethanolamide; Tritiated Anandamide**Mol. formula :** C₂₂H₃₇NO₂**FW :** 347.54 **DEA schedule :** 0**Catalog number :** NOCD-078**Name :** Tritium-labeled Arachidonylethanolamide; Tritiated Anandamide**Mol. formula :** C₂₂H₃₇NO₂**FW :** 347.54 **DEA schedule :** 0**Notes :** *Cannabinoid CB1 and CB2 receptor radioligand.***Cannabinoids: Fatty Acid Derivatives (Arachidonyl amides)****Catalog number :** NOCD-097**CASRN :** 401941-73-9**Name :** N-Arachidonyl-L-alanine**Mol. formula :** C₂₃H₃₇NO₃**FW :** 375 **DEA schedule :** 0**Catalog number :** NOCD-096**CASRN :** 179113-91-8**Name :** N-Arachidonylglycine; NAGly**Mol. formula :** C₂₂H₃₅NO₃**FW :** 361 **DEA schedule :** 0**Notes :** *Endogenous anandamide-like compound with analgesic properties (although it lacks CB1 receptor and anandamide transporter affinity).***References :** Sheskin, T; et al. *J Med Chem* 1997, 40, 659-67.Huang, SM; et al. *J Biol Chem* 2001, 276, 42639-44.**Cannabinoids: Fatty Acid Derivatives (Arachidonyl esters)****Catalog number :** NOCD-089**CASRN :** 53847-30-6**Name :** 2-Arachidonylglycerol; 2-AG**Mol. formula :** C₂₃H₃₈O₄**FW :** 378.5 **DEA schedule :** 0**Notes :** *Cannabinoid CB1 receptor agonist.***References :** Stella, N; Schweitzer P; Piomelli D. *Nature* 1997, 388, 773-8.

Catalog number : NOCD-018**Name :** [³H]-2-Arachidonylglycerol; [³H]-2-AG**Mol. formula :** C₂₃H₃₈O₄**FW :** 378.55 **DEA schedule :** 0**Notes :** Cannabinoid CB1 receptor agonist (tritium-labeled).**References :** Stella, N; Schweitzer, P; Piomelli, D *Nature* 1997, 388, 773–8.**Catalog number :** NOCD-035**Name :** Tritium-labeled 1-Arachidonylglycerol; [³H]-1-AG**Mol. formula :** C₂₃H₃₈O₄**FW :** 378.55 **DEA schedule :** 0**Notes :** Cannabinoid CB1 receptor agonist (tritium-labeled).**References :** Stella, N; Schweitzer, P; Piomelli D *Nature* 1997, 388, 773–8.**Catalog number :** NOCD-095**CASRN :** 222723-55-9**Name :** Noladin**Mol. formula :** C₂₃H₄₀O₃**FW :** 364.57 **DEA schedule :** 0**Notes :** Cannabinoid CB1 receptor agonist**References :** Hanus, L; et al. *Proc Natl Acad Sci USA* 2001, 98, 3662–5.**Cannabinoids: Fatty Acid Derivatives (Palmitoyl amides)****Catalog number :** NOCD-002**CASRN :** 544-31-0**Name :** Palmitoyl ethanolamide; Palmidrol**Mol. formula :** C₁₈H₃₇NO₂**FW :** 299.49 **DEA schedule :** 0**Notes :** Cannabinoid CB2 receptor agonist.**References :** Hanus, L; et al. *J Med Chem* 1993, 36, 3032–4.**Catalog number :** NOCD-005**Name :** Palmitoyl[1-¹⁴C]ethanolamide**Mol. formula :** C₁₈H₃₇NO₂**FW :** 299.49 **DEA schedule :** 0**Notes :** Cannabinoid CB2 receptor agonist (carbon-labeled).**Cannabinoids: Indole Analogs & Related Compounds****Catalog number :** 7694-001**CASRN :** 335161-03-0**Name :** AM-694**Mol. formula :** C₂₀H₁₉FINO**FW :** 435.27 **DEA schedule :** 1**References :** Logan BK, Reinhold LE, Xu A, Diamond FX, *J Forensic Sci*, 2012, 57(5), 1168–1180.

2 – Cannabinoids and Cannabinoid-related

Catalog number : 7203-001	CASRN : 864445-54-5	
Name : JWH-203		
Mol. formula : C ₂₁ H ₂₂ CINO	FW : 339.86 DEA schedule : 1	
Notes : Non-selective CB ₁ /CB ₂ cannabinoid receptor agonist.		
References : Huffman JW, Szklenik PV, Almond A, Bushell K, Selley DE, He H, Cassidy MP, Wiley JL, et al., <i>Bioorg Med Chem Letters</i> , 2005, 15(18), 4110–4113.		
Catalog number : 6250-001	CASRN : 864445-43-2	
Name : JWH-250		
Mol. formula : C ₂₂ H ₂₅ NO ₂	FW : 335.44 DEA schedule : 1	
Notes : Non-selective CB ₁ /CB ₂ cannabinoid receptor agonist.		
References : Huffman JW, Szklenik PV, Almond A, Bushell K, Selley DE, He H, Cassidy MP, Wiley JL, et al., <i>Bioorg Med Chem Letters</i> , 2005, 15(18), 4110–4113.		
Catalog number : NOCD-036	CASRN : 432047-72-8	
Name : Naphthalen-1-yl-(4-pentyloxynaphthalen-1-yl)methanone		
Mol. formula : C ₂₆ H ₂₄ O ₂	FW : 368.47 DEA schedule : 0	
Notes : Orally bioavailable human CB ₁ /CB ₂ dual agonist with antihyperalgesic properties and limited CNS penetration.		
References : Dziadulewicz, EK; et al. <i>J Med Chem</i> 2007, 50, 3851–6.		
Cannabinoids: Indole, Alkylnaphthoyl Class		
Catalog number : 7201-001	CASRN : 335161-24-5	
Name : AM-2201		
Mol. formula : C ₂₄ H ₂₂ FNO	FW : 359.44 DEA schedule : 1	
Notes : Non-selective CB ₁ /CB ₂ cannabinoid receptor agonist.		
References : Makriyannis, A; Deng H, Cannabimimetic Indole Derivatives, US Patent 7,241,799 B2 (2007).		
Catalog number : 7007-001	CASRN : 155471-10-6	
Name : JWH-007		
Mol. formula : C ₂₅ H ₂₅ NO	FW : 355.47 DEA schedule : 1	
References : Huffman JW, Dong D, <i>Bioorg Med Chem Letters</i> , 1994, 4(4), 563–566.		
Catalog number : 7118-001	CASRN : 209414-07-3	
Name : JWH-018		
Mol. formula : C ₂₄ H ₂₃ NO	FW : 341.45 DEA schedule : 1	
Notes : Non-selective CB ₁ /CB ₂ cannabinoid receptor agonist.		
References : Huffman J, Zengin G, Wu M, Lu J, Hynd G, Bushell K, Thompson A, Bushell S, et al., <i>Bioorg Med Chem</i> , 2005, 13(1), 89–112.		

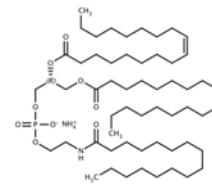
Catalog number : 7019-001**CASRN :** 209414-08-4**Name :** JWH-019**Mol. formula :** C₂₅H₂₅NO**FW :** 355.47 **DEA schedule :** 1**References :** Poso A, Huffman JW, *Br J Pharm*, 2008, 153(2), 335–346.**Catalog number :** 7173-001**CASRN :** 208987-48-8**Name :** JWH-073**Mol. formula :** C₂₃H₂₁NO**FW :** 327.42 **DEA schedule :** 1**Notes :** Non-selective CB₁/CB₂ cannabinoid receptor agonist.**References :** Wiley, JL; et al., *J Pharmacol Exp Ther* 1998, 285, 995–1004.**Catalog number :** 7081-001**CASRN :** 210179-46-7**Name :** JWH-081**Mol. formula :** C₂₅H₂₅NO₂**FW :** 371.47 **DEA schedule :** 1**Notes :** Non-selective CB₁/CB₂ cannabinoid receptor agonist.**References :** Huffman J, Zengin G, Wu M, Lu J, Hynd G, Bushell K, Thompson A, Bushell S, et al., *Bioorg Med Chem*, 2005, 13(1), 89–112.**Catalog number :** 7098-001**CASRN :** 316189-74-9**Name :** JWH-098**Mol. formula :** C₂₆H₂₇NO₂**FW :** 385.50 **DEA schedule :** 1**References :** Huffman JW, Zengin G, Wu MJ, Lu J, Hynd G, Bushell K, Thompson ALS, Bushell S, Tartal C, Hurst DP, Reggio PH, Selley DE, Cassidy MP, Wiley JL, Martin BR, *Bioorg Med Chem*, 2005, 13, 89–112.**Catalog number :** 7122-001**CASRN :** 619294-47-2**Name :** JWH-122**Mol. formula :** C₂₅H₂₅NO**FW :** 355.47 **DEA schedule :** 1**References :** Huffman J, Zengin G, Wu M, Lu J, Hynd G, Bushell K, Thompson A, Bushell S, et al., *Bioorg Med Chem*, 2005, 13(1), 89–112.**Catalog number :** 7149-001**CASRN :** 548461-82-1**Name :** JWH-149**Mol. formula :** C₂₆H₂₇NO**FW :** 369.50 **DEA schedule :** 1**References :** Huffman J, Zengin G, Wu M, Lu J, Hynd G, Bushell K, Thompson A, Bushell S, et al., *Bioorg Med Chem*, 2005, 13(1), 89–112.**Catalog number :** 7181-001**CASRN :** 824960-03-4**Name :** JWH-181**Mol. formula :** C₂₈H₃₁NO**FW :** 397.55 **DEA schedule :** 1**Notes :** Non-selective CB₁/CB₂ cannabinoid receptor agonist.**References :** Huffman J, Zengin G, Wu M, Lu J, Hynd G, Bushell K, Thompson A, Bushell S, et al., *Bioorg Med Chem*, 2005, 13(1), 89–112.

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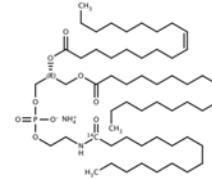
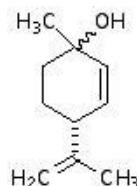
Catalog number : 7182-001	CASRN : 824960-02-3	
Name : JWH-182		
Mol. formula : C ₂₇ H ₂₉ NO	FW : 383.53	DEA schedule : 1
Notes : Non-selective CB ₁ /CB ₂ cannabinoid receptor agonist.		
References : Huffman J, Zengin G, Wu M, Lu J, Hynd G, Bushell K, Thompson A, Bushell S, et al., <i>Bioorg Med Chem</i> , 2005, 13(1), 89–112.		
Catalog number : 7193-001	CASRN : 133438-58-1	
Name : JWH-193		
Mol. formula : C ₂₆ H ₂₆ N ₂ O ₂	FW : 398.50	DEA schedule : 1
References : Huffman JW, Padgett LW, Curr Med Chem, 2005, 12, 1395–1411.		
Catalog number : 7200-001	CASRN : 103610-04-4	
Name : JWH-200; WIN 55,225		
Mol. formula : C ₂₄ H ₂₅ N ₂ O ₂	FW : 384.47	DEA schedule : 1
References : Dutta AK, Ryan W, Thomas BF, Singer M, Compton DR, Martin BR, Razdan RK, <i>Bioorg Med Chem</i> , 1997, 5(8), 1591–1600.		
Catalog number : 7210-001	CASRN : 824959-81-1	
Name : JWH-210		
Mol. formula : C ₂₆ H ₂₇ NO	FW : 369.51	DEA schedule : 1
Notes : Non-selective CB ₁ /CB ₂ cannabinoid receptor agonist.		
References : Huffman J, Zengin G, Wu M, Lu J, Hynd G, Bushell K, Thompson A, Bushell S, et al., <i>Bioorg Med Chem</i> , 2005, 13(1), 89–112.		
Catalog number : 7213-001	CASRN : 824959-83-3	
Name : JWH-213		
Mol. formula : C ₂₇ H ₂₉ NO	FW : 383.53	DEA schedule : 1
Notes : Non-selective CB ₁ /CB ₂ cannabinoid receptor agonist.		
References : Huffman J, Zengin G, Wu M, Lu J, Hynd G, Bushell K, Thompson A, Bushell S, et al., <i>Bioorg Med Chem</i> , 2005, 13(1), 89–112.		
Cannabinoids: Precursors, Biosynthetic		
Catalog number : NOCD-009		
Name : N-[1- ¹⁴ C]-Arachidonyl-1,2-dioleoyl-sn-glycero-3-phosphoethanolamine ammonium salt		
Mol. formula : C ₆₁ H ₁₁₁ N ₂ O ₉ P	FW : 1047.52	DEA schedule : 0
Notes : Putative biosynthetic precursor of endogenous cannabinoids (carbon-labeled).		
References : Morishita, J; et al., <i>J Neurochem</i> 2005, 94, 753–62.		

Catalog number : NOCD-066

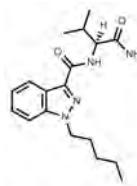
Name : N-Palmitoyl-1,2-dioleoyl-*sn*-glycero-3-phosphoethanolamine ammonium salt

Mol. formula : C₅₇H₁₁₁N₂O₉P**FW :** 999.47 **DEA schedule :** 0**Notes :** Putative biosynthetic precursor of endogenous cannabinoids.**References :** Morishita, J; et al. *J Neurochem* 2005, 94, 753–62.**Catalog number :** NOCD-000

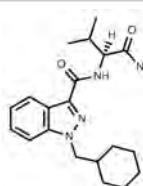
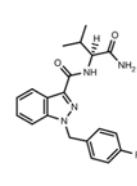
Name : N-[1-¹⁴C]-Palmitoyl-1,2-dioleoyl-*sn*-glycero-3-phosphoethanolamine ammonium salt

Mol. formula : C₅₇H₁₁₁N₂O₉P**FW :** 999.47 **DEA schedule :** 0**Notes :** Putative biosynthetic precursor of endogenous cannabinoids (carbon-labeled).**References :** Morishita J; et al. *J Neurochem* 2005, 94, 753–62.**Cannabinoids: Precursors, Synthetic****Catalog number :** NOCD-093**CASRN :** 52154-82-2**Name :** *p*-Mentha-2,8-dien-1-ol**Mol. formula :** C₁₀H₁₆O**FW :** 152.2 **DEA schedule :** 0**Notes :** Cannabinoid synthetic precursor**References :** Razdan RK; Dalzell HC; Handrick GR *J Am Chem Soc* 1974, 96, 5860–5865.**Cannabinoids: Pyrazole Class****Catalog number :** 7023-001

new

CASRN : 1445752-09-9**Name :** AB-PINACA**Mol. formula :** C₁₈H₂₆N₄O₂**FW :** 330.43 **DEA schedule :** 1**References :** Uchiyama, N., Matsuda, S., Wakana, D., et al., *Forensic Toxicol.*, 2012, 31(1), 93–100.**Catalog number :** 7031-001

new

CASRN : 1185887-21-1**Name :** AB-CHMINACA**Mol. formula :** C₂₀H₂₈N₄O₂**FW :** 356.47 **DEA schedule :** 1**References :** Uchiyama, N., Matsuda, S., Wakana, D., et al., *Forensic Toxicol.*, 2012, 31(1), 93–100.**Catalog number :** 7012-001**CASRN :** 1185282-01-2**Name :** AB-FUBINACA**Mol. formula :** C₂₀H₂₁N₄O₂**FW :** 368.41 **DEA schedule :** 1**Notes :** CB₁ receptor agonist.**References :** Uchiyama, N., Matsuda, S., Wakana, D., et al., *Forensic Toxicol.*, 2012, 31(1), 93–100.

2 – Cannabinoids and Cannabinoid-related

Catalog number : NOCD-082

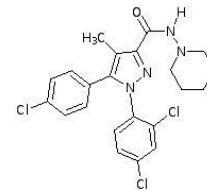
Name : SR141716

Mol. formula : C₂₂H₂₁Cl₃N₄O

FW : 463.78 **DEA schedule :** 0

Notes : *Cannabinoid CB1 receptor ligand*

References : Seltzman, H; et al. *J Chem Soc, Chem Commun* 1995, 1549–1550.



Catalog number : NOCD-083

CASRN : 170937-38-9

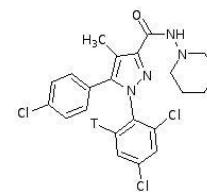
Name : [2,4-Dichlorophenyl-6-³H]SR141716

Mol. formula : C₂₂H₂₁Cl₃N₄O

FW : 465.80 **DEA schedule :** 0

Notes : *Cannabinoid CB1 receptor radioligand (tritium-labeled)*

References : Seltzman, H; et al. *J Chem Soc, Chem Commun* 1995, 1549–1550.



Catalog number : NOCD-085

CASRN : 192703-06-3

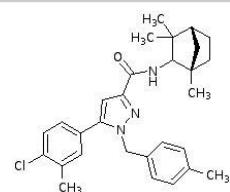
Name : SR144528

Mol. formula : C₂₉H₃₄ClN₃O

FW : 476.05 **DEA schedule :** 0

Notes : *Cannabinoid CB2 receptor antagonist*

References : Portier, M; et al. *J Pharmacol Exp Ther* 1999, 288, 582–9.
Rinaldi-Carmona, M; et al. *J Pharmacol Exp Ther* 1998, 284, 644–50.



Catalog number : NOCD-086

CASRN : 475471-24-0

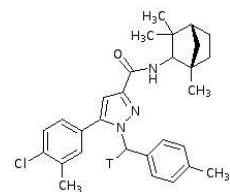
Name : Tritium-labeled SR144528

Mol. formula : C₂₉H₃₄N₃OCl

FW : 478.05 **DEA schedule :** 0

Notes : *Cannabinoid CB2 receptor radioligand (tritium-labeled)*

References : Portier, M; et al. *J Pharmacol Exp Ther* 1999, 288, 582–9.
Rinaldi-Carmona, M; et al. *J Pharmacol Exp Ther* 1998, 284, 644–50.



Cannabinoids: Tetrahydrocannabinol Class

Catalog number : 7370-042

CASRN : 137945-48-3

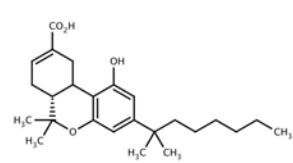
Name : Ajulemic acid; IP-751

Mol. formula : C₂₅H₃₆O₄

FW : 400.55 **DEA schedule :** 1

Notes : *Non-psych psychoactive cannabinoid*.

References : Burstein, S *AAPS J* 2005, 7, E143–8.
Wiley, JL *IDrugs* 2005, 8, 1002–11.



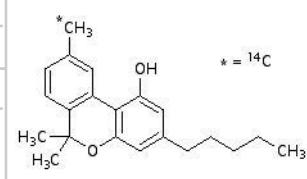
Catalog number : 7360-015

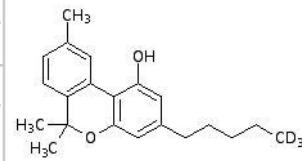
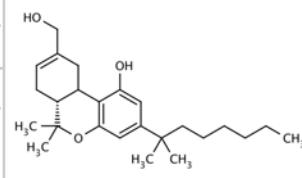
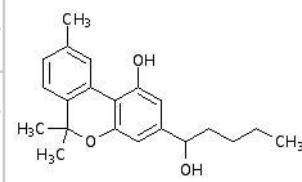
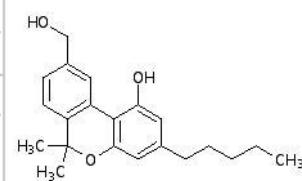
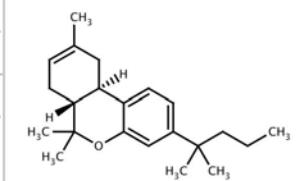
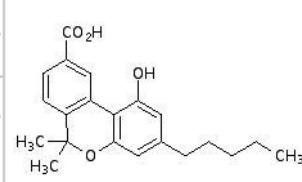
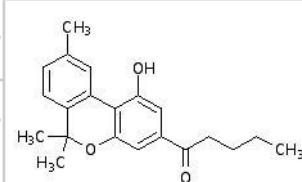
Name : [11-¹⁴C]Cannabinol

Mol. formula : C₂₀H₂₆O₂

FW : 310.43 **DEA schedule :** 1

Notes : *Inactive constituent of cannabis (carbon-labeled)*



Catalog number : 7360-014**Name :** [5'-²H₃]Cannabinol**Mol. formula :** C₂₁H₂₆O₂**FW :** 310.43 **DEA schedule :** 1**Notes :** Inactive constituent of cannabis (deuterium-labeled).**Catalog number :** 7370-057**CASRN :** 112830-95-2**Name :** HU 210**Mol. formula :** C₂₅H₃₈O₃**FW :** 386.57 **DEA schedule :** 1**References :** Ottani, A; Giuliani, D *CNS Drug Rev* 2001, 7, 131–45.**Catalog number :** 7360-017**Name :** 1'-Hydroxycannabinol**Mol. formula :** C₂₁H₂₆O₃**FW :** 326.43 **DEA schedule :** 1**Catalog number :** 7360-016**Name :** 11-Hydroxycannabinol**Mol. formula :** C₂₁H₂₆O₃**FW :** 326.43 **DEA schedule :** 1**Catalog number :** 7133-001**CASRN :** 259869-55-1**Name :** JWH-133**Mol. formula :** C₂₂H₃₂O**FW :** 312.49 **DEA schedule :** 1**Notes :** Selective CB₂ receptor agonist.**References :** Huffman, JW; et al., *Bioorganic & Medicinal Chemistry* 1999, 7, 2905–2914.**Catalog number :** 7360-019**CASRN :** 53989-32-5**Name :** 9-Carboxy-11-norcannabinol**Mol. formula :** C₂₁H₂₄O₄**FW :** 340.41 **DEA schedule :** 1**Catalog number :** 7360-018**Name :** 1'-Oxocannabinol**Mol. formula :** C₂₁H₂₄O₃**FW :** 324.41 **DEA schedule :** 1

2 – Cannabinoids and Cannabinoid-related

Catalog number : 7370-023

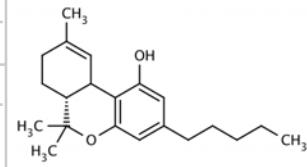
CASRN : 1972-08-3

Name : $(-)\Delta^9$ -Tetrahydrocannabinol; THC

Mol. formula : $C_{21}H_{30}O_2$

FW : 314.46 **DEA schedule :** 1

References : *Merck Index*, 14th ed., Monograph 9209.



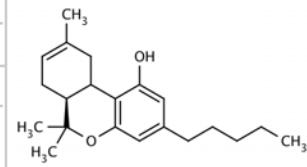
Catalog number : 7370-053

CASRN : 81586-39-2

Name : $(+)-\Delta^8$ -THC

Mol. formula : $C_{21}H_{30}O_2$

FW : 314.46 **DEA schedule :** 1



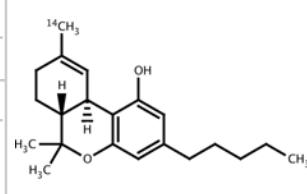
Catalog number : 7370-033

Name : $[11-^{14}C]\Delta^9$ -THC

Mol. formula : $C_{21}H_{30}O_2$

FW : 316.45 **DEA schedule :** 1

Notes : Hallucinogen; psychotropic; analgesic (carbon-labeled).



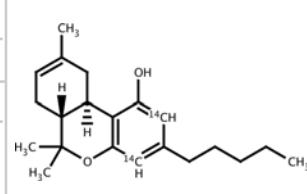
Catalog number : 7370-032

Name : $[2,4-^{14}C]\Delta^8$ -THC

Mol. formula : $C_{21}H_{30}O_2$

FW : 318.45 **DEA schedule :** 1

Notes : Hallucinogen; psychotropic; analgesic (carbon-labeled).

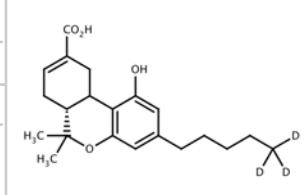


Catalog number : 7370-038

Name : $[5'-^2H_3]-11-nor-\Delta^8$ -THC-9-carboxylic acid

Mol. formula : $C_{21}H_{28}O_4$

FW : 347.46 **DEA schedule :** 1

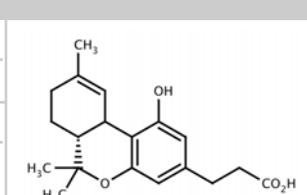


Catalog number : 7370-028

Name : 2'-Carboxy-3',4',5'-*trinor*- Δ^9 -THC

Mol. formula : $C_{19}H_{24}O_4$

FW : 316.39 **DEA schedule :** 1

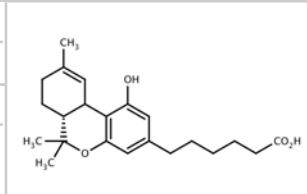


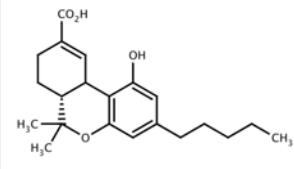
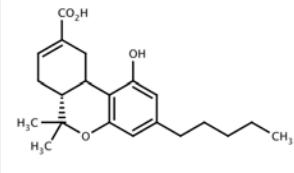
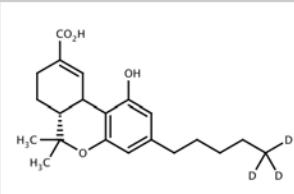
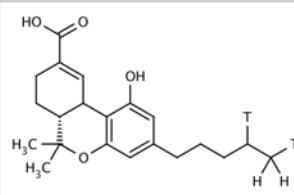
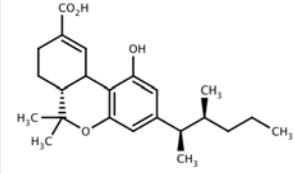
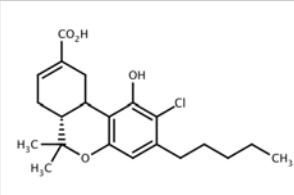
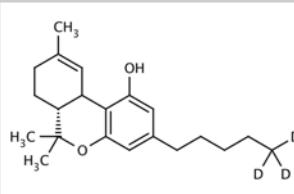
Catalog number : 7370-055

Name : 5'-Carboxy- Δ^9 -THC

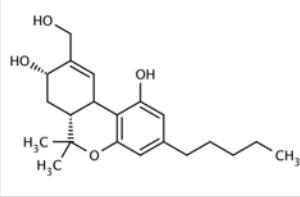
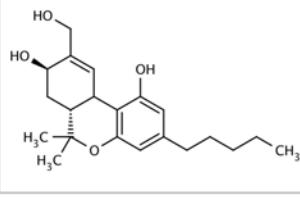
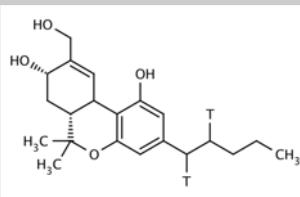
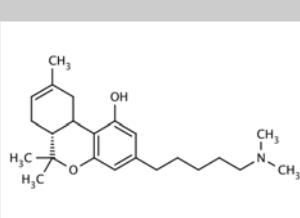
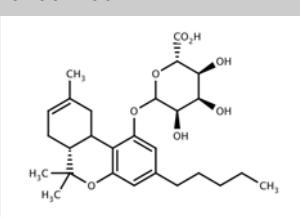
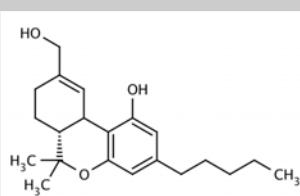
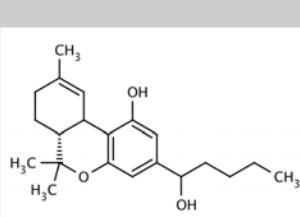
Mol. formula : $C_{22}H_{30}O_4$

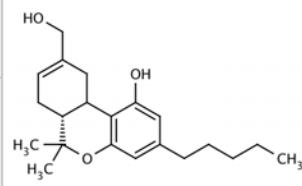
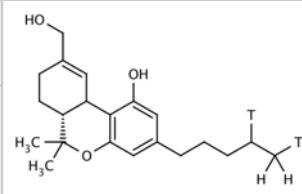
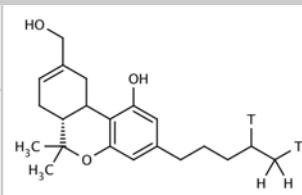
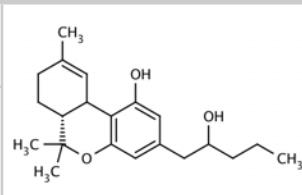
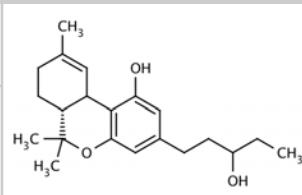
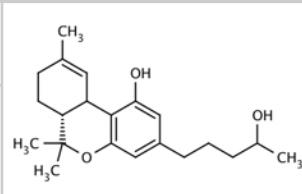
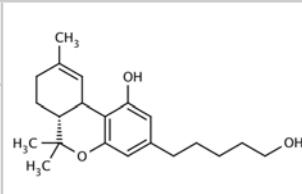
FW : 358.47 **DEA schedule :** 1



Catalog number : 7370-020**CASRN :** 56354-06-4**Name :** 9-Carboxy-11-nor- Δ^9 -THC**Mol. formula :** C₂₁H₂₈O₄**FW :** 344**DEA schedule :** 1**Notes :** Urinary metabolite of THC**Catalog number :** 7370-037**CASRN :** 39690-06-7**Name :** 9-Carboxy-11-nor- Δ^8 -THC**Mol. formula :** C₂₁H₂₈O₄**FW :** 344**DEA schedule :** 1**Catalog number :** 7370-003**CASRN :** 113269-48-0**Name :** [5'-²H₃]9-Carboxy-11-nor- Δ^9 -THC**Mol. formula :** C₂₁H₂₈O₄**FW :** 347.46**DEA schedule :** 1**Notes :** Urinary metabolite of THC (deuterium-labeled).**Catalog number :** 7370-017**Name :** [4',5'-³H₂]9-Carboxy-11-nor- Δ^9 -THC**Mol. formula :** C₂₁H₂₈O₄**FW :** 348.46**DEA schedule :** 1**Notes :** Urinary metabolite of THC (tritium-labeled).**Catalog number :** 7370-040**CASRN :** 199388-13-1**Name :** (6aR,10aR)-3-[(1S,2R)-1,2-Dimethylheptyl]-6a,7,10,10a-tetrahydro-6,6,9-trimethyl-6H-dibenzo[b,d]pyran-1-ol**Mol. formula :** C₂₃H₃₂O₄**FW :** 372.50**DEA schedule :** 1**Catalog number :** 7370-048**Name :** 9-Carboxy-11-nor-(2 or 4)-chloro- Δ^8 -THC**Mol. formula :** C₂₁H₂₇ClO₄**FW :** 378.89**DEA schedule :** 1**Catalog number :** 7370-005**CASRN :** 81586-39-2**Name :** Deuterium-labeled Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 317**DEA schedule :** 1**Notes :** Hallucinogen; psychotropic; analgesic (deuterium-labeled).

2 – Cannabinoids and Cannabinoid-related

Catalog number : 7370-018	CASRN : 36913-21-0	
Name : 8α,11-Dihydroxy-Δ⁹-THC		
Mol. formula : C ₂₁ H ₃₀ O ₄	FW : 346.47	DEA schedule : 1
Catalog number : 7370-022	CASRN : 57030-51-0	
Name : 8β,11-Dihydroxy-Δ⁹-THC		
Mol. formula : C ₂₁ H ₃₀ O ₄	FW : 346.47	DEA schedule : 1
Catalog number : 7370-026		
Name : [1',2'-³H₂]-8α,11-Dihydroxy-Δ⁹-THC		
Mol. formula : C ₂₁ H ₃₀ O ₄	FW : 350.48	DEA schedule : 1
Catalog number : 7370-050		
Name : 5'-Dimethylamino-Δ⁸-THC		
Mol. formula : C ₂₃ H ₃₅ NO ₂	FW : 357.53	DEA schedule : 1
Catalog number : 7370-049	CASRN : 62667-60-1	
Name : Δ⁸-THC-O-glucuronide		
Mol. formula : C ₂₇ H ₃₈ O ₈	FW : 490.6	DEA schedule : 1
Catalog number : 7370-008	CASRN : 36557-05-08	
Name : 11-Hydroxy-Δ⁹-THC		
Mol. formula : C ₂₁ H ₃₀ O ₃	FW : 330.47	DEA schedule : 1
Catalog number : 7370-011		
Name : 1'-Hydroxy-Δ⁹-THC (Isomer B)		
Mol. formula : C ₂₁ H ₃₀ O ₃	FW : 330.46	DEA schedule : 1

Catalog number : 7370-034**Name :** 11-Hydroxy- Δ^8 -THC**Mol. formula :** C₂₁H₃₀O₃**FW :** 330.46 **DEA schedule :** 1**Catalog number :** 7370-024**Name :** [4',5'-³H₂]-11-Hydroxy- Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.48 **DEA schedule :** 1**Catalog number :** 7370-035**Name :** [4',5'-³H₂]-11-Hydroxy- Δ^8 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.48 **DEA schedule :** 1**Catalog number :** 7370-012**Name :** 2'-Hydroxy- Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₃**FW :** 330.46 **DEA schedule :** 1**Catalog number :** 7370-013**Name :** 3'-Hydroxy- Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₃**FW :** 330.46 **DEA schedule :** 1**Catalog number :** 7370-014**Name :** 4'-Hydroxy- Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₃**FW :** 330.46 **DEA schedule :** 1**Catalog number :** 7370-015**Name :** 5'-Hydroxy- Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₃**FW :** 330.46 **DEA schedule :** 1

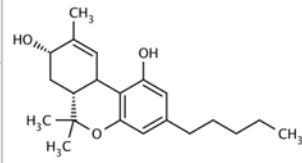
2 – Cannabinoids and Cannabinoid-related

Catalog number : 7370-016

Name : 8 α -Hydroxy- Δ^9 -THC

Mol. formula : C₂₁H₃₀O₃

FW : 330.46 **DEA schedule :** 1



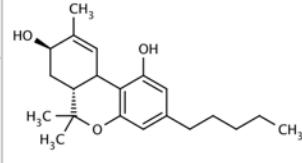
Catalog number : 7370-019

CASRN : 34984-78-6

Name : 8 β -Hydroxy- Δ^9 -THC

Mol. formula : C₂₁H₃₀O₃

FW : 330 **DEA schedule :** 1

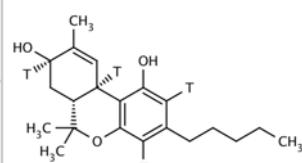


Catalog number : 7370-021

Name : [2,4,8,10a-³H₄]-8 β -Hydroxy- Δ^9 -THC

Mol. formula : C₂₁H₃₀O₃

FW : 338.49 **DEA schedule :** 1



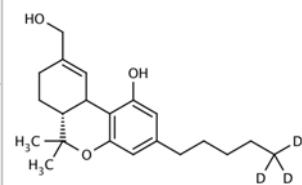
Catalog number : 7370-010

CASRN : 130410-26-3

Name : [5'-²H₃]-11-Hydroxy- Δ^9 -THC

Mol. formula : C₂₁H₃₀O₃

FW : 333 **DEA schedule :** 1

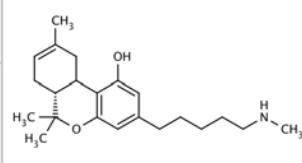


Catalog number : 7370-054

Name : 5'-Methylamino- Δ^8 -THC

Mol. formula : C₂₂H₃₂NO₂

FW : 342.49 **DEA schedule :** 1

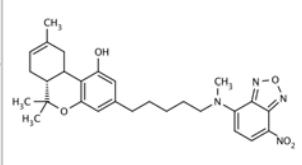


Catalog number : 7370-036

Name : 5'-N-Methyl-N-4-(7-nitrobenzofurazano)amino- Δ^8 -THC

Mol. formula : C₂₈H₃₄N₄O₅

FW : 506.59 **DEA schedule :** 1



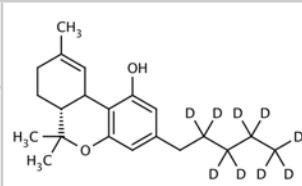
Catalog number : 7370-025

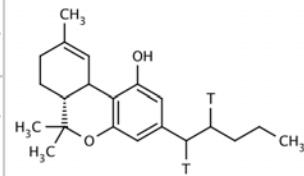
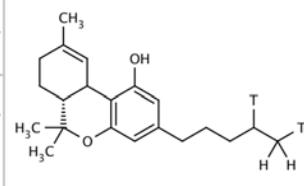
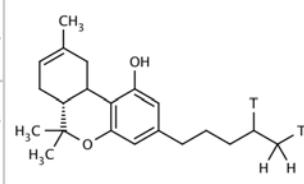
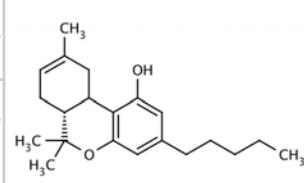
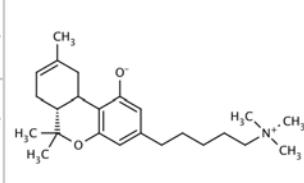
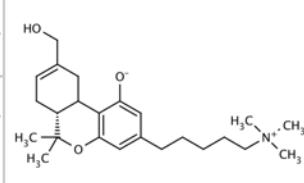
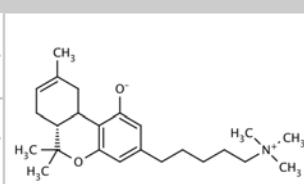
Name : [2',2',3',3',4',4',5',5',5'-²H₉]- Δ^9 -THC

Mol. formula : C₂₁H₃₀O₂

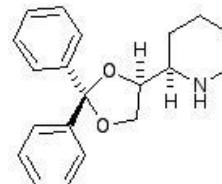
FW : 323.52 **DEA schedule :** 1

Notes : Hallucinogen; psychotropic; analgesic (deuterium-labeled).

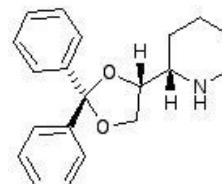


Catalog number : 7370-004**Name :** [1',2'-³H] Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.48 **DEA schedule :** 1**Notes :** Hallucinogen; psychotropic; analgesic (tritium-labeled).**Catalog number :** 7370-009**Name :** [4',5'-³H] Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.48 **DEA schedule :** 1**Catalog number :** 7370-027**Name :** [4',5'-³H] Δ^8 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.48 **DEA schedule :** 1**Notes :** Hallucinogen; psychotropic; analgesic (tritium-labeled).**Catalog number :** 7370-030**CASRN :** 5957-75-5**Name :** (-)-*trans*- Δ^8 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 314.46 **DEA schedule :** 1**Catalog number :** 7370-051**Name :** 5'-Trimethylammonium- Δ^8 -THC phenolate**Mol. formula :** C₂₄H₃₇NO₂**FW :** 371.56 **DEA schedule :** 1**Catalog number :** 7370-056**Name :** 5'-Trimethylammonium-11-hydroxy- Δ^8 -THC phenolate**Mol. formula :** C₂₄H₃₇NO₃**FW :** 388.56 **DEA schedule :** 1**Catalog number :** 7370-052**Name :** [³H]-5'-Trimethylammonium- Δ^8 -THC phenolate**Mol. formula :** C₂₄H₃₇NO₂**FW :** 371.56 **DEA schedule :** 1

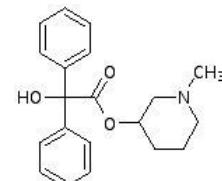
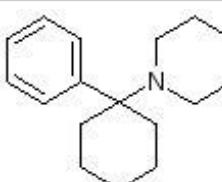
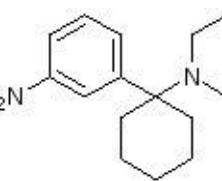
Cannabinoids (dosage form): Stock Solutions**Catalog number :** 7360-009**CASRN :** 21366-63-2**Name :** Cannabicyclol ampuls (1.0 mg/mL in absolute ethanol)**Mol. formula :** C₂₁H₃₀O₂**FW :** 314.46 **DEA schedule :** 1**Notes :** Non-psychoactive constituent of cannabis. Photodegradation product of cannabichromene.**References :** Burstein, S; Hunter, SA; Renzulli, LJ *Pharmacol Exp Ther* 1985, 235, 87-91.**Dosage
Form****Catalog number :** 7370-001**Name :** Δ⁹-THC ampuls (various concentrations in 95% ethanol)**DEA schedule :** 1**Dosage
Form**

Dissociatives: Dexoxadrol Class**Catalog number :** NOCD-068**CASRN :** 631-06-1**Name :** α -(+)-Dexoxadrol hydrochloride**Mol. formula :** $C_{20}H_{24}ClNO_2$ **FW :** 345.87**DEA schedule :** 0

References : Thurkauf, A; et al. *J Med Chem* 1988, 31, 2257-63.
Sax, M; Wunsch, B *Curr Top Med Chem* 2006, 6, 723-32.

Catalog number : NOCD-069**CASRN :** 4792-18-1**Name :** α -(-)-Levoxadrol hydrochloride**Mol. formula :** $C_{20}H_{24}ClNO_2$ **FW :** 345.87**DEA schedule :** 0

References : Thurkauf, A; et al. *J Med Chem* 1988, 31, 2257-63.
Sax, M; Wunsch, B *Curr Top Med Chem* 2006, 6, 723-32.

Catalog number : 7484-001**Name :** N-Methyl-3-piperidylbenzilate hydrochloride**Mol. formula :** $C_{20}H_{24}ClNO_3$ **FW :** 361.86**DEA schedule :** 1**Dissociatives: Phencyclidine Class****Catalog number :** 7471-004**CASRN :** 956-90-1**Name :** Phencyclidine hydrochloride; PCP HCl**Mol. formula :** $C_{17}H_{26}ClN$ **FW :** 284.36**DEA schedule :** 2**Notes :** CNS depressant; anesthetic, psychostimulant**References :** Merck Index, 14th ed., Monograph 7219.**Catalog number :** 7471-041**Name :** 1-[*m*-Aminophenyl]cyclohexylpiperidine lactate**Mol. formula :** $C_{20}H_{32}N_2O_3$ **FW :** 348.48**DEA schedule :** 1

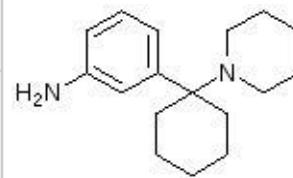
3 – Dissociatives

Catalog number : 7471-042

Name : 1-[1-(*m*-Aminophenyl)cyclohexyl]piperidine

Mol. formula : C₁₇H₂₆N₂

FW : 258.39 DEA schedule : 1

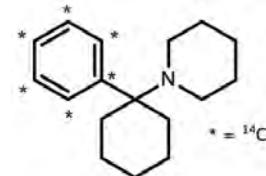


Catalog number : 7471-002

Name : [¹⁴C]Phencyclidine HBr; [¹⁴C]PCP

Mol. formula : C₁₈H₂₈BrN

FW : 523.53 DEA schedule : 2

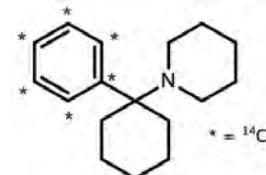


Catalog number : 7471-007

Name : [Phenyl-U-¹⁴C]-1-(1-Phenylcyclohexyl)Piperidine; [Phenyl-U-¹⁴C]PCP

Mol. formula : C₁₇H₂₅N

FW : 245.39 DEA schedule : 2

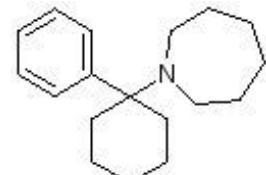


Catalog number : 7471-023

Name : 1-(1-Phenylcyclohexyl)hexamethyleneimine hydrochloride

Mol. formula : C₁₈H₂₈CIN

FW : 293.88 DEA schedule : 1

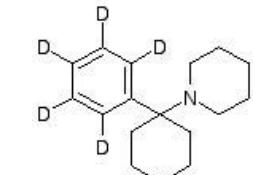


Catalog number : 7471-003

Name : [Phenyl-2,3,4,5,6-²H₅]Phencyclidine; [Phenyl-2,3,4,5,6-²H₅]PCP

Mol. formula : C₁₇H₂₅N

FW : 248.42 DEA schedule : 2

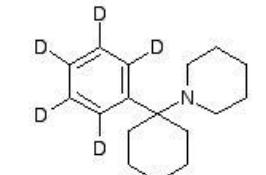


Catalog number : 7471-006

Name : [Phenyl-²H₅]Phencyclidine hydrochloride; [Phenyl-²H₅]PCP HCl

Mol. formula : C₁₇H₂₆CIN

FW : 284.89 DEA schedule : 2

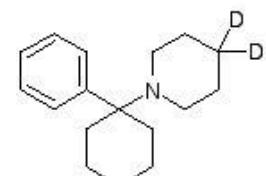


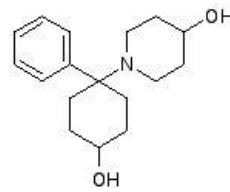
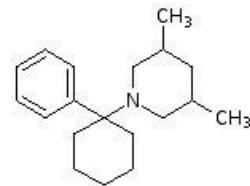
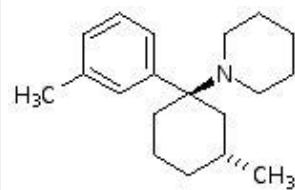
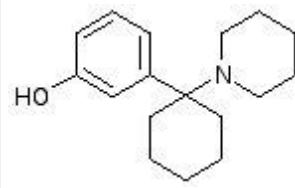
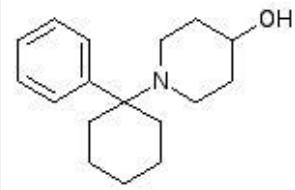
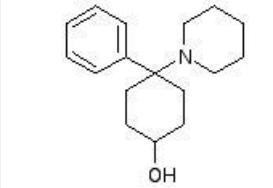
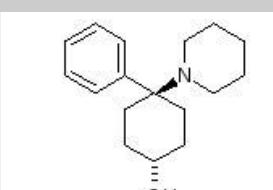
Catalog number : 7471-008

Name : [Piperidino-4,4-²H₂]Phencyclidine hydrochloride;
[Piperidino-4,4-²H₂]PCP

Mol. formula : C₁₇H₂₆CIN

FW : 281.86 DEA schedule : 2



Catalog number : 7471-087**Name :** 4-(4'-Hydroxypiperidino)-4-phenylcyclohexanol (*cis/trans*)**Mol. formula :** C₁₇H₂₅NO₂**FW :** 275.40 **DEA schedule :** 1**Catalog number :** 7471-019**Name :** 1-(1-Phenylcyclohexyl)-3,5-dimethylpiperidine hydrochloride**Mol. formula :** C₁₉H₃₀ClN**FW :** 307.91 **DEA schedule :** 1**Catalog number :** 7471-034**Name :** (±)-*cis*-1-[1-(*m*-Tolyl)-3-methylcyclohexyl]piperidine hydrochloride**Mol. formula :** C₁₉H₃₀ClN**FW :** 307.91 **DEA schedule :** 1**Catalog number :** 7471-033**CASRN :** 79787-43-2**Name :** 1-[1-(*m*-Hydroxyphenyl)cyclohexyl]piperidine hydrochloride**Mol. formula :** C₁₇H₂₆ClNO**FW :** 295.85 **DEA schedule :** 1**References :** Itzhak, Y; Kalir A; Sarne Y *Eur J Pharmacol* 1981, 73, 229-33.**Catalog number :** 7471-081**Name :** 1-(1-Phenylcyclohexyl)-4-hydroxypiperidine**Mol. formula :** C₁₇H₂₅NO**FW :** 259.40 **DEA schedule :** 1**Catalog number :** 7471-083**CASRN :** 60756-83-4**Name :** 1-(1-Phenyl-4-hydroxycyclohexyl)piperidine (*cis/trans*)**Mol. formula :** C₁₇H₂₅NO**FW :** 259.35 **DEA schedule :** 1**References :** Carroll, FI; et al. *J Med Chem* 1981, 24, 1047-51.
Martin, BR; Vincek WC; Balster RL *Subst Alcohol Actions Misuse* 1981, 2, 143-7.**Catalog number :** 7471-085**Name :** *cis*-1-(1-Phenyl-4-hydroxycyclohexyl)piperidine hydrochloride**Mol. formula :** C₁₇H₂₆ClNO**FW :** 295.87 **DEA schedule :** 1**References :** Carroll, FI; et al. *J Med Chem* 1981, 24, 1047-51.

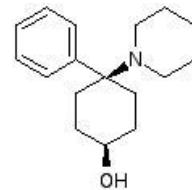
3 – Dissociatives

Catalog number : 7471-086

Name : *trans*-1-(1-Phenyl-4-hydroxycyclohexyl)piperidine hydrochloride

Mol. formula : C₁₇H₂₆CINO

FW : 295.87 DEA schedule : 1



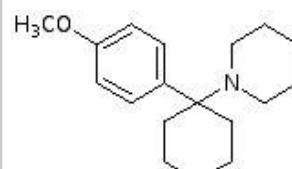
References : Carroll, FI; et al. *J Med Chem* 1981, 24, 1047-51.

Catalog number : 7471-028

Name : 1-[1-(*p*-Methoxyphenyl)cyclohexyl]piperidine hydrochloride

Mol. formula : C₁₈H₂₈CINO

FW : 309.88 DEA schedule : 1

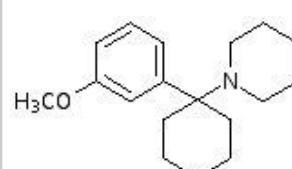


Catalog number : 7471-039

Name : 1-[1-(*m*-Methoxyphenyl)cyclohexyl]piperidine hydrochloride

Mol. formula : C₁₈H₂₈CINO

FW : 309.88 DEA schedule : 1

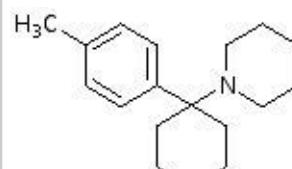


Catalog number : 7471-022

Name : 1-[1-(*p*-Tolyl)cyclohexyl]piperidine hydrochloride

Mol. formula : C₁₈H₂₈CIN

FW : 293.88 DEA schedule : 1



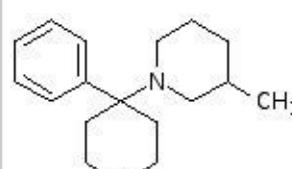
Catalog number : 7471-024

CASRN : 2201-41-4

Name : (\pm)-1-(1-Phenylcyclohexyl)-3-methylpiperidine hydrochloride

Mol. formula : C₁₈H₂₈CIN

FW : 293.88 DEA schedule : 1



References : Berry, SC; et al. *Eur J Pharmacol* 1983, 96, 261-7.

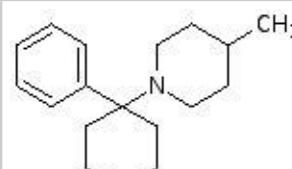
Catalog number : 7471-025

CASRN : 1934-50-5

Name : 1-(1-Phenylcyclohexyl)-4-methylpiperidine hydrochloride

Mol. formula : C₁₈H₂₈CIN

FW : 293.88 DEA schedule : 1

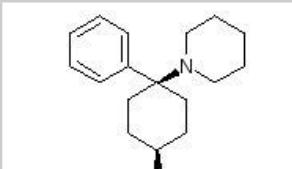


Catalog number : 7471-029

Name : *trans*-1-(1-Phenyl-4-methylcyclohexyl)piperidine hydrochloride

Mol. formula : C₁₈H₂₈CIN

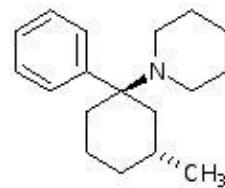
FW : 293.88 DEA schedule : 1



Catalog number : 7471-030

Name : (\pm)-*cis*-(1-Phenyl-3-methylcyclohexyl)piperidine hydrochlorideMol. formula : C₁₈H₂₈CIN

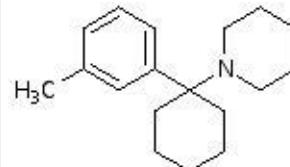
FW : 293.88 DEA schedule : 1



Catalog number : 7471-035

Name : 1-[1-(*m*-Tolyl)cyclohexyl]piperidine hydrochlorideMol. formula : C₁₈H₂₈CIN

FW : 293.88 DEA schedule : 1

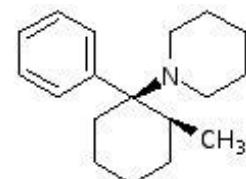


Catalog number : 7471-036

CASRN : 59397-29-4

Name : (\pm)-*trans*-1-(1-Phenyl-2-methylcyclohexyl)piperidine hydrochloride; 1-PMCPPMol. formula : C₁₈H₂₈CIN

FW : 293.88 DEA schedule : 1

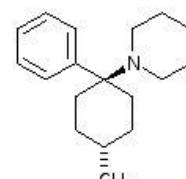
References : Iorio, MA; et al. *J Med Chem* 1991, 34, 2615–23.

Catalog number : 7471-040

CASRN : 21602-54-0

Name : *cis*-1-(1-Phenyl-4-methylcyclohexyl)piperidine hydrochlorideMol. formula : C₁₈H₂₈CIN

FW : 293.88 DEA schedule : 1



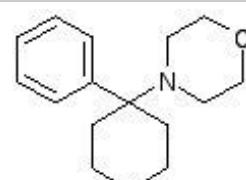
Catalog number : 7471-010

CASRN : 2201-40-3

Name : 1-(1-Phenylcyclohexyl)morpholine hydrochloride; PCM hydrochloride

Mol. formula : C₁₆H₂₄CINO

FW : 281.83 DEA schedule : 1

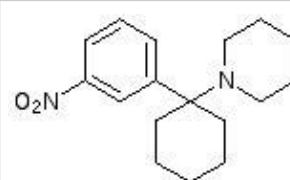
References : Brine, GA; et al. *J Heterocyclic Chem* 1979, 16, 1425.

Catalog number : 7471-026

CASRN : 70227-29-1

Name : 1-[1-(*m*-Nitrophenyl)cyclohexyl]piperidineMol. formula : C₁₇H₂₄N₂O₂

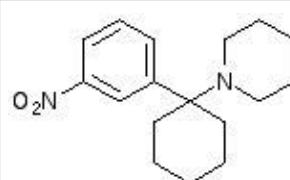
FW : 288.38 DEA schedule : 1

References : Ramoa, AS; Albuquerque EX *FEBS Lett* 1988, 235, 156–62.

Catalog number : 7471-027

Name : 1-[1-(*m*-Nitrophenyl)cyclohexyl]piperidine hydrochlorideMol. formula : C₁₇H₂₅CIN₂O₂

FW : 324.85 DEA schedule : 1

References : Ramoa, AS; Albuquerque EX *FEBS Lett* 1988, 235, 156–62.

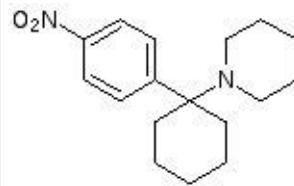
3 – Dissociatives

Catalog number : 7471-031

Name : 1-[1-(*p*-Nitrophenyl)cyclohexyl]piperidine

Mol. formula : C₁₇H₂₄N₂O₂

FW : 288.38 DEA schedule : 1

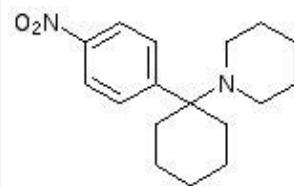


Catalog number : 7471-032

Name : 1-[1-(*p*-Nitrophenyl)cyclohexyl]piperidine hydrochloride

Mol. formula : C₁₇H₂₅ClN₂O₂

FW : 324.85 DEA schedule : 1

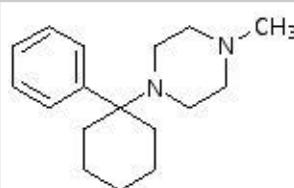


Catalog number : 7471-021

Name : 1-(1-Phenylcyclohexyl)-N'-methylpiperazine

Mol. formula : C₁₇H₂₆N₂

FW : 258.41 DEA schedule : 1

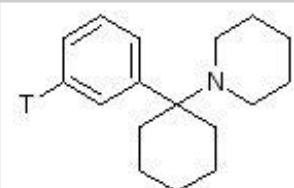


Catalog number : 7471-001

Name : [Phenyl-3-³H(n)]Phencyclidine; [Phenyl-3-³H(n)]PCP

Mol. formula : C₁₇H₂₅N

FW : 245.40 DEA schedule : 2

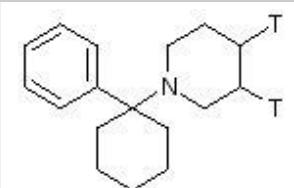


Catalog number : 7471-043

Name : [3',4'-³H]Phencyclidine; [3,4-³H]PCP

Mol. formula : C₁₇H₂₅N

FW : 247.4 DEA schedule : 2

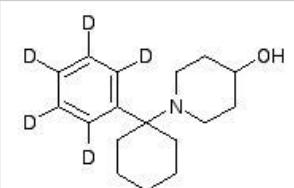


Catalog number : 7471-082

Name : [Phenyl-²H₅]-1-(1-Phenylcyclohexyl)-4-hydroxypiperidine

Mol. formula : C₁₇H₂₅NO

FW : 264.43 DEA schedule : 1

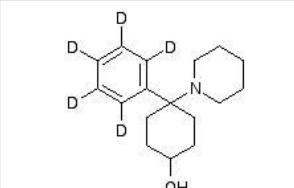


Catalog number : 7471-084

Name : [Phenyl-²H₅]-1-(1-Phenyl-4-hydroxycyclohexyl)piperidine (*cis/trans*)

Mol. formula : C₁₇H₂₅NO

FW : 264.43 DEA schedule : 1



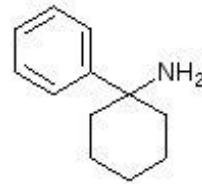
Catalog number : 7460-001

CASRN : 1934-71-0

Name : 1-Phenylcyclohexylamine hydrochloride; PCA hydrochloride

Mol. formula : C₁₂H₁₈CIN

FW : 211.74 DEA schedule : 1

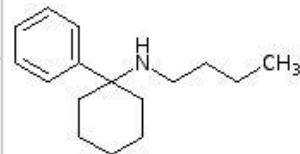
References : Brine, GA; et al. *J Heterocyclic Chem* 1979, 16, 1425.Blake, PA; Yamaguchi, S; Thurkauf, A; Rogawski, MA *Epilepsia* 1992, 33, 188-94.

Catalog number : 7460-005

Name : N-Butyl-1-phenylcyclohexylamine hydrochloride

Mol. formula : C₁₆H₂₆CIN

FW : 267.85 DEA schedule : 1

References : Brine, GA; et al. *J Heterocyclic Chem* 1979, 16, 1425.

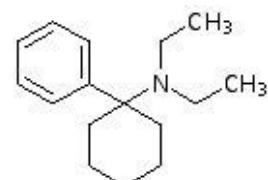
Catalog number : 7455-003

CASRN : 2201-19-6

Name : N,N-Diethyl-1-phenylcyclohexylamine hydrochloride; PCDE hydrochloride

Mol. formula : C₁₆H₂₆CIN

FW : 267.84 DEA schedule : 1

References : Brine, GA; et al. *J Heterocyclic Chem* 1979, 16, 1425.

Catalog number : 7455-002

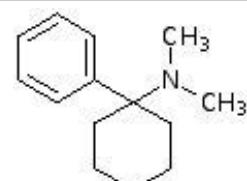
CASRN : 2201-17-4

Name : N,N-Dimethyl-1-phenylcyclohexylamine hydrochloride

Mol. formula : C₁₄H₂₂CIN

FW : 239.82 DEA schedule : 1

Notes : Current designer drug?

References : Brine, GA; et al. *J Heterocyclic Chem* 1979, 16, 1425.

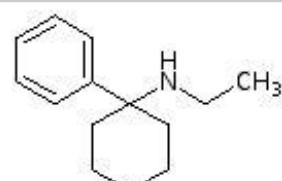
Catalog number : 7455-001

CASRN : 2201-15-2

Name : N-Ethyl-1-phenylcyclohexylamine hydrochloride; PCD hydrochloride

Mol. formula : C₁₄H₂₂CIN

FW : 239.79 DEA schedule : 1

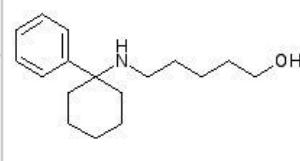
References : Brine, GA; et al. *J Heterocyclic Chem* 1979, 16, 1425.
Brady, KT; Balster, RL; Meltzer, LT; Schwertz, D *Pharmacol Biochem Behav* 1980, 12, 67-71.

Catalog number : 7471-089

Name : N-(5-Hydroxypentyl)-1-phenylcyclohexylamine hydrochloride

Mol. formula : C₁₇H₂₈CINO

FW : 297.87 DEA schedule : 0



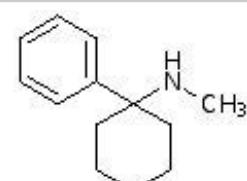
Catalog number : 7460-002

CASRN : 2201-16-3

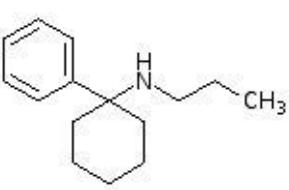
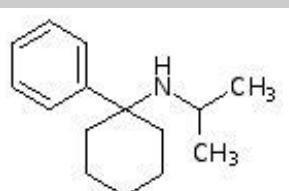
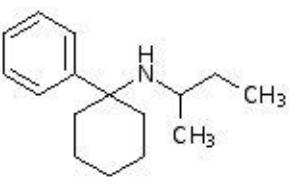
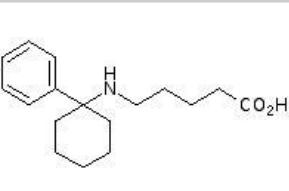
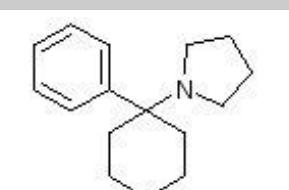
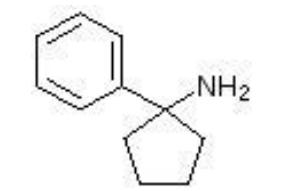
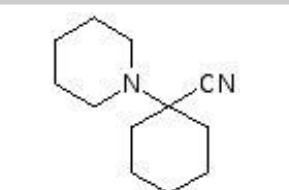
Name : N-Methyl-1-phenylcyclohexylamine hydrochloride

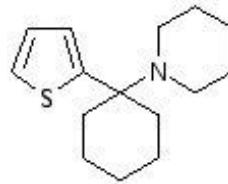
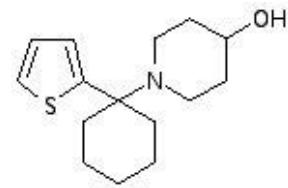
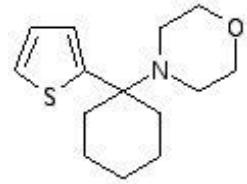
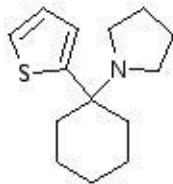
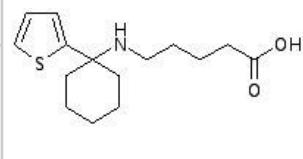
Mol. formula : C₁₃H₂₀CIN

FW : 225.76 DEA schedule : 1

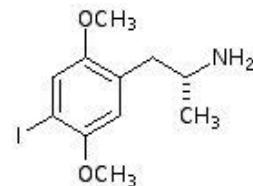
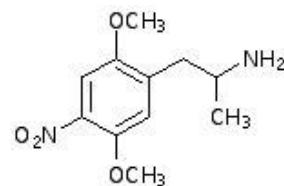
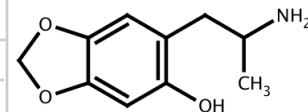
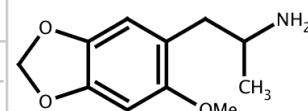
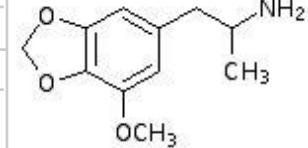
References : Brine, GA; et al. *J Heterocyclic Chem* 1979, 16, 1425.

3 – Dissociatives

Catalog number : 7460-003	CASRN : 18949-81-0	
Name : N-Propyl-1-phenylcyclohexylamine hydrochloride		
Mol. formula : C ₁₅ H ₂₄ ClN	FW : 253.82	DEA schedule : 1
		
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1979, 16, 1425.		
Catalog number : 7460-004		
Name : N-(<i>i</i> -Propyl)-1-phenylcyclohexylamine hydrochloride		
Mol. formula : C ₁₅ H ₂₄ ClN	FW : 253.82	DEA schedule : 1
		
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1979, 16, 1425.		
Catalog number : 7460-006		
Name : N-(<i>s</i> -Butyl)-1-phenylcyclohexylamine hydrochloride		
Mol. formula : C ₁₆ H ₂₆ ClN	FW : 267.85	DEA schedule : 1
		
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1979, 16, 1425.		
Catalog number : 7471-088		
Name : 5-[N-(1'-Phenylcyclohexyl)amino]pentanoic acid hydrochloride		
Mol. formula : C ₁₇ H ₂₆ ClNO ₂	FW : 311.86	DEA schedule : 0
		
Catalog number : 7458-001	CASRN : 2201-39-0	
Name : 1-(1-Phenylcyclohexyl)pyrrolidine hydrochloride; PCPy; Rolicyclidine		
Mol. formula : C ₁₆ H ₂₄ ClN	FW : 265.83	DEA schedule : 1
		
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1979, 16, 1425. Budd, RD <i>N Engl J Med</i> 1980, 303, 588.		
Catalog number : 7460-012	CASRN : 17380-74-4	
Name : 1-Phenylcyclopentylamine hydrochloride; PPA hydrochloride		
Mol. formula : C ₁₁ H ₁₆ ClN	FW : 197.71	DEA schedule : 1
		
References : Blake, PA; Yamaguchi, S; Thurkauf, A; Rogawski, MA <i>Epilepsia</i> 1992, 33, 188-94.		
Catalog number : 8603-001		
Name : 1-Piperidinocyclohexanecarbonitrile		
Mol. formula : C ₁₂ H ₂₀ N ₂	FW : 192.31	DEA schedule : 1
		

Catalog number : 7470-001**CASRN :** 1867-65-8**Name :** 1-[1-(2-Thienyl)cyclohexyl]piperidine hydrochloride; TCP hydrochloride**Mol. formula :** C₁₅H₂₄CINS**FW :** 285.88 **DEA schedule :** 1**References :** Brine, GA; et al. *J Heterocyclic Chem* 1979, 16, 1425.Vignon, J; et al. *Brain Res* 1983, 280, 194-7.**Catalog number :** 7470-002**Name :** 1-[1-(2-Thienyl)cyclohexyl]-4-hydroxypiperidine**Mol. formula :** C₁₅H₂₃NOS**FW :** 265.42 **DEA schedule :** 1**Catalog number :** 7471-013**Name :** 1-[1-(2-Thienyl)cyclohexyl]morpholine hydrochloride**Mol. formula :** C₁₄H₂₂CINOS**FW :** 287.85 **DEA schedule :** 1**References :** Brine, GA; et al. *J Heterocyclic Chem* 1979, 16, 1425.**Catalog number :** 7473-015**Name :** 1-[1-(2-Thienyl)cyclohexyl]pyrrolidine hydrochloride**Mol. formula :** C₁₄H₂₂CINS**FW :** 271.86 **DEA schedule :** 1**Catalog number :** 7470-003**Name :** 5-[N-[1'-(2-Thienyl)cyclohexyl]amino]pentanoic acid hydrochloride**Mol. formula :** C₁₅H₂₄CINO₂S**FW :** 317.88 **DEA schedule :** 0



Hallucinogens: Amphetamine Class**Catalog number :** 7396-011**CASRN :** 82864-06-0**Name :** (-)-2,5-Dimethoxy-4-iodoamphetamine hydrochloride**Mol. formula :** C₁₁H₁₇ClINO₂**FW :** 357.62 **DEA schedule :** 0**Catalog number :** 7396-009**Name :** (±)-2,5-Dimethoxy-4-nitroamphetamine hydrochloride**Mol. formula :** C₁₁H₁₇CIN₂O₄**FW :** 276.75 **DEA schedule :** 0**Catalog number :** 7402-002**CASRN :** 145284-65-7**Name :** (±)-6-Hydroxy-3,4-methylenedioxymethylamphetamine hydrochloride**Mol. formula :** C₁₀H₁₄CINO₃**FW :** 238.89 **DEA schedule :** 1**Notes :** CNS stimulant**Catalog number :** 7401-001**Name :** (±)-2-Methoxy-4,5-methylenedioxymethylamphetamine hydrochloride**Mol. formula :** C₁₁H₁₆CINO₃**FW :** 245.71 **DEA schedule :** 1**Catalog number :** 7401-003**Name :** (±)-3-Methoxy-4,5-methylenedioxymethylamphetamine hydrochloride**Mol. formula :** C₁₁H₁₆CINO₃**FW :** 245.71 **DEA schedule :** 1

4 – Hallucinogens

Catalog number : 7402-001

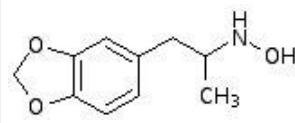
CASRN : 4764-17-4

Name : (\pm)-N-Hydroxy-3,4-methylenedioxymphetamine hydrochloride

Mol. formula : C₁₀H₁₄ClNO₃

FW : 231.68 DEA schedule : 1

Notes : CNS stimulant

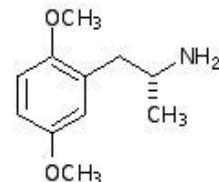


Catalog number : 7396-002

Name : (-)-2,5-Dimethoxyamphetamine hydrochloride; (-)-DMA

Mol. formula : C₁₁H₁₈ClNO₂

FW : 231.72 DEA schedule : 1

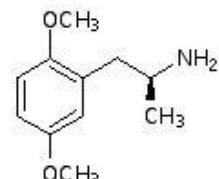


Catalog number : 7396-003

Name : (+)-2,5-Dimethoxyamphetamine hydrochloride; (+)-DMA

Mol. formula : C₁₁H₁₈ClNO₂

FW : 231.72 DEA schedule : 1

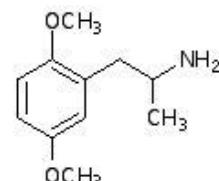


Catalog number : 7396-001

Name : (\pm)-2,5-Dimethoxyamphetamine hydrochloride; (\pm)-DMA

Mol. formula : C₁₁H₁₈ClNO₂

FW : 231.72 DEA schedule : 1

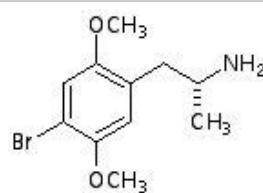


Catalog number : 7391-003

Name : (-)-4-Bromo-2,5-dimethoxyamphetamine hydrochloride; (-)-DOB

Mol. formula : C₁₁H₁₇BrClNO₂

FW : 310.62 DEA schedule : 1

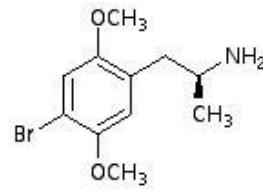


Catalog number : 7391-002

Name : (+)-4-Bromo-2,5-dimethoxyamphetamine hydrochloride; (+)-DOB

Mol. formula : C₁₁H₁₇BrClNO₂

FW : 310.62 DEA schedule : 1

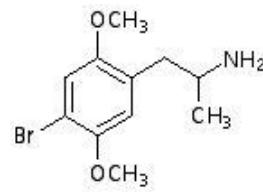


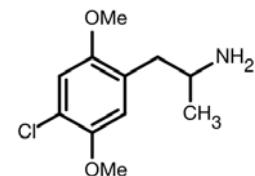
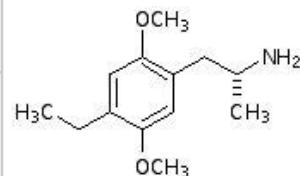
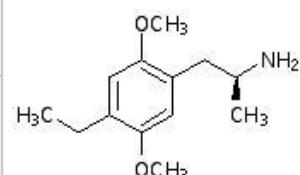
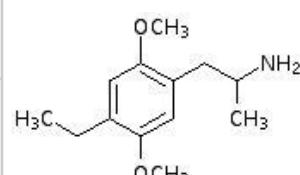
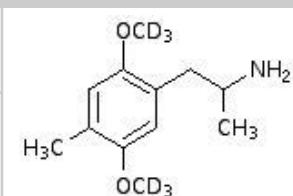
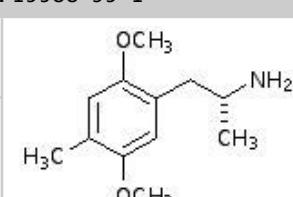
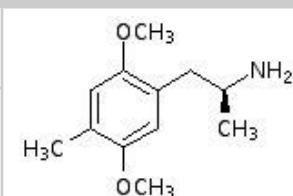
Catalog number : 7391-001

Name : (\pm)-4-Bromo-2,5-dimethoxyamphetamine hydrochloride; (\pm)-DOB

Mol. formula : C₁₁H₁₇BrClNO₂

FW : 310.62 DEA schedule : 1



Catalog number : 7391-004**CASRN :** 123431-31-2**Name :** (\pm)-4-Chloro-2,5-dimethoxyamphetamine hydrochloride; (\pm)-DOC**Mol. formula :** C₁₁H₁₆ClNO₂ • HCl**FW :** 266.166 **DEA schedule :** 1
3**Catalog number :** 7399-007**Name :** (-)-2,5-Dimethoxy-4-ethylamphetamine hydrochloride; (-)-DOET**Mol. formula :** C₁₃H₂₂ClNO₂**FW :** 259.78 **DEA schedule :** 1**Catalog number :** 7399-005**Name :** (+)-2,5-Dimethoxy-4-ethylamphetamine hydrochloride; (+)-DOET**Mol. formula :** C₁₃H₂₂ClNO₂**FW :** 259.78 **DEA schedule :** 1**Catalog number :** 7399-004**Name :** (\pm)-2,5-Dimethoxy-4-ethylamphetamine hydrochloride; (\pm)-DOET**Mol. formula :** C₁₃H₂₂ClNO₂**FW :** 259.78 **DEA schedule :** 1**Catalog number :** 7395-004**Name :** [OC²H₃]-2,5-Dimethoxy-4-methylamphetamine HCl; [OC²H₃]DOM**Mol. formula :** C₁₂H₂₀ClNO₂**FW :** 251.78 **DEA schedule :** 1**Catalog number :** 7395-002**CASRN :** 15588-95-1**Name :** (-)-2,5-Dimethoxy-4-methylamphetamine hydrochloride; (-)-DOM**Mol. formula :** C₁₂H₂₀ClNO₂**FW :** 245.75 **DEA schedule :** 1**Catalog number :** 7395-003**Name :** (+)-2,5-Dimethoxy-4-methylamphetamine hydrochloride; (+)-DOM**Mol. formula :** C₁₂H₂₀ClNO₂**FW :** 245.75 **DEA schedule :** 1

4 – Hallucinogens

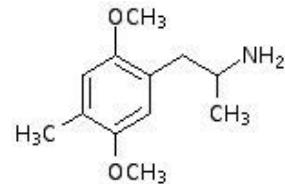
Catalog number : 7395-001

CASRN : 15588-95-1

Name : (\pm)-2,5-Dimethoxy-4-methylamphetamine hydrochloride; (\pm)-DOM

Mol. formula : C₁₂H₂₀ClNO₂

FW : 245.75 **DEA schedule :** 1



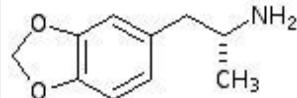
Catalog number : 7400-003

CASRN : 61614-60-6

Name : (–)-3,4-Methylenedioxymphetamine hydrochloride; (–)-MDA HCl

Mol. formula : C₁₀H₁₄ClNO₂

FW : 215.68 **DEA schedule :** 1



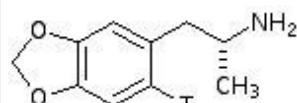
Catalog number : 7400-006

CASRN : 6292-91-7

Name : (–)-[6'-³H(n)]-3',4'-Methylenedioxymamphetamine hydrochloride;
(–)-[6'-³H(n)]MDA

Mol. formula : C₁₀H₁₄ClNO₂

FW : 217.68 **DEA schedule :** 1



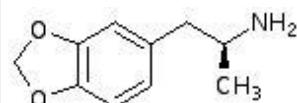
Catalog number : 7400-002

CASRN : 64057-70-1

Name : (+)-3,4-Methylenedioxymamphetamine hydrochloride; (+)-MDA HCl

Mol. formula : C₁₀H₁₄ClNO₂

FW : 215.68 **DEA schedule :** 1



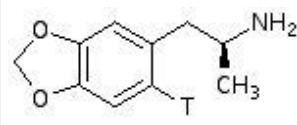
Catalog number : 7400-004

Name : (+)-[6'-³H(n)-3',4'-Methylenedioxymamphetamine hydrochloride;
(+)-[6'-³H(n)]MDA

Mol. formula : C₁₀H₁₄ClNO₂

FW : 215.68 **DEA schedule :** 1

Notes : Hallucinogen (tritium-labeled).



Catalog number : 7400-001

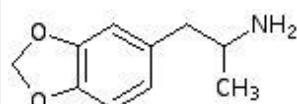
CASRN : 6292-91-7

Name : (\pm)-3,4-Methylenedioxymamphetamine hydrochloride; MDA HCl

Mol. formula : C₁₀H₁₄ClNO₂

FW : 215.68 **DEA schedule :** 1

References : Merck Index, 14th ed., Monograph 5765.

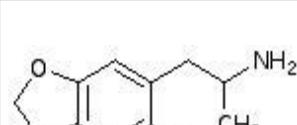


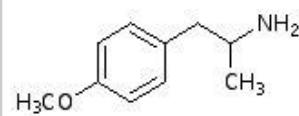
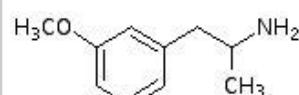
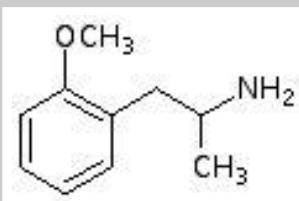
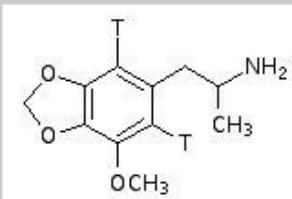
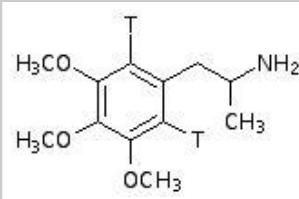
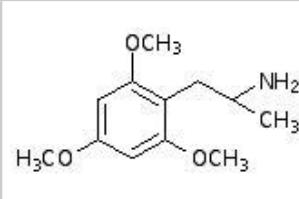
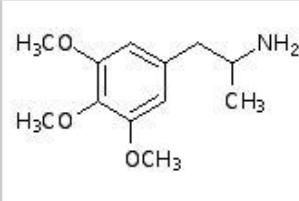
Catalog number : 7400-005

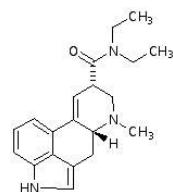
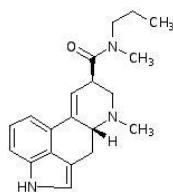
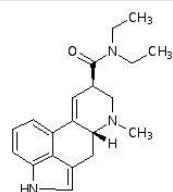
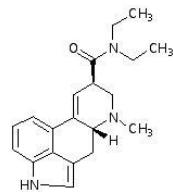
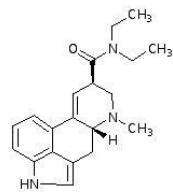
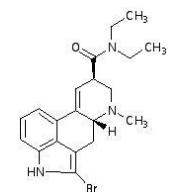
Name : [6'-³H(n)]-3,4-Methylenedioxymamphetamine hydrochloride;
[6'-³H(n)]MDA

Mol. formula : C₁₀H₁₄ClNO₂

FW : 217.68 **DEA schedule :** 1



Catalog number : 7411-001**CASRN :** 3706-26-1**Name :** (\pm)-4-Methoxyamphetamine hydrochloride**Mol. formula :** C₁₀H₁₆CINO**FW :** 201.70 **DEA schedule :** 1**Catalog number :** 7411-010**CASRN :** 17862-85-0**Name :** (\pm)-3-Methoxyamphetamine hydrochloride**Mol. formula :** C₁₀H₁₆CINO**FW :** 201.70 **DEA schedule :** 1**Catalog number :** 7411-011**CASRN :** 15402-84-3**Name :** (\pm)-2-Methoxyamphetamine hydrochloride; NDMP**Mol. formula :** C₁₀H₁₆CINO**FW :** 201.70 **DEA schedule :** 1**Catalog number :** 7401-002**Name :** [2',6'-³H(n)]-3-Methoxy-4,5-methylenedioxyamphetamine hydrochloride; [2,6-³H(n)]MMDA**Mol. formula :** C₁₁H₁₆CINO₃**FW :** 249.72 **DEA schedule :** 1**Catalog number :** 7390-002**Name :** (\pm)-[2,6-³H₂(n)]-3,4,5-Trimethoxyamphetamine hydrochloride**Mol. formula :** C₁₂H₂₀CINO₃**FW :** 265.76 **DEA schedule :** 1**Catalog number :** 7390-003**Name :** (\pm)-2,4,6-Trimethoxyamphetamine hydrochloride**Mol. formula :** C₁₂H₂₀CINO₃**FW :** 261.47 **DEA schedule :** 1**Catalog number :** 7390-001**CASRN :** 5688-80-2**Name :** (\pm)-3,4,5-Trimethoxyamphetamine hydrochloride; TMA**Mol. formula :** C₁₂H₂₀CINO₃**FW :** 261.47 **DEA schedule :** 1

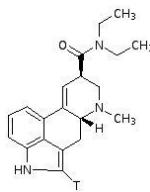
Hallucinogens: Ergot alkaloids**Catalog number :** 7315-014**CASRN :** 2126-78-5**Name :** (+)-Isolysergic acid diethylamide**Mol. formula :** C₂₀H₂₅N₃O**FW :** 323.20**DEA schedule :** 1**Catalog number :** 7315-013**CASRN :** 101692-69-7**Name :** (+)-Lysergic acid methyl-n-propylamide tartrate; LAMPA**Mol. formula :** (C₂₀H₂₅N₃O)₂**FW :** 796.33**DEA schedule :** 1**Catalog number :** 7315-004**CASRN :** 17676-08-3**Name :** (+)-Lysergic acid diethylamide (+)-tartrate (2:1); (+)-LSD (+)-tartrate**Mol. formula :** (C₂₀H₂₅N₃O)₂**FW :** 796.93**DEA schedule :** 1**References :** Merck Index, 14th ed., Monograph 5634.**Catalog number :** 7315-008**CASRN :** 50-37-3**Name :** (+)-Lysergic acid diethylamide; (+)-LSD**Mol. formula :** C₂₀H₂₅N₃O**FW :** 323.42**DEA schedule :** 1**References :** Merck Index, 14th ed., Monograph 5634.**Catalog number :** 7315-009**CASRN :** 24656-41-5**Name :** (+)-Lysergic acid diethylamide hydrogen maleate**Mol. formula :** C₂₀H₂₅N₃O₂**FW :** 439.49**DEA schedule :** 1**Catalog number :** 7315-010**CASRN :** 4004-43-7**Name :** (+)-2-Bromo-LSD (+)-hydrogen tartrate; BOL-148**Mol. formula :** C₂₀H₂₄BrN₃O**FW :** 552.44**DEA schedule :** 1**Notes :** Serotonin antagonist without the hallucinogenic activity of LSD**References :** Merck Index, 14th ed., Monograph 1423.

Catalog number : 7315-007

CASRN : 377756-22-4

Name : [2-³H]-(+)-Lysergic acid diethylamide; (+)-[3H]-LSDMol. formula : C₂₀H₂₅N₃O

FW : 323.42 DEA schedule : 1



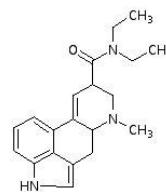
Catalog number : 7315-006

CASRN : 51064-36-9

Name : (±)-Lysergic acid diethylamide; (±)-LSD

Mol. formula : C₂₀H₂₅N₃O

FW : 323.42 DEA schedule : 1

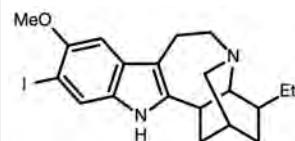
References : *Merck Index*, 14th ed., Monograph 5634.***Hallucinogens:*** Ibogaine Class

Catalog number : 7260-003

Name : 11-Iodoibogaine

Mol. formula : C₂₀H₂₅IN₂O

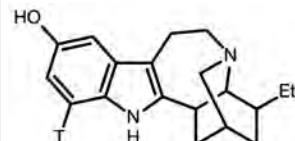
FW : 436.33 DEA schedule : 1



Catalog number : 7260-005

Name : [12-³H]-NoribogaineMol. formula : C₁₉H₂₄N₂O

FW : 296.41 DEA schedule : 0

Notes : *Ibogaine-like effect without tremors (tritium-labeled).*

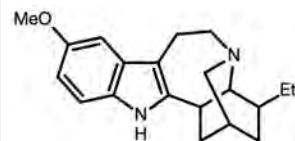
Catalog number : 7260-001

CASRN : 36415-61-9

Name : Ibogaine hydrochloride

Mol. formula : C₂₀H₂₇ClN₂O

FW : 346.89 DEA schedule : 1

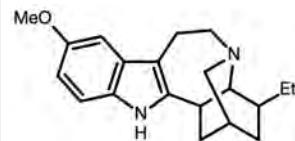
References : *Merck Index*, 14th ed., Monograph 4875.

Catalog number : 7260-002

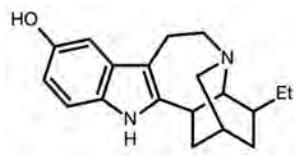
CASRN : 146560-35-2

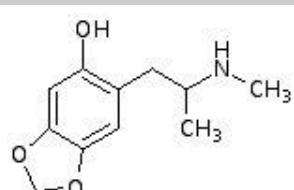
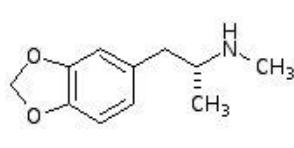
Name : Tritium-labeled Ibogaine; [12-³H]IbogaineMol. formula : C₂₀H₂₆N₂O

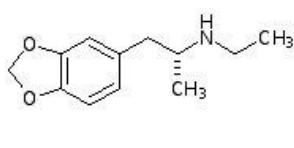
FW : 312.44 DEA schedule : 1

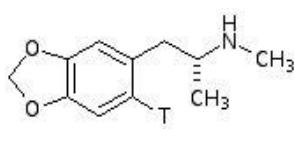
Notes : *Hallucinogen (tritium-labeled).*

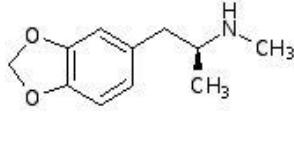
4 – Hallucinogens

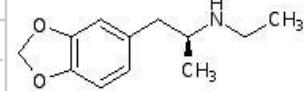
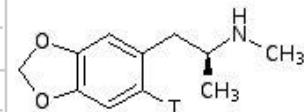
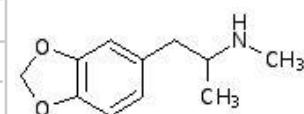
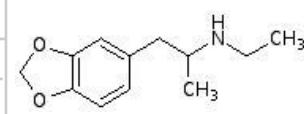
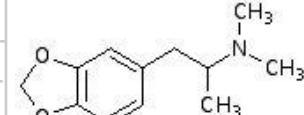
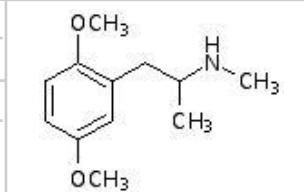
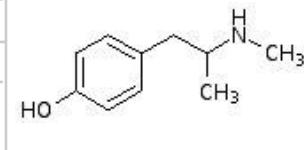
Catalog number : 7260-006	CASRN : 481-88-9		
Name : Noribogaine			
Mol. formula : C ₁₉ H ₂₄ N ₂ O	FW : 292.42	DEA schedule :	0
Notes : <i>Ibogaine-like effect without tremors</i>			
References : Glick, SD; Pearl, SM; Cai, J; Maisonneuve, IM <i>Brain Res</i> 1996, 713, 294–7.			

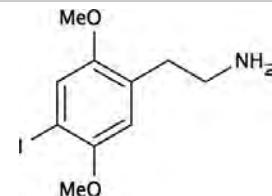
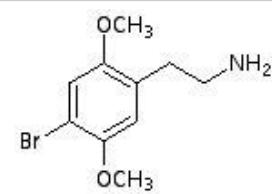
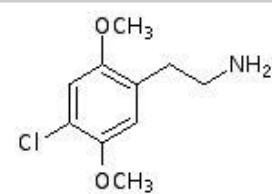
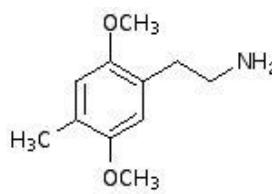
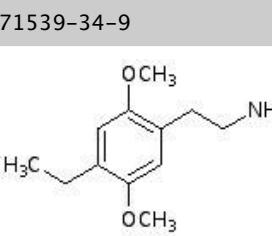
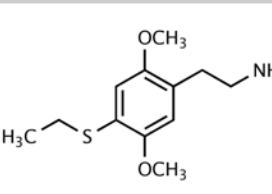
Hallucinogens: Methamphetamine Class			
Catalog number : 7401-004			
Name : (±)-6-Hydroxy-3,4-methylenedioxymethamphetamine fumarate			
Mol. formula : C ₂₆ H ₃₄ N ₂ O ₁₀	FW : 534.56	DEA schedule :	1
			
Catalog number : 7405-003	CASRN : 69558-31-2		
Name : (–)-3,4-Methylenedioxymethamphetamine hydrochloride; (–)-MDMA			
Mol. formula : C ₁₁ H ₁₆ CINO ₂	FW : 229.71	DEA schedule :	1
			

Catalog number : 7404-003	CASRN : 82801-81-8		
Name : (–)N-Ethyl-3,4-methylenedioxymethamphetamine hydrochloride			
Mol. formula : C ₁₂ H ₁₈ CINO ₂	FW : 243.73	DEA schedule :	1
Notes : CNS stimulant			
			

Catalog number : 7405-005	CASRN : 4764-17-4		
Name : (–)[6'- ³ H(n)]-3',4'-Methylenedioxymethamphetamine hydrochloride; (–)[³ H]MDMA			
Mol. formula : C ₁₁ H ₁₆ CINO ₂	FW : 231.71	DEA schedule :	1
Notes : CNS stimulant; hallucinogen (tritium-labeled).			
			

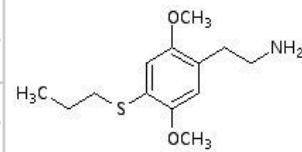
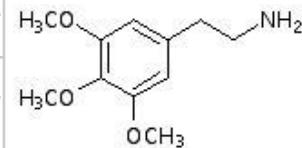
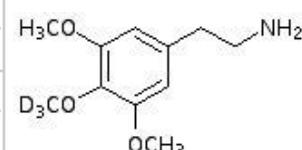
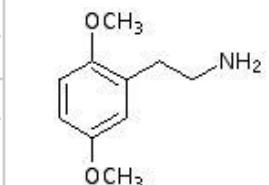
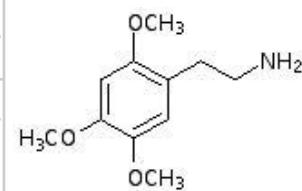
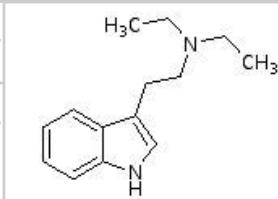
Catalog number : 7405-002	CASRN : 69558-32-3		
Name : (+)-3,4-Methylenedioxymethamphetamine hydrochloride; (+)-MDMA			
Mol. formula : C ₁₁ H ₁₆ CINO ₂	FW : 229.71	DEA schedule :	1
			

Catalog number : 7404-002**CASRN :** 82801-81-8**Name :** (+)-N-Ethyl-3,4-methylenedioxymphetamine hydrochloride**Mol. formula :** C₁₂H₁₈CINO₂**FW :** 243.73 **DEA schedule :** 1**Notes :** CNS stimulant**Catalog number :** 7405-006**Name :** (+)-[6'-³H(n)]-3',4'-Methylenedioxymethamphetamine hydrochloride; (+)-[³H]MDMA**Mol. formula :** C₁₁H₁₆CINO₂**FW :** 231.71 **DEA schedule :** 1**Notes :** CNS stimulant; hallucinogen (tritium-labeled).**Catalog number :** 7405-001**CASRN :** 64057-70-1**Name :** (±)-3,4-Methylenedioxymethamphetamine hydrochloride; MDMA**Mol. formula :** C₁₁H₁₆CINO₂**FW :** 229.71 **DEA schedule :** 1**References :** Merck Index, 14th ed., Monograph 5767.**Catalog number :** 7404-001**CASRN :** 82801-81-8**Name :** (±)-N-Ethyl-3,4-methylenedioxymamphetamine hydrochloride**Mol. formula :** C₁₂H₁₈CINO₂**FW :** 243.73 **DEA schedule :** 1**Notes :** CNS stimulant**Catalog number :** 7404-004**Name :** (±)-N,N-Dimethyl-3,4-methylenedioxymamphetamine hydrochloride**Mol. formula :** C₁₂H₁₈CINO₂**FW :** 243.73 **DEA schedule :** 1**Notes :** CNS stimulant**Catalog number :** 7395-005**Name :** (±)-2,5-Dimethoxymethamphetamine hydrochloride**Mol. formula :** C₁₂H₂₀CINO₂**FW :** 245.75 **DEA schedule :** 1**Catalog number :** 7411-012**CASRN :** 370-14-9**Name :** (±)-4-Hydroxymethamphetamine hydrochloride; Pholedrine**Mol. formula :** C₁₀H₁₆CINO**FW :** 201.70 **DEA schedule :** 2**References :** Merck Index, 14th ed., Monograph 4810.

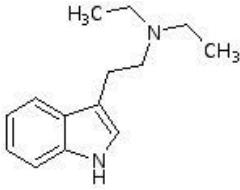
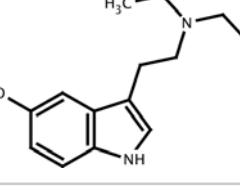
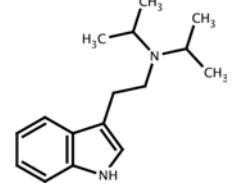
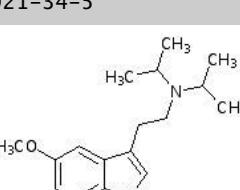
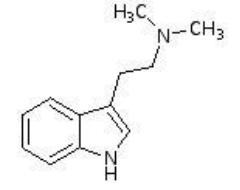
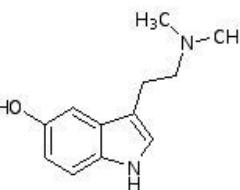
Hallucinogens: Phenethylamine Class**Catalog number :** 7518-001**CASRN :** 69587-11-7 (base)**Name :** 2,5-Dimethoxy-4-iodo-β-phenethylamine hydrochloride; 2C-I**Mol. formula :** C₁₀H₁₅ClINO₂**FW :** 343.59 **DEA schedule :** 1**Catalog number :** 7392-012**CASRN :** 56281-37-9**Name :** 4-Bromo-2,5-dimethoxy-β-phenethylamine hydrochloride; 2C-B**Mol. formula :** C₁₀H₁₅BrClNO₂**FW :** 296.59 **DEA schedule :** 1**References :** Merck Index, 14th ed., Monograph 1908.**Catalog number :** 7519-001**CASRN :** 88441-14-9**Name :** 4-Chloro-2,5-dimethoxy-β-phenethylamine hydrochloride; 2C-C**Mol. formula :** C₁₀H₁₅Cl₂NO₂**FW :** 252.14 **DEA schedule :** 0**References :** Jacob, P., 3rd; Shulgin AT *NIDA Res Monogr* 1994, 146, 74-91.**Catalog number :** 7508-001**CASRN :** 25505-65-1**Name :** 2,5-Dimethoxy-4-methyl-β-phenethylamine hydrochloride; 2C-D**Mol. formula :** C₁₁H₁₈ClNO₂**FW :** 231.72 **DEA schedule :** 0**References :** Jacob, P., 3rd; Shulgin AT *NIDA Res Monogr* 1994, 146, 74-91.**Catalog number :** 7509-001**CASRN :** 71539-34-9**Name :** 2,5-Dimethoxy-4-ethyl-β-phenethylamine hydrochloride; 2C-E**Mol. formula :** C₁₂H₂₀ClNO₂**FW :** 245.75 **DEA schedule :** 0**References :** Jacob, P., 3rd; Shulgin AT *NIDA Res Monogr* 1994, 146, 74-91.**Catalog number :** 7385-001**CASRN :** 207740-24-7 (base)**Name :** 2,5-Dimethoxy-4-ethylthio-β-phenethylamine hydrochloride; 2C-T-2**Mol. formula :** C₁₂H₂₀ClNO₂S**FW :** 277.81 **DEA schedule :** 0**References :** Fantegrossi, WE; et al. *Psychopharmacology (Berl)* 2005, 181, 496-503.

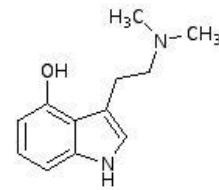
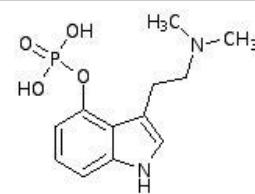
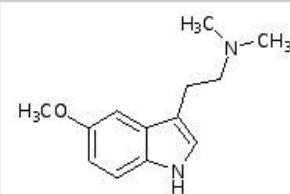
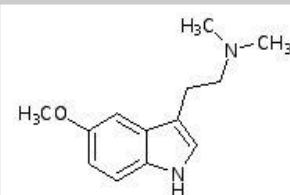
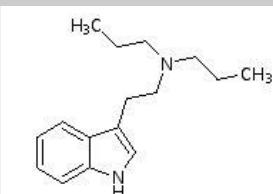
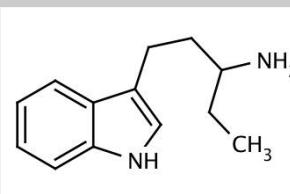
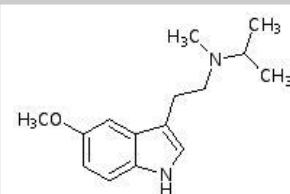
Catalog number : 7348-001**CASRN :** 207740-26-9

Name : 2,5-Dimethoxy-4-n-propylthio- β -phenethylamine hydrochloride;
2C-T-7

Mol. formula : C₁₃H₂₂CINO₂S**FW :** 291.84 **DEA schedule :** 1**References :** Fantegrossi, WE; et al. *Psychopharmacology (Berl)* 2005, 181, 496–503.**Catalog number :** 7381-001**CASRN :** 832-92-8**Name :** Mescaline hydrochloride**Mol. formula :** C₁₁H₁₈CINO₃**FW :** 247.72 **DEA schedule :** 1**References :** Merck Index, 14th ed., Monograph 5905.**Catalog number :** 7381-002**Name :** [4'-OC²H₃]Mescaline hydrochloride**Mol. formula :** C₁₁H₁₈CINO₃**FW :** 247.72 **DEA schedule :** 1**Catalog number :** 7381-011**CASRN :** 3166-74-3**Name :** 2,5-Dimethoxy- β -phenethylamine hydrochloride**Mol. formula :** C₁₀H₁₆CINO₂**FW :** 217.70 **DEA schedule :** 0**Catalog number :** 7381-010**CASRN :** 15394-83-9**Name :** 2,4,5-Trimethoxy- β -phenethylamine hydrochloride**Mol. formula :** C₁₁H₁₈CINO₃**FW :** 247.72 **DEA schedule :** 1**Hallucinogens: Tryptamine Class****Catalog number :** 7434-001**CASRN :** 20671-78-7**Name :** N,N-Diethyltryptamine fumarate; DET fumarate**Mol. formula :** C₁₈H₂₄N₂O₄**FW :** 332.40 **DEA schedule :** 1**References :** Heinze, WJ; Schlemmer, RF; Tyler, CB; Davis, JM *Biol Psychiatry* 1983, 18, 829–36.

4 – Hallucinogens

Catalog number : 7434-002	CASRN : 61-51-8	
Name : N,N-Diethyltryptamine (base); DET		
Mol. formula : C ₁₄ H ₂₀ N ₂	FW : 216.33	DEA schedule : 1
References : Heinze, WJ; Schlemmer, RF, Jr; Tyler, CB; Davis JM <i>JM Biol Psychiatry</i> 1983, 18, 829–36.		
Catalog number : 7525-000	CASRN : 14780-24-6	
Name : 5-Methoxy-N,N-diethyltryptamine, 5-MeO-DET		
Mol. formula : C ₁₅ H ₂₂ N ₂ O	FW : 246.35	DEA schedule : 1
References : Fumiko N, Ryouchi N, Kanak S, Hisashi K, <i>Eur J Pharmacol</i> , 2007, 559 (2–3), 132–137.		
Catalog number : 7522-000	CASRN : 63065-90-7	
Name : N,N-Diisopropyltryptamine hydrochloride, DiPT		
Mol. formula : C ₁₆ H ₂₄ N ₂	FW : 244.38	DEA schedule : 1
Notes : <i>Psilocin analog.</i>		
References : Pichini, S; et al., <i>J Pharm Biomed Analysis</i> 2008, 47, 335–342.		
Catalog number : 7439-001	CASRN : 4021-34-5	
Name : 5-Methoxy-N,N-diisopropyltryptamine hydrochloride; 5-MeO-DiPT; FOXY		
Mol. formula : C ₁₇ H ₂₇ ClN ₂ O	FW : 310.87	DEA schedule : 1
References : Shulgin, AT; Carter MF <i>Commun Psychopharmacol</i> 1980, 4, 363–9.		
Catalog number : 7435-001	CASRN : 69321-46-6	
Name : N,N-Dimethyltryptamine fumarate; DMT fumarate		
Mol. formula : C ₁₆ H ₂₀ N ₂ O ₄	FW : 304.35	DEA schedule : 1
References : Barker, SA; Monti, JA; Christian, ST <i>Int Rev Neurobiol</i> 1981, 22, 83–110.		
Catalog number : 7433-001	CASRN : 487-93-4	
Name : Bufotenine (base); 5-Hydroxy-N,N-dimethyltryptamine; 5-OH-DMT		
Mol. formula : C ₁₂ H ₁₆ N ₂ O	FW : 204.26	DEA schedule : 1
References : Emanuele, E; et al., <i>Neuro Endocrinol Lett</i> 2010, 31, 117–21.		

Catalog number : 7438-001**CASRN :** 520-53-6**Name :** Psilocin**Mol. formula :** C₁₂H₁₆N₂O**FW :** 204.27 **DEA schedule :** 1**References :** *Merck Index*, 14th ed., Monograph 7925.**Catalog number :** 7437-001**CASRN :** 520-52-5**Name :** Psilocybin**Mol. formula :** C₁₂H₁₇N₂O₄**FW :** 284.25 **DEA schedule :** 1**References :** *Merck Index*, 14th ed., Monograph 7926.**Catalog number :** 7431-001**CASRN :** 1019-45-0**Name :** 5-Methoxy-N,N-dimethyltryptamine (base); 5-MeO-DMT**Mol. formula :** C₁₃H₁₈N₂O**FW :** 218.29 **DEA schedule :** 1**Catalog number :** 7431-002**CASRN :** 1019-45-0**Name :** 5-Methoxy-N,N-dimethyltryptamine fumarate; 5-MeO-DMT fumarate**Mol. formula :** C₃₀H₄₀N₄O₆**FW :** 552.67 **DEA schedule :** 1**Catalog number :** 7436-001**CASRN :** 7558-73-8**Name :** N,N-Dipropyltryptamine hydrochloride**Mol. formula :** C₁₆H₂₅ClN₂**FW :** 280.84 **DEA schedule :** 0**Catalog number :** 7249-001**CASRN :** 2235-90-7**Name :** α-Ethyltryptamine acetate; AET**Mol. formula :** C₁₄H₂₀N₂O₂**FW :** 248.33 **DEA schedule :** 1**References :** *Merck Index*, 13th ed., Monograph 3924.**Catalog number :** NOCD-001**CASRN :** 96096-54-7**Name :** 5-Methoxy-N-isopropyl-N-methyltryptamine hydrochloride; 5-MeO-MiPT**Mol. formula :** C₁₅H₂₂N₂O • HCl**FW :** 282.81 **DEA schedule :** 1**References :** Repke, DB; Grotjahn DB; Shulgin AT, *J Med Chem* 1985, 28, 892-6; Nagai, F; et al., *Eur J Pharmacol* 2007, 559(2-3), 132-137.

4 – Hallucinogens

Catalog number : 7432-001

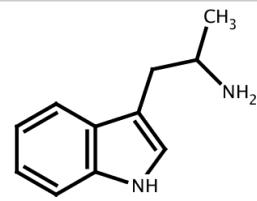
CASRN : 299-26-3

Name : α -Methyltryptamine; AMT

Mol. formula : $C_{11}H_{14}N_2$

FW : 174.24

DEA schedule : 1



Catalog number : 7435-002

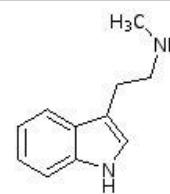
CASRN : 61-49-4

Name : N-Methyltryptamine fumarate

Mol. formula : $C_{15}H_{18}N_2O_4$

FW : 290.31

DEA schedule : 0



Catalog number : 7506-001

CASRN : 1137-04-8

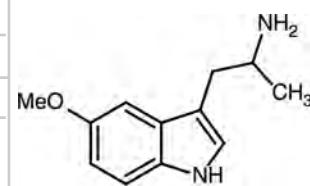
Name : 5-Methoxy- α -methyltryptamine; 5-MeO-AMT

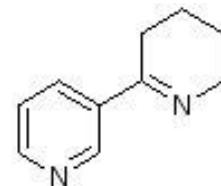
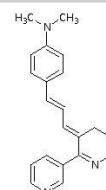
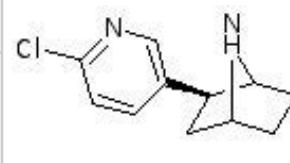
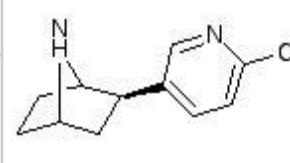
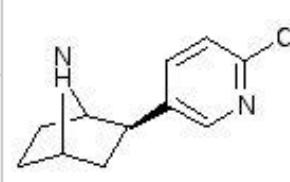
Mol. formula : $C_{12}H_{16}N_2O$

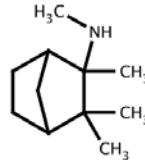
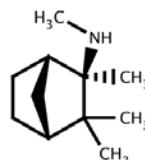
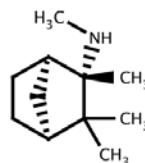
FW : 204.27

DEA schedule : 0

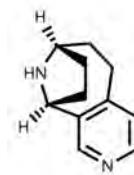
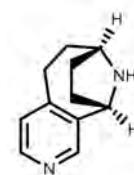
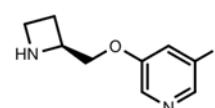
References : Jacob, P; Shulgin AT *NIDA Res Monogr* 1994, 146, 74-91.

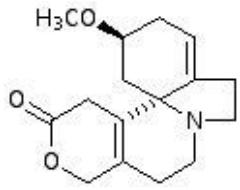
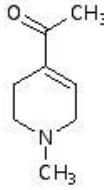
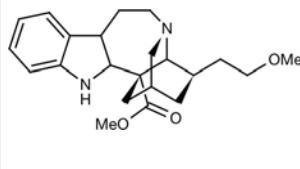
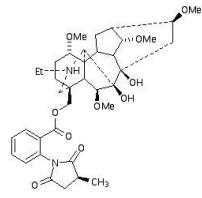
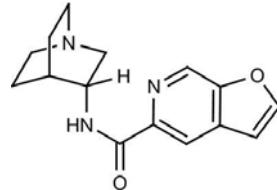
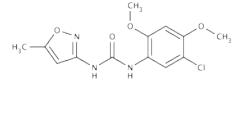
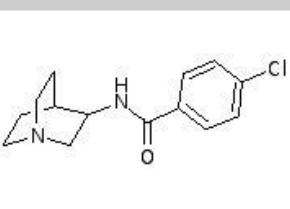


Nicotinics: Anabaseine Class**Catalog number :** NICT-011**CASRN :** 3471-05-4**Name :** Anabaseine dihydrochloride**Mol. formula :** C₁₀H₁₄Cl₂N₂**FW :** 233.14 **DEA schedule :** 0**Notes :** Nicotinic $\alpha 7$ receptor agonist.**References :** Wheeler, JW; et al. *Science* 1981, 211, 1051–1052.Kem WR; Mahnir VM; Papke RL; Lingle CJ *J Pharmacol Exp Ther* 1997, 283, 979–92.**Catalog number :** NICT-012**Name :** (E,E)-3-(4'-Dimethylaminocinnamylidene)anabaseine dihydrochloride; DMAC**Mol. formula :** C₂₁H₂₅Cl₂N₃**FW :** 390.36 **DEA schedule :** 0**Notes :** Selective $\alpha 7$ nicotinic acetylcholine receptor agonist.**References :** de Fiebre, C; et al. *Mol Pharmacol* 1995, 47, 164–171.**Nicotinics: Epibatidine Class****Catalog number :** NICT-002**CASRN :** 152378-30-8**Name :** (–)-Epibatidine hydrochloride**Mol. formula :** C₁₁H₁₄Cl₂N₂**FW :** 245.15 **DEA schedule :** 0**References :** Qian C; Li T; Shen TY; Libertine–Garahan L; Eckman J; Biftu T; Ip S *Eur J Pharmacol* 1993, 250, R13–4.Badio B; Daly JW *Mol Pharmacol* 1994, 45, 563–9.**Catalog number :** NICT-003**CASRN :** 140111-52-0**Name :** (+)-Epibatidine hydrochloride**Mol. formula :** C₁₁H₁₄Cl₂N₂**FW :** 245.15 **DEA schedule :** 0**Notes :** Epibatidine natural isomer.**References :** Qian C; Li T; Shen TY; Libertine–Garahan L; Eckman J; Biftu T; Ip S *Eur J Pharmacol* 1993, 250, R13–4.Badio B; Daly JW *Mol Pharmacol* 1994, 45, 563–9.**Catalog number :** NICT-001**Name :** (±)-Epibatidine hydrochloride**Mol. formula :** C₁₁H₁₄Cl₂N₂**FW :** 245.15 **DEA schedule :** 0**Notes :** Potent nicotinic receptor agonist and non-opioid analgesic.**References :** Sullivan, JP; et al. *J Pharmacol Exp Ther* 1994, 271, 624–31.

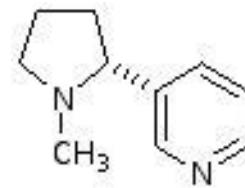
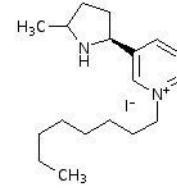
Nicotinics: Mecamylamine Class**Catalog number :** NICT-004**CASRN :** 826-39-1**Name :** (±)-Mecamylamine HCl**Mol. formula :** C₁₁H₂₂CIN**FW :** 203.75 **DEA schedule :** 0**Notes :** Non-competitive nicotinic acetylcholine receptor antagonist.**References :** Sanberg, PR; et al. *Int J Neurosci* 2001, 109, 81-90.**Catalog number :** NOCD-123**CASRN :** 107596-31-6**Name :** (−)-(R)-Mecamylamine HCl**Mol. formula :** C₁₁H₂₂CIN**FW :** 203.75 **DEA schedule :** 0**Notes :** Non-competitive nicotinic acetylcholine receptor antagonist.**References :** Suchocki, JA, *Journal of Medicinal Chemistry* 1991, 34(3), 1003-10.
Schoenenberger, B, *Helvetica Chimica Acta* 1986, 69(2), 283-7.**Catalog number :** NICT-018**CASRN :** 107596-30-5**Name :** (+)-(S)-Mecamylamine HCl**Mol. formula :** C₁₁H₂₂CIN**FW :** 203.75 **DEA schedule :** 0**Notes :** Non-competitive nicotinic acetylcholine receptor antagonist.**References :** Stone, CA; et al., *J. Med. Chem.*, 34, 1003 (1991).**Nicotinics: Miscellaneous****Catalog number :** NOCD-142

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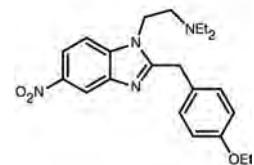
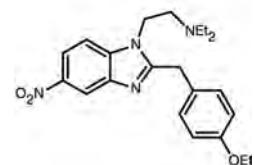
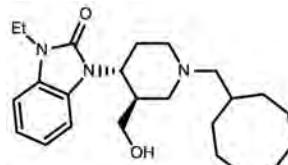
CASRN : 895518-77-1**Name :** (+)-PHT dihydrochloride**Mol. formula :** C₁₁H₁₆Cl₂N₂**FW :** 247.17 **DEA schedule :** 0**Notes :** Conformationally rigid nicotine analog.**References :** Kanne, D. B.; Ashworth, D. J.; Cheng, M. T.: Mutter, L. C.; Abood, L. G. *J. Am. Chem. Soc.* 1986, 108, 7864.
Carroll FI, Hu X, Navarro HA, Deschamps J, Abdurakhmanova GR, Damaj MI, Martin BR *J. Med. Chem.* 2006, 49, 3244.**Catalog number :** NOCD-122**CASRN :** 895518-76-0**Name :** (−)-PHT dihydrochloride**Mol. formula :** C₁₁H₁₆Cl₂N₂**FW :** 247.17 **DEA schedule :** 0**Notes :** Conformationally rigid nicotine analog.**References :** Kanne, D. B.; Ashworth, D. J.; Cheng, M. T.: Mutter, L. C.; Abood, L. G. *J. Am. Chem. Soc.* 1986, 108, 7864.
Carroll FI, Hu X, Navarro HA, Deschamps J, Abdurakhmanova GR, Damaj MI, Martin BR *J. Med. Chem.* 2006, 49, 3244.**Catalog number :** NICT-019**CASRN :** 1217837-17-6**Name :** 5-I-A85380**Mol. formula :** C₉H₁₁IN₂O • 2 HCl**FW :** 363.02 **DEA schedule :** 0**Notes :** α4β2-selective nAChR ligand**References :** Koren, et al., *J Med Chem* 1998, 41, 3690. Muhkin, et al., *Mol Pharmacol* 2000, 57, 642.

Catalog number : NICT-006	CASRN : 29734-68-7	
Name : Dihydro-β-erythroidine hydrobromide		
Mol. formula : C ₁₆ H ₂₂ BrNO ₃	FW : 356.26	DEA schedule : 0
References : <i>Merck Index</i> , 14th ed., Monograph 3175.		
Catalog number : NICT-005	CASRN : 100752-88-3	
Name : Isoarecolone hydrochloride		
Mol. formula : C ₈ H ₁₄ ClNO	FW : 175.66	DEA schedule : 0
Notes : Nicotinic receptor agonist.		
References : Mirza, NR; et al. <i>Eur J Pharmacol</i> 1996, 295, 207–10. Shoaib, M <i>Psychopharmacology (Berl)</i> 2006, 188, 252–7.		
Catalog number : NOCD-012	CASRN : 308123-60-6	
Name : 18-Methoxycoronaridine; 18-MC		
Mol. formula : C ₂₂ H ₃₀ N ₂ O ₃	FW : 370.49	DEA schedule : 0
Notes : Selective α3β4 nicotinic antagonist.		
References : Glick, SD; et al. <i>Brain Res</i> 1996, 719, 29–35.		
Catalog number : NICT-007	CASRN : 21019-30-7	
Name : Methyllycaconitine citrate; MLA		
Mol. formula : C ₄₃ H ₅₈ N ₂ O ₁₇	FW : 928.98	DEA schedule : 0
Notes : Nicotinic receptor antagonist.		
References : Ward, JM; et al. <i>FEBS Lett</i> 1990, 270, 45–8. Alkondon, M; et al. <i>Mol Pharmacol</i> 1992, 41, 802–8.		
Catalog number : NOCD-013	CASRN : 478149-53-0	
Name : PHA-543613		
Mol. formula : C ₁₅ H ₁₉ Cl ₂ N ₃ O ₂	FW : 344.24	DEA schedule : 0
Notes : α7 nicotinic acetylcholine receptor agonist		
References : Wishka, DG; et al. <i>J Med Chem</i> 2006, 49, 4425–36.		
Catalog number : NICT-017	CASRN : 501925-31-1	
Name : PNU-120596		
Mol. formula : C ₁₃ H ₁₄ ClN ₃ O ₄	FW : 311.72	
Notes : Positive allosteric modulator of the α ₇ neuronal nicotinic acetylcholine receptor.		
References : Hurst, RS; et al. <i>J Neurosci</i> 2005, 25, 4396–405.		
Catalog number : NICT-016	CASRN : 123464-89-1	
Name : PNU-282987		
Mol. formula : C ₁₄ H ₁₀ Cl ₂ N ₂₀	FW : 310.21	DEA schedule : 0
Notes : Selective α7 nicotinic acetylcholine receptor agonist.		
References : Hajos, M; et al. <i>J Pharmacol Exp Ther</i> 2005, 312, 1213–22.		

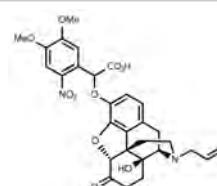
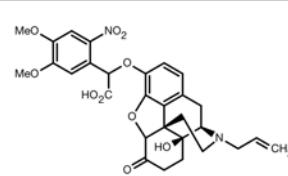
Catalog number : MEDD-031	CASRN : 820231-95-6	
Name : Sazetidine hydrochloride		
Mol. formula : C ₁₅ H ₂₂ ClN ₂ O ₂	FW : 333.26	DEA schedule : 0
Notes : Selective $\alpha_4\beta_2$ nicotinic receptor agonist.		
References : Xiao, et al., <i>Mol Pharmacol</i> , 2006, 70, 1454. Zwart, et al., <i>Mol Pharmacol</i> , 2008, 73 1843.		
Catalog number : NICT-013	CASRN :	
Name : Trimethaphan bromide		
Mol. formula : C ₂₂ H ₂₅ BrN ₂ O ₃	FW : 445.43	DEA schedule : 0
Notes : Nicotinic antagonist; antihypertensive.		
References : <i>Merck Index</i> , 14th ed., Monograph 9707.		
Catalog number : NICT-014	CASRN :	
Name : 2,3,6-Trimethyl-1,4-naphthoquinone		
Mol. formula : C ₁₃ H ₁₂ O ₂	FW : 200.24	DEA schedule : 0
Catalog number : NOCD-046	CASRN : 249296-44-4	
Name : Varenicline dihydrochloride		
Mol. formula : C ₁₃ H ₁₅ Cl ₂ N ₃	FW : 288.685	DEA schedule : 0
Notes : $\alpha 4\beta 2$ nicotinic receptor partial agonist.		
References : Coe, JW; et al. <i>J Med Chem</i> 2005, 48, 3474-7.		
Nicotinics: Nicotine Class		
Catalog number : NICT-020	new	CASRN : 65-31-6
Name : (-)-Nicotine ditartrate		
Mol. formula : C ₁₈ H ₂₆ N ₂ O ₁₂	FW : 462.41	DEA schedule : 0
Catalog number : NICT-009		
Name : (-)-N-Decylnicotinium iodide (-)-di-p-toluoyl-L-tartrate; NDNI		
Mol. formula : C ₄₀ H ₅₃ IN ₂ O ₈	FW : 816.77	DEA schedule : 0

Catalog number : NICT-015**CASRN :** 121949-85-7**Name :** (-)-Nicotine-4,4-d₂ (-)-di-p-toluoyl-L-tartrate**Mol. formula :** C₃₀H₃₂N₂O₈**FW :** 550.60 **DEA schedule :** 0**References :** Jacob, PJ *Labelled Comp Radiopharm* **1988**, *25*, 1117-28.**Catalog number :** NICT-008**CASRN :** 163804-20-4**Name :** (+)-Nicotine (+)-di-p-toluoyl-D-tartrate**Mol. formula :** C₃₀H₃₂N₂O₈**FW :** 548.58 **DEA schedule :** 0**Catalog number :** NICT-010**Name :** (-)-N-Octylnicotinium iodide (-)-di-p-toluoyl-L-tartrate; NONI**Mol. formula :** C₃₈H₄₉IN₂O₈**FW :** 788.72 **DEA schedule :** 0



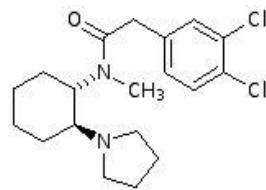
Opioids: Benzodiazole Class**Catalog number :** 9624-002**Name :** Etonitazene**Mol. formula :** C₂₂H₂₈N₄O₃**FW :** 396.48 **DEA schedule :** 1**Notes :** Narcotic analgesic; potent μ opiate receptor agonist**References :** Merck Index, 14th ed., Monograph 3883.**Catalog number :** 9624-001**CASRN :** 911-65-9**Name :** Etonitazene hydrochloride**Mol. formula :** C₂₂H₂₉ClN₄O₃**FW :** 432.95 **DEA schedule :** 1**Notes :** Narcotic analgesic; potent μ opiate receptor agonist**References :** Merck Index, 14th ed., Monograph 3883.**Catalog number :** NOCD-057**CASRN :** 361343-48-8**Name :** (+)-J-113397**Mol. formula :** C₂₄H₃₈ClN₃O₂**FW :** 454.05 **DEA schedule :** 0**Notes :** Nociceptin/orphanin FQ (NOP) receptor antagonist.**References :** Kawamoto H; et al. *Tetrahedron* 2001, 57, 981–986.
Kawamoto H; et al. *J Med Chem* 1999, 42, 5061–3.**Opioids: Caged****Catalog number :** 9652-071

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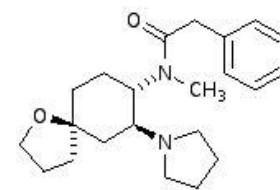
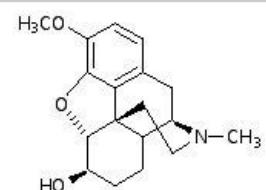
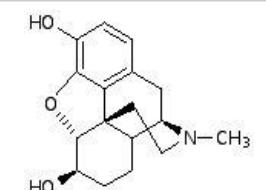
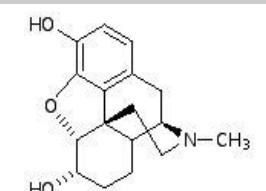
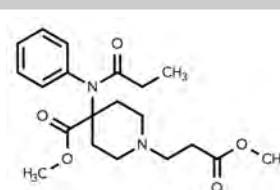
CASRN : n/a**Name :** CNV-NLX**Mol. formula :** C₂₉H₃₀N₂O₁₀**FW :** 566.56**Notes :** "Caged" opioid antagonist.**References :** M. R. Banghart, J. T. Williams, R. C. Shah, L. D. Lavis, B. L. Sabatini, Mol Pharmacol. 84, 687–695 (2013); M. R. Bruchas, B. L. Roth, Trends in Pharmacological Sciences. 37, 279–289 (2016).**Catalog number :** MPSP-120**Name :** (Carboxynitroveratryl)naloxone; CNV-NLX**Mol. formula :** C₂₉H₃₀N₂O₁₀**FW :** 566.47 **DEA schedule :** 0**Notes :** Photoactivatable analog of naloxone.**References :** Banghart, MR; et al., *Mol Pharmacol* 2013, 84, 687–95.

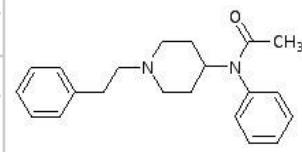
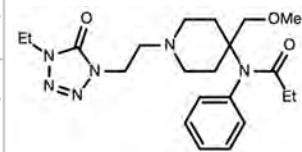
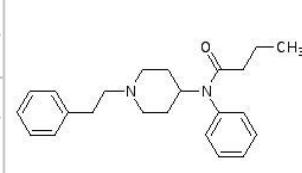
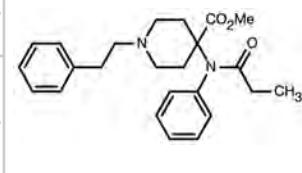
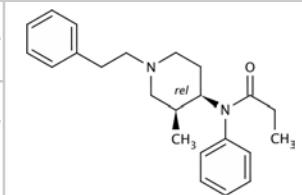
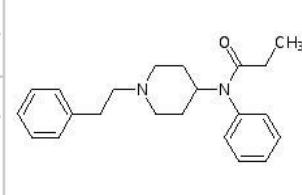
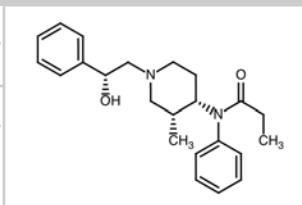
Opioids: Cyclohexyldiamine Class**Catalog number :** NOCD-064**CASRN :** 67198-13-4**Name :** U50,488H**Mol. formula :** C₂₀H₃₀Cl₂N₂O₄S**FW :** 465.44 **DEA schedule :** 0**Notes :** Kappa-opioid receptor agonist.

References : Vonvoigtlander, PF; et al. *J Pharmacol Exp Ther* 1983, 224, 7-12.
 Negus, SS; et al. *J Pharmacol Exp Ther* 1997, 282, 44-55.
 Taylor, CC; et al. *J Pharmacol Exp Ther* 1997, 280, 416-21.

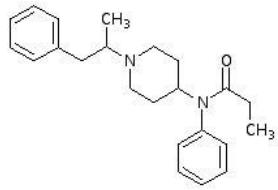
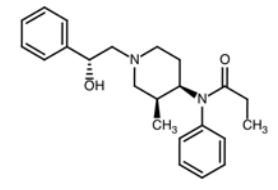
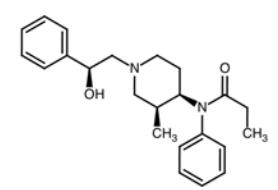
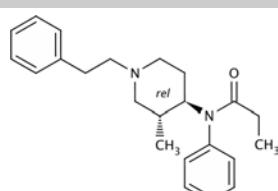
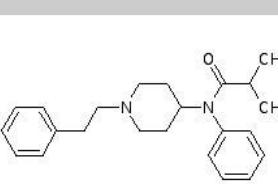
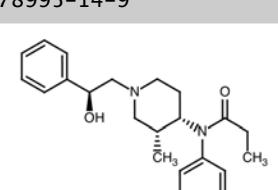
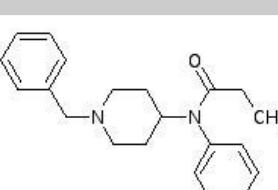
**Catalog number :** NOCD-065**CASRN :** 96744-75-1**Name :** U69,593**Mol. formula :** C₂₂H₃₂N₂O₂**FW :** 356.51 **DEA schedule :** 0**Notes :** Selective κ-opioid agonist.

References : Lahti, RA; et al. *Eur J Pharmacol* 1985, 109, 281-4.
 Puig-Ramos, A; et al. *Behav Neurosci* 2008, 122, 151-60.

**Opioids: Dihydromorphine Class****Catalog number :** 9120-002**Name :** Dihydroisocodeine**Mol. formula :** C₁₈H₂₃NO₃**FW :** 301.39 **DEA schedule :** 2**References :** *Merck Index*, 14th ed., Monograph 3176.**Catalog number :** 9145-002**Name :** Dihydroisomorphone**Mol. formula :** C₁₇H₂₁NO₃**FW :** 287.36 **DEA schedule :** 1**Catalog number :** 9145-001**CASRN :** 509-60-4**Name :** Dihydromorphine**Mol. formula :** C₁₇H₂₁NO₃**FW :** 287.36 **DEA schedule :** 1**References :** *Merck Index*, 14th ed., Monograph 3177.**Opioids: Fentanyl Class****Catalog number :** 9739-001**new****CASRN :** 132539-07-2**Name :** Remifentanil hydrochloride**Mol. formula :** C₂₀H₂₉ClN₂O₅**FW :** 412.91 **DEA schedule :** 2**Notes :** Short-acting fentanyl analog.

Catalog number : 9821-001**CASRN :** 117332-89-5**Name :** Acetylfentanyl hydrochloride**Mol. formula :** C₂₁H₂₇ClN₂O**FW :** 358.91 **DEA schedule :** 2**References :** Brine, GA; et al. *J Heterocyclic Chem* 1989, 26, 677.**Catalog number :** 9737-001**CASRN :** 69049-06-5**Name :** Alfentanil hydrochloride**Mol. formula :** C₂₁H₃₃ClN₆O₃**FW :** 452.98 **DEA schedule :** 2**Notes :** Analgesic, μ -Opioid agonist**References :** Merck Index, 14th ed., Monograph 236.**Catalog number :** 9822-001**CASRN :** 1443-52-3**Name :** Butyrylfentanyl hydrochloride**Mol. formula :** C₂₃H₃₁ClN₂O**FW :** 386.97 **DEA schedule :** 1**References :** Brine, GA; et al. *J Heterocyclic Chem* 1989, 26, 677.**Catalog number :** 9743-001**CASRN :** 59708-52-0 (parent)**Name :** Carfentanil hydrochloride**Mol. formula :** C₂₄H₃₁ClN₂O₃**FW :** 430.98 **DEA schedule :** 2**Notes :** Analgesic, μ -Opioid agonist.
(see footnotes 1 & 2 in Section C)**Catalog number :** 9813-001**CASRN :** 42045-86-3**Name :** (\pm)-*cis*-3-Methylfentanyl hydrochloride**Mol. formula :** C₂₃H₃₁ClN₂O**FW :** 386.96 **DEA schedule :** 1**Notes :** Analgesic, μ -Opioid agonist**References :** Brine, GA; et al. *J Heterocyclic Chem* 1989, 26, 677.**Catalog number :** 9801-001**CASRN :** 1443-54-5**Name :** Fentanyl hydrochloride**Mol. formula :** C₂₂H₂₉ClN₂O**FW :** 372.92 **DEA schedule :** 2**Notes :** Analgesic, μ -Opioid agonist**References :** Merck Index, 14th ed., Monograph 4001.
Brine, GA; et al. *J Heterocyclic Chem* 1989, 26, 677.**Catalog number :** 9831-003**CASRN :** 155168-97-1**Name :** (–)(*β*R,*3R*,*4S*)- β -Hydroxy-3-methylfentanyl oxalate**Mol. formula :** C₂₅H₃₂N₂O₆**FW :** 456.55 **DEA schedule :** 1**Notes :** Analgesic, μ -Opioid agonist**References :** Brine, GA; et al. *J Med Chem* 1995, 38, 1547-57.

6 – Opioids and Opioid-related

Catalog number : 9814-001	CASRN : 79704-88-4	
Name : α -Methylfentanyl hydrochloride		
Mol. formula : $C_{23}H_{31}ClN_2O$	FW : 386.97	DEA schedule : 1
Notes : Analgesic, μ -Opioid agonist		
References : Ayres, WA; et al. <i>J Psychoactive Drugs</i> 1981, 13, 91–3. Kram, TC; et al. <i>Anal Chem</i> 1981, 53, 1379A–1386A. Brine, GA; et al. <i>J Heterocyclic Chem</i> 1989, 26, 677.		
Catalog number : 9831-004	CASRN : 143553-99-5	
Name : $(-)(\beta R,3S,4R)$ - β -Hydroxy-3-methylfentanyl hydrochloride		
Mol. formula : $C_{23}H_{31}ClN_2O_2$	FW : 402.97	DEA schedule : 1
Notes : Analgesic, μ -Opioid agonist		
References : Brine, GA; et al. <i>J Med Chem</i> 1995, 38, 1547–57.		
Catalog number : 9831-005	CASRN : 155169-00-9	
Name : $(+)(\beta S,3S,4R)$ - β -Hydroxy-3-methylfentanyl oxalate		
Mol. formula : $C_{25}H_{32}N_2O_6$	FW : 456.55	DEA schedule : 1
Notes : Analgesic, μ -Opioid agonist		
References : Brine, GA; et al. <i>J Med Chem</i> 1995, 38, 1547–57.		
Catalog number : 9813-002	CASRN : 42045-87-4	
Name : (\pm) - <i>trans</i> -3-Methylfentanyl hydrochloride		
Mol. formula : $C_{23}H_{31}ClN_2O$	FW : 386.96	DEA schedule : 1
Notes : Analgesic, μ -Opioid agonist		
Catalog number : 9801-006	CASRN : 117332-90-8	
Name : Isopropylfentanyl hydrochloride		
Mol. formula : $C_{23}H_{31}ClN_2O$	FW : 386.96	DEA schedule : 2
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1989, 26, 677.		
Catalog number : 9831-002	CASRN : 78995-14-9	
Name : $(+)(\beta S,3R,4S)$ - β -Hydroxy-3-methylfentanyl hydrochloride; Ohmefentanyl; F 7302		
Mol. formula : $C_{23}H_{31}ClN_2O_2$	FW : 402.97	DEA schedule : 1
Notes : Analgesic, μ -Opioid agonist		
References : Xu, H; Chen J; Chi, ZQ <i>Sci Sin [B]</i> 1985, 28, 504–11. Brine, GA; et al. <i>J Med Chem</i> 1995, 38, 1547–57.		
Catalog number : 9818-002	CASRN : 1474-02-8	
Name : Benzylfentanyl hydrochloride		
Mol. formula : $C_{21}H_{27}ClN_2O$	FW : 358.91	DEA schedule : 0
Notes : Pharmacologically inert fentanyl analog.		
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1989, 26, 677. <i>Federal Register</i> 2010, 75(124), 37300.		

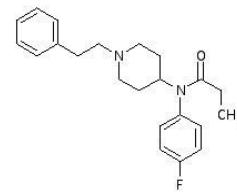
Catalog number : 9812-001

CASRN : 90736-23-5

Name : p-Fluorofentanyl hydrochloride

Mol. formula : C₂₂H₂₈ClFN₂O

FW : 390.93 DEA schedule : 1



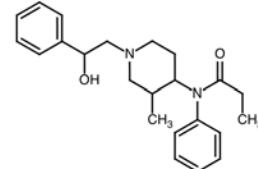
Catalog number : 9831-001

CASRN : 78995-14-9

Name : (±)-cis-β-Hydroxy-3-methylfentanyl hydrochloride

Mol. formula : C₂₃H₃₁CIN₂O₂

FW : 402.97 DEA schedule : 1

Notes : μ -Opioid agonistReferences : Brine, GA; et al. *J Med Chem* 1995, 38, 1547-57.

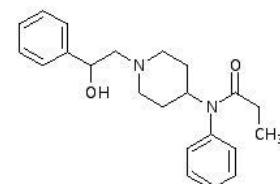
Catalog number : 9830-001

CASRN : 78995-10-5

Name : β-Hydroxyfentanyl hydrochloride

Mol. formula : C₂₂H₂₉CIN₂O₂

FW : 388.93 DEA schedule : 1

Notes : μ -Opioid agonist

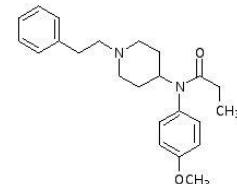
Catalog number : 9801-011

CASRN : 23609-41-8

Name : p-Anisoylfentanyl hydrochloride

Mol. formula : C₂₃H₃₁CIN₂O₂

FW : 402.97 DEA schedule : 2

References : Brine, GA; et al. *J Heterocyclic Chem* 1989, 26, 677.

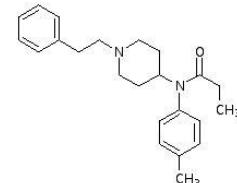
Catalog number : 9801-010

CASRN : 1807-12-1

Name : p-Tolylfentanyl hydrochloride

Mol. formula : C₂₃H₃₁CIN₂O₂

FW : 386.97 DEA schedule : 2

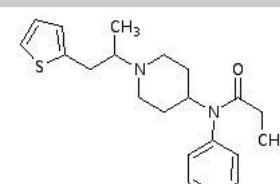


Catalog number : 9832-001

CASRN : 117332-94-2

Name : α -Methylthiofentanyl hydrochlorideMol. formula : C₂₁H₂₉CIN₂OS

FW : 392.99 DEA schedule : 1

Notes : μ -Opioid agonistReferences : Brine, GA; et al. *J Heterocyclic Chem* 1989, 26, 677.

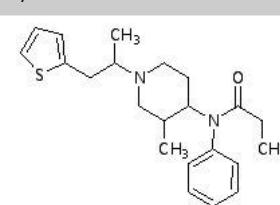
Catalog number : 9833-001

CASRN : n/a

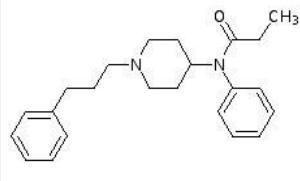
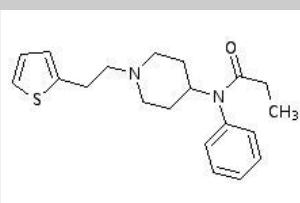
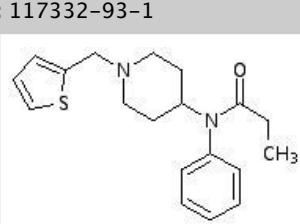
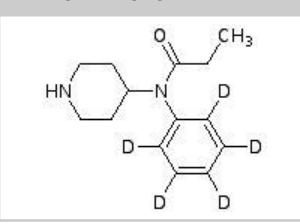
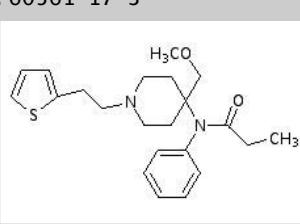
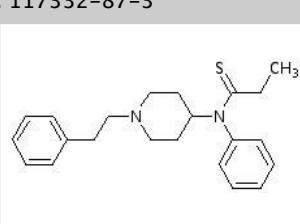
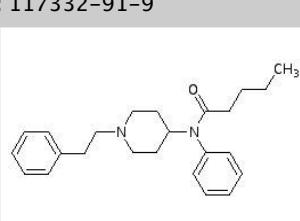
Name : 3-Methylthiofentanyl hydrochloride

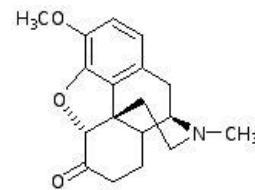
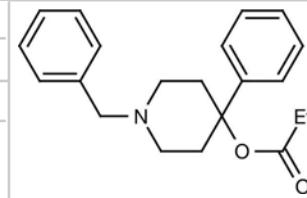
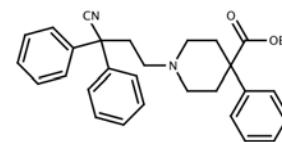
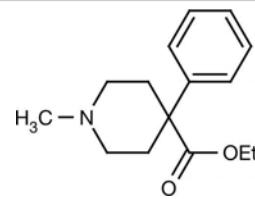
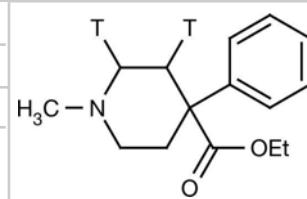
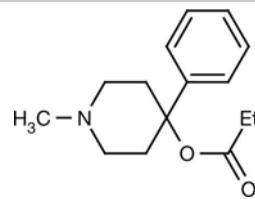
Mol. formula : C₂₁H₂₉CIN₂OS

FW : 392.99 DEA schedule : 1

Notes : μ -Opioid agonistReferences : Brine, GA; et al. *J Heterocyclic Chem* 1989, 26, 677.

6 – Opioids and Opioid-related

Catalog number : 9801-003	CASRN : 117332-88-4	
Name : N-[1-(3-Phenylpropyl)-4-piperidyl]-N-phenylpropanamide hydrochloride		
Mol. formula : C ₂₃ H ₃₁ ClN ₂ O	FW : 386.96	DEA schedule : 2
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1989, 26, 677.		
Catalog number : 9835-001	CASRN : 79278-88-9	
Name : Thiofentanyl hydrochloride		
Mol. formula : C ₂₀ H ₂₇ ClN ₂ O ₅	FW : 378.97	DEA schedule : 1
Notes : μ -Opioid agonist		
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1989, 26, 677.		
Catalog number : 9834-001	CASRN : 117332-93-1	
Name : Thenylfentanyl hydrochloride		
Mol. formula : C ₁₉ H ₂₅ ClN ₂ O ₅	FW : 364.94	DEA schedule : 0
Notes : Pharmacologically inert fentanyl analog.		
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1989, 26, 677. <i>Federal Register</i> 2010, 75(124), 37300.		
Catalog number : 9801-008	CASRN : 1211527-23-9	
Name : [Phenyl- ² H ₅]N-4-Piperidyl-N-phenylpropanamide		
Mol. formula : C ₁₄ H ₂₀ N ₂ O ₅	FW : 237.36	DEA schedule : 2
Notes : Precursor for the synthesis of mass-labeled fentanyl analogs.		
Catalog number : 9740-001	CASRN : 60561-17-3	
Name : Sufentanil citrate		
Mol. formula : C ₂₈ H ₃₈ N ₂ O ₉ S	FW : 578.67	DEA schedule : 2
Notes : Analgesic, μ -Opioid agonist		
References : <i>Merck Index</i> , 14th ed., Monograph 8887. Brine, GA; et al. <i>J Heterocyclic Chem</i> 1989, 26, 677.		
Catalog number : 9801-002	CASRN : 117332-87-3	
Name : N-[1-(2-Phenylethyl)-4-piperidyl]-N-phenylthiopropanamide hydrochloride		
Mol. formula : C ₂₂ H ₂₉ ClN ₂ S	FW : 389.01	DEA schedule : 2
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1989, 26, 677.		
Catalog number : 9801-007	CASRN : 117332-91-9	
Name : Valerylfentanyl hydrochloride		
Mol. formula : C ₂₄ H ₃₃ ClN ₂ O	FW : 400.99	DEA schedule : 2
References : Brine, GA; et al. <i>J Heterocyclic Chem</i> 1989, 26, 677.		

Opioids: Hydrocodone Class**Catalog number :** 9193-002**CASRN :** 34195-34-1**Name :** Dihydrocodeinone tartrate; hydrocodone bitartrate**Mol. formula :** C₂₂H₂₇NO₉**FW :** 449.46 **DEA schedule :** 2**Notes :** Narcotic analgesic; antitussive**References :** Merck Index, 14th ed., Monograph 4785.**Opioids: Meperidine Class****Catalog number :** 9661-002**CASRN :** 63916-24-5**Name :** 1-Benzyl-4-phenyl-4-propionoxypiperidine hydrochloride**Mol. formula :** C₂₁H₂₆CINO₂**FW :** 359.90 **DEA schedule :** 0**Catalog number :** 9170-001**CASRN :** 3810-80-8**Name :** Diphenoxylate hydrochloride**Mol. formula :** C₃₀H₃₃CINO₂**FW :** 489.04 **DEA schedule :** 2**Notes :** Antiperistaltic; antidiarrheal**References :** Merck Index, 14th ed., Monograph 3313.**Catalog number :** 9230-002**CASRN :** 50-13-5**Name :** Meperidine hydrochloride; Demerol hydrochloride**Mol. formula :** C₁₅H₂₂CINO₂**FW :** 283.80 **DEA schedule :** 2**Notes :** Narcotic analgesic; sedative; anesthetic**References :** Merck Index, 14th ed., Monograph 5849.**Catalog number :** 9230-003**Name :** [2,3-³H₂]Meperidine hydrochloride**Mol. formula :** C₁₅H₂₂CINO₂**FW :** 287.81 **DEA schedule :** 2**Catalog number :** 9661-001**CASRN :** 13147-09-6**Name :** 1-Methyl-4-phenyl-4-propionoxypiperidine hydrochloride; Desmethylprodine; MPPP HCl**Mol. formula :** C₁₅H₂₂CINO₂**FW :** 283.80 **DEA schedule :** 1**Notes :** μ -Opioid agonist; narcotic analgesic**References :** Johannessen, JN; Markey SP *Drug Alcohol Depend* 1984, 13, 367-74.

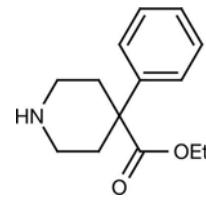
Catalog number : 9230-001

CASRN : 24465-45-0

Name : Normeperidine hydrochloride

Mol. formula : C₁₄H₂₀CINO₂

FW : 269.77 DEA schedule : 0



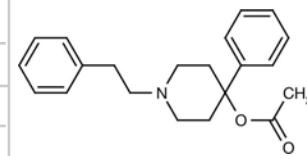
Catalog number : 9663-001

CASRN : 94-30-4

Name : 1-(2-Phenylethyl)-4-phenyl-4-acetoxypiperidine hydrochloride; PEPAP HCl

Mol. formula : C₂₁H₂₆CINO₂

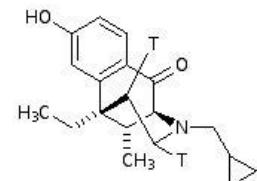
FW : 359.90 DEA schedule : 1

References : Pritzker, D; et al. *J Clin Psychopharmacol* 2002, 22, 330-2.**Opioids: Metazocine Class**

Catalog number : 9240-003

Name : [11,12-³H₂](-)-EthylketazocineMol. formula : C₁₉H₂₅NO₂

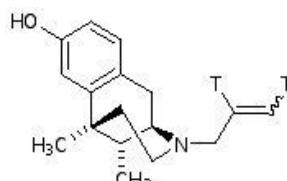
FW : 303.42 DEA schedule : 1



Catalog number : 9240-019

Name : (-)-[17,18-³H]N-AllylnormetazocineMol. formula : C₁₇H₂₃NO

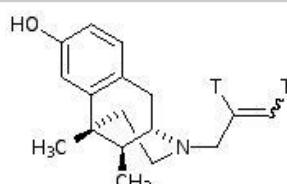
FW : 261.39 DEA schedule : 0



Catalog number : 9240-040

Name : (+)-[17,18-³H]N-AllylnormetazocineMol. formula : C₁₇H₂₃NO

FW : 261.39 DEA schedule : 0



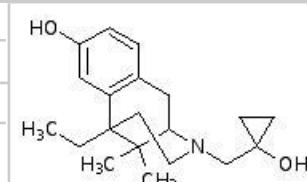
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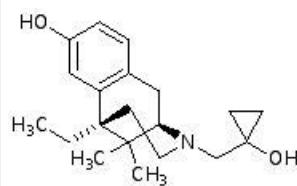
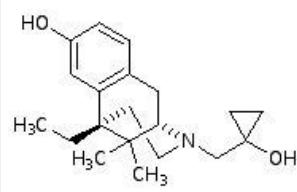
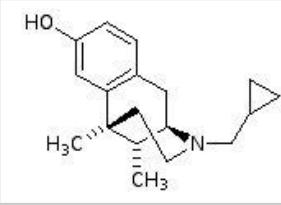
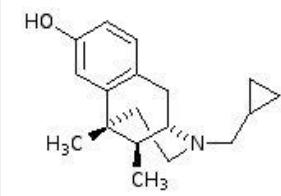
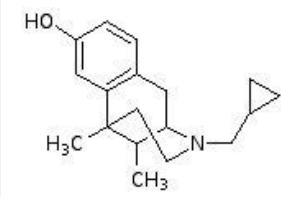
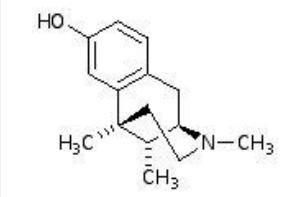
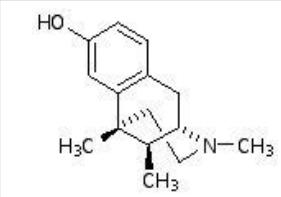
CASRN : 71990-00-6

Name : (\pm)-Bremazocine hydrochlorideMol. formula : C₂₀H₃₀CINO₂

FW : 351.92 DEA schedule : 0

Notes : Kappa-opioid receptor standard ligand.

References : Dortch-Carnes, J; Potter, DE *CNS Drug Rev* 2005, 11, 195-212.

Catalog number : 9240-031**CASRN :** 75684-07-0**Name :** (-)-Bremazocine hydrochloride**Mol. formula :** C₂₀H₃₀CINO₂**FW :** 351.92 **DEA schedule :** 0**Notes :** Kappa-opioid receptor standard ligand.**References :** Dortch-Carnes, J; Potter, DE *CNS Drug Rev* 2005, 11, 195–212.**Catalog number :** 9240-030**Name :** (+)-Bremazocine hydrochloride**Mol. formula :** C₂₀H₃₀CINO₂**FW :** 351.92 **DEA schedule :** 0**Notes :** Kappa-opioid receptor standard ligand.**References :** Dortch-Carnes, J; Potter, DE *CNS Drug Rev* 2005, 11, 195–212.**Catalog number :** 9240-011**Name :** (-)-Cyclazocine**Mol. formula :** C₁₈H₂₅NO**FW :** 271.39 **DEA schedule :** 0**Notes :** Narcotic antagonist**Catalog number :** 9240-012**Name :** (+)-Cyclazocine**Mol. formula :** C₁₈H₂₅NO**FW :** 271.39 **DEA schedule :** 0**Notes :** Narcotic antagonist**References :** Carroll, FI; et al. *J Med Chem* 1992, 35, 2812–8.**Catalog number :** 9240-010**CASRN :** 3572-80-3**Name :** (±)-Cyclazocine**Mol. formula :** C₁₈H₂₅NO**FW :** 271.39 **DEA schedule :** 0**Notes :** Narcotic antagonist**References :** Merck Index, 14th ed., Monograph 2705.
Carroll, FI; et al. *J Med Chem* 1992, 35, 2812–8.**Catalog number :** 9240-008**CASRN :** 133005-39-7**Name :** (-)-Metazocine fumarate**Mol. formula :** C₃₄H₄₆N₂O₆**FW :** 578.75 **DEA schedule :** 2**References :** Carroll, FI; et al. *J Med Chem* 1992, 35, 2812–8.**Catalog number :** 9240-007**CASRN :** 133005-40-0**Name :** (+)-Metazocine fumarate**Mol. formula :** C₃₄H₄₆N₂O₆**FW :** 578.75 **DEA schedule :** 2**References :** Carroll, FI; et al. *J Med Chem* 1992, 35, 2812–8.

6 – Opioids and Opioid-related

Catalog number : 9240-018	CASRN : 14198-28-8	
Name : (-)-N-Allylnormetazocine hydrochloride, (-)-NANM; (-)-SKF-10,047		
Mol. formula : C ₁₇ H ₂₄ CINO	FW : 293.85	DEA schedule : 0
Notes : Sigma receptor standard ligand.		
References : Carroll, FI; et al. <i>J Med Chem</i> 1992, 35, 2812-8.		
Catalog number : 9240-017	CASRN : 133005-41-1	
Name : (+)-N-Allylnormetazocine hydrochloride, (+)-NANM; (+)-SKF-10,047		
Mol. formula : C ₁₇ H ₂₄ CINO	FW : 293.85	DEA schedule : 0
Notes : Sigma receptor standard ligand.		
References : Carroll, FI; et al. <i>J Med Chem</i> 1992, 35, 2812-8.		
Catalog number : 9240-016	CASRN : 7619-35-4	
Name : (±)-N-Allylnormetazocine hydrochloride; (±)-NANM; (±)-SKF-10,047; Alazocine		
Mol. formula : C ₁₇ H ₂₄ CINO	FW : 293.85	DEA schedule : 0
Notes : Sigma receptor standard ligand.		
Catalog number : 9240-004	CASRN : 16603-67-1	
Name : (-)- <i>cis</i> -Normetazocine		
Mol. formula : C ₁₄ H ₁₉ NO	FW : 217.31	DEA schedule : 0
References : Carroll, FI; et al. <i>J Med Chem</i> 1992, 35, 2812-8.		
Catalog number : 9240-005	CASRN : 16670-83-0	
Name : (+)- <i>cis</i> -Normetazocine		
Mol. formula : C ₁₄ H ₁₉ NO	FW : 217.31	DEA schedule : 0
References : Carroll, FI; et al. <i>J Med Chem</i> 1992, 35, 2812-8.		
Catalog number : 9240-002	CASRN : 16808-63-2	
Name : (±)-Normetazocine		
Mol. formula : C ₁₄ H ₁₉ NO	FW : 217.31	DEA schedule : 0
References : Carroll, FI; et al. <i>J Med Chem</i> 1992, 35, 2812-8.		
Catalog number : 9709-003	CASRN : 124819-25-6	
Name : (-)-Pentazocine succinate		
Mol. formula : C ₂₃ H ₃₃ NO ₅	FW : 403.52	DEA schedule : 4
Notes : Sigma receptor standard ligand.		
References : Merck Index, 14th ed., Monograph 7121. Brogden, RN; Speight, TM; Avery, GS <i>Drugs</i> 1973, 5, 6-91. Carroll, FI; et al. <i>J Med Chem</i> 1992, 35, 2812-8.		

Catalog number : 9709-002

CASRN : 124819-26-7

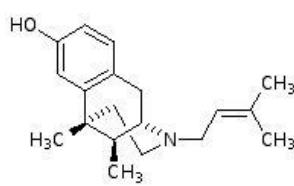
Name : (+)-Pentazocine succinate

Mol. formula : C₂₃H₃₃NO₅

FW : 403.52 DEA schedule : 4

Notes : Sigma receptor standard ligand.

References : Merck Index, 14th ed., Monograph 7121.

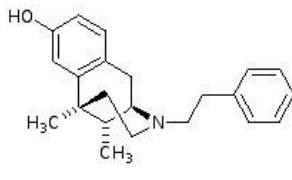
Brogden, RN; Speight, TM; Avery, GS *Drugs* 1973, 5, 6–91.Carroll, FI; et al. *J Med Chem* 1992, 35, 2812–8.

Catalog number : 9715-003

Name : (-)-Phenazocine hydrobromide

Mol. formula : C₂₂H₂₈BrNO

FW : 402.39 DEA schedule : 2

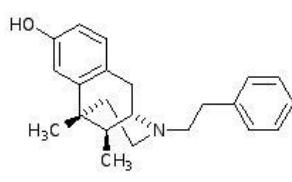


Catalog number : 9715-002

Name : (+)-Phenazocine hydrobromide

Mol. formula : C₂₂H₂₈BrNO

FW : 402.39 DEA schedule : 2



Catalog number : 9715-001

CASRN : 1239-04-9

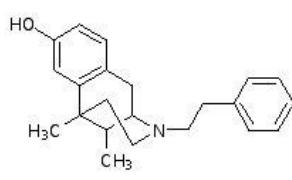
Name : (±)-Phenazocine hydrobromide

Mol. formula : C₂₂H₂₈BrNO

FW : 402.39 DEA schedule : 2

Notes : Narcotic analgesic, σ receptor ligand

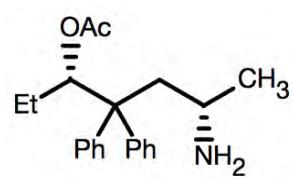
References : Merck Index, 14th ed., Monograph 7218.

**Opioids: Methadone Class**

Catalog number : 9633-010

Name : (-)- α -Acetyl-N,N-dinormethadol hydrochlorideMol. formula : C₂₁H₂₈ClNO₂

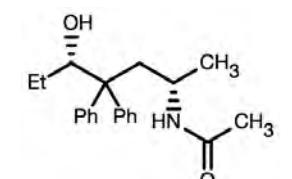
FW : 361.92 DEA schedule : 0

References : Carroll, FI; et al. *J Org Chem* 1976, 41, 3521–4.

Catalog number : 9633-050

Name : (-)- α -N-Acetyl-N,N-dinormethadolMol. formula : C₂₁H₂₇NO₂

FW : 325.45 DEA schedule : 1

References : Carroll, FI; et al. *J Org Chem* 1976, 41, 3521–4.

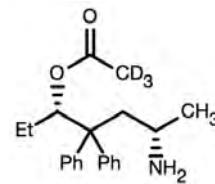
6 – Opioids and Opioid-related

Catalog number : 9633-012

Name : (–)-[Acetyl-²H₃] α -Acetyl-N,N-dinormethadol hydrochloride

Mol. formula : C₂₁H₂₈ClNO₂

FW : 361.92 **DEA schedule :** 0

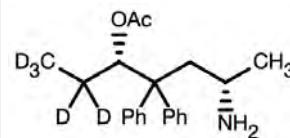


Catalog number : 9633-014

Name : (–)-[1,1,1,2,2,3-²H₆] α -Acetyl-N,N-dinormethadol hydrochloride

Mol. formula : C₂₁H₂₈ClNO₂

FW : 361.92 **DEA schedule :** 0

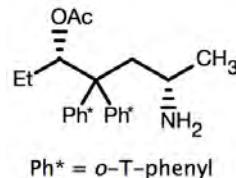


Catalog number : 9633-011

Name : (–)-[o,o'-³H₂(n)]- α -Acetyl-N,N-dinormethadol hydrochloride

Mol. formula : C₂₁H₂₈ClNO₂

FW : 389.97 **DEA schedule :** 0

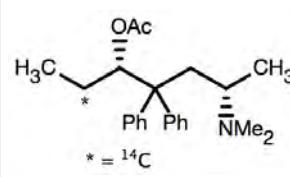


Catalog number : 9648-013

Name : (–)-[2-¹⁴C]- α -Acetylmethadol hydrochloride

Mol. formula : C₂₃H₃₂ClNO₂

FW : 353.50 **DEA schedule :** 2

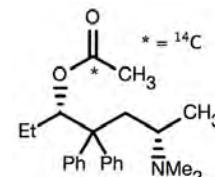


Catalog number : 9648-014

Name : (–)-[Acetyl-¹⁴C]- α -Acetylmethadol

Mol. formula : C₂₃H₃₂ClNO₂

FW : 353.50 **DEA schedule :** 2

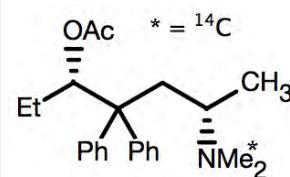


Catalog number : 9648-015

Name : (–)-[N-¹⁴CH₃]- α -Acetylmethadol hydrochloride

Mol. formula : C₂₃H₃₁NO₂

FW : 353.50 **DEA schedule :** 2

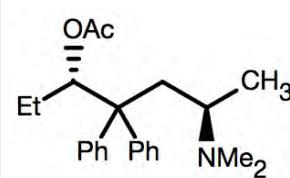


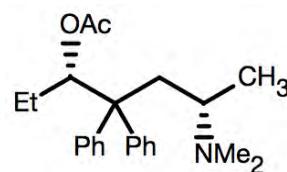
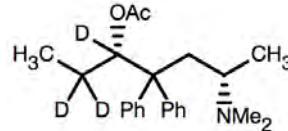
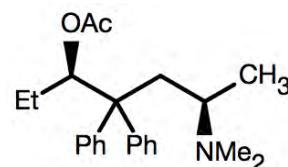
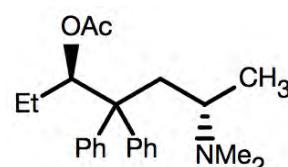
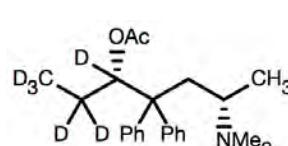
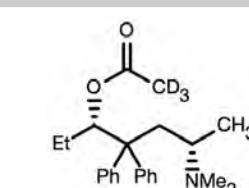
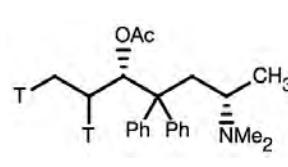
Catalog number : 9607-001

Name : (–)- β -Acetylmethadol hydrochloride

Mol. formula : C₂₃H₃₂ClNO₂

FW : 389.97 **DEA schedule :** 1



Catalog number : 9648-001**CASRN :** 43033-72-3**Name :** (–)- α -Acetylmethadol hydrochloride; LAAM hydrochloride**Mol. formula :** C₂₃H₃₂CINO₂**FW :** 389.97 **DEA schedule :** 2**Catalog number :** 9648-016**Name :** (–)[2,2,3–²H₃] α -Acetylmethadol hydrochloride**Mol. formula :** C₂₃H₃₂CINO₂**FW :** 353.50 **DEA schedule :** 2**Catalog number :** 9603-003**Name :** (+)- α -Acetylmethadol hydrochloride**Mol. formula :** C₂₃H₃₂CINO₂**FW :** 389.97 **DEA schedule :** 1**Catalog number :** 9607-002**Name :** (+)- β -Acetylmethadol hydrochloride**Mol. formula :** C₂₃H₃₂CINO₂**FW :** 389.97 **DEA schedule :** 1**Catalog number :** 9648-010**Name :** (–)[1,1,1,2,2,3–²H₆] α -Acetylmethadol hydrochloride**Mol. formula :** C₂₄H₃₄CINO**FW :** 389.97 **DEA schedule :** 2**Catalog number :** 9648-011**Name :** (–)[Acetyl–²H₃] α -acetylmethadol hydrochloride**Mol. formula :** C₂₁H₂₈CINO₂**FW :** 389.97 **DEA schedule :** 2**Catalog number :** 9648-012**Name :** (–)[1,2–³H₂] α -Acetylmethadol**Mol. formula :** C₂₃H₃₂CINO₂**FW :** 353.50 **DEA schedule :** 2

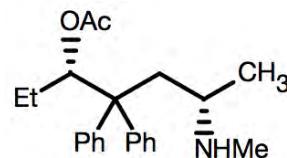
6 – Opioids and Opioid-related

Catalog number : 9633-001

Name : (-)- α -Acetylnormethadol hydrochloride

Mol. formula : C₂₂H₃₀ClNO₂

FW : 375.94 **DEA schedule :** 1



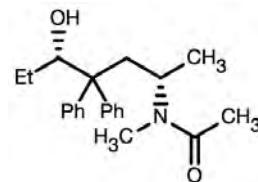
References : Carroll, FI; et al. *J Org Chem* 1976, 41, 3521-4.

Catalog number : 9633-040

Name : (-)- α -N-Acetylnormethadol

Mol. formula : C₂₂H₂₉NO₂

FW : 339.48 **DEA schedule :** 1

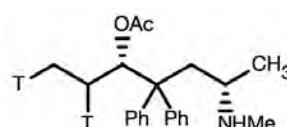


Catalog number : 9633-006

Name : (-)-[1,2-³H] α -Acetylnormethadol hydrochloride

Mol. formula : C₂₃H₃₂ClNO₂

FW : 375.94 **DEA schedule :** 1

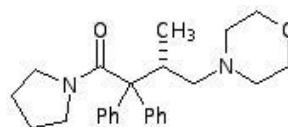


Catalog number : 9613-001

Name : Dextromoramide tartrate

Mol. formula : C₂₅H₃₂N₂O₂

FW : 542.60 **DEA schedule :** 1

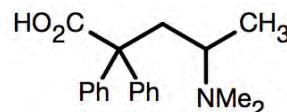


Catalog number : 9254-004

Name : (\pm)-4-Dimethylamino-2,2-diphenylvaleric acid

Mol. formula : C₁₉H₂₃NO₂

FW : 297.40 **DEA schedule :** 0

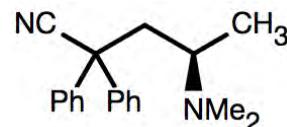


Catalog number : 9254-002

Name : (-)-(R)-4-Dimethylamino-2,2-diphenylvaleronitrile

Mol. formula : C₁₉H₂₂N₂

FW : 278.40 **DEA schedule :** 2

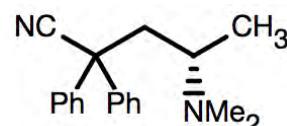


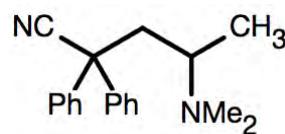
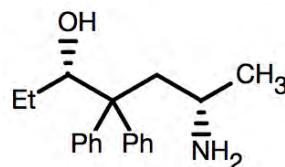
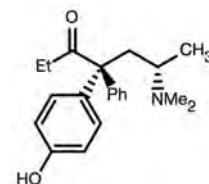
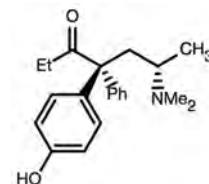
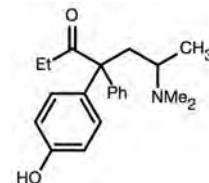
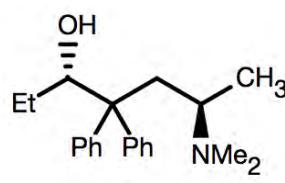
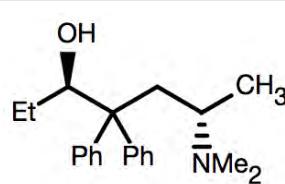
Catalog number : 9254-001

Name : (+)-(S)-4-Dimethylamino-2,2-diphenylvaleronitrile

Mol. formula : C₁₉H₂₂N₂

FW : 278.40 **DEA schedule :** 2



Catalog number : 9254-003**Name :** (\pm)-4-Dimethylamino-2,2-diphenylvaleronitrile**Mol. formula :** C₁₉H₂₂N₂**FW :** 278.40 **DEA schedule :** 2**Catalog number :** 9605-030**Name :** (-)- α -N,N-Dinormethadol maleate**Mol. formula :** C₂₃H₂₉NO₅**FW :** 399.49 **DEA schedule :** 1**References :** Carroll, FI; et al. *J Org Chem* 1976, 41, 3521–4.**Catalog number :** 9250-061**Name :** (4R,6S)-*p*-Hydroxymethadone hydrochloride**Mol. formula :** C₂₁H₂₈ClNO₂**FW :** 361.91 **DEA schedule :** 2**Catalog number :** 9250-062**Name :** (4S,6S)-*p*-Hydroxymethadone hydrochloride**Mol. formula :** C₂₁H₂₈ClNO₂**FW :** 361.91 **DEA schedule :** 2**Catalog number :** 9250-060**Name :** (4RS,6RS)-*p*-Hydroxymethadone hydrochloride**Mol. formula :** C₂₁H₂₈ClNO₂**FW :** 361.91 **DEA schedule :** 2**Catalog number :** 9609-001**CASRN :** 17199-55-2**Name :** (-)- β -Methadol**Mol. formula :** C₂₁H₂₉NO**FW :** 311.47 **DEA schedule :** 1**References :** Portoghese, PS; Williams DA *J Med Chem* 1969, 12, 839–44.**Catalog number :** 9609-002**Name :** (+)- β -Methadol**Mol. formula :** C₂₁H₂₉NO**FW :** 311.47 **DEA schedule :** 1**References :** Portoghese, PS; Williams DA *J Med Chem* 1969, 12, 839–44.

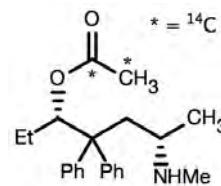
6 – Opioids and Opioid-related

Catalog number : 9633-005

Name : (–)[Acetyl- $^{14}\text{C}_2$]- α -Acetylnormethadol hydrochloride

Mol. formula : $\text{C}_{22}\text{H}_{30}\text{ClNO}_2$

FW : 375.94 **DEA schedule :** 1



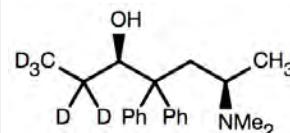
Catalog number : 9605-010

Name : [1,1,1,2,2- $^2\text{H}_5$] α -Methadol hydrochloride

Mol. formula : $\text{C}_{21}\text{H}_{30}\text{ClNO}$

FW : 347.93 **DEA schedule :** 1

References : Portoghese, PS; Williams DA *J Med Chem* 1969, 12, 839–44.

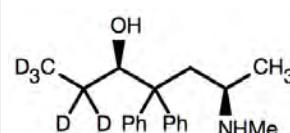


Catalog number : 9605-021

Name : [1,1,1,2,2- $^2\text{H}_5$] α -Normethadol perchlorate

Mol. formula : $\text{C}_{20}\text{H}_{28}\text{ClNO}_5$

FW : 397.90 **DEA schedule :** 1

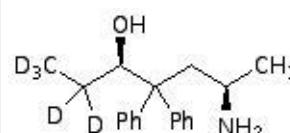


Catalog number : 9605-031

Name : [1,1,1,2,2- $^2\text{H}_5$] α -N,N-Dinormethadol maleate

Mol. formula : $\text{C}_{19}\text{H}_{25}\text{NO}$

FW : 399.49 **DEA schedule :** 1

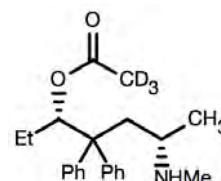


Catalog number : 9633-003

Name : (–)[Acetyl- $^2\text{H}_3$] α -Acetylnormethadol hydrochloride

Mol. formula : $\text{C}_{22}\text{H}_{30}\text{ClNO}_2$

FW : 375.94 **DEA schedule :** 1

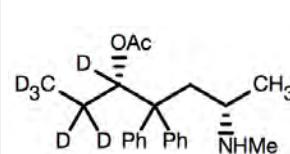


Catalog number : 9633-004

Name : (–)[1,1,1,2,2,3- $^2\text{H}_6$] α -Acetylnormethadol hydrochloride

Mol. formula : $\text{C}_{23}\text{H}_{22}\text{ClNO}_2$

FW : 375.94 **DEA schedule :** 1



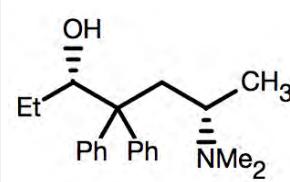
Catalog number : 9605-001

Name : (–) α -Methadol hydrochloride

Mol. formula : $\text{C}_{21}\text{H}_{30}\text{ClNO}$

FW : 347.93 **DEA schedule :** 1

References : Portoghese, PS; Williams DA *J Med Chem* 1969, 12, 839–44.

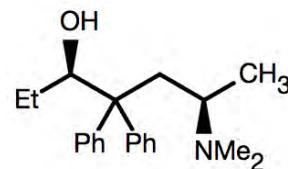


Catalog number : 9605-002

CASRN : 17199-54-1

Name : (+)- α -Methadol hydrochlorideMol. formula : C₂₁H₃₀CINO

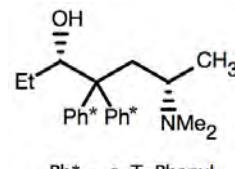
FW : 347.93 DEA schedule : 1

References : Portoghesi, PS; Williams DA *J Med Chem* 1969, 12, 839–44.

Catalog number : 9605-003

Name : (−)-[o,o'−³H₂(n)] α -Methadol hydrochlorideMol. formula : C₂₁H₂₉NO

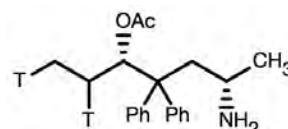
FW : 347.93 DEA schedule : 1

References : Portoghesi, PS; Williams DA *J Med Chem* 1969, 12, 839–44.

Catalog number : 9633-015

Name : (−)-[1,2−³H₂] α -Acetyl-N,N-dinormethadol hydrochlorideMol. formula : C₂₁H₂₈CINO₂

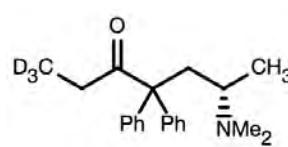
FW : 361.9–2 DEA schedule : 0



Catalog number : 9250-011

Name : (+)-[1,1,1−²H₃]Methadone hydrochlorideMol. formula : C₂₁H₂₈CINO

FW : 348.92 DEA schedule : 2



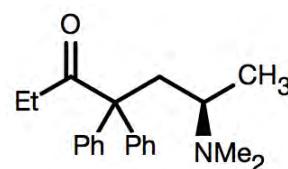
Catalog number : 9250-002

CASRN : 125-58-6

Name : (−)-(R)-Methadone

Mol. formula : C₂₁H₂₇NO

FW : 309.46 DEA schedule : 2

References : *Merck Index*, 14th ed., Monograph 5944.

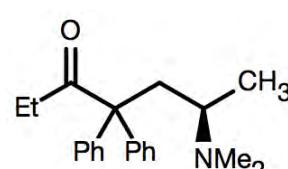
Catalog number : 9250-005

CASRN : 125-58-6

Name : (−)-((R)-Methadone hydrochloride

Mol. formula : C₂₁H₂₈CINO

FW : 345.92 DEA schedule : 2



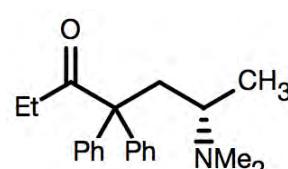
Catalog number : 9250-001

CASRN : 5653-80-5

Name : (+)-(S)-Methadone

Mol. formula : C₂₁H₂₇NO

FW : 309.46 DEA schedule : 2



6 – Opioids and Opioid-related

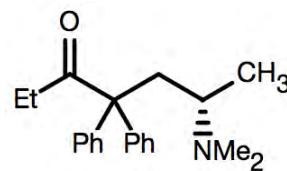
Catalog number : 9250-004

CASRN : 5653-80-5

Name : (+)-(S)-Methadone hydrochloride

Mol. formula : C₂₁H₂₈ClNO

FW : 345.92 **DEA schedule :** 2



Catalog number : 9250-003

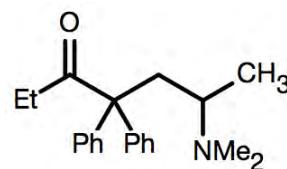
CASRN : 1095-90-5

Name : (±)-Methadone hydrochloride

Mol. formula : C₂₁H₂₈ClNO

FW : 345.92 **DEA schedule :** 2

References : Merck Index, 14th ed., Monograph 5944.

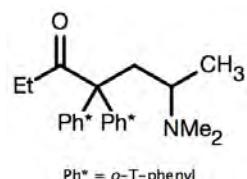


Catalog number : 9250-006

Name : (±)-[o,o'-³H₂(n)]Methadone

Mol. formula : C₂₁H₂₇NO

FW : 313.46 **DEA schedule :** 2



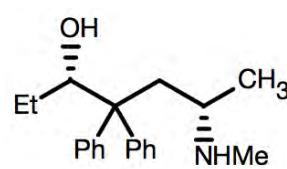
Catalog number : 9605-020

Name : (−)-α-Normethadol perchlorate

Mol. formula : C₂₀H₂₈ClNO₅

FW : 397.90 **DEA schedule :** 1

References : Carroll, FI; et al. *J Org Chem* 1976, 41, 3521-4.

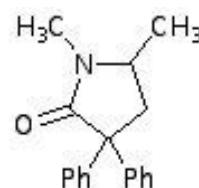


Catalog number : 9254-005

Name : (±)-1,5-Dimethyl-3,3-diphenyl-2-pyrrolidone

Mol. formula : C₁₈H₁₉NO

FW : 265.36 **DEA schedule :** 0

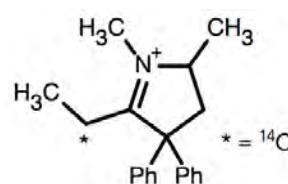


Catalog number : 9250-026

Name : [1'-¹⁴C]-2-Ethyl-1,5-dimethyl-3,3-diphenylpyrrolinium perchlorate

Mol. formula : C₂₀H₂₅ClNO₄

FW : 536.01 **DEA schedule :** 0

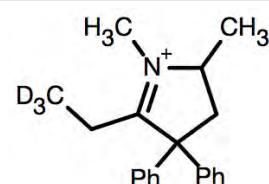


Catalog number : 9250-027

Name : [Ethyl-2',2',2'-²H₃]-1,5-Dimethyl-3,3-diphenyl-2-ethylpyrrolinium perchlorate

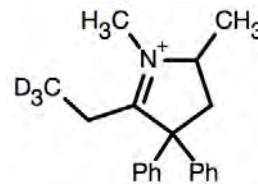
Mol. formula : C₂₀H₂₅ClNO₄

FW : 384.91 **DEA schedule :** 0

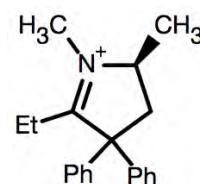


Catalog number : 9250-032

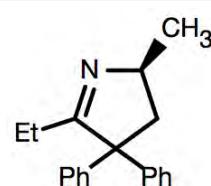
Name : [Ethyl-2',2',2'-²H₃]-3,3-Diphenyl-2-ethyl-5-methyl-1-pyrroline hydrochloride

Mol. formula : C₂₀H₂₅CIN**FW :** 302.86 **DEA schedule :** 0**Catalog number :** 9250-024

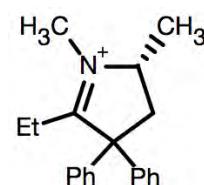
Name : (–)-(S)-2-Ethyl-1,5-dimethyl-3,3-diphenylpyrrolinium perchlorate

Mol. formula : C₂₀H₂₄CINO₄**FW :** 377.88 **DEA schedule :** 0**Catalog number :** 9250-033

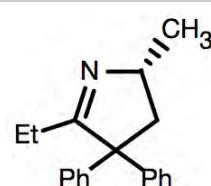
Name : (–)-(S)-2-Ethyl-5-methyl-3,3-diphenyl-1-pyrroline hydrochloride

Mol. formula : C₁₉H₂₂CIN**FW :** 299.85 **DEA schedule :** 0**Catalog number :** 9250-025

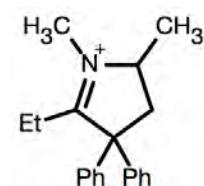
Name : (+)-(R)-2-Ethyl-1,5-dimethyl-3,3-diphenylpyrrolinium perchlorate

Mol. formula : C₂₀H₂₄CINO₄**FW :** 377.88 **DEA schedule :** 0**Catalog number :** 9250-034

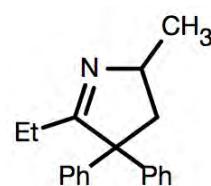
Name : (+)-(R)-2-Ethyl-5-methyl-3,3-diphenyl-1-pyrroline hydrochloride

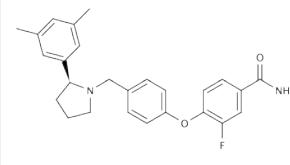
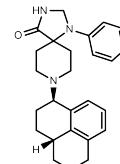
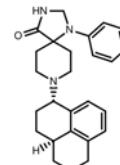
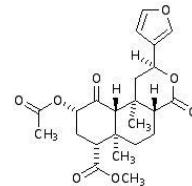
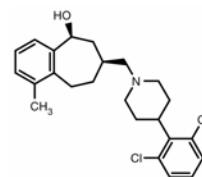
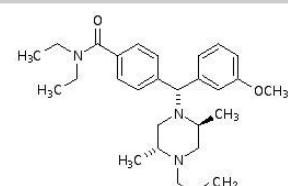
Mol. formula : C₁₉H₂₂CIN**FW :** 299.85 **DEA schedule :** 0**Catalog number :** 9250-021

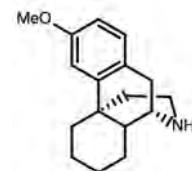
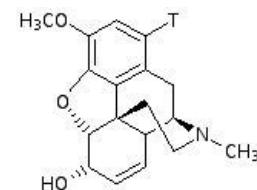
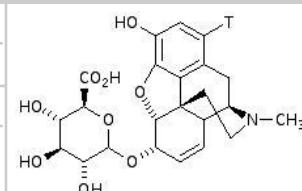
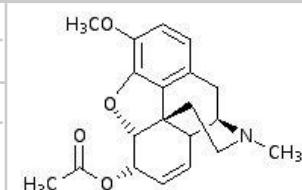
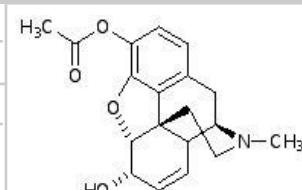
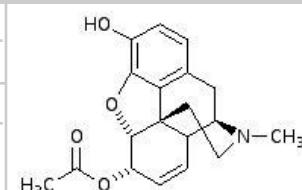
Name : (±)-2-Ethyl-1,5-dimethyl-3,3-diphenylpyrrolinium perchlorate

Mol. formula : C₂₀H₂₄CINO₄**FW :** 377.86 **DEA schedule :** 0**Catalog number :** 9250-031

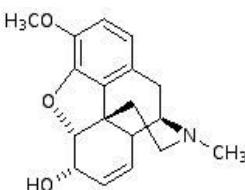
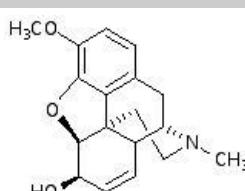
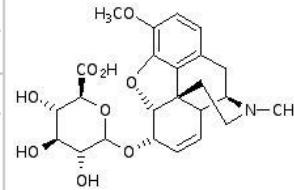
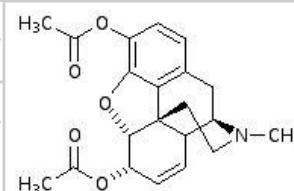
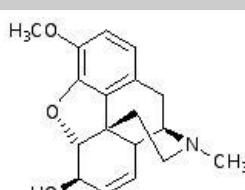
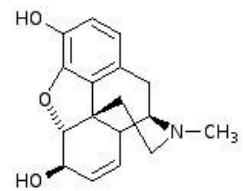
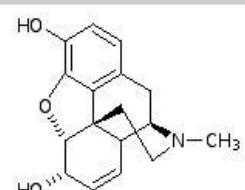
Name : (±)-2-Ethyl-5-methyl-3,3-diphenyl-1-pyrroline hydrochloride

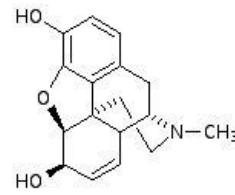
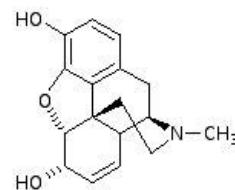
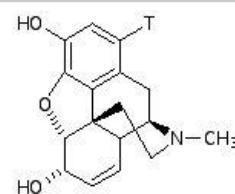
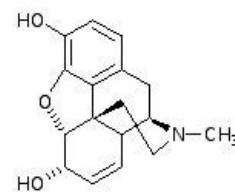
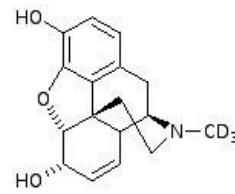
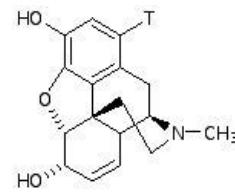
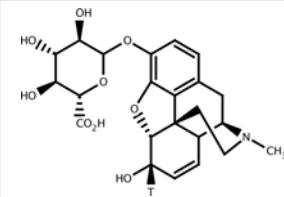
Mol. formula : C₁₉H₂₂CIN**FW :** 299.85 **DEA schedule :** 0

Opioids: Miscellaneous**Catalog number :** NOCD-107**CASRN :** 1174130-61-0**Name :** LY-2456302**Mol. formula :** C₂₆H₂₈ClFN₂O₂**FW :** 454.97 **DEA schedule :** 0**Notes :** κ -Opioid antagonist.**References :** Peters, MF; et al., *Eur J Pharmacol* 2011, 661, 27–34.**Catalog number :** NOCD-138**CASRN :** 309254-79-3**Name :** Ro 64-6198 hydrochloride**Mol. formula :** C₂₆H₃₁N₃O • HCl**FW :** 438.02 **DEA schedule :** 0**Notes :** Potent nociceptin opioid receptor agonist.**References :** Wichmann, J; et al., *Eur. J. Med. Chem.* 2000, 35, 839–851.**Catalog number :** NOCD-139**Name :** Ro 64-6198 analog [(+)-isomer of Ro 64-6198]**Mol. formula :** C₂₆H₃₁N₃O • HCl**FW :** 438.02 **DEA schedule :** 0**Notes :** Nociceptin opioid receptor agonist. 32-fold less potent (+)-enantiomer of Ro 64-6198.**References :** Jenck, F; et al., *Proc Natl Acad Sci USA* 2000, 97, 4938–4943.**Catalog number :** NOCD-099**CASRN :** 83729-01-5**Name :** Salvinorin A**Mol. formula :** C₂₃H₂₈O₈**FW :** 432.47 **DEA schedule :** 0**Notes :** Hallucinogen; κ -opioid agonist.**References :** Roth, BL; et al. *Proc Natl Acad Sci USA* 2002, 99, 11934–9.**Catalog number :** NOCD-134**CASRN :** 371980-98-2**Name :** SB-612111 hydrochloride**Mol. formula :** C₂₄H₂₉Cl₂NO • HCl**FW :** 454.87 **DEA schedule :** 0**Notes :** Nociceptin/orphanin FQ (NOP) receptor antagonist.**References :** Zaratin, PF; et al., *J Pharmacol Exp Ther* 2004, 308, 454–61.**Catalog number :** NOCD-055**CASRN :** 156727-74-1**Name :** SNC 80**Mol. formula :** C₂₈H₃₉N₃O₂**FW :** 449.64 **DEA schedule :** 0**Notes :** Highly-selective δ -opioid receptor agonist.**References :** Do Carmo, GP; et al. *Eur J Pharmacol* 2006, 547, 92–100.

Opioids: Morphinan Class**Catalog number :** 9732-001**CASRN :** 1087-69-0**Name :** (+)-3-Methoxymorphinan hydrochloride**Mol. formula :** C₁₇H₂₃ClNO**FW :** 293.84**DEA schedule :** 0**Opioids: Morphine Class****Catalog number :** 9050-011**Name :** [1-³H]Codeine**Mol. formula :** C₁₈H₂₁NO₃**FW :** 301.37**DEA schedule :** 2**Catalog number :** 9300-014**Name :** [1-³H]Morphine-6-glucuronide**Mol. formula :** C₂₃H₂₇NO₉**FW :** 463.47**DEA schedule :** 2**Catalog number :** 9050-004**CASRN :** 6703-27-1**Name :** 6-Acetylcodeine**Mol. formula :** C₂₀H₂₃NO**FW :** 341.41**DEA schedule :** 2**Catalog number :** 9300-010**Name :** 3-Acetylmorphine sulfamate**Mol. formula :** C₁₉H₂₄N₂O₇S**FW :** 424.47**DEA schedule :** 2**Catalog number :** 9300-004**Name :** 6-Acetylmorphine**Mol. formula :** C₁₉H₂₁NO₄**FW :** 327.38**DEA schedule :** 2

6 – Opioids and Opioid-related

Catalog number : 9050-001		CASRN : 1422-07-7
Name : Codeine hydrochloride		
Mol. formula : C ₁₈ H ₂₂ ClNO ₃	FW : 335.84	DEA schedule : 2
Notes : Narcotic analgesic; antitussive		
References : Merck Index, 14th ed., Monograph 2455.		
		
Catalog number : 9050-010		CASRN : 76-57-3
Name : (+)-Codeine base		
Mol. formula : C ₁₈ H ₂₁ NO ₃	FW : 299.36	DEA schedule : 2
		
Catalog number : 9050-009		
Name : Codeine-6-β-D-glucuronide		
Mol. formula : C ₂₄ H ₂₉ NO ₉	FW : 475.50	DEA schedule : 2
		
Catalog number : 9200-001		CASRN : 1502-95-0
Name : 3,6-Diacetylmorphine hydrochloride; Diamorphine HCl		
Mol. formula : C ₂₁ H ₂₄ ClNO ₅	FW : 405.88	DEA schedule : 1
Notes : Narcotic analgesic		
References : Merck Index, 14th ed., Monograph 2968.		
		
Catalog number : 9050-006		CASRN : 509-64-8
Name : Isocodeine		
Mol. formula : C ₁₈ H ₂₁ NO ₃	FW : 299.35	DEA schedule : 2
		
Catalog number : 9300-020		CASRN : 143-70-4
Name : α-Isomorphine		
Mol. formula : C ₁₇ H ₁₉ NO ₃	FW : 285.35	DEA schedule : 2
		
Catalog number : 9300-007		CASRN : 57-27-2
Name : Morphine base		
Mol. formula : C ₁₇ H ₁₉ NO ₃	FW : 285.33	DEA schedule : 2
Notes : Narcotic analgesic; antitussive; antiperistaltic		
References : Merck Index, 14th ed., Monograph 6276.		
		

Catalog number : 9300-012**CASRN :** 65165-99-3**Name :** (+)-Morphine base**Mol. formula :** C₁₇H₁₉NO₃**FW :** 285.35 **DEA schedule :** 2**Catalog number :** 9300-016**CASRN :** 52-26-6**Name :** Morphine hydrochloride monohydrate**Mol. formula :** C₁₇H₂₀ClNO₃**FW :** 321.81 **DEA schedule :** 2**References :** *Merck Index*, 14th ed., Monograph 6276.**Catalog number :** 9300-002**Name :** Tritium-labeled Morphine sulfate**Mol. formula :** C₁₇H₁₉NO₃**FW :** 475.50 **DEA schedule :** 2**Catalog number :** 9300-001**CASRN :** 6211-15-0**Name :** Morphine sulfate pentahydrate**Mol. formula :** C₃₄H₄₀N₂O₁₀S**FW :** 758.83 **DEA schedule :** 2**Notes :** Narcotic analgesic; prototypic μ opioid receptor agonist; sedative**References :** *Merck Index*, 14th ed., Monograph 6276.**Catalog number :** 9300-005**CASRN :** 67293-88-3**Name :** [N-C²H₃]Morphine**Mol. formula :** C₁₇H₁₉NO₃**FW :** 288.36 **DEA schedule :** 2**Catalog number :** 9300-006**CASRN :** 80573-75-7**Name :** [1-³H(n)]Morphine**Mol. formula :** C₁₇H₁₉NO₃**FW :** 287.35 **DEA schedule :** 2**Catalog number :** 9300-011**Name :** Morphine-(6-³H)-3-glucuronide**Mol. formula :** C₂₃H₂₇NO₉**FW :** 463.47 **DEA schedule :** 2

6 – Opioids and Opioid-related

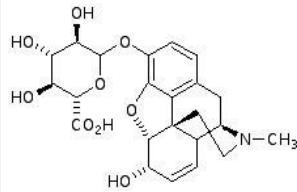
Catalog number : 9300-003

Name : Morphine-3- β -D-glucuronide

Mol. formula : C₂₃H₂₇NO₉

FW : 461.44 DEA schedule : 2

References : Berrang, B; et al. *Synthetic Communications* 1975, 5, 231–236.

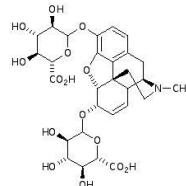


Catalog number : 9300-015

Name : Morphine-3,6-di- β -D-glucuronide, monolithium salt

Mol. formula : C₂₉H₃₄LiNO₁₅

FW : 643.52 DEA schedule : 2



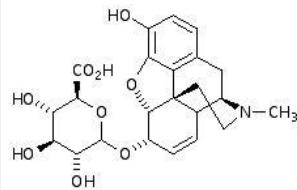
Catalog number : 9300-013

CASRN : 20290-10-2

Name : Morphine-6- β -D-glucuronide

Mol. formula : C₂₃H₂₇NO₉

FW : 461.47 DEA schedule : 2



Catalog number : 9400-001

CASRN : 62-67-9

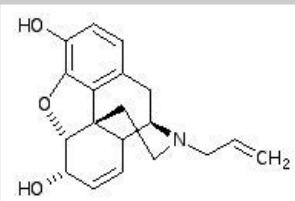
Name : Nalorphine hydrochloride

Mol. formula : C₁₉H₂₂ClNO₃

FW : 347.84 DEA schedule : 3

Notes : Narcotic antagonist

References : *Merck Index*, 14th ed., Monograph 6361.

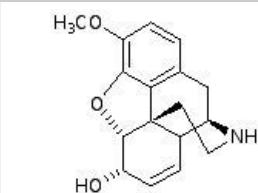


Catalog number : 9050-002

Name : Norcodeine hydrochloride

Mol. formula : C₁₇H₂₀ClNO₃

FW : 321.81 DEA schedule : 2



Catalog number : 9313-001

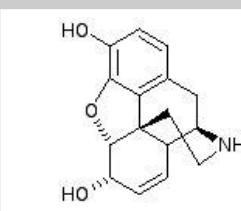
CASRN : 3372-02-9

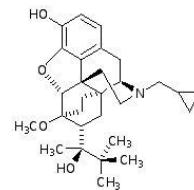
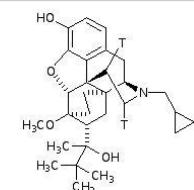
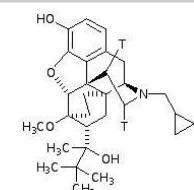
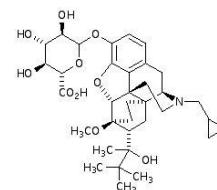
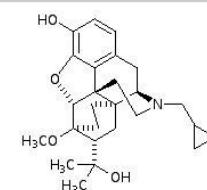
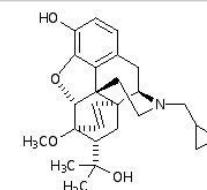
Name : Normorphine hydrochloride

Mol. formula : C₁₆H₁₈ClNO₃

FW : 307.78 DEA schedule : 1

References : *Merck Index*, 14th ed., Monograph 6711.



Opioids: Orvinol Class**Catalog number :** 9064-001**CASRN :** 53152-21-9**Name :** Buprenorphine hydrochloride**Mol. formula :** C₂₉H₄₂ClNO₄**FW :** 504.11**DEA schedule :** 3**Notes :** Narcotic Analgesic**References :** Robinson, SE *CNS Drug Rev* 2002, 8, 377–90.**Catalog number :** 9064-002**CASRN :** 161772-95-8**Name :** [15,16-³H₂]Buprenorphine hydrochloride**Mol. formula :** C₂₆H₃₈ClNO₄**FW :** 508.11**DEA schedule :** 5**Notes :** Narcotic analgesic (tritium-labeled).**References :** Robinson, SE *CNS Drug Rev* 2002, 8, 377–90.**Catalog number :** 9064-003**CASRN :** 161772-95-8**Name :** [15,16-³H₂]Buprenorphine**Mol. formula :** C₂₉H₄₁NO₄**FW :** 508.11**DEA schedule :** 3**Notes :** Narcotic analgesic (tritium-labeled).**References :** Robinson, SE *CNS Drug Rev* 2002, 8, 377–90.**Catalog number :** 9064-005**CASRN :** 101224-22-0**Name :** Buprenorphine-3-β-D-glucuronide**Mol. formula :** C₃₅H₄₉NO₁₀**FW :** 654.59**DEA schedule :** 5**Catalog number :** 9058-003**CASRN :** 14357-78-9**Name :** Diprenorphine**Mol. formula :** C₂₆H₃₅NO₄**FW :** 425.54**DEA schedule :** 2**Notes :** (see footnotes 1 & 2 in Section C)**References :** Merck Index, 14th ed., Monograph 3340.**Catalog number :** 9058-001**CASRN :** 14357-78-9**Name :** Diprenorphine hydrochloride**Mol. formula :** C₂₆H₃₆ClNO₄**FW :** 462.04**DEA schedule :** 2**Notes :** (see footnotes 1 & 2 in Section C)**References :** Merck Index, 14th ed., Monograph 3340.

6 – Opioids and Opioid-related

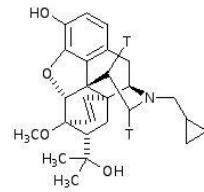
Catalog number : 9058-002

Name : [15,16- $^3\text{H}_2$]Diprenorphine

Mol. formula : $\text{C}_{26}\text{H}_{36}\text{ClNO}_4$

FW : 429.58 **DEA schedule :** 2

Notes : (see footnotes 1 & 2 in Section C)



Catalog number : 9059-001

CASRN : 14521-96-1

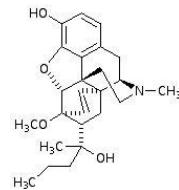
Name : Etorphine hydrochloride

Mol. formula : $\text{C}_{25}\text{H}_{34}\text{ClNO}_4$

FW : 448.01 **DEA schedule :** 2

Notes : (see footnotes 1 & 2 in Section C)

References : Merck Index, 14th ed., Monograph 3888.



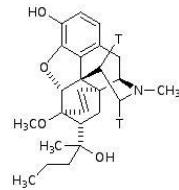
Catalog number : 9056-001

Name : [15,16- $^3\text{H}_2$]Etorphine

Mol. formula : $\text{C}_{25}\text{H}_{33}\text{NO}_4$

FW : 429.58 **DEA schedule :** 1

Notes : (see footnotes 1 & 2 in Section C)



Catalog number : 9333-013

CASRN : 78715-23-8

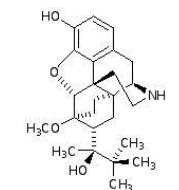
Name : Norbuprenorphine base

Mol. formula : $\text{C}_{25}\text{H}_{35}\text{NO}_4$

FW : 413.55 **DEA schedule :** 2

Notes : Buprenorphine metabolite.

References : Robinson, SE *CNS Drug Rev* 2002, 8, 377–90.



Opioids: Oxymorphone Class

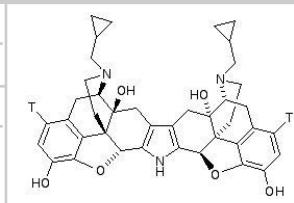
Catalog number : NOCD-084

Name : [1,1'- $^3\text{H}(n)$]Norbinaltorphimine; [^3H]norBNI

Mol. formula : $\text{C}_{40}\text{H}_{43}\text{N}_3\text{O}_6$

FW : 665.80 **DEA schedule :** 0

References : Birch, PJ; et al. *Eur J Pharmacol* 1987, 144, 405–8.
Portoghese, PS; Lipkowski, AW; Takemori, AE *Life Sci* 1987, 40, 1287–92.

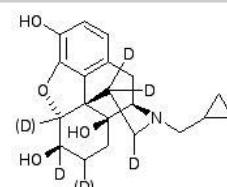


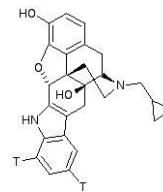
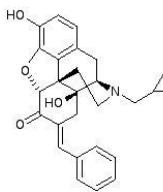
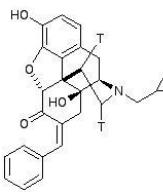
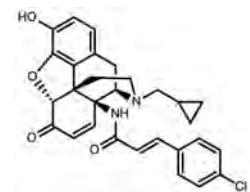
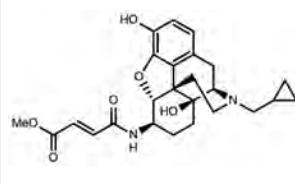
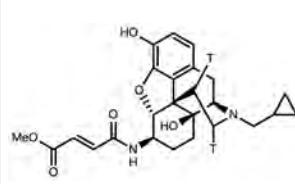
Catalog number : 9652-057

Name : [5,6,7,15,15,16- $^2\text{H}_6$]-6 α -Naltrexol

Mol. formula : $\text{C}_{20}\text{H}_{25}\text{NO}_4$

FW : 349.45 **DEA schedule :** 0



Catalog number : 9652-061**Name :** [5',7'-³H]Naltrindole**Mol. formula :** C₂₆H₂₆N₂O₃**FW :** 418.51 **DEA schedule :** 0**Catalog number :** 9652-067**CASRN :** 173556-52-0**Name :** 7-Benzylidene-7-dehydronaltrexone (BNTX) hydrochloride**Mol. formula :** C₂₇H₂₈CINO₄**FW :** 465.97 **DEA schedule :** 0**Catalog number :** 9652-068**Name :** [15,16-³H]-7-Benzylidine-7-dehydronaltrexone; [³H]BNTX**Mol. formula :** C₂₇H₂₈CINO₄**FW :** 433.52 **DEA schedule :** 0**Catalog number :** NOCD-100**CASRN :** 117332-69-1**Name :** Clocinnamox mesylate**Mol. formula :** C₃₀H₃₁CIN₂O₇S**FW :** 599.09**Notes :** Irreversible μ -opioid receptor antagonist.**References :** Comer, SD; et al. *J Pharmacol Exp Ther* 1992, 262, 1051-6.**Catalog number :** NOCD-079**Name :** Tritium-labeled Clocinnamox**Mol. formula :** C₂₉H₂₉CIN₂O₄**FW :** 507.01 **DEA schedule :** 0**References :** Comer, SD; Burke, TF; Lewis, JW; Woods, JH *J Pharmacol Exp Ther* 1992, 262, 1051-6.**Catalog number :** NOCD-141**CASRN :** 72786-10-8**Name :** β -Funaltrexamine hydrochloride; β -FNA hydrochloride**Mol. formula :** C₂₅H₃₁CIN₂O₆**FW :** 490.99 **DEA schedule :** 2**Notes :** Irreversible binding antagonist for the μ -opioid receptor.**References :** Portoghese, PS; el Kouhen, R; Law, PY; Loh, HH; Le Bourdonnec, B *Farmaco* 2001, 56, 191-6.**Catalog number :** 9333-012**Name :** [15,16-³H]-6 β -Funaltrexamine**Mol. formula :** C₂₆H₃₃CIN₂O₆**FW :** 458.53 **DEA schedule :** 2**Notes :** Irreversible μ -opioid receptor antagonist (tritium-labeled).**References :** Portoghese, PS; el Kouhen, R; Law, PY; Loh, HH; Le Bourdonnec, B *Farmaco* 2001, 56, 191-6.

6 – Opioids and Opioid-related

Catalog number : 9652-062

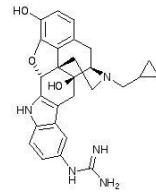
Name : 5'-GNTI trihydrochloride

Mol. formula : C₂₇H₃₂Cl₃N₅O₃

FW : 580.94 **DEA schedule :** 0

Notes : Selective κ-opioid receptor antagonist.

References : Jones, RM; Portoghesi, PS *Eur J Pharmacol* 2000, 396, 49-52.



Catalog number : 9652-063

CASRN : 219655-56-8

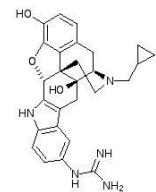
Name : 5'-GNTI dihydrochloride

Mol. formula : C₂₇H₃₁Cl₂N₅O₃

FW : 544.48 **DEA schedule :** 0

Notes : Selective κ-opioid receptor antagonist.

References : Jones, RM; Portoghesi, PS *Eur J Pharmacol* 2000, 396, 49-52.



Catalog number : 9652-069

CASRN : 350693-06-0

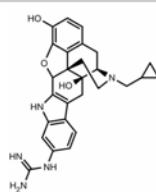
Name : 6'-GNTI dihydrochloride

Mol. formula : C₂₇H₂₉N₅O₃ • 2HCl

FW : 544.48 **DEA schedule :** 0

Notes : Kappa-opioid receptor agonist that inhibits arrestin recruitment.

References : Rives, ML; et al., *J Biol Chem* 2012, 287, 27050-4.



Catalog number : 9652-070

CASRN : 1314879-44-1

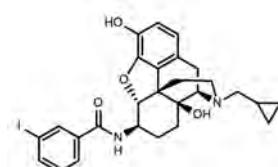
Name : IBNtxA

Mol. formula : C₂₇H₂₉IN₂O₄ • HCl

FW : 608.90 **DEA schedule :** 0

Notes : Kappa opioid receptor agonist.

References : Majumdar, S; et al., *J Med Chem* 2012, 55, 6352-62.



Catalog number : 9668-001

CASRN : 152658-17-8

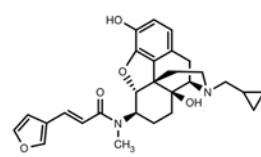
Name : Nalfurafine

Mol. formula : C₂₈H₃₂N₂O₅ • HCl

FW : 513.03 **DEA schedule :** 2

Notes : Kappa opioid receptor agonist.

References : Kawai, K., et al., *Bioorg Med Chem* (2008) 16, 9188.

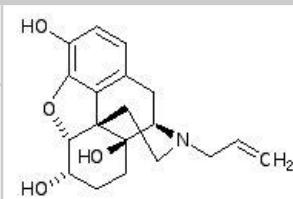


Catalog number : 9652-040

Name : 6α-Naloxol hydrochloride

Mol. formula : C₁₉H₂₄CINO₄

FW : 365.86 **DEA schedule :** 0

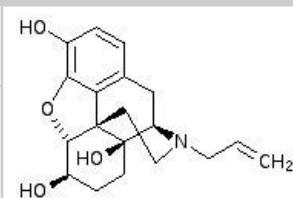


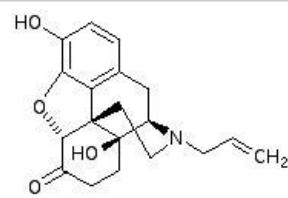
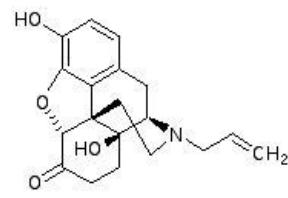
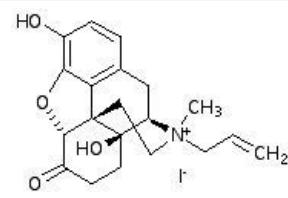
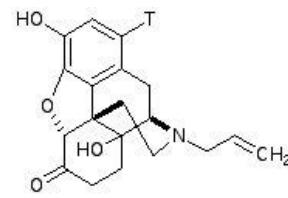
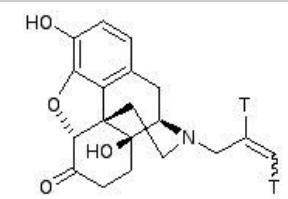
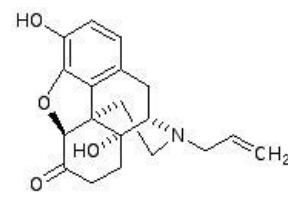
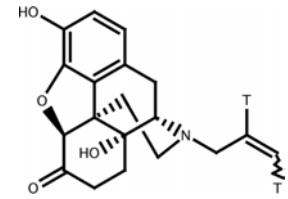
Catalog number : 9652-041

Name : 6β-Naloxol hydrochloride

Mol. formula : C₁₉H₂₄CINO₄

FW : 365.86 **DEA schedule :** 0



Catalog number : 9652-030**CASRN :** 51481-60-8**Name :** Naloxone hydrochloride**Mol. formula :** C₁₉H₂₂ClNO₄**FW :** 363.84 **DEA schedule :** 0**Notes :** Narcotic antagonist**References :** Merck Index, 14th ed., Monograph 6362.**Catalog number :** 9652-032**Name :** Naloxone pamoate**Mol. formula :** C₆₁H₅₈N₂O₁₄**FW :** 1043.14 **DEA schedule :** 0**Catalog number :** 9652-037**Name :** Naloxone methiodide**Mol. formula :** C₂₀H₂₄INO₄**FW :** 469.32 **DEA schedule :** 0**Catalog number :** 9652-031**Name :** (–)-[1-³H(n)]Naloxone**Mol. formula :** C₁₉H₂₁NO₄**FW :** 329.38 **DEA schedule :** 0**Catalog number :** 9652-035**Name :** (–)-[19,20-³H₂]Naloxone**Mol. formula :** C₁₉H₂₁NO₄**FW :** 331.39 **DEA schedule :** 0**Catalog number :** 9652-038**CASRN :** 357-08-4**Name :** (+)-Naloxone hydrochloride**Mol. formula :** C₁₉H₂₂ClNO₄**FW :** 363.85 **DEA schedule :** 0**Notes :** Unnatural isomer of naloxone. Selective Toll-like receptor 4 antagonist.**References :** Hutchinson, MR; et al., *Brain Behav Immun* 2010, 24, 83–95.; Watkins, LR; et al., *Trends Pharmacol Sci* 2009, 30, 581–91.**Catalog number :** 9652-039**Name :** (+)-[19,20-³H₂]Naloxone**Mol. formula :** C₁₉H₂₁NO₄**FW :** 331.39 **DEA schedule :** 0

6 – Opioids and Opioid-related

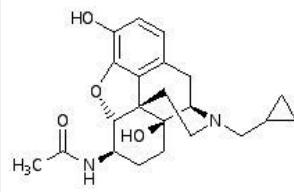
Catalog number : 9333-017

CASRN : 360770-17-8

Name : 6 β -Naltrexamide

Mol. formula : C₂₂H₂₈N₂O₄

FW : 388.975 DEA schedule : 2

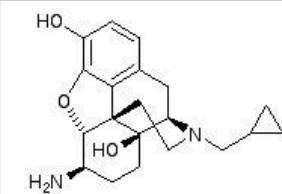


Catalog number : 9333-010

Name : 6 β -Naltrexamine dihydrochloride

Mol. formula : C₂₀H₂₈Cl₂N₂O₃

FW : 415.36 DEA schedule : 2



Catalog number : 9333-009

Name : [15,16-³H]-6 β -Naltrexamine

Mol. formula : C₂₀H₂₆N₂O₃

FW : 346.45 DEA schedule : 2



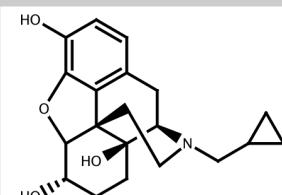
Catalog number : 9652-050

Name : 6 α -Naltrexol hydrochloride

Mol. formula : C₂₀H₂₆CINO₄

FW : 379.89 DEA schedule : 0

Notes : Naltrexone metabolite.



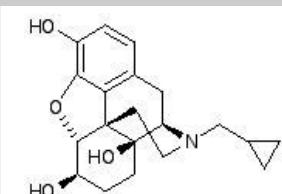
Catalog number : 9652-051

CASRN : 49625-89-0

Name : 6 β -Naltrexol hydrochloride

Mol. formula : C₂₀H₂₆CINO₄

FW : 393.41 DEA schedule : 0

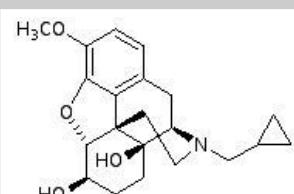


Catalog number : 9652-055

Name : 3-O-Methyl-6 β -naltrexol

Mol. formula : C₂₁H₂₇NO₄

FW : 357.45 DEA schedule : 0

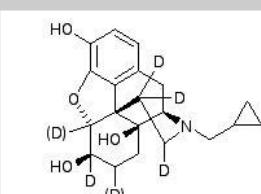


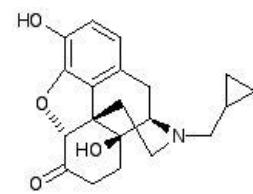
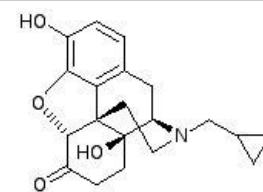
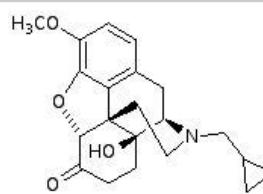
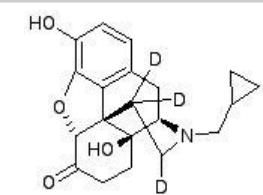
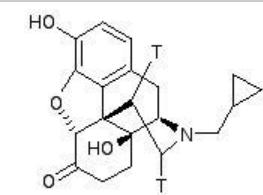
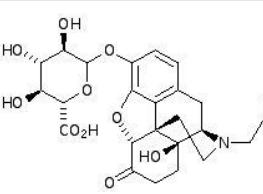
Catalog number : 9652-052

Name : [5,6,7,15,15,16-²H₆]-6 β -Naltrexol hydrochloride

Mol. formula : C₂₀H₂₆CINO₄

FW : 385.91 DEA schedule : 0



Catalog number : 9652-019**CASRN :** 16676-29-2**Name :** Naltrexone hydrochloride**Mol. formula :** C₂₀H₂₄ClNO₄**FW :** 377.88 **DEA schedule :** 0**Notes :** Narcotic antagonist**References :** Merck Index, 14th ed., Monograph 6363.**Catalog number :** 9652-020**CASRN :** 16590-41-3**Name :** Naltrexone base**Mol. formula :** C₂₀H₂₃NO₄**FW :** 341.42 **DEA schedule :** 0**Notes :** Narcotic antagonist**References :** Merck Index, 14th ed., Monograph 6363.**Catalog number :** 9652-027**Name :** 3-O-MethylNaltrexone**Mol. formula :** C₂₁H₂₅NO₄**FW :** 355.44 **DEA schedule :** 0**Catalog number :** 9652-014**Name :** [15,15,16-²H₃]Naltrexone**Mol. formula :** C₂₀H₂₃NO₄**FW :** 344.42 **DEA schedule :** 0**Catalog number :** 9652-012**Name :** [15,16-³H₂]Naltrexone**Mol. formula :** C₂₀H₂₃NO₄**FW :** 345.42 **DEA schedule :** 0**Catalog number :** 9652-013**Name :** [15,16-³H₂]Naltrexone hydrochloride**Mol. formula :** C₂₀H₂₄ClNO₄**FW :** 381.88 **DEA schedule :** 0**Catalog number :** 9652-025**Name :** Naltrexone-3-β-D-glucuronide**Mol. formula :** C₂₆H₃₁NO₁₀**FW :** 517.53 **DEA schedule :** 0

6 – Opioids and Opioid-related

Catalog number : 9652-064

CASRN : 122517-78-6

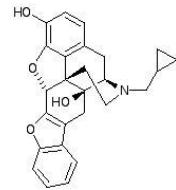
Name : Naltriben methanesulfonate; NTB methanesulfonate

Mol. formula : C₂₇H₂₉NO₇S

FW : 511.60 **DEA schedule :** 0

Notes : Selective δ-opioid receptor antagonist.

References : Sofuooglu, M; Portoghese, PS; Takemori, AE *J Pharmacol Exp Ther* **1991**, 257, 676–80.



Catalog number : 9652-065

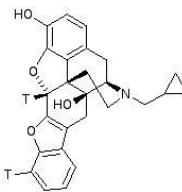
Name : Tritium-labeled Naltriben

Mol. formula : C₂₆H₂₅NO₄

FW : 419.50 **DEA schedule :** 0

Notes : Selective δ-opioid receptor antagonist (tritium-labeled).

References : Sofuooglu, M; Portoghese, PS; Takemori, AE *J Pharmacol Exp Ther* **1991**, 257, 676–80.



Catalog number : 9652-060

CASRN : 111469-81-9

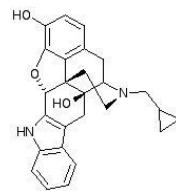
Name : Naltrindole hydrochloride; NTI hydrochloride

Mol. formula : C₂₆H₂₇ClN₂O₃

FW : 450.96 **DEA schedule :** 0

Notes : Selective δ-opioid receptor antagonist.

References : Portoghese, PS *Trends Pharmacol Sci* **1989**, 10, 230–5.



Catalog number : NOCD-067

CASRN : 113158-35-3

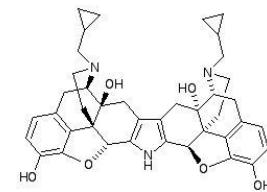
Name : Norbinaltorphimine dihydrochloride; norBNI

Mol. formula : C₄₀H₄₅Cl₂N₃O

FW : 734.73 **DEA schedule :** 0

Notes : Selective κ-opioid antagonist.

References : Birch, PJ; et al. *Eur J Pharmacol* **1987**, 144, 405–8.
Portoghese, PS; Lipkowski, AW; Takemori, AE *Life Sci* **1987**, 40, 1287–92.

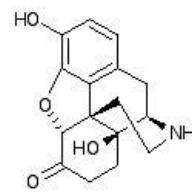


Catalog number : 9652-006

Name : (–)-Noroxyphorphone

Mol. formula : C₁₆H₁₇NO₄

FW : 287.30 **DEA schedule :** 2

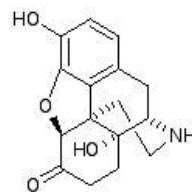


Catalog number : 9652-007

Name : (+)-Noroxyphorphone; (+)-14-Hydroxydihydronormorphinone

Mol. formula : C₁₆H₁₇NO₄

FW : 287.30 **DEA schedule :** 2



Catalog number : 9143-002

CASRN : 124-90-3

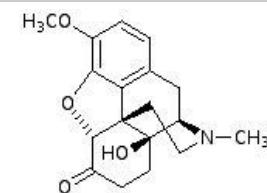
Name : Oxycodone hydrochloride; Dihydrohydroxycodeinone hydrochloride

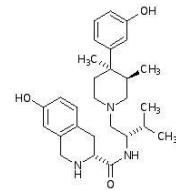
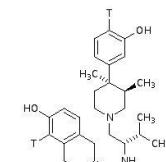
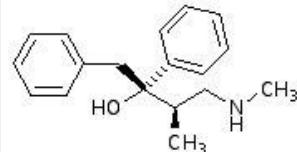
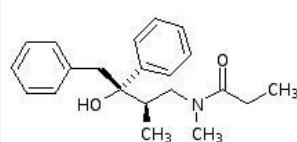
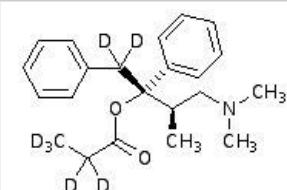
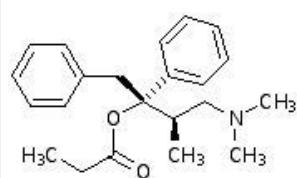
Mol. formula : C₁₈H₂₂ClNO₄

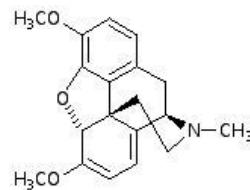
FW : 351.84 **DEA schedule :** 2

Notes : Narcotic analgesic

References : Merck Index, 14th ed., Monograph 6959.



Opioids: Phenylpiperidine Class**Catalog number :** NOCD-056**CASRN :** 785835-79-2**Name :** JDTic dihydrochloride**Mol. formula :** C₂₈H₄₁Cl₂N₃O₃**FW :** 538.56 **DEA schedule :** 0**Notes :** Selective kappa-opioid antagonist.**References :** Thomas, JB; Atkinson, RN; Rothman, RB; Fix, SE; Mascarella, SW; Vinson, NA; Xu, H; Dersch, CM; Lu, Y; Cantrell, BE; Zimmerman, DM; Carroll, FI *J Med Chem* 2001, 44, 2687–90.**Catalog number :** NOCD-059**Name :** [³H]JDTic**Mol. formula :** C₂₈H₃₉N₃O₃**FW :** 469.64 **DEA schedule :** 0**Notes :** Selective kappa-opioid antagonist (tritium-labeled).**References :** Thomas, JB; et al. *J Med Chem* 2001, 44, 2687–90.**Opioids: Propoxyphene Class****Catalog number :** 9273-011**Name :** (+)- α -Norpropoxyphene maleate**Mol. formula :** C₂₅H₃₁NO₆**FW :** 441.53 **DEA schedule :** 0**Catalog number :** 9273-010**Name :** (+)- α -N-Propionylnorpropoxyphene**Mol. formula :** C₂₁H₂₇NO₂**FW :** 325.45 **DEA schedule :** 0**Catalog number :** 9273-004**Name :** [²H₇]Propoxyphene hydrochloride**Mol. formula :** C₂₂H₂₉NO₂**FW :** 346.51 **DEA schedule :** 2**Catalog number :** 9273-002**Name :** (+)- α -Propoxyphene hydrochloride; Darvon hydrochloride**Mol. formula :** C₂₂H₃₀ClNO₂**FW :** 375.94 **DEA schedule :** 2**Notes :** Narcotic analgesic**References :** Miller, RR *Am J Hosp Pharm* 1977, 34, 413–23.
Barkin, RL; Barkin SJ; Barkin DS *Am J Ther* 2006, 13, 534–42.

Opioids: Thebaine Class**Catalog number :** 9333-002**CASRN :** 115-37-7**Name :** Thebaine base**Mol. formula :** C₁₉H₂₁NO₃**FW :** 311.36 **DEA schedule :** 2**Notes :** Opiate; synthetic starting material for many opiate compounds**References :** Merck Index, 14th ed., Monograph 9276.**Opioids (dosage form): Implantable****Catalog number :** 9250-014**Name :** Methadone HCl implant pellets**DEA schedule :** 2Dosage
Form**Catalog number :** 9300-008**Name :** Morphine base implant pellets**DEA schedule :** 2Dosage
Form**Catalog number :** 9300-009**Name :** Placebo for morphine base implant pellets**DEA schedule :** 0Dosage
Form**Catalog number :** 9652-021**Name :** Naltrexone base implant pellets**DEA schedule :** 0Dosage
Form**Catalog number :** 9652-022**Name :** Placebo for naltrexone implant pellets**DEA schedule :** 0Dosage
Form

(All peptides prepared using solution-phase methods.)

Peptides: Caged**Catalog number :** MPSP-118**Name :** CNB-Y-DYN8 (CYD8)**Sequence :** H-(Carboxynitrobenzyl)Tyr-Gly-Gly-Phe-Leu- Arg-Arg-Ile-OH**Mol. Formula :** C₅₄H₇₇N₁₅O₁₄**FW :** 1159.31**Note :** Photoactivatable analog of Dynorphin A (1–8).**Reference:** Banghart, MR; Sabatini BL, *Neuron* 2012, 73, 249–59.**Catalog number :** MPSP-117**Name :** CNB-Y-LE (CYLE)**Sequence :** H-(Carboxynitrobenzyl)Tyr-Gly-Gly-Phe-Leu-OH**Mol. Formula :** C₃₆H₄₂N₆O₁₁**FW :** 734.76**Note :** Photoactivatable analog of Leu-enkephalin.**Reference:** Banghart, MR; Sabatini BL, *Neuron* 2012, 73, 249–59.**Catalog number :** MPSP-119**Name :** CNV-Y-LE**Sequence :** H-(Carboxynitroveratryl)Tyr-Gly-Gly-Phe-Leu-OH**Mol. Formula :** C₃₈H₄₆N₆O₁₃**FW :** 794.70**Note :** Photoactivatable analog of Leu-enkephalin.**Reference:** Russell, AG; et al., *J Org Chem* 2010, 75, 4648–4651.
Russell, AG; et al., *Photochem Photobiol Sci* 2012, 11, 556–563.**Peptides: Cannabinoid-related****Catalog number :** MPSP-090**Name :** Hemopressin**Sequence :** H-Pro-Val-Asn-Phe-Lys-Phe-Leu-Ser-His-OH**Mol. Formula :** C₅₃H₇₇N₁₃O₁₂**FW :** 1088.3**Note :** CB₁ cannabinoid receptor inverse agonist. Analgesic.**Reference :** Heimann, AS; et al. *Proc Natl Acad Sci USA* 2007, 104, 20588–93.

7 – Peptides

Catalog number : PEPT-053

Name : Hemopressin *tris*(trifluoroacetate)

Sequence : H-Pro-Val-Asp-Phe-Lys-Phe-Leu-Ser-His

Mol. Formula : C₅₉H₈₁F₉N₁₃O₁₈

FW : 1431.36

Note : *CB₁* cannabinoid receptor inverse agonist. Analgesic.

Reference : Heimann, AS; et al. *Proc Natl Acad Sci USA* 2007, 104, 20588–93.

Catalog number : MPSP-096

Name : [³H₂]Hemopressin

Sequence : H-Pro-Val-Asn-[³H]Phe-Lys-[³H]Phe-Leu-Ser-His-OH

Mol. Formula : C₅₃H₇₇N₁₃O₁₂

FW : 1092.3

Note : Radiolabeled hemopressin (MPSP-90; PEPT-053).

Catalog number : MPSP-091

Name : [³H₄]Hemopressin

Sequence : H-[³H₂]Pro-Val-Asn-[³H]Phe-Lys-[³H]Phe-Leu-Ser-His-OH

Mol. Formula : C₅₃H₇₇N₁₃O₁₂

FW : 1096.3

Note : Radiolabeled hemopressin (MPSP-90; PEPT-053).

Catalog number : MPSP-098

Name : ABz-Hemopressin

Sequence : 2-ABz-Pro-Val-Asn-Phe-Lys-Phe-Leu-Ser-His-OH

Mol. Formula : C₆₀H₈₂N₁₄O₁₃

FW : 1207.1

Note : Hemopressin analog (MPSP-90; PEPT-053).

Catalog number : MPSP-099

Name : K⁵-ABz-Hemopressin

Sequence : H-Pro-Val-Asn-Phe-Lys(2-ABz)-Phe-Leu-Ser-His-OH

Mol. Formula : C₆₀H₈₂N₁₄O₁₃

FW : 1207.1

Note : Hemopressin analog (MPSP-90; PEPT-053).

Catalog number : MPSP-100

Name : Asp³-Hemopressin

Sequence : H-Pro-Val-Asp-Phe-Lys-Phe-Leu-Ser-His-OH

Mol. Formula : C₅₃H₇₆N₁₂O₁₃

FW : 1089.3

Note : Hemopressin analog (MPSP-90; PEPT-053).

Catalog number : MPSP-094**Name :** DIT⁷-Hemopressin (1-6)**Sequence :** H-Pro-Val-Asn-Phe-Lys-Phe-Tyr(3,5-diido)-OH**Mol. Formula :** C₄₇H₆₁N₉O₁₀I₂**FW :** 1165.9**Note :** *Hemopressin analog (MPSP-90; PEPT-053).***Catalog number :** MPSP-101**Name :** D-His⁹-Hemopressin**Sequence :** H-Pro-Val-Asn-Phe-Lys-Phe-Leu-Ser-D-His-OH**Mol. Formula :** C₅₃H₇₇N₁₃O₁₂**FW :** 1088.3**Note :** *Hemopressin analog (MPSP-90; PEPT-053).***Catalog number :** PEPT-054**Name :** Leu⁶-Hemopressin**Sequence :** H-Pro-Val-Asn-Phe-Lys-Leu-Leu-Ser-His-OH**Mol. Formula :** C₅₀H₇₉N₁₃O₁₂**FW :** 1054.26**Note :** *CB₁ cannabinoid receptor antagonist.***Reference :** Gomes et al., The FASEB Journal, 23, 1 (2009)**Catalog number :** MPSP-097**Name :** [³H]Leu⁶-Hemopressin**Sequence :** H-Pro-Val-Asn-[³H]Phe-Lys-Leu-Leu-Ser-His-OH**Mol. Formula :** C₅₀H₇₉N₁₃O₁₂**FW :** 1056.3**Note :** *Radiolabeled Leu⁶-Hemopressin (PEPT-054).***Catalog number :** MPSP-105**Name :** RVD-Hemopressin**Sequence :** H-Arg-Val-Asp-Pro-Val-Asn-Phe-Lys-Phe-Leu-Ser-His-OH**Mol. Formula :** C₆₈H₁₀₃N₁₉O₁₇**FW :** 1458.7**Note :** *Agonist of CB₁ cannabinoid receptors.***Reference :** Gomes et al., FASEB J. 2009; 23(9): 3020-9**Catalog number :** PEPT-057**Name :** RVD-Leu⁶-Hemopressin tetrakis(trifluoroacetate)**Sequence :** H-Arg-Val-Asp-Pro-Val-Asn-Phe-Lys-Leu-Leu-Ser-His-OH**Mol. Formula :** C₆₅H₁₀₅N₁₉O₁₇ • 4 CF₃CO₂H**FW :** 1880.75**Note :** *N-terminal extended hemopressin, CB1 receptor agonist.***Reference :** Dale, C. S.; Pagano Rde, L.; Rioli, V.; Hyslop, S.; Giorgi, R.; Ferro, E. S., *Peptides*, 26, 431 (2005). Gomes, I.; Grushko, J. S.; Golebiewska, U.; Hoogendoorn, S.; Gupta, A.; Heimann, A. S.; Ferro, E. S.; Scarlata, S.; Fricker, L. D.; Devi, L. A., *FASEB J.*, 23, 3020 (2009).

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Catalog number : MPSP-107

Name : RVD–Hemopressin (1–7)

Sequence : H–Arg–Val–Asp–Pro–Val–Asn–Phe–Lys–Phe–Leu–OH

Mol. Formula : C₅₉H₉₁N₁₅O₁₄

FW : 1234.5

Note : *RVD–Hemopressin analog (MPSP-105; PEPT-057).*

Catalog number : MPSP-092

Name : Tyr⁷–Hemopressin (1–6)

Sequence : H–Pro–Val–Asn–Phe–Lys–Phe–Tyr–OH

Mol. Formula : C₄₇H₆₃N₉O₁₀

FW : 914.1

Note : *Analog of MPSP-090.*

Catalog number : MPSP-093

Name : [³H₂]Tyr⁷–Hemopressin (1–6)

Sequence : H–Pro–Val–Asn–Phe–Lys–Phe–[³H₂]Tyr–OH

Mol. Formula : C₄₇H₆₁N₉O₁₀₃H₂

FW : 918.1

Note : *Radioligand for MPSP-092.*

Catalog number : MPSP-103

Name : VD–Hemopressin

Sequence : H–Val–Asp–Pro–Val–Asn–Phe–Lys–Phe–Leu–Ser–His–OH

Mol. Formula : C₆₂H₉₁N₁₅O₁₆

FW : 1302.5

Note : *Agonist of CB₁ cannabinoid receptors.*

Reference : Gomes et al., FASEB J. 2009; 23(9): 3020–9.

Catalog number : PEPT-055

Name : VD–Leu⁶–Hemopressin

Sequence : H–Val–Asp–Pro–Val–Asn–Phe–Lys–Leu–Leu–Ser–His–OH

Mol. Formula : C₅₉H₉₃N₁₅O₁₆

FW : 1268.1

Note : *CB₁ cannabinoid receptor antagonist.*

Reference : Gelman et al., J. Neurochem., 113, 871 (2010)

Peptides: Nicotinic Class**Catalog number :** PEPT-058**Name :** α -Conotoxin MII *tris*(trifluoroacetate)**Sequence :** H-Gly-Cys-Cys-Ser-Asn-Pro-Val-Cys-His-Leu-Glu-His-Ser-Asn-Leu-Cys-NH₂ (disulfide bonds Cys2-Cys8, Cys3-Cys16)**Mol. Formula :** C₆₇H₁₀₃N₂₃O₂₂S₄ • 3 CF₃CO₂H **FW :** 2053.02**Note :** Highly potent and selective $\alpha 3\beta 2$ nicotinic receptor antagonist.**Reference :** Cartier, G.; Yoshikami, D.; Gray, W.; Luo, S.; Olivera, B.; McIntosh, J., *Journal of Biological Chemistry*, 271, 7522 (1966). Hill, J. M.; Oomen, C. J.; Miranda, L. P.; Bingham, J. P.; Alewood, P. F.; Craik, D. J., *Biochemistry*, 37, 15621 (1998).**Catalog number :** MPSP-109**Name :** ω -Conotoxin GVIA**Sequence :** H-Cys-Lys-Ser-Hyp-Gly-Ser-Ser-Cys-Ser-Hyp-Thr-Ser-Tyr-Asn-Cys-Cys-Arg-Ser-Cys-Asn-Hyp-Tyr-Thr-Lys-Arg-Cys-Tyr-NH₂ (disulfide bonds Cys1-Cys16, Cys8-Cys19, Cys15-Cys26)**Mol. Formula :** C₁₂₀H₁₈₂N₃₈O₄₃S₆ **FW :** 3037.4**Note :** ω -Conotoxin GVIA, extracted from the venom of the fish-hunting cone snail *Conus geographus* is a neuronal calcium channel blocker.**Reference :** B.M. Olivera et al., *Biochemistry* 26, 2086 (1987); D.R. Hillyard et al., *Neuron* 9, 69 (1992); K. Sato et al., *Biochem. Biophys. Res. Commun.* 194, 1292 (1993); J.R. Abbott et al., *Int. J. Devl Neurosci.* 12, 43 (1994)**Catalog number :** MPSP-111**Name :** ω -Conotoxin MVIIA**Sequence :** H-Cys-Lys-Gly-Lys-Gly-Ala-Lys-Cys-Ser-Arg-Leu-Met-Tyr-Asp-Cys-Cys-Thr-Gly-Ser-Cys-Arg-Ser-Gly-Lys-Cys-NH₂ (disulfide bonds Cys1-Cys16, Cys8-Cys20, Cys15-Cys25)**Mol. Formula :** C₁₀₂H₁₇₂N₃₆O₃₂S₇ **FW :** 2639.2**Note :** ω -Conotoxin MVIIA, extracted from the venom of the fish-hunting cone snail *Conus magus* is a neuronal calcium channel blocker.**Reference :** B.M. Olivera et al., *Biochemistry* 26, 2086 (1987); D.R. Hillyard et al., *Neuron* 9, 69 (1992); J.R. Abbott et al., *Int. J. Devl Neurosci.* 12, 43 (1994); R. Newcomb et al., *Brain Res.* 638, 95 (1994)**Catalog number :** MPSP-113**Name :** ω -Conotoxin MVIIC**Sequence :** H-Cys-Lys-Gly-Lys-Gly-Ala-Pro-Cys-Arg-Lys-Thr-Met-Tyr-Asp-Cys-Cys-Ser-Gly-Ser-Cys-Gly-Arg-Arg-Gly-Lys-Cys-NH₂ (disulfide bonds Cys1-Cys16, Cys8-Cys20, Cys15-Cys26)**Mol. Formula :** C₁₀₆H₁₇₈N₄₀O₃₂S₇ **FW :** 2749.3**Note :** ω -Conotoxin MVIIC, extracted from the venom of the fish-hunting cone snail *Conus magus* is a neuronal calcium channel blocker.**Reference :** D.R. Hillyard et al., *Neuron* 9, 69 (1992); R. Newcomb et al., *Brain Res.* 638, 95 (1994)

Peptides: Opioid**Catalog number :** PEPT-042**Name :** Biphalin trifluoroacetate/acetate**Sequence :** H-Tyr-D-Ala-Gly-Phe-NH-Tyr-D-Ala-Gly-Phe-NH₂**Mol. Formula :** C₅₀H₅₈•₃F₅•₇N₁₀O₁₄**FW :** 1120.26**Note :** Synthetic dimeric enkephalin having potent agonist activity at μ and δ receptors.**Reference :** Lipkowski, AW; et al. *Life Sci* 1987, 40, 2283-8.
Li, G; et al. *Bioorg Med Chem Lett* 1998, 8, 555-60.**Catalog number :** PEPT-015**Name :** N-(tert-Butyloxycarbonyl)tyrosyl-D-alanylglycyl-N α -methylphenylalanylglycine methyl ester**Sequence :** Boc-Tyr-D-Ala-N α -Me-Phe-Gly-OCH₃**Mol. Formula :** C₃₂H₄₃N₅O₉**FW :** 641.72**Catalog number :** PEPT-022**Name :** N-(tert-Butyloxycarbonyl)tyrosyl-O-tert-butyl-D-threonylglycylphenylalanylleucine hydrazide**Sequence :** Boc-Tyr(Bu^t)-D-Thr-Gly-Phe-Leu-NH-NH₂**Mol. Formula :** C₃₉H₅₉N₇O₉**FW :** 769.95**Catalog number :** PEPT-060**Name :** CJ-15,208**Sequence :** [c(Phe¹-D-Pro²-Phe³-Trp⁴)]**Mol. Formula :** C₃₄H₃₅N₅O₄ • CF₃CO₂H**FW :** 691.70**Note :** Kappa-opioid receptor antagonist. Peptide content 63%.**Reference:** Ross, NC; et al., *Tetrahedron Lett.* 2010, 51, 5020-5023.**Catalog number :** PEPT-061**Name :** D-Trp⁴-CJ-15,208**Sequence :** [c(Phe¹-D-Pro²-Phe³-D-Trp⁴)]**Mol. Formula :** C₃₄H₃₅N₅O₄ • CF₃CO₂H**FW :** 691.70**Note :** Kappa-opioid receptor antagonist. Peptide content 54%.**Reference:** Ross, NC; et al., *Tetrahedron Lett.* 2010, 51, 5020-5023.

Catalog number : MPSP-013**Name :** CTAP**Sequence :** H-D-Phe-c[Cys-Tyr-D-Trp-Arg-Thr-Pen]-Thr-NH₂**Mol. Formula :** C₅₁H₆₉N₁₃O₁₁S₂**FW :** 1103.0**Note :** *Cyclic somatostatin analog with high μ selectivity.***Reference :** Pelton, JT; et al. *J Med Chem* 1986, 29, 2370–5.
Bilsky, EJ; et al. *J Pharmacol Exp Ther* 1996, 277, 484–90.**Catalog number :** MPSP-014**Name :** [³H]CTAP**Sequence :** H-D-Phe-[³H]-c[Cys-Tyr-D-Trp-Arg-Thr-Pen]-Thr-NH₂**Mol. Formula :** C₅₁H₆₈N₁₃O₁₁S₂₃H**FW :** 1105.0**Note :** *Radioactive ligand for MPSP-013***Reference :** Abbruscato, TJ; et al. *J Pharmacol Exp Ther* 1997, 280, 402–9.**Catalog number :** MPSP-054**Name :** CTOP**Sequence :** H-D-Phe-c[Cys-Tyr-D-Trp-Orn-Thr-Pen]-Thr-NH₂**Mol. Formula :** C₅₀H₆₇N₁₁O₁₁S₂**FW :** 1062.4**Note :** *Cyclic somatin analog with high μ selectivity.***Reference :** Gulya, K; et al. *Life Sci* 1986, 38, 2221–9.
Pelton, JT; et al. *J Med Chem* 1986, 29, 2370–5.**Catalog number :** MPSP-006**Name :** DADLE**Sequence :** H-Tyr-D-Ala-Gly-Phe-D-Leu-OH**Mol. Formula :** C₂₉H₃₉N₅O₇**FW :** 569.7**Note :** *Synthetic enkephalin agonist having improved biological stability and some selectivity for the δ receptor over the μ receptor. It shows roughly the same anti-nociceptive activity as β-endorphin.***Reference :** Chang, KJ; Cuatrecasas P. *J Biol Chem* 1979, 254, 2610–8.
Gorin, FA; et al. *J Med Chem* 1980, 23, 1113–22.
Knapp, RJ; Yamamura HI. *Biochem Pharmacol* 1992, 44, 1687–95.**Catalog number :** PEPT-001**Name :** DADLE hydrochloride; [D-Ala²,D-Leu⁵]Enkephalin**Sequence :** H-Tyr-D-Ala-Gly-Phe-D-Leu-OH**Mol. Formula :** C₂₉H₄₀CIN₅O₇**FW :** 606.12**Note :** *Synthetic enkephalin agonist having improved biological stability and some selectivity for the δ receptor over the μ receptor.***Reference :** Chang, KJ; Cuatrecasas, PJ. *J Biol Chem* 1979, 254, 2610.
Gorin, FA; et al. *J Med Chem* 1980, 23, 1113.
Knapp, RJ; Yamamura, HI. *Biochem Pharmacol* 1992, 44, 1687.

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Catalog number : MPSP-007

Name : [$^3\text{H}_2$]DADLE

Sequence : H-Tyr[$^3\text{H}_2$]-D-Ala-Gly-Phe-D-Leu-OH

Mol. Formula : $\text{C}_{29}\text{H}_{37}\text{N}_5\text{O}_{7.5}\text{H}_2$

FW : 573.7

Note : Tritiated DADLE

Reference : Luciano, MG; et al. *Brain Res Bull* 1981, 7, 677-82.

Catalog number : PEPT-008

Name : DALA hydrochloride

Sequence : H-Tyr-D-Ala-Gly-Phe-Met-NH₂

Mol. Formula : $\text{C}_{28}\text{H}_{39}\text{ClN}_6\text{O}_6\text{S}$

FW : 623.17

Note : Potent, long-lasting analog of [Met5]enkephalin.

Reference : Pert, CB; et al. *Science* 1976, 194, 330.

Catalog number : MPSP-034

Name : DALCE

Sequence : H-Tyr-D-Ala-Gly-Phe-Leu-Cys-OH

Mol. Formula : $\text{C}_{32}\text{H}_{43}\text{N}_6\text{O}_8\text{S}$

FW : 673.0

Note : Irreversible antagonist for the $\delta 1$ receptor.

Reference : Bowen, WD; et al. *J Biol Chem* 1987, 262, 13434-9.
Traynor, JR; Elliott, J *TIPS* 1993, 14, 84.

Catalog number : MPSP-035

Name : [$^3\text{H}_2$]DALCE

Sequence : H-Tyr[$^3\text{H}_2$]-D-Ala-Gly-Phe-Leu-Cys-OH

Mol. Formula : $\text{C}_{33}\text{H}_{44}\text{N}_6\text{O}_{10}\text{S}_3\text{H}_2$

FW : 677.0

Note : Radioactive ligand for MPSP-034

Catalog number : PEPT-035

Name : DALCE trifluoroacetate

Sequence : Tyr-D-Ala-Gly-Phe-Leu-Cys-OH

Mol. Formula : $\text{C}_{34}\text{H}_{45}\text{F}_3\text{N}_6\text{O}_{10}\text{S}$

FW : 786.82

Note : Irreversible antagonist for the $\delta 1$ receptor.

Reference : Bowen, WD; et al. *J Biol Chem* 1987, 262, 13434-9.
Traynor, JR; Elliott, J *TIPS* 1993, 14, 84.

Catalog number : MPSP-027**Name :** DALDA**Sequence :** H-Tyr-D-Arg-Phe-Lys-NH₂**Mol. Formula :** C₃₀H₄₅N₉O₅**FW :** 611.8**Note :** *Synthetic tetrapeptide agonist highly selective for the μ receptor.***Reference :** Schiller, PW; et al. *J Med Chem* 1989, 32, 698–703.**Catalog number :** MPSP-084**Name :** (DMT)¹-DALDA**Sequence :** H-DMT-D-Arg-Phe-Lys-NH₂**Mol. Formula :** C₃₂H₄₉N₉O₅**FW :** 639.6**Note :** *Super DALDA***Reference :** Zhao, K; et al. *J Biol Chem* 2004, 279, 34682–90.**Catalog number :** PEPT-038**Name :** DALDA *tris(trifluoroacetate)***Sequence :** H-Tyr-D-Arg-Phe-Lys-NH₂**Mol. Formula :** C₃₆H₄₈F₉N₉O₁₁**FW :** 611.75**Note :** *Synthetic tetrapeptide agonist highly selective for the μ receptor.***Reference :** Schiller, PW; et al. *J Med Chem* 1989, 32, 698.**Catalog number :** PEPT-047**Name :** Super DALDA; SS-02; [Dmt1]DALDA *tris(trifluoroacetate)***Sequence :** Dmt-D-Arg-Phe-Lys-NH₂**Mol. Formula :** C₃₈H₅₂F₉N₉O₁₁**FW :** 981.354**Reference :** Zhao, K; et al. *J Biol Chem* 2004, 279, 34682–90.
Schiller, PW; et al. *Eur J Med Chem* 2000, 35, 895–901.**Catalog number :** MPSP-028**Name :** [³H₂]DALDA**Sequence :** H-Tyr[³H₂]-D-Arg-Phe-Lys-NH₂**Mol. Formula :** C₃₀H₄₃N₉O₅₃H₂**FW :** 615.8**Note :** *Radioactive ligand for MPSP-027*

Catalog number : MPSP-011**Name :** DAMGO**Sequence :** H-Tyr-D-Ala-Gly-N^α-Me-Phe-Gly-ol**Mol. Formula :** C₂₆H₃₅N₅O₆**FW :** 513.7**Note :** Synthetic enkephalin agonist highly selective for the μ receptor.**Reference :** Handa, BK; et al. *Eur J Pharmacol* 1981, 70, 531-40.
Kosterlitz, HW; Paterson SJ; Robson LE *Br J Pharmacol* 1981, 73, 939-49.
Reddy, PA; et al. *Org Prep Proc Internat* 1995, 27, 469.**Catalog number :** PEPT-014**Name :** DAMGO trifluoroacetate**Sequence :** Tyr-D-Ala-Gly-N^α-Me-Phe-Gly-ol**Mol. Formula :** C₂₈H₃₆F₃N₅O₈**FW :** 627.62**Note :** Synthetic enkephalin agonist highly selective for the μ receptor.**Reference :** Handa, BK; et al. *Eur J Pharmacol* 1981, 70, 531.
Kosterlitz, HW; Paterson, SJ *Br J Pharmacol* 1981, 73, 299P.
Reddy, PA; et al. *Org Prep Proc Internat* 1995, 27, 469.**Catalog number :** MPSP-012**Name :** [³H₂]DAMGO**Sequence :** H-Tyr[³H₂]-D-Ala-Gly-N^α-Me-Phe-Gly-ol**Mol. Formula :** C₂₆H₃₃N₅O₆H₂**FW :** 517.7**Note :** Radioactive ligand for MPSP-011**Reference :** Zajac, JM; Roques BP *Life Sci* 1983, 33 Suppl 1, 155-8.**Catalog number :** MPSP-030**Name :** Deltorphin I**Sequence :** H-Tyr-D-Ala-Phe-Asp-Val-Val-Gly-NH₂**Mol. Formula :** C₃₇H₅₂N₈O₁₀**FW :** 768.9**Note :** Selective δ receptor agonist isolated from the skin of *Phyllomedusa bicolor***Reference :** Amodeo, P; et al. *Pept Res* 1992, 5, 48-55.
Melchiorri, P; et al. *Eur J Pharmacol* 1991, 195, 201-7.
Salvadori, S; et al. *J Med Chem* 1991, 34, 1656-61.
Lazarus, LH; et al. *J Med Chem* 1991, 34, 1350-5.**Catalog number :** PEPT-033**Name :** Deltorphin I hemitrifluoroacetate**Sequence :** Tyr-D-Ala-Phe-Asp-Val-Val-Gly-NH₂**Mol. Formula :** C₃₇H₅₂N₈O₁₀**FW :** 768.87**Note :** Selective δ receptor agonist isolated from the skin of *Phyllomedusa bicolor*.**Reference :** Ersperger, V; et al. *Proc Natl Acad Sci USA* 1989, 86, 5188.
Salvadori, S; et al. *J Med Chem* 1991, 34, 1656.
DLHeyl and SESchullery, *CurrMedChem.*, 4, 117 (1997).

Catalog number : MPSP-031**Name :** [³H₂]Deltorphin I**Sequence :** H-Tyr[³H₂]-D-Ala-Phe-Asp-Val-Val-Gly-NH₂**Mol. Formula :** C₃₇H₅₀N₈O₁₀H₂**FW :** 772.9**Note :** Radioactive ligand for MPSP-030**Reference :** Erspamer, V; et al. *Proc Natl Acad Sci USA* 1989, 86, 5188.**Catalog number :** MPSP-036**Name :** Deltorphin II**Sequence :** H-Tyr-D-Ala-Phe-Glu-Val-Val-Gly-NH₂**Mol. Formula :** C₃₈H₅₄N₈O₁₀**FW :** 782.9**Note :** A selective δ-opioid receptor agonist isolated from the skin of *Phyllomedusa bicolor***Reference :** Amodeo, P; et al. *Pept Res* 1992, 5, 48-55.
Melchiorri, P; et al. *Eur J Pharmacol* 1991, 195, 201-7.
Salvadori, S; et al. *J Med Chem* 1991, 34, 1656-61.
Lazarus, LH; et al. *J Med Chem* 1991, 34, 1350-5.**Catalog number :** PEPT-034**Name :** Deltorphin II trifluoroacetate**Sequence :** Tyr-D-Ala-Phe-Glu-Val-Val-Gly-NH₂**Mol. Formula :** C₄₀H₅₅F₃N₈O₁₂**FW :** 896.92**Note :** Selective δ2 receptor agonist isolated from the skin of *Phyllomedusa bicolor*.**Reference :** Erspamer, V; et al. *Proc Natl Acad Sci USA* 1989, 86, 5188.
Salvadori, S; et al. *J Med Chem* 1991, 34, 1656.
Traynor, JR; Elliott, J *TIPS* 1993, 14, 84.**Catalog number :** MPSP-037**Name :** [³H₂]Deltorphin II**Sequence :** H-Tyr[³H₂]-D-Ala-Phe-Glu-Val-Val-Gly-NH₂**Mol. Formula :** C₃₈H₅₂N₈O₁₀H₂**FW :** 786.9**Note :** Radioactive ligand for MPSP-036**Reference :** Erspamer, V; et al. *Proc Natl Acad Sci USA* 1989, 86, 5188.**Catalog number :** PEPT-037**Name :** Dermorphin trifluoroacetate**Sequence :** Tyr-D-Ala-Phe-Gly-Tyr-Pro-Ser-NH₂**Mol. Formula :** C₄₂H₅₁F₃N₈O₁₂**FW :** 916.91**Note :** Selective μ receptor agonist isolated from the skin of *Phyllomedusa sauvagei*.**Reference :** Montecucchi, PC; et al. *Int J Peptide Protein Res* 1981, 17, 275.
Melchiorri, P; Negri, L *Gen Pharmac* 1996, 27, 1099.

7 – Peptides

Catalog number : PEPT-049

Name : 2',6'-Dimethyltyrosyl-Ng-nitro-d-arginylphenylalanyl-Ne-2-Cl-Cbz-lysinamide trifluoroacetate

Sequence : Dmt-d-Arg(NO₂)-Phe-Lys(2-Cl-Cbz)-NH₂

Mol. Formula : C₄₂H₅₅ClF₃N₁₀O₁₁

FW : 949.376

Reference : Zhao, K; et al. *J Biol Chem* 2004, 279, 34682-90.

Catalog number : PEPT-048

Name : 2',6'-Dimethyltyrosyl-Ng-nitro-d-arginylphenylalanyllysinamide trifluoroacetate

Sequence : Dmt-d-Arg(NO₂)-Phe-Lys-NH₂

Mol. Formula : C₃₆H₅₀F₆N₁₀O₁₁

FW : 944.2

Reference : Zhao, K; et al. *J Biol Chem* 2004, 279, 34682-90.

Catalog number : PEPT-046

Name : DIPP-NH₂[Ψ] bis(Trifluoroacetate)

Sequence : 2TFA •Dmt-TicΨ[CH₂NH]Phe-Phe-NH₂

Mol. Formula : C₄₃H₄₇F₆N₅O₈

FW : 875.85

Note : Synthetic high potency pseudopeptide having balanced μ agonist/δ antagonist properties.

Reference : Schiller, PW; et al. *J Med Chem* 1999, 42, 3520-6.

Catalog number : MPSP-001

Name : DPDPE

Sequence : H-Tyr-c[D-Pen-Gly-Phe-d-Pen]-OH

Mol. Formula : C₃₀H₃₉N₇O₅S₂

FW : 645.8

Note : Synthetic, conformationally restricted enkephalin agonist selective for the δ1 receptor. Minimal cross-reactivity with μ- and κ- receptors.

Reference : Mosberg, HI; et al. *Proc Natl Acad Sci USA* 1983, 80, 5871-4.

Akiyama, K; et al. *Proc Natl Acad Sci USA* 1985, 82, 2543-7.

Mosberg, HI; Omnaas JR; Goldstein A *Mol Pharmacol* 1987, 31, 599-602.

Knapp, RJ; Yamamura HI *Biochem Pharmacol* 1992, 44, 1687-95.

Catalog number : MPSP-044

Name : p-Cl-Phe-DPDPE

Sequence : H-Tyr-c[D-Pen-Gly-4-Cl-Phe-d-Pen]-OH

Mol. Formula : C₃₀H₃₈ClN₅O₇S₂

FW : 680.4

Note : Analog of DPDPE. δ-selectivity 5 times higher than DPDPE due to a 5-fold increased δ-receptor affinity.

Reference : Toth, G; et al. *J Med Chem* 1990, 33, 249-53.
Vaughn, LK; et al. *Life Sci* 1989, 45, 1001-8.

Catalog number : PEPT-010**Name :** DPDPE trifluoroacetate; [D-Pen²,D-Pen⁵]Enkephalin**Sequence :** H-Tyr-c[D-Pen-Gly-Phe-D-Pen]-OH**Mol. Formula :** C₃₂H₄₀F₃N₅O₉S₂**FW :** 759.82**Note :** Synthetic, conformationally restricted enkephalin agonist selective for the δ1 receptor.**Reference :** Mosberg, HI; et al. *Proc Natl Acad Sci USA* 1983, 80, 5871–4.
Knapp, RJ; Yamamura, HI *Biochem Pharmacol* 1992, 44, 1687.
Traynor, JR; Elliott, J *TIPS* 1993, 14, 84.**Catalog number :** MPSP-002**Name :** [³H₂]DPDPE**Sequence :** H-Tyr[³H₂]-c[D-Pen-Gly-Phe-D-Pen]-OH**Mol. Formula :** C₃₀H₃₇N₇O₅S₂₃H₂**FW :** 649.8**Note :** Radioactive ligand for MPSP-001**Reference :** Cotton, R; et al. *Eur J Pharmacol* 1984, 97, 331–2.**Catalog number :** MPSP-045**Name :** [³H₂]p-Cl-Phe-DPDPE**Sequence :** H-Tyr[³H₂]-c[D-Pen-Gly-4-Cl-Phe-D-Pen]-OH**Mol. Formula :** C₃₀H₃₆ClN₅O₇S₂₃H₂**FW :** 684.4**Note :** Radioactive ligand for MPSP-044**Catalog number :** MPSP-086**Name :** DV1.2.DA1.5.Lan.Enk.(DPDPE lanthionine analog)**Sequence :** H-Tyr-[D-S(Val L-Gly-Phe-D-Ala) L]-OH**Mol. Formula :** C₂₈H₃₅N₅O₇S**FW :** 584.7**Note :** DPDPE lanthionine analog.**Catalog number :** MPSP-003**Name :** DSLET**Sequence :** H-Tyr-D-Ser-Gly-Phe-Leu-Thr-OH**Mol. Formula :** C₃₃H₄₆N₆O₁₀**FW :** 686.9**Note :** Synthetic enkephalin agonist selective for the δ2 receptor.**Reference :** Fournie-Zaluski, MC; et al. *Mol Pharmacol* 1981, 20, 484–91.
Traynor, JR; Elliott J *Trends Pharmacol Sci* 1993, 14, 84–6.

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Catalog number : PEPT-039

Name : DSLET trifluoroacetate

Sequence : Tyr-D-Ser-Gly-Phe-Leu-Thr-OH

Mol. Formula : C₃₅H₄₇F₃N₆O₁₂

FW : 800.79

Note : Synthetic enkephalin agonist selective for the δ2 receptor.

Reference : Gacel, G; et al. *FEBS Lett* 1980, 118, 245.
Gacel, G; et al. *J Med Chem* 1981, 24, 1119.
Traynor, JR; Elliott, J *TIPS* 1993, 14, 84.

Catalog number : MPSP-004

Name : [³H₂]DSLET

Sequence : H-Tyr[³H₂]-D-Ser-Gly-Phe-Leu-Thr-OH

Mol. Formula : C₃₃H₄₄N₆O₁₀H₂

FW : 690.9

Note : Radioactive ligand for MPSP-003

Reference : David, M; et al. *Eur J Pharmacol* 1982, 78, 385-7.

Catalog number : MPSP-032

Name : DSTBULET

Sequence : H-Tyr-D-Ser(Bu^t)-Gly-Phe-Leu-Thr-OH

Mol. Formula : C₃₇H₄₄N₆O₁₀

FW : 742.9

Note : Synthetic enkephalin agonist selective for the δ receptor

Reference : Delay-Goyet, P; et al. *J Biol Chem* 1988, 263, 4124-30.
Gacel, G; et al. *J Med Chem* 1988, 31, 1891-7.

Catalog number : PEPT-043

Name : DSTBULET acetate

Sequence : H-Tyr-D-Ser(Bu^t)-Gly-Phe-Leu-Thr-OH

Mol. Formula : C₃₉H₅₈N₆O₁₂

FW : 802.92

Note : Synthetic enkephalin agonist selective for the δ receptor.

Reference : Delay-Goyet, P; et al. *J Biol Chem* 1988, 263, 4124-30.
Gacel, G; et al. *J Med Chem* 1988, 31, 1891-7.

Catalog number : MPSP-033

Name : [³H₂]DSTBULET

Sequence : H-Tyr[³H₂]-D-Ser(Bu^t)-Gly-Phe-Leu-Thr-OH

Mol. Formula : C₃₃H₄₄N₆O₁₀H₂

FW : 746.9

Note : Radioactive ligand for MPSP-032

Reference : Delay-Goyet, P *NIDA Res Mono* 1986, 75, 197-200.

Catalog number : MPSP-005**Name :** DTLET**Sequence :** H-Tyr-D-Ser-Gly-Phe-Leu-Thr-OH**Mol. Formula :** C₃₄H₄₈N₆O₁₀**FW :** 700.9**Note :** Synthetic enkephalin agonist moderately selective for the δ receptor.**Reference :** Zajac, JM; et al. *Biochem Biophys Res Commun* 1983, 111, 390–7.
Delay-Goyet, P; et al. *FEBS Lett* 1985, 183, 439–43.**Catalog number :** PEPT-021**Name :** DTLET trifluoroacetate**Sequence :** Tyr-D-Thr-Gly-Phe-Leu-Thr-OH**Mol. Formula :** C₃₆H₄₉F₃N₆O₁₂**FW :** 814.81**Reference :** Zajac, JM; et al. *Biochem Biophys Res Commun* 1983, 111, 390.
Delay-Goyet, P; et al. *FEBS Lett* 1985, 183, 439.**Catalog number :** MPSP-088**Name :** [(2S)-Mdp]¹-Dyn NH₂ (1–11); Dynantin**Sequence :** [(2S)-Mdp]¹-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-NH₂**Mol. Formula :** C₆₆H₁₀₉N₂₁O₁₂**FW :** 1388.4**Note :** Highly and potent selective kappa antagonist.**Reference :** Lu, Y et al. *J Med Chem* 2001, 44, 3048–53.**Catalog number :** PEPT-056**Name :** Dynantin**Sequence :** [2S]-Mdp-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-NH₂**Mol. Formula :** C₆₆H₁₀₉N₂₁O₁₂ • 4 CF₃CO₂H**FW :** 1844.41**Note :** κ-Opioid antagonist.**Reference :** (1) Lu, Y.; et al., *J Med Chem*, 2001, 44, 3048. (2) Schiller, P. W.; et al., *Life Sci*, 2003, 73, 691.**Catalog number :** MPSP-019**Name :** Dynorphin (1–11)**Sequence :** H-Tyr-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-OH**Mol. Formula :** C₆₃H₁₀₃N₂₁O₁₃**FW :** 1362.7**Note :** Truncated analog of dynorphin A with similar potency

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Catalog number : MPSP-020

Name : [³H₂]Dynorphin (1-11)

Sequence : H-Tyr[³H₂]-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-OH

Mol. Formula : C₆₃H₁₀₁N₂₁O₁₃₃H₂

FW : 1366.7

Note : *Radioactive ligand for MPSP-019*

Reference : Chavkin, C; Goldstein, A *Proc Natl Acad Sci USA* **1981**, *78*, 6543-7.

Catalog number : MPSP-016

Name : Dynorphin (1-13)

Sequence : H-Tyr-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-Leu-Lys-OH

Mol. Formula : C₇₅H₁₂₆N₂₄O₁₅

FW : 1604.2

Note : *Truncated analog of dynorphin A with similar potency*

Reference : Goldstein, A; et al. *Proc Natl Acad Sci USA* **1979**, *78*, 6666-70.

Catalog number : MPSP-017

Name : Dynorphin (1-13) amide

Sequence : H-Tyr-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-Leu-Lys-NH₂

Mol. Formula : C₇₅H₁₂₇N₂₅O₁₄

FW : 1603.2

Note : *Truncated amide analog of dynorphin A with similar potency*

Reference : Chavkin, C; Goldstein, A *Proc Natl Acad Sci USA* **1981**, *78*, 6543-7.

Catalog number : MPSP-018

Name : [³H₂]Dynorphin (1-13) amide

Sequence : H-Tyr[³H₂]-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-Leu-Lys-NH₂

Mol. Formula : C₇₅H₁₂₅N₂₅O₁₄₃H₂

FW : 1607.2

Note : *Radioactive ligand for MPSP-017*

Catalog number : MPSP-021

Name : Dynorphin (1-9)

Sequence : H-Tyr-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-OH

Mol. Formula : C₅₂H₈₄N₁₈O₁₁

FW : 1137.5

Note : *Selective ligand for κ-binding sites.*

Reference : Corbett, AD; et al. *Nature* **1982**, *299*, 79-81.
Yoshimura, K; et al. *J Pharmacol Exp Ther* **1982**, *222*, 71-9.

Catalog number : MPSP-022**Name :** [³H₂]Dynorphin (1-9)**Sequence :** H-Tyr[³H₂]-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-OH**Mol. Formula :** C₅₂H₈₂N₁₈O₁₁₃H₂**FW :** 1141.5**Note :** Radioactive ligand for MPSP-021**Reference :** Robson, LE; et al. *Life Sci* 1983, 33 Suppl 1, 283-6.
Wood, MS; Rodriguez FD; Traynor JR *Neuropharmacology* 1989, 28, 1041-6.**Catalog number :** MPSP-048**Name :** Dynorphin (2-11)**Sequence :** H-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-OH**Mol. Formula :** C₃₃H₄₄N₆O₁₀**FW :** 1199.7**Note :** Dynorphin truncation which does not bind to opioid receptors**Reference :** Takemori, AE; Loh HH; Lee NM *J Pharmacol Exp Ther* 1993, 266, 121-4.
Meyer, ME *Pharmacol Biochem Behav* 1993, 44, 329-32.**Catalog number :** MPSP-046**Name :** Dynorphin (2-13)**Sequence :** H-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-Leu-Lys-OH**Mol. Formula :** C₆₆H₁₁₇N₂₃O₁₃**FW :** 1441.1**Note :** Dynorphin truncation that does not bind to opioid receptors. Highly immunoreactive as the first tyrosine appears not to be essential for immunoreactivity.**Reference :** Young, EA; et al. *Peptides* 1987, 8, 701-7.
Walker, JM; et al. *Eur J Pharmacol* 1982, 85, 121-2.**Catalog number :** MPSP-040**Name :** Dynorphin (2-17)**Sequence :** H-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-Leu-Lys-Trp-Asp-Asn-Gln-OH**Mol. Formula :** C₉₀H₁₅₆N₃₀O₂₁**FW :** 1984.8**Note :** Dynorphin truncation that does not bind to opioid receptors.**Reference :** Chavkin, C; Goldstein, A *Proc Natl Acad Sci USA* 1981, 78, 6543-6547.**Catalog number :** MPSP-041**Name :** [³H]Dynorphin (2-17)**Sequence :** H-Gly-Gly-Phe[³H]-Leu-Arg-Arg-Ile-Arg-Pro-Lys-Leu-Lys-Trp-Asp-Asn-Gln-OH**Mol. Formula :** C₉₀H₁₅₅N₃₀O₂₁₃H**FW :** 1986.8**Note :** Radioactive ligand for MPSP-040

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Catalog number : MPSP-050

Name : Dynorphin A (1-8)

Sequence : H-Tyr-Gly-Gly-Phe-Leu-Arg-Arg-Ile-OH

Mol. Formula : C₄₆H₇₂N₁₄O₁₀

FW : 981.4

Note : *Endogenous peptide agonist for opioid receptors.*

Reference : Nakao, K; et al. *Biochem Biophys Res Commun* 1983, 117, 695-701.
Seizinger, BR; et al. *J Neurochem* 1984, 42, 447-57.
Bell, KM; Traynor JR *Can J Physiol Pharmacol* 1998, 76, 325-33.

Catalog number : MPSP-066

Name : [³Ψ⁴,D-Leu⁸]Dynorphin A (1-8)amide

Sequence : H-Tyr-Gly-Gly-Ψ(CH₂-NH)-Phe-Leu-Arg-Arg-D-Leu-NH₂

Mol. Formula : C₃₃H₄₄N₆O₁₀

FW : 967.4

Note : *Kappa-selective dynorphin A analog*

Reference : Ambo, A; et al. *Chem Pharm Bull* 1995, 43, 1547-50.

Catalog number : MPSP-052

Name : Dynorphin A (2-8)

Sequence : H-Gly-Gly-Phe-Leu-Arg-Arg-Ile-OH

Mol. Formula : C₃₇H₆₃N₁₃O₈

FW : 818.2

Note : *Truncated analog of dynorphin A which does not bind to opioid receptors*

Reference : Takemori, AE; Loh HH; Lee NM *J Pharmacol Exp Ther* 1993, 266, 121-4.
Meyer, ME *Pharmacol Biochem Behav* 1993, 44, 329-32.

Catalog number : PEPT-040

Name : Dynorphin A 1-8 tris(trifluoroacetate)

Sequence : H-Tyr-Gly-Gly-Phe-Leu-Arg-Arg-Ile-OH

Mol. Formula : C₅₂H₇₅F₉N₁₄O₁₆

FW : 1323.24

Note : *Endogenous peptide agonist for opioid receptors.*

Reference : Weber, E; et al. *Nature* 1982, 299, 77.
Corbett, A; et al. *Nature* 1982, 299, 79.
Bell, KM; Traynor, JR *Can J Physiol Pharmacol* 1998, 76, 325.

Catalog number : PEPT-041

Name : Dynorphin A 2-8 tris(trifluoroacetate)

Sequence : H-Gly-Gly-Phe-Leu-Arg-Arg-Ile-OH

Mol. Formula : C₄₃H₆₆F₉N₁₃O₁₄

FW : 1159.69

Note : *Des-Tyr1 analog of dynorphin A (1-8).*

Reference : Takemori, AE; Loh HH; Lee NM *J Pharmacol Exp Ther* 1993, 266, 121-4.
Meyer, ME *Pharmacol Biochem Behav* 1993, 44, 329-32.

Catalog number : MPSP-023**Name :** Dynorphin B**Sequence :** H-Tyr-Gly-Gly-Phe-Leu-Arg-Arg-Gln-Phe-Lys-Val-Val-Thr-OH**Mol. Formula :** C₇₄H₁₁₅N₂₁O₁₇**FW :** 1571.1**Note :** *N-terminal part of leumorphin***Reference :** Kilpatrick, DL; et al. *Proc Natl Acad Sci USA* 1982, 79, 6480–83.
Seizinger, BR; et al. *J Neurochem* 1984, 42, 447–457.**Catalog number :** MPSP-015**Name :** Dynorphin(1-17)**Sequence :** H-Tyr-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-Leu-Lys-Trp-Asp-Asn-Gln-OH**Mol. Formula :** C₉₉H₁₅₅N₃₁O₂₃**FW :** 2147.8**Note :** *Considered the endogenous ligand for kappa receptors.***Reference :** Goldstein, A; et al. *Proc Natl Acad Sci USA* 1981, 78, 7219–23.
Cox, BM; et al. *Life Sci* 1975, 16, 1777–1782.
Goldstein, A; et al. *Proc Natl Acad Sci USA* 1979, 76, 6666–70.**Catalog number :** MPSP-070**Name :** Endomorphin-1**Sequence :** H-Tyr-Pro-Trp-Phe-NH₂**Mol. Formula :** C₃₄H₃₈N₆O₅**FW :** 610.7**Note :** *Endogenous tetrapeptide agonist selective for the μ receptor.***Reference :** Zadina, JE; et al. *Nature* 1997, 386, 499.
Goldberg, IE; et al. *J Pharmacol Exp Ther* 1998, 286, 1007.
McConalogue, K; et al. *Neuroscience* 1999, 90, 1051.**Catalog number :** PEPT-044**Name :** Endomorphin-1 Trifluoroacetate/Acetate**Sequence :** TFA/HOAc •Tyr-Pro-Trp-Phe-NH₂**Mol. Formula :** C₃₆H₃₉₋₁₅F₂₋₈₅N₆O₇**FW :** 722.04**Note :** *Endogenous tetrapeptide agonist selective for the μ receptor.***Reference :** JEZadina et al., *Nature*, 386, 499 (1997)
IEGoldberg et al., *JPharmacolExpTher.*, 286, 1007 (1998)
KMcConalogue et al., *Neuroscience*, 90, 1051 (1999)**Catalog number :** MPSP-073**Name :** Endomorphin-2**Sequence :** H-Tyr-Pro-Phe-Phe-NH₂**Mol. Formula :** C₃₂H₃₇N₅O₅**FW :** 571.7**Note :** *Endogenous tetrapeptide agonist selective for the μ receptor.***Reference :** Zadina, JE; et al. *Nature* 1997, 386, 499–502.
Goldberg, IE; et al. *J Pharmacol Exp Ther* 1998, 286, 1007–13.
McConalogue, K; et al. *Neuroscience* 1999, 90, 1051–9.

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Catalog number : PEPT-045

Name : Endomorphin-2 Trifluoroacetate/Acetate

Sequence : TFA/HOAc •Tyr-Pro-Phe-Phe-NH₂

Mol. Formula : C₃₄H₃₈F₂N₅O₇

FW : 680.30

Note : *Endogenous tetrapeptide agonist selective for the μ receptor.*

Reference : Zadina, JE; et al. *Nature* 1997, 386, 499-502.
Goldberg, IE; et al. *J Pharmacol Exp Ther* 1998, 286, 1007-13.
McConalogue, K; et al. *Neuroscience* 1999, 90, 1051-9.

Catalog number : MPSP-074

Name : [³H₂]Endomorphin-2

Sequence : H-Tyr[³H₂]-Pro-Phe-Phe-NH₂

Mol. Formula : C₃₂H₃₅N₅O₅H₂

FW : 575.7

Note : *Radioactive ligand for MPSP-073*

Reference : Spetea, M; et al. *Biochem Biophys Res Commun* 1998, 250, 720-5.

Catalog number : MPSP-024

Name : β-Endorphin

Sequence : H-Tyr-Gly-Gly-Phe-Met-Thr-Ser-Glu-Lys-Ser-Gln-Thr-Pro-Leu-Val-Thr-Leu-Phe-Lys-Asn-Ala-Ile-IIe-Lys-Asn-Ala-Tyr-Lys-Lys-Gly-Glu-OH

Mol. Formula : C₁₅₈H₂₅₁N₃₉O₄₆S

FW : 3465.6

Note : *Endogenous opioid peptide. Contains enkephalin sequence.*

Reference : Li, CH; et al. *Biochem Biophys Res Commun* 1976, 72, 1542-1547.
Cox, BM; et al. *Proc Natl Acad Sci USA* 1976, 73, 1821-1823.
Bradbury, AF; et al. *Nature* 1976, 260, 165-166.

Catalog number : MPSP-026

Name : [¹²⁵I]β-Endorphin

Sequence : H-YGGFMTSEKSQTPLVTLFKNAlIKNAYKY[¹²⁵I₂]GE-OH

Mol. Formula : C₁₅₈H₂₄₉N₃₉O₄₆S₁₂₅I₂

FW : 3589.6

Note : *Radioactive ligand for MPSP-024*

Reference : Deby-Dupont, G; et al. *C R Seances Soc Biol Fil* 1983, 177, 259-68.
Schweigerer, L; et al. *J Biol Chem* 1983, 258, 12287-92.

Catalog number : MPSP-025

Name : [³H₂]β-Endorphin

Sequence : H-Tyr-Gly-Gly-Phe-Met-Thr-Ser-Glu-Lys-Ser-Gln-Thr-Pro-Leu-Val-Thr-Leu-Phe-Lys-Asn-Ala-Ile-IIe-Lys-Asn-Ala-[³H₂]Tyr-Lys-Lys-Gly-Glu-OH

Mol. Formula : C₁₅₈H₂₄₉N₃₉O₄₆S₃H₂

FW : 3469.6

Note : *Radioactive ligand for MPSP-024*

Reference : Li, CH; et al. *Proc Natl Acad Sci USA* 1980, 77, 2303-4.

Catalog number : MPSP-081**Name :** [$^3\text{H}_2$]Leu-Enkephalin-acid**Sequence :** H-[$^3\text{H}_2$]Tyr-Gly-Gly-Phe-Leu-OH**Mol. Formula :** C₂₈H₃₇N₅O₇**FW :** 559.3**Note :** Radioactive ligand for MPSP-080.**Catalog number :** PEPT-036**Name :** FMRF amide bis(trifluoroacetate)**Sequence :** H-Phe-Met-Arg-Phe-NH₂**Mol. Formula :** C₃₃H₄₄F₆N₈O₈S**FW :** 826.82**Note :** Molluscan cardioexcitatory neuropeptide.**Reference :** Price, DA; Greenberg, MJ *Science* **1997**, *197*, 670.
Tang, J; et al. *Proc Natl Acad Sci USA* **1984**, *81*, 5002.**Catalog number :** MPSP-008**Name :** ICI 174,864**Sequence :** N,N-Diallyl-Tyr-Aib-Aib-Phe-Leu-OH**Mol. Formula :** C₃₈H₅₃N₅O₇**FW :** 691.87**Note :** Synthetic enkephalin antagonist selective for the δ receptor.**Reference :** Cotton, R; et al. US Patent 4,474,767 (1984).
Cotton, R; et al. *Eur J Pharmacol* **1984**, *97*, 331-2.
Dray, A; Nunan L *Peptides* **1984**, *5*, 1015-6.
Smith, CB; Bennett-Kelly L; Woods JH *Neuropeptides* **1984**, *5*, 161-4.**Catalog number :** PEPT-017**Name :** ICI 174864 hydrochloride**Sequence :** N,N-Diallyl-Tyr-Aib-Aib-Phe-Leu-OH**Mol. Formula :** C₃₈H₅₄ClN₅O₇**FW :** 728.34**Note :** Synthetic enkephalin antagonist selective for the δ receptor.**Reference :** Cotton, R; et al. US Patent 4,474,767 (1984).
Cotton, R; et al. *Eur J Pharmacol* **1984**, *97*, 331.
Dray, D; Nunan, L *Peptides* **1984**, *5*, 1015.**Catalog number :** MPSP-078**Name :** Kaffiralin-1**Sequence :** H-D-Phe-D-Phe-D-Ile-D-Arg-NH₂**Mol. Formula :** C₃₀H₄₄N₈O₄**FW :** 580.7**Note :** Kappa-selective agonist identified from a combinatorial library**Reference :** Dooley, CT; et al. *J Biol Chem* **1998**, *273*, 18848-56.

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Catalog number : MPSP-079

Name : [³H₂]Kaffiralin-1

Sequence : H-[³H]D-Phe-[³H]D-Phe-D-Ile-D-Arg-NH₂

Mol. Formula : C₃₀H₄₂N₈O₄₃H₂

FW : 584.7

Note : Radioactive ligand for MPSP-076

Reference : Dooley, CT; et al. *J Biol Chem* 1998, 273, 18848-56.

Catalog number : MPSP-076

Name : Kaffiralin-2

Sequence : H-D-Phe-D-Phe-D-Nle-D-Arg-NH₂

Mol. Formula : C₃₀H₄₄N₈O₄

FW : 580.7

Note : Kappa-selective agonist identified from a combinatorial library.

Reference : Dooly, CT; et al. *J Bio Chem* 1998, 273, 18848-18856.

Catalog number : MPSP-077

Name : [³H₂]Kaffiralin-2

Sequence : H-[³H]D-Phe-[³H]D-Phe-D-Nle-D-Arg-NH₂

Mol. Formula : C₃₀H₄₂N₈O₄₃H₂

FW : 584.7

Note : Radioactive ligand for MPSP-076.

Catalog number : PEPT-003

Name : [Leu⁵]Enkephalin trifluoroacetate

Sequence : H-Tyr-Gly-Gly-Phe-Leu-OH

Mol. Formula : C₃₀H₃₈F₃N₅O₉

FW : 669.62

Note : Endogenous enkephalin.

Reference : Hughes, J; et al. *Nature* 1975, 258, 577.
Jones, D *Tetrahedron Lett* 1977, 2853.
Vilkas, E; et al. *Int J Peptide Protein Res* 1980, 15, 29.

Catalog number : MPSP-080

Name : Leu-Enkephalin-acid

Sequence : H-Tyr-Gly-Gly-Phe-Leu-OH

Mol. Formula : C₂₈H₃₇N₅O₇

FW : 555.3

Note : Endogenous enkephalin.

Reference : Hughes, J; et al. *Nature* 1975, 258, 577.
Jones, D *Tetrahedron Lett* 1977, 2853.
Vilkas, E; et al. *Int J Peptide Protein Res* 1980, 15, 29.

Catalog number : PEPT-007**Name :** [Met⁵]Enkephalinamide hydrochloride**Sequence :** H-Tyr-Gly-Gly-Phe-Met-NH₂**Mol. Formula :** C₂₇H₃₇ClN₆O₆S**FW :** 609.14**Note :** Amide analog of [Met5]enkephalin.**Reference :** Puig, MM; et al. *Arch Int Pharmacol Ther* 1979, 226, 69.
Berger, E; et al. *Pharmazie* 1979, 34, 349.**Catalog number :** MPSP-064**Name :** Met-Enkephalin-acid**Sequence :** H-Tyr-Gly-Gly-Phe-Met-OH**Mol. Formula :** C₂₈H₃₇N₅O₇**FW :** 573.7**Note :** Endogenous opioid peptide**Catalog number :** MPSP-029**Name :** Met-Enkephalin-amide**Sequence :** H-Tyr-Gly-Gly-Phe-Met-NH₂**Mol. Formula :** C₂₇H₃₆N₆O₆S**FW :** 572.8**Note :** Amide analog of [Met5]enkephalin, endogenous opioid peptide.**Reference :** Hughes, J; et al. *Nature* 1975, 258, 577–80.
Vavrek, RJ; et al. *Peptides* 1981, 2, 303–8.**Catalog number :** PEPT-020**Name :** Methyl N,N-Diallyl-O-tert-butyltyrosyl- α -aminoisobutyryl- α -aminoisobutyrate**Sequence :** N,N-Dallyl-Tyr(Bu^t)-Aib-Aib-OMe**Mol. Formula :** C₂₈H₄₃N₃O₅**FW :** 501.67**Catalog number :** MPSP-042**Name :** Metkephamid**Sequence :** H-Tyr-Ala-Gly-N ^{α} -Me-Phe-Met-NH₂**Mol. Formula :** C₂₉H₄₀N₆O₆S**FW :** 600.8**Note :** Stable, systemically active analog of [Met5]enkephalin.**Reference :** RCFredrickson, et al., *Science*, 211, 603–605(1981)**Catalog number :** PEPT-009**Name :** Metkephamide hydrochloride**Sequence :** Tyr-D-Ala-Gly-N ^{α} -Me-Phe-Met-NH₂**Mol. Formula :** C₂₉H₄₁ClN₆O₆S**FW :** 637.19**Note :** Stable, systemically active analog of [Met⁵]enkephalin.**Reference :** Frederickson, RCA; et al. *Science* 1981, 211, 603.

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Catalog number : PEPT-004

Name : N-(*tert*-Butyloxycarbonyl)tyrosylglycylglycylphenylalanylleucine methyl ester

Sequence : Boc-Tyr-Gly-Phe-Leu-OMe

Mol. Formula : C₃₄H₄₇N₅O₉

FW : 669.77

Note : [Leu⁵]enkephalin synthetic precursor.

Reference : Vilkas, E; et al. *Int J Peptide Protein Res* 1980, 15, 29.

Catalog number : PEPT-006

Name : N-(*tert*-Butyloxycarbonyl)tyrosylglycylglycylphenylalanylmetionine methyl ester

Sequence : Boc-Tyr-Gly-Gly-Phe-Met-OMe

Mol. Formula : C₃₃H₄₅N₅O₉S

FW : 687.81

Note : [Met⁵]enkephalin synthetic precursor.

Reference : Bhotre, BJ; et al. *J Indian Chem Soc* 1978, 55, 1128.

Catalog number : PEPT-018

Name : N,N-Diallyl-O-*tert*-butyltyrosyl- α -aminoisobutyryl- α -aminoisobutyrylphenylalanylleucine

Sequence : N,N-Diallyl-Tyr(Bu^t)-Aib-Aib-Phe-Leu-OH

Mol. Formula : C₄₂H₆₁N₅O₇

FW : 747.98

Reference : Cotton, R.; et al. *Eur J Pharmacol* 1984, 97, 331-2.

Catalog number : PEPT-019

Name : N,N-Diallyl-O-*tert*-butyltyrosyl- α -aminoisobutyryl- α -aminoisobutyrylphenylalanylleucine methyl ester

Sequence : N,N-Diallyl-Tyr(Bu^t)-Aib-Aib-Phe-Leu-OMe

Mol. Formula : C₄₃H₆₃N₅O₇

FW : 762.01

Note : ICI 174864 synthetic intermediate.

Reference : Flippin-Anderson, JL; et al. *Lett Pept Sci* 1994, 1, 107.

Catalog number : MPSP-059

Name : [³H₂]Orphanin FQ; [³H₂]Nociceptin

Sequence : H-Phe[³H]-Gly-Gly-Phe[³H]-Thr-Gly-Ala-Arg-Lys-Ser-Ala-Arg-Lys-Leu-Ala-Asn-Gln-OH

Mol. Formula : C₇₉H₁₂₇N₂₇O₂₂₃H₂

FW : 1813.4

Note : Radioactive ligand for MPSP-058

Reference : Dooley, CT; Houghten RA *Life Sci* 1996, 59, PL23-9.

Catalog number : MPSP-082**Name :** [³H]⁴Tyr¹⁴Orphanin FQ**Sequence :** H-Phe[³H]-Gly-Gly-Phe[³H]-Thr-Gly-Ala-Arg-Lys-Ser-Ala-Arg-Lys-Tyr[³H₂]-Ala-Asn-Gln-OH**Mol. Formula :** C₈₂H₁₂₃N₂₇O₂₃H₄**FW :** 1867.1**Note :** *Radioactive ligand for MPSP-058***Reference :** Dooley, CT; et al. *Life Science* 1996, 59, PL23–PL29.**Catalog number :** MPSP-095**Name :** [³H]⁴Orphanin FQ**Sequence :** H-[³H]Phe-Gly-Gly-[³H]Phe-Thr-Gly-Ala-Arg-Lys-Ser-Ala-Arg-Lys-[³H₂]Leu-Ala-Asn-Gln-OH**Mol. Formula :** C₇₉H₁₂₅N₂₇O₂₂H₄**FW :** 1817.4**Note :** *Radioactive ligand for MPSP-058.***Catalog number :** MPSP-083**Name :** (Tyr¹⁴)Orphanin FQ analog**Sequence :** H-Phe-Gly-Gly-Phe-Thr-Gly-Ala-Arg-Lys-Ser-Ala-Arg-Lys-Tyr-Ala-Asn-Gln-OH**Mol. Formula :** C₈₂H₁₂₇N₂₇O₂₃**FW :** 1859.1**Note :** *Control ligand for MPSP-082.***Catalog number :** MPSP-058**Name :** Orphanin FQ; Nociceptin**Sequence :** H-Phe-Gly-Gly-Phe-Thr-Gly-Ala-Arg-Lys-Ser-Ala-Arg-Lys-Leu-Ala-Asn-Gln-OH**Mol. Formula :** C₇₉H₁₂₉N₂₇O₂₂**FW :** 1809.4**Note :** *Endogenous ligand for orphan opioid like receptor (ORL)***Reference :** Reinscheid, RK; et al. *Science* 1995, 270, 792–4.
Meunier, JC; et al. *Nature* 1995, 377, 532–5.**Catalog number :** PEPT-023**Name :** PL017 trifluoroacetate**Sequence :** Tyr-Pro-N^α-Me-Phe-D-Pro-NH₂**Mol. Formula :** C₃₁H₃₈F₃N₅O₇**FW :** 649.67**Note :** *Synthetic morphiceptin agonist selective for the μ receptor.***Reference :** Chang, KJ; et al. *Science* 1981, 212, 75.

7 – Peptides

Catalog number : MPSP-009

Name : PLO17

Sequence : H-Tyr-Pro-N^α-Me-Phe-D-Pro-NH₂

Mol. Formula : C₂₉H₃₇N₅O₅

FW : 535.65

Note : Synthetic morphiceptin agonist selective for the μ receptor. Casomorphin analog.

Reference : Chang, KJ; et al. *J Pharmacol Exp Ther* 1983, 227, 403–8.

Catalog number : MPSP-010

Name : [³H₂]PLO17

Sequence : H-Tyr[³H₂]-Pro-N^α-Me-Phe-D-Pro-NH₂

Mol. Formula : C₂₉H₃₅N₅O₅H₂

FW : 539.6

Note : Radioactive ligand for MPSP-009.

Reference : Hawkins, KN; et al. *Eur J Pharmacol* 1987, 133, 351–2.

Catalog number : MPSP-060

Name : Dansyl-PQR-amide

Sequence : Dansyl-Pro-Gln-Arg-NH₂

Mol. Formula : C₂₈H₄₁N₉O₆S

FW : 631.5

Note : Fluorescent truncation analog of neuropeptide FF(NPFF) an anti-opiate peptide

Reference : Malin, DH; et al. *Life Sci* 1993, 53, PL261–6.

Catalog number : MPSP-062

Name : Benzyl-PQR-amide

Sequence : Benzyl-Pro-Gln-Arg-NH₂

Mol. Formula : C₂₃H₃₆N₈O₄

FW : 488.6

Note : Truncation analog of neuropeptide FF(NPFF) an anti-opiate peptide

Catalog number : PEPT-051

Name : SS-20; Phenylalanyl-D-arginylphenylalanyllysineamide tris(trifluoroacetate)

Sequence : Phe-D-Arg-Phe-Lys-NH₂

Mol. Formula : C₃₆H₄₈F₉N₈O₉

FW : 938.9

Reference : Zhao, K; et al. *J Biol Chem* 2004, 279, 34682-90.

Catalog number : PEPT-050**Name :** SS-31; D-Arginyl-(2',6'-dimethyl)tyrosyllysylphenylalaninamide *tris*(Trifluoroacetate)**Sequence :** D-Arg-Dmt-Lys-Phe-NH₂**Mol. Formula :** C₃₈H₅₂F₉N₉O₁₁**FW :** 981.354**Reference :** Zhao, K; et al. *J Biol Chem* 2004, 279, 34682-90.**Catalog number :** MPSP-068**Name :** TICP[Ψ]**Sequence :** H-Tyr-TicΨ[CH₂-NH]Cha-Phe-OH**Mol. Formula :** C₃₄H₃₈N₆O₅**FW :** 610.7**Note :** Very stable potent δ antagonist**Reference :** Schiller, PW; et al. *J Signal Transduction Research* 1999, 19, 573-88.
Szatmari, I; et al. *Peptides* 1999, 20, 1079-83.**Catalog number :** MPSP-056**Name :** TIPP[Ψ]**Sequence :** H-Tyr-TicΨ[CH₂-NH]Phe-Phe-OH**Mol. Formula :** C₃₇H₄₀N₄O₅**FW :** 620.8**Note :** Highly potent and stable δ receptor antagonist with extraordinary δ selectivity.**Reference :** Schiller, PW; et al. *J Med Chem* 1993, 36, 3182-7.
Visconti, LM; et al. *Neurosci Lett* 1994, 181, 47-9.**Catalog number :** PEPT-030**Name :** TIPP trifluoroacetate**Sequence :** H-Tyr-Tic-Phe-Phe-OH**Mol. Formula :** C₃₉H₃₉F₃N₄O₈**FW :** 748.76**Note :** Highly potent and selective δ receptor antagonist.**Reference :** Schiller, PW; et al. *Proc Natl Acad Sci USA* 1992, 89, 11871.**Catalog number :** PEPT-032**Name :** TIPP[Ψ] acetate/trifluoroacetate**Sequence :** H-Tyr-TicΨ[CH₂NH]Phe-Phe-OH**Mol. Formula :** C₄₁H₄₅F₃N₄O₉**FW :** 794.82**Note :** Highly potent and stable δ receptor antagonist with extraordinary δ selectivity.**Reference :** Schiller, PW; et al. *J Med Chem* 1993, 36, 3182.
Visconti, LM; et al. *Neurosci Lett* 1994, 181, 47.

7 – Peptides

Catalog number : PEPT-052

Name : TIPP[Ψ] bis(trifluoroacetate)

Sequence : Tyr-Tic-Ψ[CH₂NH]Phe-Phe-OH

Mol. Formula : C₄₁H₄₂F₆N₄O₉

FW : 848.80

Note : Selective delta opioid receptor antagonist.

Reference : Schiller, PW; et al. *J Med Chem* 1993, 36, 3182–7.

Catalog number : MPSP-038

Name : TIPP-Enkephalin (TIPP)

Sequence : H-Tyr-Tic-Phe-Phe-OH

Mol. Formula : C₃₇H₃₈N₄O₆

FW : 634.8

Note : Highly potent and selective δ-opioid receptor antagonist

Reference : Schiller, PW; et al., *Proc Natl Acad Sci USA* 1992, 89, 11871–11875.
Flippen-Anderson, JL; et al. *Lett Pept Sci* 1994, 1, 107.

Catalog number : MPSP-039

Name : [³H₂]Tipp-Enkephalin

Sequence : H-Tyr[³H₂]-Tic-Phe-Phe-OH

Mol. Formula : C₃₇H₃₆N₄O₆₃H₂

FW : 638.8

Note : Radioactive ligand for MPSP-038

Reference : Nevin, ST; et al. *Life Sci* 1993, 53, PL57–62.

Catalog number : PEPT-031

Name : D-TIPP-NH₂ trifluoroacetate/acetate

Sequence : Tyr-D-Tic-Phe-Phe-NH₂

Mol. Formula : C₃₉H₄₀F₂N₅O₇

FW : 731.58

Note : Potent and highly selective μ receptor agonist.

Reference : Schiller, PW; et al. *Proc Natl Acad Sci USA* 1992, 89, 11871.
Flippen-Anderson, JL; et al. *J Peptide Res* 1997, 49, 384.

Catalog number : MPSP-115

Name : Orexin A (human, bovine, rat, mouse), Hypocretin-1

Sequence : pGlu-Pro-Leu-Pro-Asp-Cys-Cys-Arg-Gln-Lys-Thr-Cys-Ser-Cys-Arg-Leu-Tyr-Glu-Leu-His-Gly-Ala-Gly-Asn-His-Ala-Ala-Gly-Ile-Leu-Thr-Leu-NH₂ (Disulfide bridges Cys6-Cys12 and Cys7-Cys14)

Mol. Formula : C₁₅₂H₂₄₃N₄₇O₄₄S

FW : 3561.2

Note : Hypothalamic neuropeptide that regulates feeding behavior.

Reference : M.W. Schwartz, *Nat. Med.* 4, 385 (1998); T. Sakurai et al., *Cell* 92, 573 (1998); M.R. Jain et al., *Regul. Peptides* 87, 19 (2000); E. Goncz et al., *Endocrinology* 149, 1618 (2008)

Catalog number : MPSP-116

Name : Orexin B (human), Hypocretin-2

Sequence : H-Arg-Ser-Gly-Pro-Pro-Gly-Leu-Gln-Gly-Arg-Leu-Gln-Arg-Leu-Leu-Gln-Ala-Ser-Gly-Asn-His-Ala-A
Ia-Gly-Ile-Leu-Thr-Met-NH₂

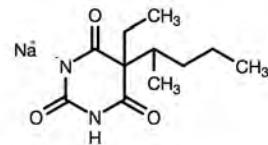
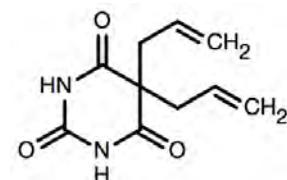
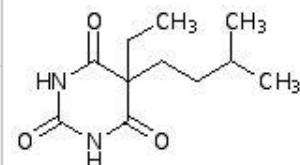
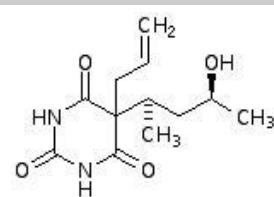
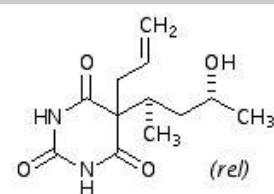
Mol. Formula : C₁₂₃H₂₁₂N₄₄O₃₅S

FW : 2899.4

Note : *Hypothalamic neuropeptide that regulates feeding behavior.*

Reference : M.W. Schwartz, Nat. Med. 4, 385 (1998); T. Sakurai et al., Cell 92, 573 (1998); M.R. Jain et al., Regul. Peptides 87, 19 (2000)



Sedatives: Barbiturate Class**Catalog number :** 2270-012**new****CASRN :** 57-33-0**Name :** Sodium pentobarbital**Mol. formula :** C₁₁H₁₇N₂NaO₃**FW :** 249.26**DEA schedule :** 2**References :** ChemSpider ID 5762.**Catalog number :** 2100-001**CASRN :** 52-43-7**Name :** 5,5-Diallylbarbituric acid; Allobarbital**Mol. formula :** C₁₀H₁₂N₂O₃**FW :** 208.21**DEA schedule :** 2**Notes :** Anti-convulsant**References :** Merck Index, 14th ed., Monograph 263.**Catalog number :** 2125-001**CASRN :** 57-43-2**Name :** (±)-Amobarbital**Mol. formula :** C₁₁H₁₈N₂O₃**FW :** 226.27**DEA schedule :** 2**Notes :** Hypnotic; sedative (but not anti-anxiety)**References :** Merck Index, 14th ed., Monograph 570.**Catalog number :** 2315-010**Name :** (1'R,S,3'SR)-5-Allyl-5-(3'-hydroxy-1'-methylbutyl)barbituric acid**Mol. formula :** C₁₂H₁₈N₂O₄**FW :** 254.28**DEA schedule :** 2**Notes :** Secobarbital metabolite**Catalog number :** 2315-011**Name :** (1'R,S,3'RS)-5-Allyl-5-(3'-hydroxy-1'-methylbutyl)barbituric acid**Mol. formula :** C₁₂H₁₈N₂O₄**FW :** 254.28**DEA schedule :** 2**Notes :** Secobarbital metabolite**References :** Carroll, FI; Mitchell, GN J Med Chem 1975, 18, 37-41.

8 – Sedatives and Hypnotics

Catalog number : 2270-010

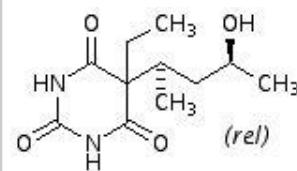
Name : (1'*RS*,3'*SR*)-5-Ethyl-5-(3'-hydroxy-1'-methylbutyl)barbituric acid

Mol. formula : C₁₁H₁₈N₂O₄

FW : 242.27 **DEA schedule :** 2

Notes : Pentobarbital metabolite

References : Carroll, FI; Mitchell, GN *J Med Chem* 1975, 18, 37–41.



Catalog number : 2270-011

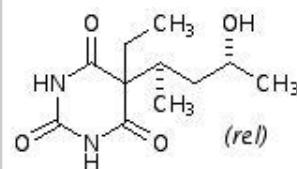
Name : (1'*RS*,3'*S*)-5-Ethyl-5-(3'-hydroxy-1'-methylbutyl)barbituric acid

Mol. formula : C₁₁H₁₈N₂O₄

FW : 242.27 **DEA schedule :** 2

Notes : Pentobarbital metabolite

References : Carroll, FI; Mitchell, GN *J Med Chem* 1975, 18, 37–41.



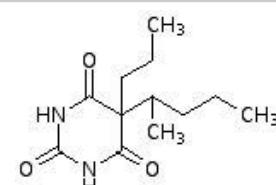
Catalog number : 2100-010

Name : (±)-5-Propyl-5-(1'-methylbutyl)barbituric acid

Mol. formula : C₁₂H₂₀N₂O₃

FW : 240.30 **DEA schedule :** 2

Notes : Sedative



Catalog number : 2100-006

CASRN : 56-29-1

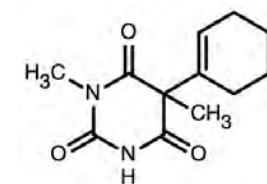
Name : (±)-Hexobarbital

Mol. formula : C₁₂H₁₆N₂O₃

FW : 236.26 **DEA schedule :** 2

Notes : Hypnotic; sedative; GABA moderator

References : Merck Index, 14th ed., Monograph 4704.



Catalog number : 2270-002

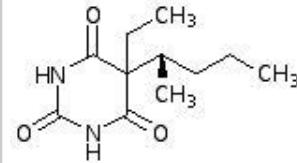
CASRN : 21642-82-0

Name : (−)-(S)-Pentobarbital

Mol. formula : C₁₁H₁₈N₂O₃

FW : 226.27 **DEA schedule :** 2

Notes : Hypnotic; sedative (but not anti-anxiety)



Catalog number : 2270-003

CASRN : 21642-83-1

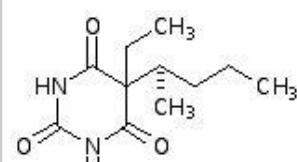
Name : (+)-(R)-Pentobarbital

Mol. formula : C₁₁H₁₈N₂O₃

FW : 226.27 **DEA schedule :** 2

Notes : Hypnotic; sedative (but not anti-anxiety)

References : Merck Index, 14th ed., Monograph 7130.



Catalog number : 2315-001

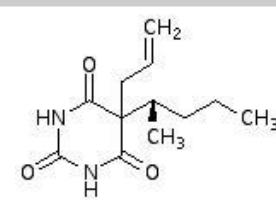
CASRN : 20224-45-7

Name : (−)-(S)-Secobarbital

Mol. formula : C₁₂H₁₈N₂O₃

FW : 238.27 **DEA schedule :** 2

Notes : Sedative (but not anti-anxiety)



Catalog number : 2315-002

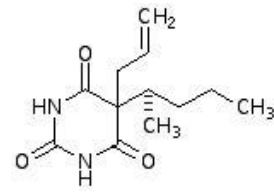
CASRN : 22328-94-5

Name : (+)-(R)-Secobarbital

Mol. formula : C₁₂H₁₈N₂O₃

FW : 238.27 DEA schedule : 2

Notes : Hypnotic; sedative (but not anti-anxiety)

References : *Merck Index*, 14th ed., Monograph 8420.Carroll, FL; Mitchell, GN *J Med Chem* 1975, 18, 37–41.**Sedatives: Benzodiazepine Class**

Catalog number : 2763-001

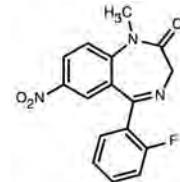
CASRN : 1622-62-4

Name : Flunitrazepam; Rohypnol

Mol. formula : C₁₆H₁₂FN₃O₃

FW : 313.28 DEA schedule : 4

Notes : Hypnotic.

References : Mattila, MA; Larni, HM *Drugs* 1980, 20, 353–74.

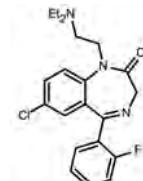
Catalog number : 2767-001

CASRN : 1172-18-5

Name : Flurazepam dihydrochloride

Mol. formula : C₂₁H₂₃ClFN₃O

FW : 460.80 DEA schedule : 4

References : *Merck Index*, 14th ed., Monograph 4198.

Catalog number : 2835-001

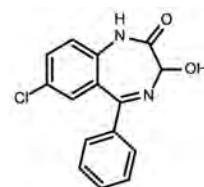
CASRN : 604-75-1

Name : (±)-Oxazepam

Mol. formula : C₁₅H₁₁ClN₂O₂

FW : 286.71 DEA schedule : 4

Notes : Anti-anxiety, anti-insomnia

References : *Merck Index*, 14th ed., Monograph 6926.**Sedatives: Butyrolactam Class**

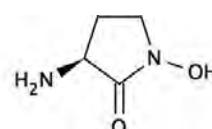
Catalog number : NOCD-071

Name : (−)-(S)-3-Amino-1-hydroxypyrrolidin-2-one; (−)-(S)-HA-966

Mol. formula : C₄H₈N₂O₂

FW : 116.12 DEA schedule : 0

Notes : Potent γ-butyrolactone-like sedative

References : Singh, L; et al. *Proc Natl Acad Sci USA* 1990, 87, 347–51.

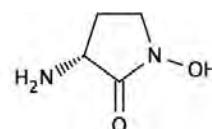
Catalog number : NOCD-072

Name : (+)-((R)-3-Amino-1-hydroxypyrrolidin-2-one; (+)-((R)-HA-966

Mol. formula : C₄H₈N₂O₂

FW : 116.12 DEA schedule : 0

Notes : Glycine/NMDA receptor antagonist

References : Singh, L; et al. *Proc Natl Acad Sci USA* 1990, 87, 347–51.

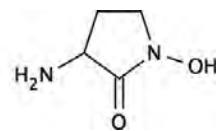
Catalog number : NOCD-070

CASRN : 1003-51-6

Name : (\pm)-3-Amino-1-hydroxypyrrolidin-2-one; (\pm)-HA-966Mol. formula : C₄H₈N₂O₂

FW : 116.12 DEA schedule : 0

Notes : Glycine/NMDA receptor antagonist

References : Leeson, PD; Iversen, LL *J Med Chem* 1994, 37, 4053-4067.**Sedatives: Methaqualone Class**

Catalog number : 2565-021

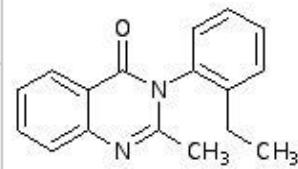
CASRN : 7432-25-9

Name : Etaqualone hydrochloride

Mol. formula : C₁₇H₁₇ClN₂O

FW : 300.79 DEA schedule : 0

Notes : CNS depressant; sedative; hypnotic

References : *Merck Index*, 14th ed., Monograph 3714.

Catalog number : 2572-001

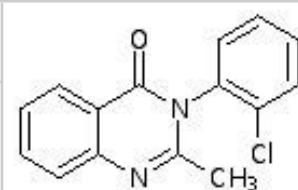
CASRN : 340-57-8

Name : Mecloqualone

Mol. formula : C₁₅H₁₁ClN₂O

FW : 270.74 DEA schedule : 1

Notes : CNS depressant; sedative; hypnotic

References : *Merck Index*, 14th ed., Monograph 5781.

Catalog number : 2565-003

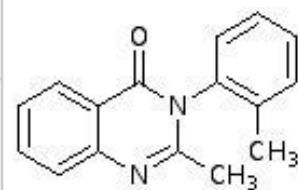
CASRN : 72-44-6

Name : Methaqualone

Mol. formula : C₁₆H₁₄N₂O

FW : 250.30 DEA schedule : 1

Notes : CNS depressant; sedative; hypnotic

References : *Merck Index*, 14th ed., Monograph 5960.

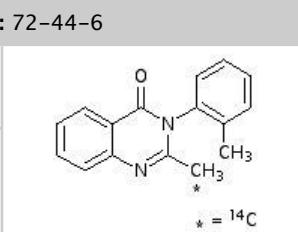
Catalog number : 2565-002

CASRN : 72-44-6

Name : [2-¹⁴C]MethaqualoneMol. formula : C₁₆H₁₄N₂O

FW : 251.30 DEA schedule : 1

Notes : Sedative; hypnotic (carbon-labeled).

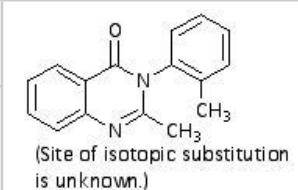
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Catalog number : 2565-001

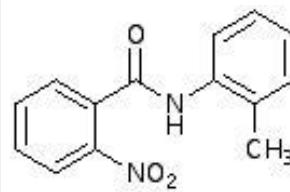
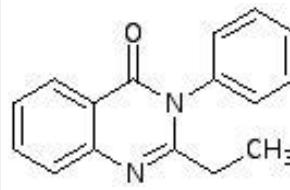
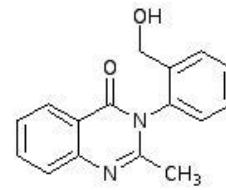
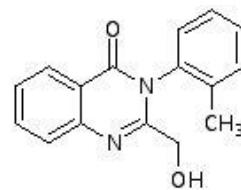
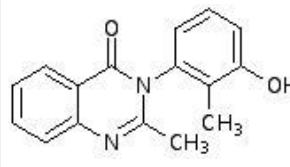
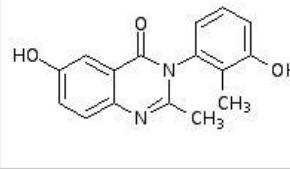
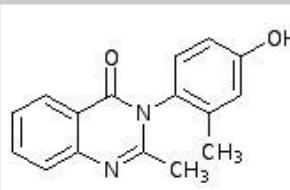
CASRN : 72-44-6

Name : [²H₄]MethaqualoneMol. formula : C₁₆H₁₄N₂O

FW : 254.32 DEA schedule : 1



(Site of isotopic substitution is unknown.)

Catalog number : 2565-012**CASRN :** 2385-25-3**Name :** 2-Nitro-*o*-benzotolidide**Mol. formula :** C₁₄H₁₂N₂O₃**FW :** 256.26 **DEA schedule :** 0**Notes :** Methaqualone urinary metabolite**References :** Murata, T; Yamamoto, I *Chem Pharm Bull (Tokyo)* 1970, 18, 133-7.**Catalog number :** 2565-020**CASRN :** 5260-41-3**Name :** 2-Ethyl-3-phenyl-4(3H)-quinazolinone**Mol. formula :** C₁₆H₁₄N₂O**FW :** 250.30 **DEA schedule :** 0**References :** Brine, GA; Coleman, ML; Carroll, FI *J Heterocyclic Chem* 1979, 16, 25.**Catalog number :** 2565-015**CASRN :** 5060-50-4**Name :** 2'-Hydroxymethaqualone**Mol. formula :** C₁₆H₁₄N₂O₂**FW :** 266.30 **DEA schedule :** 0**References :** Brine, GA; Coleman, ML; Carroll, FI *J Heterocyclic Chem* 1979, 16, 25.**Catalog number :** 2565-019**CASRN :** 5060-49-1**Name :** 2-Hydroxymethaqualone**Mol. formula :** C₁₆H₁₄N₂O₂**FW :** 266.30 **DEA schedule :** 0**References :** Brine, GA; Coleman, ML; Carroll, FI *J Heterocyclic Chem* 1979, 16, 25.**Catalog number :** 2565-014**CASRN :** 5060-63-9**Name :** 3'-Hydroxymethaqualone**Mol. formula :** C₁₆H₁₄N₂O₂**FW :** 266.30 **DEA schedule :** 0**References :** Brine, GA; Coleman, ML; Carroll, FI *J Heterocyclic Chem* 1979, 16, 25.**Catalog number :** 2565-017**CASRN :** 29541-82-0**Name :** 6, 3'-Dihydroxymethaqualone**Mol. formula :** C₁₀H₁₄N₂O₃**FW :** 282.29 **DEA schedule :** 0**References :** Brine, GA; Coleman, ML; Carroll, FI *J Heterocyclic Chem* 1979, 16, 25.**Catalog number :** 2565-013**CASRN :** 5060-52-6**Name :** 4'-Hydroxymethaqualone**Mol. formula :** C₁₆H₁₄N₂O₂**FW :** 266.30 **DEA schedule :** 0**References :** Brine, GA; Coleman, ML; Carroll, FI *J Heterocyclic Chem* 1979, 16, 25.

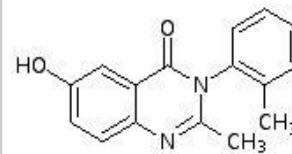
Catalog number : 2565-016

CASRN : 5060-51-5

Name : 6-Hydroxymethaqualone

Mol. formula : C₁₆H₁₄N₂O₂

FW : 266.30 DEA schedule : 0

References : Brine, GA; Coleman, ML; Carroll, FI *J Heterocyclic Chem* 1979, 16, 25.

Catalog number : 2565-018

CASRN : 5060-53-7

Name : 8-Hydroxymethaqualone

Mol. formula : C₁₆H₁₄N₂O₂

FW : 266.30 DEA schedule : 0

References : Brine, GA; Coleman, ML; Carroll, FI *J Heterocyclic Chem* 1979, 16, 25.**Sedatives: Miscellaneous**

Catalog number : 2010-001

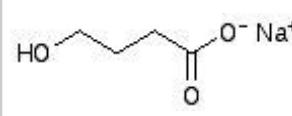
CASRN : 502-85-2

Name : γ-Hydroxybutyric acid, sodium salt; GHB

Mol. formula : C₄H₇O₃Na

FW : 126.09 DEA schedule : 1

Notes : CNS depressant; analgesic

References : *Merck Index*, 14th ed., Monograph 4815.

Catalog number : 2605-001

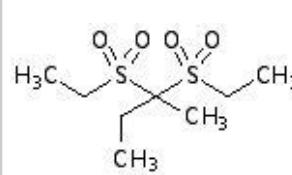
CASRN : 76-20-0

Name : Sulfonethylmethane; Ethylsulfonal

Mol. formula : C₈H₁₈O₄S₄

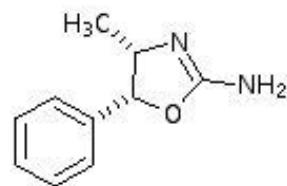
FW : 242.36 DEA schedule : 3

Notes : Hypnotic

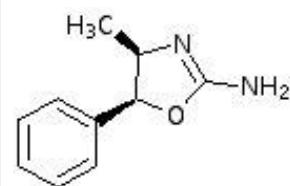
References : *Merck Index*, 14th ed., Monograph 8958.

Stimulants: Aminorex Class**Catalog number :** 1590-005**CASRN :** 133633-24-6

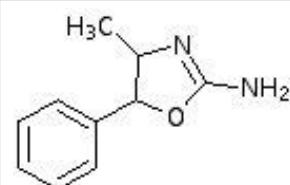
Name : $(-)$ -*cis*-(4*S*,5*R*)-2-Amino-4-methyl-5-phenyl-2-oxazoline hydrochloride; $(-)$ -*cis*-4-Methylaminorex HCl

Mol. formula : C₁₀H₁₃ClN₂O**FW :** 212.68**DEA schedule :** 1**Notes :** CNS stimulant**References :** Merck Index, 14th ed., Monograph 6018.**Catalog number :** 1590-006**CASRN :** 933777-34-5

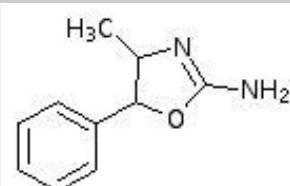
Name : $(+)$ -*cis*-(4*R*,5*S*)-2-Amino-4-methyl-5-phenyl-2-oxazoline hydrochloride; $(+)$ -*cis*-4-Methylaminorex HCl

Mol. formula : C₁₀H₁₃ClN₂O**FW :** 212.68**DEA schedule :** 1**Notes :** CNS stimulant**References :** Merck Index, 14th ed., Monograph 6018.**Catalog number :** 1590-001**CASRN :** 3568-94-3

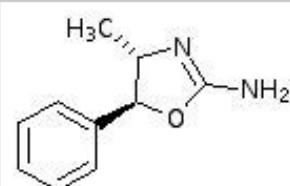
Name : (\pm) -*cis*-(4*R*,5*S*)-2-Amino-4-methyl-5-phenyl-2-oxazoline; (\pm) -*cis*-4-Methylaminorex

Mol. formula : C₁₀H₁₂N₂O**FW :** 176.22**DEA schedule :** 1**Notes :** CNS stimulant**References :** Merck Index, 14th ed., Monograph 6018.**Catalog number :** 1590-002**CASRN :** 3568-94-3

Name : (\pm) -*cis*-(4*R*,5*S*)-2-Amino-4-methyl-5-phenyl-2-oxazoline hydrochloride; (\pm) -*cis*-4-Methylaminorex HCl

Mol. formula : C₁₀H₁₃ClN₂O**FW :** 212.68**DEA schedule :** 1**Notes :** CNS stimulant**References :** Merck Index, 14th ed., Monograph 6018.**Catalog number :** 1590-003**CASRN :** 933777-34-5

Name : $(-)$ -*trans*-(4*S*,5*S*)-2-Amino-4-methyl-5-phenyl-2-oxazoline hydrochloride; $(-)$ -*trans*-4-Methylaminorex HCl

Mol. formula : C₁₀H₁₃ClN₂O**FW :** 212.68**DEA schedule :** 1**Notes :** CNS stimulant**References :** Merck Index, 14th ed., Monograph 6018.

Catalog number : 1590-004

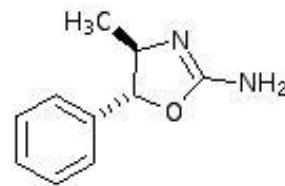
CASRN : 933777-37-8

Name : (+)-*trans*-(4*R*,5*R*)-2-Amino-4-methyl-5-phenyl-2-oxazoline hydrochloride; (+)-*trans*-4-Methylaminorex HCl

Mol. formula : C₁₀H₁₃ClN₂O

FW : 212.68 DEA schedule : 1

Notes : CNS stimulant

References : *Merck Index*, 14th ed., Monograph 6018.

Catalog number : 1590-007

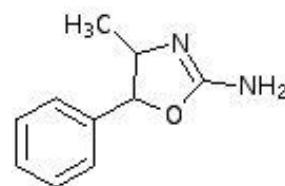
CASRN : 2077-59-0

Name : (±)-*trans*-(4*RS*,5*RS*)-2-Amino-4-methyl-5-phenyl-2-oxazoline hydrochloride; (±)-*trans*-4-Methylaminorex HCl

Mol. formula : C₁₀H₁₃ClN₂O

FW : 212.68 DEA schedule : 1

Notes : CNS stimulant

References : *Merck Index*, 14th ed., Monograph 6018.**Stimulants: Amphetamine Class**

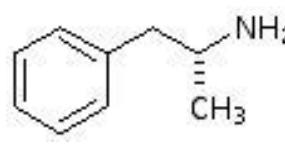
Catalog number : 1100-003

CASRN : 51-62-7

Name : (-)-Amphetamine sulfate; Levamphetamine

Mol. formula : C₁₈H₂₈N₂O₄S

FW : 368.50 DEA schedule : 2

References : *Merck Index*, 14th ed., Monograph 582.

Catalog number : 1100-006

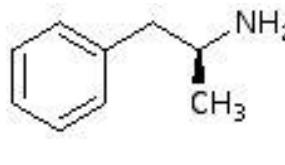
CASRN : 60-13-9

Name : (+)-Amphetamine sulfate; Dextroamphetamine sulfate

Mol. formula : C₁₈H₂₈N₂O₄S

FW : 368.50 DEA schedule : 2

Notes : CNS stimulant; sympathomimetic; anorexic

References : *Merck Index*, 14th ed., Monograph 2954.

Catalog number : 1100-007

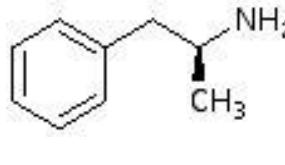
CASRN : 1462-73-3

Name : (+)-Amphetamine hydrochloride; Dextroamphetamine hydrochloride

Mol. formula : C₉H₁₄ClN

FW : 171.70 DEA schedule : 2

Notes : CNS stimulant

References : *Merck Index*, 14th ed., Monograph 2954.

Catalog number : 1100-010

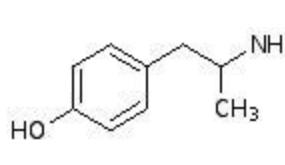
CASRN : 306-21-8

Name : (±)-4-Hydroxyamphetamine hydrobromide

Mol. formula : C₉H₁₄BrNO

FW : 232.12 DEA schedule : 2

Notes : CNS stimulant; sympathomimetic; mydriatic

References : *Merck Index*, 14th ed., Monograph 4810.

Catalog number : 1100-011

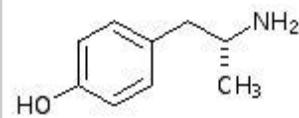
CASRN : 41509-97-1

Name : (-)-4-Hydroxyamphetamine hydrobromide

Mol. formula : C₉H₁₄BrNO

FW : 232.12 DEA schedule : 2

Notes : CNS stimulant; sympathomimetic; mydriatic

References : *Merck Index*, 14th ed., Monograph 4810.

Catalog number : 1100-012

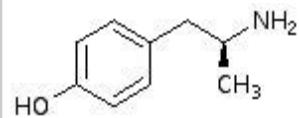
CASRN : 1693-66-9

Name : (+)-4-Hydroxyamphetamine hydrobromide

Mol. formula : C₉H₁₄BrNO

FW : 232.12 DEA schedule : 2

Notes : CNS stimulant; sympathomimetic; mydriatic

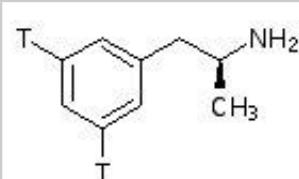
References : *Merck Index*, 14th ed., Monograph 4810.

Catalog number : 1100-009

Name : (+)-(S)-[3,5-³H(n)]AmphetamineMol. formula : C₉H₁₃N

FW : 139.22 DEA schedule : 2

Notes : CNS stimulant (tritium-labeled).

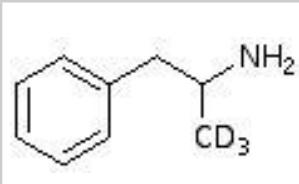


Catalog number : 1100-004

CASRN : 38875-35-3

Name : (±)-[1,1,1-²H₃]Amphetamine sulfateMol. formula : C₉H₁₃N

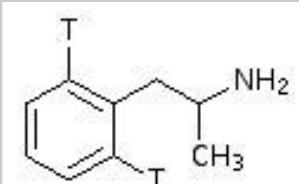
FW : 374.53 DEA schedule : 2

References : Cho, AK; et al., *Anal Chem* 1973, 45, 570–4.
Valtier, S; Cody, JT *J Anal Toxicol* 1995, 19, 375–80.

Catalog number : 1100-005

Name : (±)-[2',6'-³H₂]Amphetamine; 2,6-TritoamphetamineMol. formula : C₉H₁₃N

FW : 135.21 DEA schedule : 2



Catalog number : 1645-001

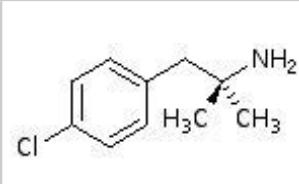
CASRN : 461-78-9

Name : Chlorphentermine; 4-Chloro-α,α-dimethyl-β-phenethylamine

Mol. formula : C₁₀H₁₄ClN

FW : 183.68 DEA schedule : 3

Notes : Anorectic

References : *Merck Index*, 14th ed., Monograph 2182.

Catalog number : 1503-001

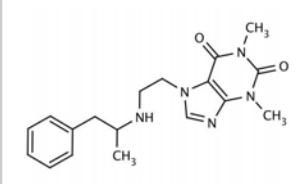
CASRN : 1892-80-4

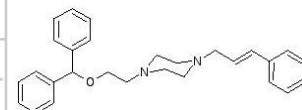
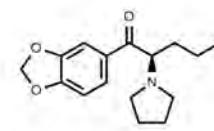
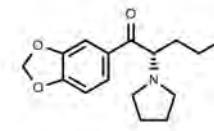
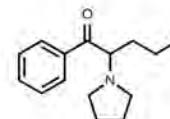
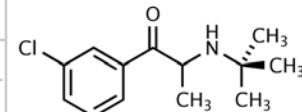
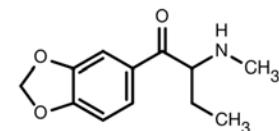
Name : Fenethylline hydrochloride; Captagon HCl

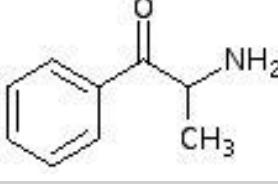
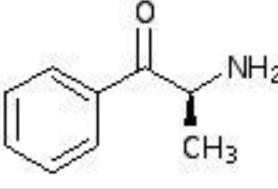
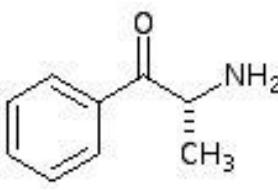
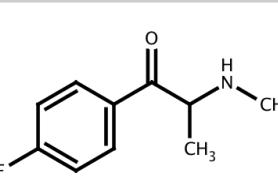
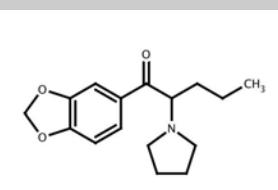
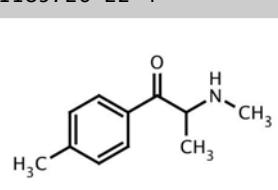
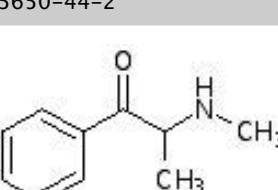
Mol. formula : C₁₈H₂₄N₅O₂Cl

FW : 341.419 DEA schedule : 2

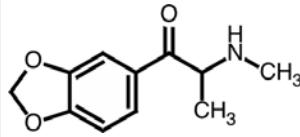
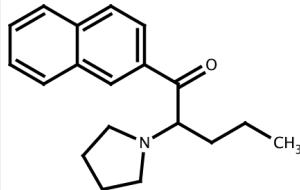
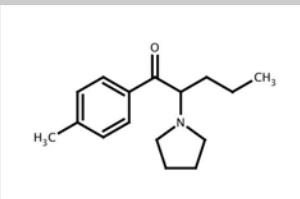
Notes : CNS stimulant

References : *Merck Index*, 14th ed., Monograph 3972.

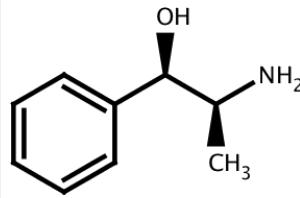
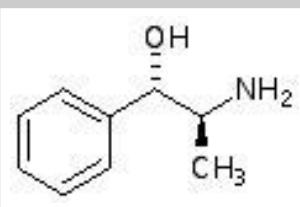
Stimulants: Benzhydrol Class**Catalog number :** NOCD-040**CASRN :** 67469-57-2**Name :** GBR 12783**Mol. formula :** C₂₈H₃₄Cl₂N₂O**FW :** 485.48 **DEA schedule :** 0**Notes :** Dopamine uptake inhibitor.**References :** Bonnet, JJ; Costentin J Eur J Pharmacol 1986, 121, 199–209.
Chagraoui, A; et al. Neurosci Lett 1987, 78, 175–9.**Stimulants: Cathinone Class****Catalog number :** 7535-002**new****CASRN :** 1669434-93-8**Name :** (R)-MDPV hydrochloride**Mol. formula :** C₁₆H₂₁NO₃**FW :** 275.35 **DEA schedule :** 1**Catalog number :** 7535-003**new****CASRN :** 1669434-94-9**Name :** (S)-MDPV hydrochloride**Mol. formula :** C₁₆H₂₁NO₃**FW :** 275.35 **DEA schedule :** 1**Catalog number :** 7545-001**new****CASRN :** 5485-65-4**Name :** α-PVP hydrochloride**Mol. formula :** C₁₅H₂₁NO**FW :** 231.34**References :** Meltzer, PC; et al., J Med Chem 2006, 49, 1420–32.**Catalog number :** 1610-002**CASRN :** 31677-93-7**Name :** (±)-Bupropion hydrochloride**Mol. formula :** C₁₃H₁₉Cl₂NO**FW :** 276.21 **DEA schedule :** 0**References :** Martin, P; Massol, J; Colin, JN; Lacomblez, L; Puech, AJ
Pharmacopsychiatry 1990, 23, 87–94.**Catalog number :** 7541-001**CASRN :** 802575-11-7**Name :** Butyline; bk-MBDB**Mol. formula :** C₁₂H₁₅NO₃**FW :** 221.26 **DEA schedule :** 1**References :** Gatch, MB; Taylor CM; Forster MJ, Behav Pharmacol 2013, 24, 437–47.

Catalog number : 1235-010	CASRN : 76333-53-4	
Name : (±)-Cathinone hydrochloride		
Mol. formula : C ₉ H ₁₂ CINO	FW : 185.66	DEA schedule : 1
Notes : CNS stimulant; psychotropic drug		
References : Kalix, P <i>Pharmacol Toxicol</i> 1992, 70, 77–86. <i>Merck Index</i> , 14th ed., Monograph 1906.		
Catalog number : 1235-011	CASRN : 71031-15-7	
Name : (−)(S)-Cathinone hydrochloride		
Mol. formula : C ₉ H ₁₂ CINO	FW : 185.66	DEA schedule : 1
Notes : CNS stimulant; psychotropic drug		
References : <i>Merck Index</i> , 14th ed., Monograph 1906.		
Catalog number : 1235-012	CASRN : 80096-54-4	
Name : (+)-(R)-Cathinone hydrochloride		
Mol. formula : C ₉ H ₁₂ CINO	FW : 185.66	DEA schedule : 1
Notes : CNS stimulant; psychotropic drug		
Catalog number : 1237-003	CASRN : 7589-35-7	
Name : 4-Fluoromethcathinone; 4-FMC; Flephedrone		
Mol. formula : C ₁₀ H ₁₂ FNO	FW : 181.21	DEA schedule : 1
References : Archer RP, <i>Forensic Sci Int</i> , 2009, 185(1–3), 10–20.		
Catalog number : 7535-001	CASRN : 24622-62-6	
Name : (±)-Methyleneidioxypyrovalerone HCl; MDPV HCl		
Mol. formula : C ₁₆ H ₂₂ CINO ₃	FW : 311.81	DEA schedule : 1
Notes : Stimulant; norepinephrine/dopamine reuptake inhibitor.		
References : Coppola, M; Mondola R, <i>Toxicol Lett</i> 2012, 208, 12–5.		
Catalog number : 1248-001	CASRN : 1189726-22-4	
Name : (±)-Mephedrone hydrochloride		
Mol. formula : C ₁₁ H ₁₆ CINO	FW : 213.70	DEA schedule : 0
References : Wood, DM; et al. <i>J Med Toxicol</i> 2010, 10.1007/s13181-010-0018-5.		
Catalog number : 1237-001	CASRN : 5650-44-2	
Name : (±)-N-Methcathinone hydrochloride		
Mol. formula : C ₁₀ H ₁₄ CINO	FW : 199.67	DEA schedule : 1
Notes : CNS stimulant; psychotropic drug		

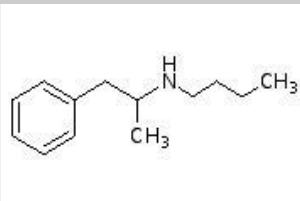
9 – Stimulants

Catalog number : 7540-001		CASRN : 186028-79-5
Name : Methylone HCl		
Mol. formula : C ₁₁ H ₁₃ NO ₃	FW : 207.23	DEA schedule : 1
References : Cozzi NV. <i>Neuropsychopharmacology</i> , 2012, 37(5), 1192–1203. Niesink RJ. <i>Addict Biol</i> , 2005, 10(4), 321–323.		
Catalog number : NOCD-126		CASRN : 850352-11-3
Name : Naphyrone; Naphthylpyrovalerone		
Mol. formula : C ₁₉ H ₂₃ NO	FW : 281.39	DEA schedule : 1
References : Meltzer PC, Butler D, Deschamps JR, Madras BK. <i>J Med Chem</i> , 2006, 49(4), 1420–1432.		
Catalog number : 1485-001		CASRN : 1147-62-2
Name : (±)-Pyrovalerone HCl		
Mol. formula : C ₁₆ H ₂₄ CINO	FW : 281.82	DEA schedule : 5
References : Stille, G; et al., <i>Arzneimittelforschung</i> 1963, 13, 871–7. Michaelis, W; Russel JH; Schindler O, <i>J Med Chem</i> 1970, 13, 497–503.		

Stimulants: Ephedrine Class

Catalog number : 1230-016		CASRN : 53643-20-2
Name : (–)-Norephedrine hydrochloride		
Mol. formula : C ₉ H ₁₄ CINO	FW : 187.68	DEA schedule : 4
Notes : Nasal decongestant; appetite suppressant		
References : <i>Merck Index</i> , 14th ed., Monograph 7307.		
Catalog number : 1230-017		CASRN : 36393-56-3
Name : (+)-Norpseudoephedrine hydrochloride		
Mol. formula : C ₉ H ₁₄ CINO	FW : 187.68	DEA schedule : 4
Notes : Nasal decongestant; appetite suppressant		
References : <i>Merck Index</i> , 14th ed., Monograph 6714.		

Stimulants: Methamphetamine Class

Catalog number : 1100-008		CASRN : 51799-33-8
Name : (±)-N-(n-Butyl)amphetamine hydrochloride		
Mol. formula : C ₁₃ H ₂₂ CIN	FW : 227.78	DEA schedule : 2
Notes : CNS stimulant		
References : Woolverton WL; Shybut, G; Johanson, CE <i>Pharmacol Biochem Behav</i> 1980, 13, 869–76.		

Catalog number : 1480-001

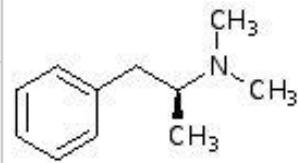
CASRN : 17279-39-9

Name : (+)-N,N-Dimethylamphetamine hydrochloride

Mol. formula : C₁₁H₁₈CIN

FW : 199.72 DEA schedule : 1

Notes : CNS stimulant

References : Ricaurte, GA; et al. *Brain Res* 1989, 490, 301–6.Katz, JL; et al. *Psychopharmacology (Berl)* 1992, 107, 315–8.

Catalog number : 1475-001

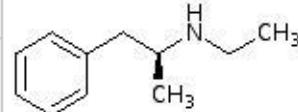
CASRN : 33817-11-7

Name : (+)-N-Ethylamphetamine hydrochloride

Mol. formula : C₁₁H₁₈CIN

FW : 199.71 DEA schedule : 1

Notes : CNS stimulant

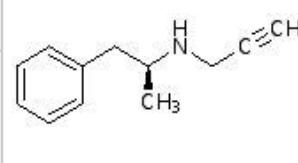
References : *Merck Index*, 14th ed., Monograph 3763.

Catalog number : MEDD-007

Name : (−)(S)-Desmethylselegiline hydrochloride

Mol. formula : C₁₂H₁₆CIN

FW : 209.72 DEA schedule : 0

References : Mytilineou C; Radcliffe PM; Olanow CW *J Neurochem* 1997, 68, 434–6.

Catalog number : 1105-001

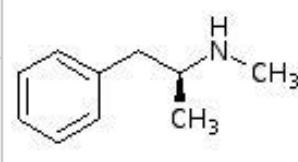
CASRN : 51-57-0

Name : (+)-(S)-Methamphetamine hydrochloride

Mol. formula : C₁₀H₁₆CIN

FW : 185.70 DEA schedule : 2

Notes : CNS stimulant; sympathomimetic; anorexic

References : *Merck Index*, 14th ed., Monograph 5948.

Catalog number : 1105-003

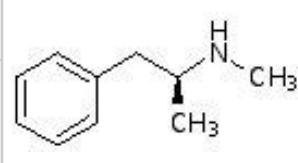
CASRN : 537-46-2

Name : (+)-Methamphetamine base

Mol. formula : C₁₀H₁₅N

FW : 149.24 DEA schedule : 2

Notes : CNS stimulant

References : *Merck Index*, 14th ed., Monograph 5948.

Catalog number : 1105-005

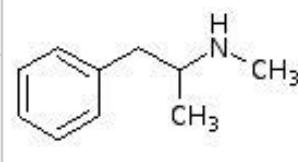
CASRN : 300-42-5

Name : (±)-Methamphetamine hydrochloride

Mol. formula : C₁₀H₁₆CIN

FW : 185.69 DEA schedule : 2

Notes : CNS stimulant

References : *Merck Index*, 14th ed., Monograph 5948.

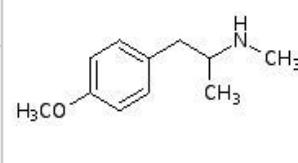
Catalog number : 1105-014

CASRN : 22331-70-0

Name : (±)-4-Methoxymethamphetamine hydrochloride; PMMA

Mol. formula : C₁₁H₁₈CINO

FW : 215.75 DEA schedule : 0

References : Glennon, RA; Ismaiel, AE; Martin, B; Poff, D; Sutton, M *Pharmacol Biochem Behav* 1988, 31, 9–13.

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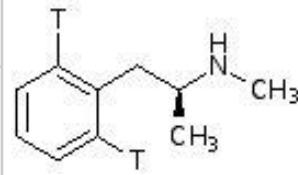
Catalog number : 1105-004

Name : (+)-[2',6'-³H(n)]Methamphetamine hydrochloride

Mol. formula : C₁₀H₁₆CIN

FW : 185.69 DEA schedule : 2

Notes : CNS stimulant (tritium-labeled).



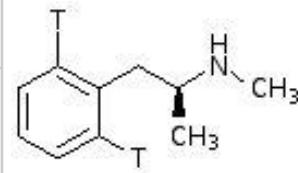
Catalog number : 1105-008

Name : (+)-[2,6-³H(n)]Methamphetamine

Mol. formula : C₁₀H₁₅N

FW : 153.25 DEA schedule : 2

Notes : CNS stimulant (tritium-labeled).

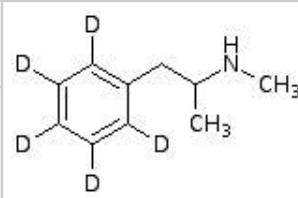


Catalog number : 1105-002

Name : (±)-[2',3',4',5',6'-²H₅]Methamphetamine hydrochloride

Mol. formula : C₁₀H₁₆CIN

FW : 190.72 DEA schedule : 2



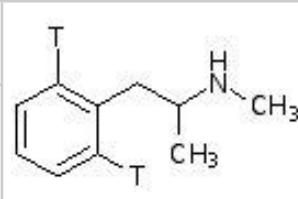
Catalog number : 1105-006

Name : (±)-[2',6'-³H(n)]Methamphetamine hydrochloride

Mol. formula : C₁₀H₁₆CIN

FW : 189.71 DEA schedule : 2

Notes : CNS stimulant (tritium-labeled).



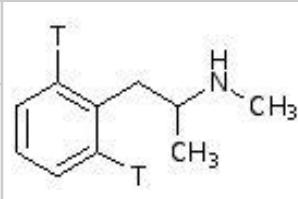
Catalog number : 1105-007

Name : (±)-[2,6-³H(n)]Methamphetamine

Mol. formula : C₁₀H₁₅N

FW : 153.25 DEA schedule : 2

Notes : CNS stimulant (tritium-labeled).

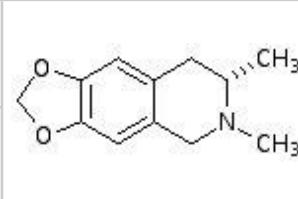


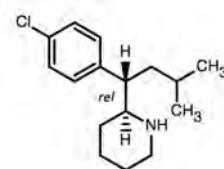
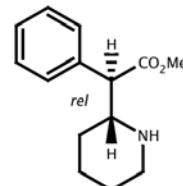
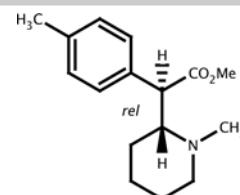
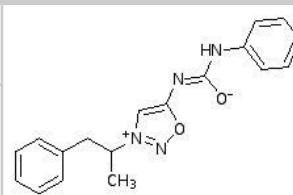
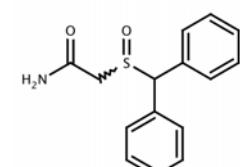
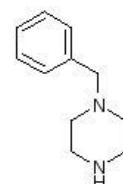
Catalog number : NOCD-074

Name : (3S)-(+)-2,3-Dimethyl-6,7-methylenedioxytetrahydroisoquinoline hydrochloride

Mol. formula : C₁₂H₁₆CINO₂

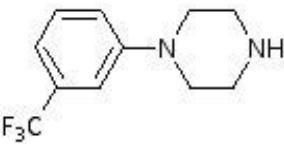
FW : 241.72 DEA schedule : 0



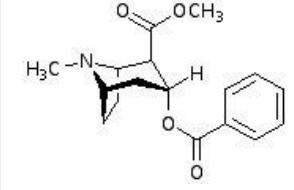
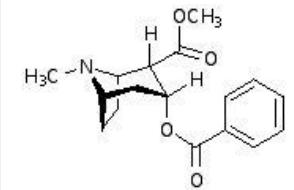
Stimulants: Methylphenidate Class**Catalog number :** NOCD-140**CASRN :** 928046-68-8**Name :** (RR/SS)-2-[1-(4-Chlorophenyl)-3-methylbutyl]piperidine hydrochloride**Mol. formula :** C₁₆H₂₄ClN**FW :** 265.83**DEA schedule :** 0**Notes :** Methylphenidate analog**References :** Froimowitz, M; et al. *J. Med. Chem.* 2007, 50, 219–232.**Catalog number :** 1724-001**CASRN :** 298-59-9**Name :** (±)-*threo*-Methylphenidate hydrochloride; Ritalin**Mol. formula :** C₁₄H₂₀ClNO₂**FW :** 269.77**DEA schedule :** 2**Notes :** Mild CNS stimulant**References :** Merck Index, 14th ed., Monograph 6110.**Catalog number :** MEDD-014**CASRN :** 204981-87-3**Name :** (±)-N-Methyl-*threo*-*p*-methylmethylphenidate hydrochloride**Mol. formula :** C₁₆H₂₄ClNO₂**FW :** 297.83**DEA schedule :** 0**Stimulants: Miscellaneous****Catalog number :** NOCD-076**CASRN :** 34262-84-5**Name :** Mesocarb; Sydnocarb**Mol. formula :** C₁₈H₁₈N₄O₂**FW :** 322.14**DEA schedule :** 3**Notes :** Stimulant; dopamine reuptake inhibitor**References :** Bashkatova V; et al. *Ann N Y Acad Sci* 2002, 965, 180–92.**Catalog number :** 1680-001**CASRN :** 112111-43-0**Name :** (±)-Modafinil**Mol. formula :** C₁₅H₁₅NO₂S**FW :** 273.35**DEA schedule :** 4**References :** Minzenberg, M; Carter, C *Neuropsychopharmacology* 2007, 33, 1477–1502.**Stimulants: Piperazine Class****Catalog number :** 7493-001**CASRN :** 2759-28-6**Name :** 1-Benzylpiperazine difumarate; BZP**Mol. formula :** C₁₉H₂₄N₂O₈**FW :** 408.41**DEA schedule :** 1**Notes :** Psychomotor stimulant; serotonin (5-HT1) agonist.**References :** Lyon, RA; et al. *J Med Chem* 1986, 29, 630–4.Staack, RF *Lancet* 2007, 369, 1411–3.Lecompte, Y; Roussel O; Perrin M *Ann Pharm Fr* 2008, 66, 85–91.

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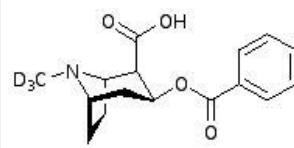
Catalog number : 7494-001	CASRN : 16015-69-3
Name : 1-(3-Trifluoromethylphenyl)piperazine hydrochloride; TFMPP	
Mol. formula : C ₁₁ H ₁₄ ClF ₃ N ₂	FW : 266.69 DEA schedule : 1
Notes : Psychomotor stimulant; serotonin (5-HT1) agonist.	
References : Glennon, RA; McKenney JD; Young R <i>Life Sci</i> 1984, 35, 1475–80. Lecompte, Y; Roussel O; Perrin M <i>Ann Pharm Fr</i> 2008, 66, 85–91.	



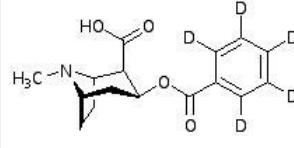
Stimulants: Tropane Class

Catalog number : 9041-010	CASRN : 518-97-8
Name : (±)-Allococaine	
Mol. formula : C ₁₇ H ₂₁ NO ₄	FW : 303.35 DEA schedule : 2
	
Catalog number : 9041-008	
Name : (±)-Allopseudoephedrine hydrochloride	
Mol. formula : C ₁₇ H ₂₂ ClNO ₄	FW : 339.81 DEA schedule : 2
	

Catalog number : 9180-003

Name : (–)-[N-C ² H ₃]Benzoyleccgonine	CASRN : 519-09-05
Mol. formula : C ₁₆ H ₁₉ NO ₄	FW : 292.34 DEA schedule : 2
	
Catalog number : 9180-011	

Catalog number : 9180-011

Name : (–)-[Phenyl- ² H ₅]Benzoyleccgonine	CASRN : 519-09-05
Mol. formula : C ₁₆ H ₁₉ NO ₄	FW : 294.36 DEA schedule : 2
	
Catalog number : 9180-001	

Catalog number : 9180-001

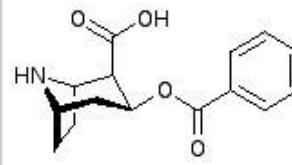
Name : (–)-Benzoyleccgonine	CASRN : 519-09-05
Mol. formula : C ₁₆ H ₁₉ NO ₄	FW : 289.34 DEA schedule : 2
Notes : Major metabolite of cocaine.	
References : Merck Index, 14th ed., Monograph 1113.	

Catalog number : 9180-004

Name : (-)-Benzoylnorecgonine hydrochloride

Mol. formula : C₁₅H₁₈ClNO₄

FW : 311.77 DEA schedule : 2



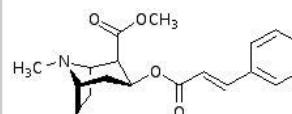
Catalog number : 9041-011

CASRN : 521-67-5

Name : (-)-trans-Cinnamoylcocaine

Mol. formula : C₁₉H₂₃NO₄

FW : 329.38 DEA schedule : 2

References : Novak, M; Salemink, CA; Khan, I *J Ethnopharmacol* 1984, 10, 261-74.

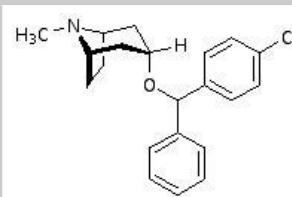
Catalog number : NOCD-004

CASRN : 5627-46-3

Name : Clobenztropine HCl

Mol. formula : C₂₁H₂₄ClNO • HCl

FW : 378.35 DEA schedule : 0



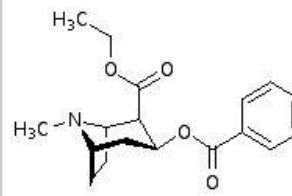
Catalog number : 9041-018

Name : (-)-Cocaethylene fumarate

Mol. formula : C₂₂H₂₇NO₈

FW : 433.45 DEA schedule : 2

Notes : Cocaine metabolite.

References : Hearn, WL; et al. *Pharmacol Biochem Behav* 1991, 39, 531-3.
Hearn, WL; et al. *J Neurochem* 1991, 56, 698-701.

Catalog number : 9041-001

CASRN : 53-21-4

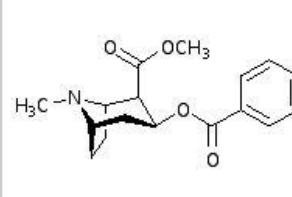
Name : (-)-Cocaine hydrochloride

Mol. formula : C₁₇H₂₂ClNO₄

FW : 339.81 DEA schedule : 2

Notes : CNS stimulant; local anesthetic

References : Merck Index, 14th ed., Monograph 2455.

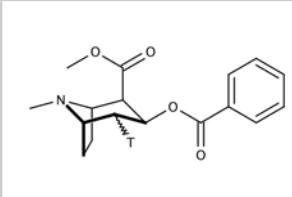


Catalog number : 9041-004

CASRN : 85438-94-4

Name : (-)-[4-³H]CocaineMol. formula : C₁₇H₂₁NO₄

FW : 303.35 DEA schedule : 2



Catalog number : 9041-012

CASRN : 50-36-2

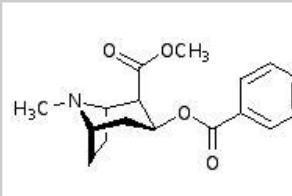
Name : (-)-Cocaine base

Mol. formula : C₁₇H₂₁NO₄

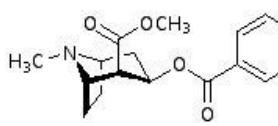
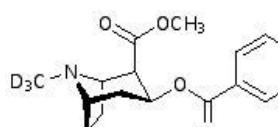
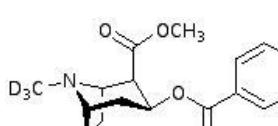
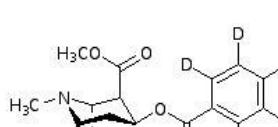
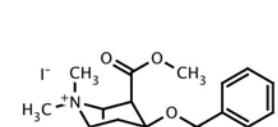
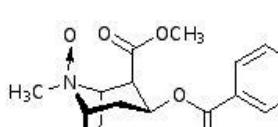
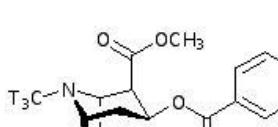
FW : 303.35 DEA schedule : 2

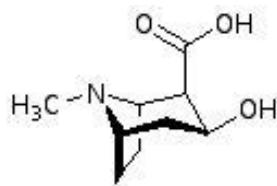
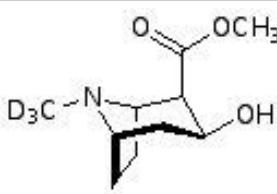
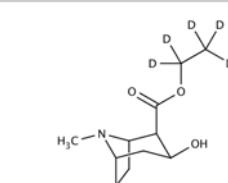
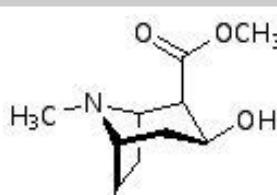
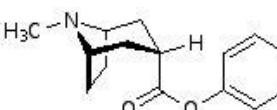
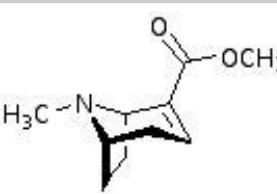
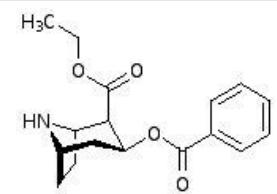
Notes : CNS stimulant; local anesthetic

References : Merck Index, 14th ed., Monograph 2455.

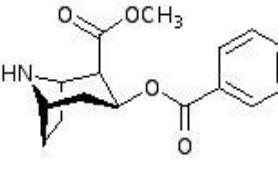
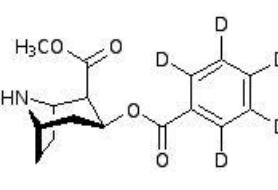
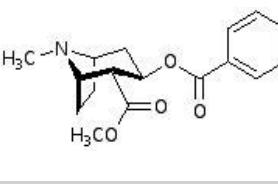
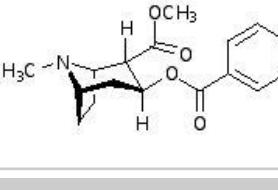
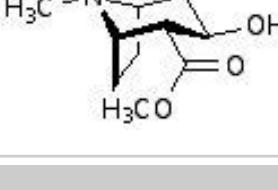
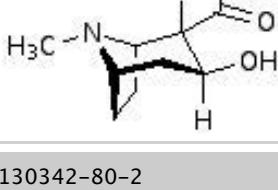
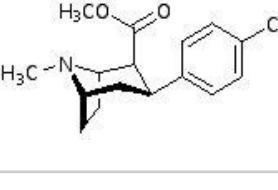


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Catalog number : 9041-020	CASRN : 47195-07-3	
Name : (+)-Cocaine base		
Mol. formula : C ₁₇ H ₂₁ NO ₄	FW : 303.35	DEA schedule : 2
		
Catalog number : 9041-006		
Name : (-)-[N-C ² H ₃]Cocaine		
Mol. formula : C ₁₇ H ₂₁ NO ₄	FW : 306.37	DEA schedule : 2
		
Catalog number : 9041-007		
Name : (-)-[N-C ² H ₃]Cocaine hydrochloride		
Mol. formula : C ₁₇ H ₂₂ ClNO ₄	FW : 342.83	DEA schedule : 2
		
Catalog number : 9041-009		
Name : (-)-[Phenyl- ² H ₅]Cocaine		
Mol. formula : C ₁₇ H ₂₁ NO ₄	FW : 344.84	DEA schedule : 2
		
Catalog number : 9041-013	CASRN : 5937-29-1	
Name : (-)-Cocaine methiodide		
Mol. formula : C ₁₈ H ₂₄ INO ₄	FW : 445.28	DEA schedule : 2
References : Abraham, P; Pitner, JB; Lewin, AH; Boja, JW; Kuhar, MJ; Carroll, FI <i>J Med Chem</i> 1992, 35, 141-4.		
Catalog number : 9041-014		
Name : (-)-Cocaine N-oxide hydrochloride		
Mol. formula : C ₁₇ H ₂₂ ClNO ₅	FW : 355.81	DEA schedule : 2
		
Catalog number : 9041-003		
Name : (-)-[N-C ³ H ₃]Cocaine		
Mol. formula : C ₁₇ H ₂₁ NO ₄	FW : 309.38	DEA schedule : 2
		

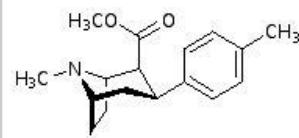
Catalog number : 9180-002**CASRN :** 5796-31-6**Name :** (–)–Ecgonine hydrochloride**Mol. formula :** C₉H₁₆ClNO₃**FW :** 221.69 **DEA schedule :** 2**Notes :** Cocaine metabolite.**References :** Merck Index, 14th ed., Monograph 3493.**Catalog number :** 9180-007**Name :** [N–C²H₃]Ecgonine methyl ester hydrochloride**Mol. formula :** C₁₀H₁₈ClNO₃**FW :** 238.73 **DEA schedule :** 2**Catalog number :** 9180-021**CASRN :** 259526-73-3**Name :** Ecgonine (1,1,2,2,2–²H₅)ethyl ester perchlorate**Mol. formula :** C₁₁H₁₄D₅NO₃ • HClO₄**FW :** 318.76 **DEA schedule :** 2**Catalog number :** 9180-006**CASRN :** 7143-09-1**Name :** (–)–Ecgonine methyl ester hydrochloride**Mol. formula :** C₁₀H₁₈ClNO₃**FW :** 235.71 **DEA schedule :** 2**Catalog number :** 9041-015**Name :** Isotropacocaine hydrochloride**Mol. formula :** C₁₅H₂₀ClNO₂**FW :** 281.79 **DEA schedule :** 2**Catalog number :** 9180-015**Name :** (–)–Anhydroecgonine methyl ester fumarate; Methyl ecgonidine**Mol. formula :** C₁₄H₁₉NO₆**FW :** 297.31 **DEA schedule :** 0**Catalog number :** 9041-019**CASRN :** 137220-02-1**Name :** (–)–Norcocacethylene fumarate**Mol. formula :** C₂₁H₂₅NO₈**FW :** 419.43 **DEA schedule :** 2

9 – Stimulants

Catalog number : 9041-002		CASRN : 18717-72-1
Name : (-)-Norcocaine		
Mol. formula : C ₁₆ H ₁₉ NO ₄	FW : 289.32	DEA schedule : 2
References : Wang, Q; Simpao, A; Sun L; Falk, JL; Lau, CE <i>Psychopharmacology (Berl)</i> 2001, 153, 341–52. Kovacic, P <i>Med Hypotheses</i> 2005, 64, 350–6.		
Catalog number : 9041-017		CASRN : 478-73-9
Name : (-)-[Phenyl- ² H ₅]Norcocaine fumarate		
Mol. formula : C ₂₀ H ₂₃ NO ₈	FW : 410.43	DEA schedule : 2
		
Catalog number : 9041-021		CASRN : 478-73-9
Name : (-)-Pseudococaine hydrochloride		
Mol. formula : C ₁₇ H ₂₂ CINO ₄	FW : 339.81	DEA schedule : 2
		
Catalog number : 9041-005		CASRN : 478-73-9
Name : (+)-Pseudococaine hydrochloride		
Mol. formula : C ₁₇ H ₂₂ CINO ₄	FW : 339.81	DEA schedule : 2
		
Catalog number : 9180-020		
Name : (-)-Pseudoecgonine methyl ester		
Mol. formula : C ₁₀ H ₁₇ NO ₃	FW : 199.25	DEA schedule : 2
		
Catalog number : 9180-005		
Name : (+)-Pseudoecgonine methyl ester		
Mol. formula : C ₁₀ H ₁₇ NO ₃	FW : 199.25	DEA schedule : 2
		
Catalog number : NOCD-020		CASRN : 130342-80-2
Name : (-)-3β-(4-Chlorophenyl)tropan-2β-carboxylic acid methyl ester tartrate salt; RTI-31		
Mol. formula : C ₂₀ H ₂₆ CINO ₈	FW : 443.87	DEA schedule : 0
Notes : Nonselective monoamine uptake inhibitor.		
References : Carroll, FI; et al. <i>J Med Chem</i> 1995, 38, 379–88.		

Catalog number : NOCD-021**CASRN :** 130342-81-3

Name : (-)-3 β -(4-Methylphenyl)tropan-2 β -carboxylic acid methyl ester tartrate; RTI-32

Mol. formula : C₂₁H₂₉NO₈**FW :** 423.46 **DEA schedule :** 0**Notes :** Nonselective monoamine uptake inhibitor.

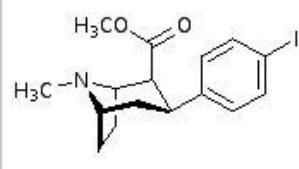
References : Boja, JW; Carroll, FI; Rahman, MA; Philip, A; Lewin, AH; Kuhar, MJ *Eur J Pharmacol* 1990, 184, 329-32.

Catalog number : NOCD-025**CASRN :** 133647-95-7

Name : (-)-3 β -(4-Iodophenyl)tropan-2 β -carboxylic acid methyl ester tartrate; RTI-55

Mol. formula : C₂₀H₂₆INO₈**FW :** 535.32 **DEA schedule :** 0**Notes :** Nonselective monoamine uptake inhibitor.

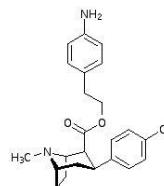
References : Boja, JW; et al. *Eur J Pharmacol* 1991, 194, 133-4.
Carroll, FI; et al. *J Med Chem* 1991, 34, 2719-25.
Carroll, FI; et al. *J Med Chem* 1995, 38, 379-88.

**Catalog number :** NOCD-029

Name : (-)-3 β -(4-Chlorophenyl)tropan-2 β -carboxylic acid (4'-aminophenyl)ethyl ester dihydrochloride; RTI-75

Mol. formula : C₂₃H₂₉Cl₃N₂O₂**FW :** 471.86 **DEA schedule :** 0

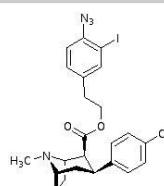
References : Carroll, FI; et al. *J Med Chem* 1992, 35, 1813-7.

**Catalog number :** NOCD-030

Name : (-)-3 β -(4-Chlorophenyl)tropan-2 β -carboxylic acid (4'-Azido-3'-iodophenyl)ethyl ester hydrochloride; RTI-82

Mol. formula : C₂₃H₂₅Cl₂IN₄O₂**FW :** 587.29 **DEA schedule :** 0**Notes :** Dopamine transporter photoaffinity ligand.

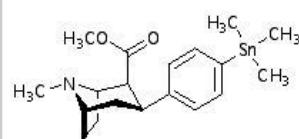
References : Carroll, FI; et al. *J Med Chem* 1992, 35, 1813-7.

**Catalog number :** NOCD-031

Name : 3 β -[4-(Trimethylstannyl)phenyl]tropan-2 β -carboxylic acid methyl ester; RTI-89

Mol. formula : C₁₉H₂₉NO₂Sn**FW :** 422.12 **DEA schedule :** 0**Notes :** Precursor for the synthesis of [¹²⁵I]- and [¹²³I]-RTI-55.

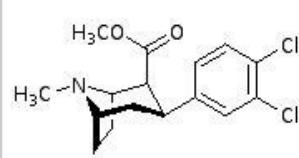
References : Carroll, FI; et al. *Med Chem Res* 1991, 1, 289-294.

**Catalog number :** NOCD-028

Name : (-)-3 β -(3,4-Dichlorophenyl)tropan-2 β -carboxylic acid methyl ester hydrochloride; RTI-111

Mol. formula : C₁₆H₂₀Cl₃NO₂**FW :** 364.70 **DEA schedule :** 0**Notes :** Nonselective monoamine uptake inhibitor.

References : Carroll, FI; et al. *J Med Chem* 1995, 38, 379-88.



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Catalog number : NOCD-026	CASRN : 316790-73-5	
Name : (-)-3β-(4-Chlorophenyl)tropan-2β-carboxylic acid phenyl ester hydrochloride; RTI-113		
Mol. formula : C ₂₁ H ₂₅ Cl ₂ NO ₃	FW : 410.32	DEA schedule : 0
Notes : Selective dopamine uptake inhibitor.		
References : Carroll, FI; et al. <i>J Med Chem</i> 1995, 38, 379-88.		
Catalog number : NOCD-027	CASRN : 146145-21-3	
Name : (-)-3β-(4-Iodophenyl)tropan-2β-carboxylic acid isopropyl ester hydrochloride; RTI-121; IPCIT		
Mol. formula : C ₁₈ H ₂₅ ClINO ₄	FW : 449.74	DEA schedule : 0
References : Carroll, FI; et al. <i>J Med Chem</i> 1995, 38, 379-88.		
Catalog number : NOCD-077		
Name : [¹²⁵ I]RTI-121		
Mol. formula : C ₁₈ H ₂₄ INO ₂	FW : 413.29	DEA schedule : 0
References : Scheffel, U; et al. <i>Neuroreport</i> 1992, 3, 969-72. Carroll, FI; et al. <i>J Med Chem</i> 1995, 38, 379-88.		
Catalog number : NOCD-032		
Name : (-)-3β-[4-(Trimethylstannyl)phenyl]tropan-2β-carboxylic acid isopropyl ester; RTI-136		
Mol. formula : C ₂₁ H ₃₃ NO ₂ Sn	FW : 450.19	DEA schedule : 0
Notes : Precursor for the synthesis of [¹²⁵ I]- and [¹²³ I]-RTI-121.		
References : Carroll, FI; et al. <i>Med Chem Res</i> 1991, 1, 289-294.		
Catalog number : 9041-016	CASRN : 637-23-0	
Name : Tropacocaine hydrochloride		
Mol. formula : C ₁₅ H ₂₀ CINO ₂	FW : 281.79	DEA schedule : 2
References : Novak, M; Saleminik, CA; Khan, I <i>J Ethnopharmacol</i> 1984, 10, 261-74. Meyer, EM; et al. <i>J Pharmacol Exp Ther</i> 1990, 254, 584-90.		
Catalog number : NOCD-023	CASRN : 50372-80-0	
Name : (-)-3β-Phenyltropan-2β-carboxylic acid methyl ester tartrate; WIN 35,065-2		
Mol. formula : C ₂₀ H ₂₂ NO ₈	FW : 409.43	DEA schedule : 0
Notes : Nonselective monoamine uptake inhibitor.		
References : Carroll, FI; et al. <i>J Med Chem</i> 1995, 38, 379-88.		
Catalog number : NOCD-024		
Name : (+)-3β-Phenyltropan-2β-carboxylic acid methyl ester tartrate; WIN 35,065-3		
Mol. formula : C ₂₀ H ₂₂ NO ₈	FW : 409.43	DEA schedule : 0

Catalog number : NOCD-022	CASRN : 50370-56-4
Name : (-)-3β-(4-Fluorophenyl)tropan-2β-carboxylic acid methyl ester tartrate; WIN 35,428	
Mol. formula : C ₂₀ H ₂₆ FNO ₈	FW : 427.42 DEA schedule : 0
Notes : <i>Nonselective monoamine uptake inhibitor.</i>	
References : Carroll, FI; et al. <i>J Med Chem</i> 1995, 38, 379-88.	

Stimulants (dosage form): Stock Solutions

Catalog number : 9041-022
Name : Injectable cocaine hydrochloride (10 mg/mL and 20 mg/mL)
DEA schedule : 2
Notes : <i>Not for human use.</i>

Dosage Form



Miscellaneous: Dopaminergic**Catalog number :** MEDD-002

Name : (\pm)-1-(2-Bromo-4,5-dimethoxybenzyl)-7-hydroxy-6-methoxy-2-methyl-1,2,3,4-tetrahydroisoquinoline hydrobromide; A69024 HBr

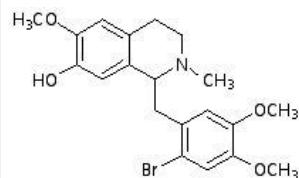
Mol. formula : C₂₀H₂₅Br₂NO₄

FW : 503.23

DEA schedule : 0

Notes : Dopamine D1 receptor antagonist

References : Caine, SB; Koob, GF., *J Pharmacol Exp Ther* 1994, 270, 209-18.

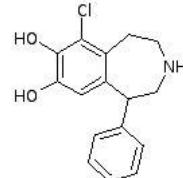
**Catalog number :** MEDD-004

Name : (\pm)-6-Chloro-7,8-dihydroxy-1-phenyl-2,3,4,5-tetrahydro[1H]-3-benzazepine hydrobromide

Mol. formula : C₁₆H₁₇BrClNO₂

FW : 370.67

DEA schedule : 0

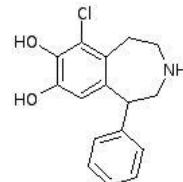
**Catalog number :** MEDD-005

Name : (\pm)-6-Chloro-7,8-dihydroxy-1-phenyl-2,3,4,5-tetrahydro[1H]-3-benzazepine hydrochloride

Mol. formula : C₁₆H₁₇Cl₂NO₂

FW : 326.22

DEA schedule : 0

**Catalog number :** NOCD-135**CASRN :** 34233-69-7

Name : Clozapine N-oxide

Mol. formula : C₁₈H₁₉ClN₄O

FW : 342.82

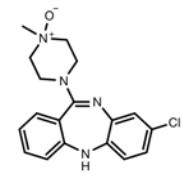
DEA schedule : 0

Notes : Clozapine metabolite.

References : Jann, MW; et al., *Clin Pharmacokinet* 1993, 24, 161-76.

Nawaratne, V; et al., *Mol Pharmacol* 2008, 74, 1119-31.

Becnel, J; et al., *Cell Rep* 2013, 4, 1049-59.

**Catalog number :** NOCD-047**CASRN :** 162408-66-4

Name : GR103691

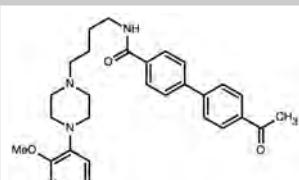
Mol. formula : C₃₀H₃₇N₃O₃

FW : 566.57

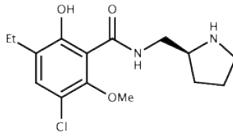
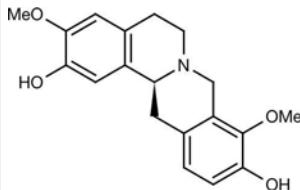
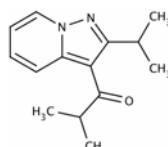
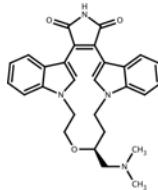
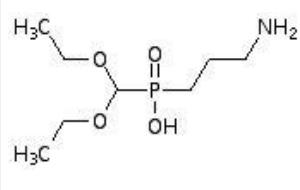
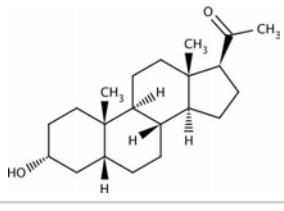
DEA schedule : 0

Notes : Selective dopamine D3 receptor antagonist

References : Audinot, V; et al. *J Pharmacol Exp Ther* 1998, 287, 187-97.



10 – Miscellaneous Compounds

Catalog number : NOCD-120	CASRN : 101536-82-7
Name : (+)-(S)-Noreticlopride HCl	
Mol. formula : C ₁₅ H ₂₂ Cl ₂ N ₂ O ₃	FW : 349.26 DEA schedule : 0
Notes : Dopamine D ₂ receptor antagonist.	
References : de Paulis, T; Hall H; Ogren SO, <i>European Journal of Medicinal Chemistry</i> 1985, 20, 273–276.	
Catalog number : NOCD-014	CASRN : 16562-13-3
Name : L-Stepholidine	
Mol. formula : C ₁₉ H ₂₁ NO ₄	FW : 327.37 DEA schedule : 0
Notes : D ₁ agonist / D ₂ antagonist	
References : Mo, J; et al. <i>Curr Med Chem</i> 2007, 14, 2996–3002. Wang, W; et al. <i>Neuropharmacology</i> 2007, 52, 355–61.	
Miscellaneous: Enzyme Inhibitors	
Catalog number : NOCD-104	CASRN : 50847-11-5
Name : Ibudilast	
Mol. formula : C ₁₄ H ₁₈ N ₂ O	FW : 230.31 DEA schedule : 0
Notes : Orally-available phosphodiesterase inhibitor. Induces cerebral vasodilation.	
References : (1) Nishino, K; et al., <i>Jpn J Pharmacol</i> 1983, 33, 267–78; (2) Souness, JE; et al., <i>Br J Pharmacol</i> 1994, 111, 1081–8.	
Catalog number : NOCD-127	CASRN : 169939-94-0
Name : Ruboxistaurin; LY-333,531	
Mol. formula : C ₂₈ H ₂₈ N ₄ O ₃	FW : 468.55 DEA schedule : 0
Notes : PKC β inhibitor.	
References : Gani O, Engh RA, <i>Nat Prod Rep</i> , 2010, 27, 489.	
Miscellaneous: GABA Receptor Related	
Catalog number : NOCD-045	CASRN : 123690-79-9
Name : (3-Aminopropyl)(diethoxymethyl)phosphinic acid; CGP-35348	
Mol. formula : C ₈ H ₂₀ NO ₄ P	FW : 225.23 DEA schedule : 0
Notes : GABAB receptor antagonist.	
References : Olpe, HR; et al. <i>Eur J Pharmacol</i> 1990, 187, 27–38.	
Catalog number : NOCD-102	CASRN : 128-20-1
Name : Pregnanolone	
Mol. formula : C ₂₁ H ₃₄ O ₂	FW : 318.49 DEA schedule : 0
Notes : Progestrone metabolite and barbiturate-like modulator of GABA _A receptors.	
References : Quinton, MS; et al. <i>Pharmacol Biochem Behav</i> 2006, 85, 385–92. Kaminski, RM; et al. <i>Eur J Pharmacol</i> 2003, 474, 217–22. Leskiewicz, M; et al. <i>Pol J Pharmacol</i> 2003, 55, 1131–6.	

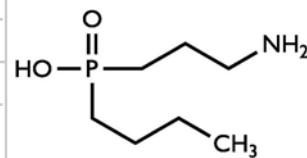
Catalog number : NOCD-103

CASRN : 145537-81-1

Name : SGS-742; CGP-36742

Mol. formula : C₇H₁₈NO₂P

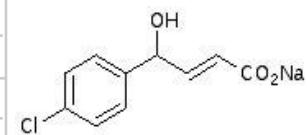
FW : 179.19 DEA schedule : 0

Notes : GABA_B receptor antagonist.References : Bullock, R *Curr Opin Investig Drugs* 2005, 6, 108–13.Froestl, W; et al. *Biochem Pharmacol* 2004, 68, 1479–87.**Miscellaneous: GHB Receptor Related**

Catalog number : NOCD-044

Name : *trans*-4-(4-Chlorophenyl)-4-hydroxy-2-butenoic acid sodium salt; NCS 356Mol. formula : C₁₀H₈ClO₃Na

FW : 234.61 DEA schedule : 0

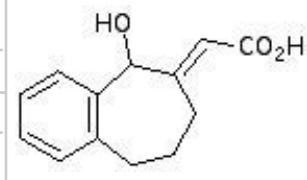
Notes : γ -Hydroxybutyrate receptor agonist.References : Gobaille, S; et al. *J Pharmacol Exp Ther* 1999, 290, 303–9.

Catalog number : NOCD-042

Name : 6,7,8,9-Tetrahydro-5-[H]benzocycloheptene-5-ol-4-ylideneacetic acid; NCS 382

Mol. formula : C₁₃H₁₄O₃

FW : 218.25 DEA schedule : 0

Notes : γ -Hydroxybutyrate receptor antagonist; anticonvulsant.References : Maitre, M; et al. *J Pharmacol Exp Ther* 1990, 255, 657–63.

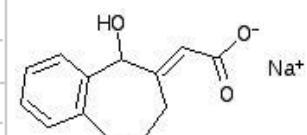
Catalog number : NOCD-043

CASRN : 131733-92-1

Name : 6,7,8,9-Tetrahydro-5-[H]benzocycloheptene-5-ol-4-ylideneacetic acid, sodium salt; NCS 382 sodium salt

Mol. formula : C₁₃H₁₃O₃Na

FW : 240.24 DEA schedule : 0

Notes : γ -Hydroxybutyrate receptor antagonist; anticonvulsant.References : Maitre, M; et al. *J Pharmacol Exp Ther* 1990, 255, 657–63.**Miscellaneous: Glutamate Receptor Related**

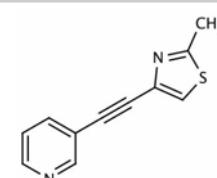
Catalog number : MEDD-026

CASRN : 329205-68-7

Name : 3-[(2-Methyl-4-thiazolyl)ethynyl]pyridine; MTEP

Mol. formula : C₁₁H₈N₂S

FW : 200.26 DEA schedule : 0

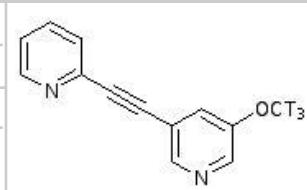
References : Bradbury, MJ; et al. *J Pharmacol Exp Ther* 2005, 313, 395–402.Kłodzinska, A; et al. *Neuropharmacology* 2004, 47, 342–50.Busse, CS; et al. *Neuropsychopharmacology* 2004, 29, 1971–9.

Catalog number : NOCD-058

Name : 3-[³H]Methoxy-5-(pyridin-2-yl-ethynyl)pyridineMol. formula : C₁₃H₁₀N₂O

FW : 216.26 DEA schedule : 0

Notes : Tritium-labeled MPEP analog



10 – Miscellaneous Compounds

Catalog number : MEDD-018

CASRN : 219911-35-0

Name : 6-Methyl-2-(phenylethynyl)pyridine hydrochloride; MPEP

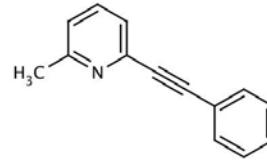
Mol. formula : C₁₄H₁₂ClN

FW : 229.71 DEA schedule : 0

Notes : *mGlu5 antagonist*

References : Gasparini, F; et al. *Neuropharmacology* 1999, 38, 1493–503.

Alagille, D; et al. *Bioorg Med Chem* 2005, 13, 197–209.



Catalog number : NOCD-033

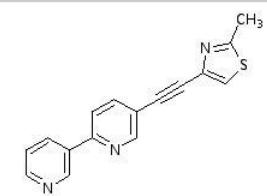
CASRN : 329204-25-3

Name : 5-[(2-Methyl-4-thiazolyl)ethynyl]-2,3'-bipyridine; MTEb

Mol. formula : C₁₆H₁₁N₃S

FW : 277.34 DEA schedule : 0

References : Roppe, JR; et al. *Bioorg Med Chem Lett* 2004, 14, 3993–3996.



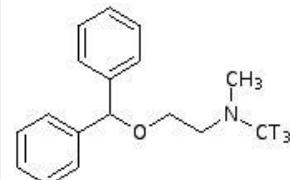
Miscellaneous: Histamine Receptor Related

Catalog number : NOCD-063

Name : [N-C³H₃]Diphenhydramine

Mol. formula : C₁₇H₂₁NO

FW : 261.38 DEA schedule : 0



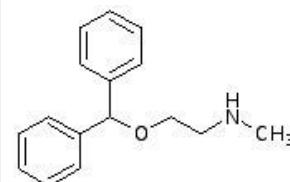
Catalog number : NOCD-060

Name : Nordiphenhydramine hydrochloride

Mol. formula : C₁₆H₂₀ClNO

FW : 277.80 DEA schedule : 0

Notes : Diphenhydramine metabolite



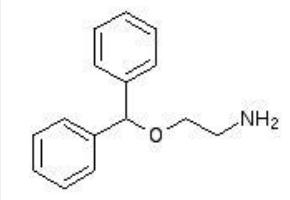
Catalog number : NOCD-061

Name : Dinordiphenhydramine hydrochloride

Mol. formula : C₁₅H₁₈ClNO

FW : 263.76 DEA schedule : 0

Notes : Diphenhydramine metabolite



Miscellaneous: NMDA Receptor Related

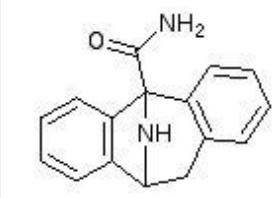
Catalog number : MEDD-003

Name : (±)-5-(Aminocarbonyl)-10,11-dihydro-5H-dibenzo[a,d]cyclohepten-5,10-imine hydrochloride; ADCI HCl

Mol. formula : C₁₆H₁₅ClN₂O

FW : 286.76 DEA schedule : 0

Notes : NMDA channel blocker



Catalog number : NOCD-019

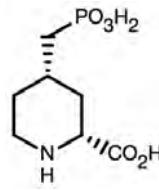
CASRN : 110347-85-8

Name : CGS 19755; Selfotel

Mol. formula : C₇H₁₄NO₅P

FW : 223.16 DEA schedule : 0

Notes : NMDA receptor antagonist.

References : Hutchison, AJ; et al. *J Med Chem* 1989, 32, 2171-8.

Catalog number : NOCD-010

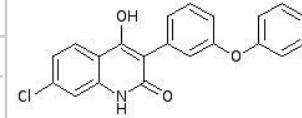
CASRN : 142326-59-8

Name : 7-Chloro-4-hydroxy-3-(3-phenoxyphenyl)-2(1H)quinoline; L 701,324

Mol. formula : C₂₁H₁₄ClNO₃

FW : 363.79 DEA schedule : 0

Notes : Glycine/NMDA receptor antagonist.

References : Bristow, LJ; et al. *Psychopharmacology (Berl)* 1995, 118, 230-2.
Bristow, LJ; et al. *J Pharmacol Exp Ther* 1996, 279, 492-501.**Miscellaneous: Orexin Receptor Related**

Catalog number : NOCD-111

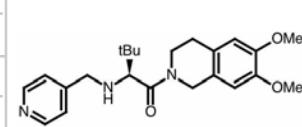
CASRN : 372523-75-6

Name : Cp-5

Mol. formula : C₂₃H₃₃ClN₃O₃

FW : 470.44 DEA schedule : 0

Notes : Non-peptidic orexin-2 receptor selective antagonist.

References : Hirose, M, et al., *Bioorg Med Chem Letters*, 2003, 13, 4497-4499.

Catalog number : NOCD-006

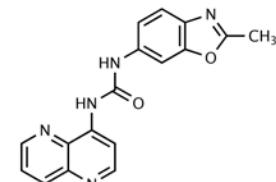
CASRN : 249889-64-3

Name : SB-334867 (base)

Mol. formula : C₁₇H₁₃N₅O₂

FW : 392.24 DEA schedule : 0

Notes : Orexin (hypocretin) antagonist

References : Coe, JW; et al., *Bioorg Med Chem Lett*, 2005, 15, 4889-4897.
McElhinny, CJ, Jr.; et al., *Bioorg Med Chem Lett*, 2012, 22, 6661-6664.**Miscellaneous: Serotonergic**

Catalog number : 1670-002

CASRN : 3616-78-2

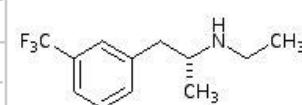
Name : (-)-Fenfluramine hydrochloride

Mol. formula : C₁₂H₁₇ClF₃N

FW : 267.72 DEA schedule : 4

Notes : CNS stimulant

References : Merck Index, 14th ed., Monograph 3973.



Catalog number : 1670-001

CASRN : 3239-45-0

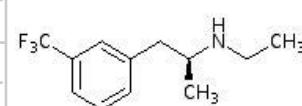
Name : (+)-Fenfluramine hydrochloride

Mol. formula : C₁₂H₁₇ClF₃N

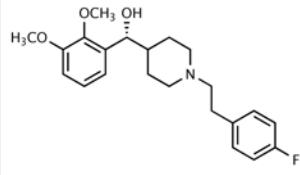
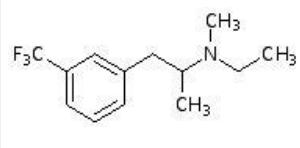
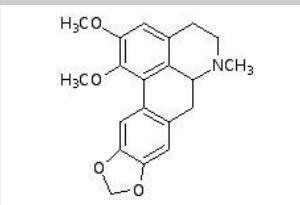
FW : 267.72 DEA schedule : 4

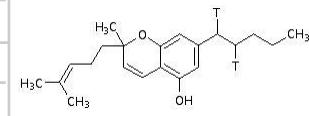
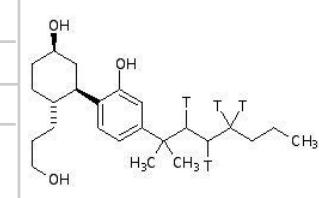
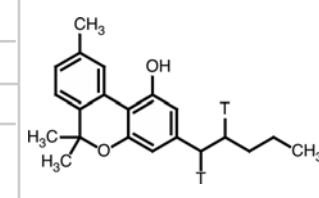
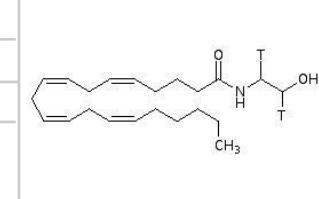
Notes : CNS stimulant; serotonin releaser

References : Merck Index, 14th ed., Monograph 3973.



10 – Miscellaneous Compounds

Catalog number : NOCD-015	CASRN : 139290-65-6
Name : MDL-100907; Volinanserin	
Mol. formula : C ₂₂ H ₂₉ ClFNO ₃	FW : 418.94 DEA schedule : 0
Notes : 5-HT _{2A} antagonist	
References : Ullrich, T; Rice KCBioorg Med Chem 2000, 8, 2427-32.	
Catalog number : 1670-003	
Name : (±)-N-Methylfenfluramine hydrochloride	
Mol. formula : C ₁₃ H ₁₉ ClF ₃ N	FW : 281.75 DEA schedule : 0
Notes : CNS stimulant	
Catalog number : NOCD-048	CASRN : 2565-01-7
Name : Nantenine	
Mol. formula : C ₂₀ H ₂₁ NO ₄	FW : 339.39 DEA schedule : 0
Notes : Serotonergic receptor antagonist.	
References : Fantegrossi, WE; et al. Psychopharmacology (Berl) 2004, 173, 270-7.	

Cannabinoids: Cannabichromene Class**Catalog number :** 7360-008**CASRN :** 20675-51-8 (parent cmpd)**Name :** [1',2'-³H₂]Cannabichromene**Mol. formula :** C₂₁H₃₀O₂**FW :** 314.46**DEA schedule :** 1**Notes :** Non-psychoactive constituent of cannabis (tritium-labeled).**References :** Instrumental Data for Drug Analysis, 2nd Ed., 1996, Volume 1, p304.**Cannabinoids: Cannabicyclohexanol Class****Catalog number :** NOCD-092**CASRN :** 119095-48-6**Name :** [2,3,4,4-³H₄](-)-CP 55,940**Mol. formula :** C₂₄H₄₀O₃**FW :** 376.58**DEA schedule :** 0**Notes :** High affinity CB1 and CB2 receptor radioligand (tritium-labeled).**Cannabinoids: Cannabinol Class****Catalog number :** 7360-023**Name :** [1',2'-³H₂]Cannabinol**Mol. formula :** C₂₁H₂₆O₂**FW :** 310.43**DEA schedule :** 1**Cannabinoids: Fatty Acid Derivatives (Anandamides)****Catalog number :** NOCD-008**Name :** Arachidonyl[1,2-³H]ethanolamide; Tritiated Anandamide**Mol. formula :** C₂₂H₃₇NO₂**FW :** 347.54**DEA schedule :** 0

11 – Tritium Labeled

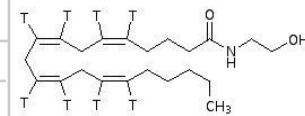
Catalog number : NOCD-078

Name : Tritium-labeled Arachidonylethanamide; Tritiated Anandamide

Mol. formula : C₂₂H₃₇NO₂

FW : 347.54 **DEA schedule :** 0

Notes : *Cannabinoid CB1 and CB2 receptor radioligand.*



Cannabinoids: Fatty Acid Derivatives (Arachidonyl esters)

Catalog number : NOCD-018

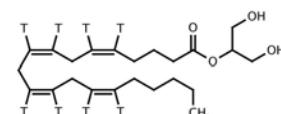
Name : [³H]-2-Arachidonylglycerol; [³H]-2-AG

Mol. formula : C₂₃H₃₈O₄

FW : 378.55 **DEA schedule :** 0

Notes : *Cannabinoid CB1 receptor agonist (tritium-labeled).*

References : Stella, N; Schweitzer, P; Piomelli, D *Nature* 1997, 388, 773–8.



Catalog number : NOCD-035

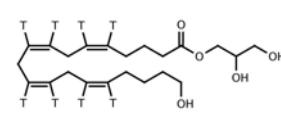
Name : Tritium-labeled 1-Arachidonylglycerol; [³H]-1-AG

Mol. formula : C₂₃H₃₈O₄

FW : 378.55 **DEA schedule :** 0

Notes : *Cannabinoid CB1 receptor agonist (tritium-labeled).*

References : Stella, N; Schweitzer, P; Piomelli D *Nature* 1997, 388, 773–8.



Cannabinoids: Pyrazole Class

Catalog number : NOCD-083

CASRN : 170937-38-9

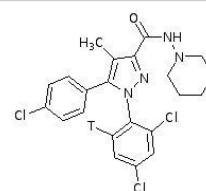
Name : [2,4-Dichlorophenyl-6-³H]SR141716

Mol. formula : C₂₂H₂₁Cl₃N₄O

FW : 465.80 **DEA schedule :** 0

Notes : *Cannabinoid CB1 receptor radioligand (tritium-labeled).*

References : Seltzman, H; et al. *J Chem Soc, Chem Commun* 1995, 1549–1550.



Catalog number : NOCD-086

CASRN : 475471-24-0

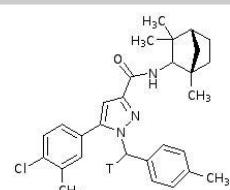
Name : Tritium-labeled SR144528

Mol. formula : C₂₉H₃₄N₃OCl

FW : 478.05 **DEA schedule :** 0

Notes : *Cannabinoid CB2 receptor radioligand (tritium-labeled).*

References : Portier, M; et al. *J Pharmacol Exp Ther* 1999, 288, 582–9.
Rinaldi-Carmona, M; et al. *J Pharmacol Exp Ther* 1998, 284, 644–50.



Cannabinoids: Tetrahydrocannabinol Class

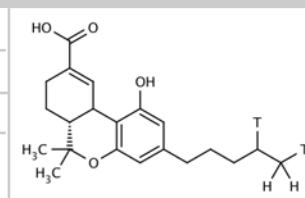
Catalog number : 7370-017

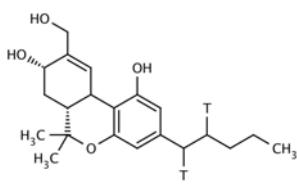
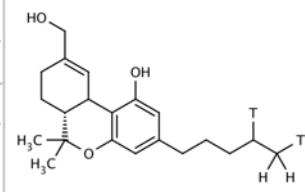
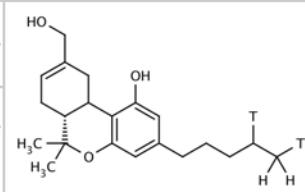
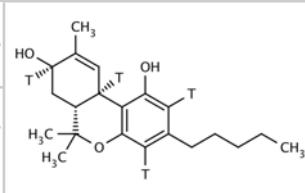
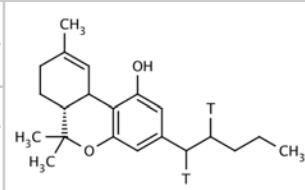
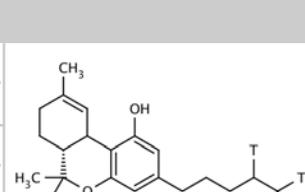
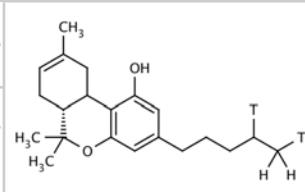
Name : [4',5'-³H₂]9-Carboxy-11-nor- Δ^9 -THC

Mol. formula : C₂₁H₂₈O₄

FW : 348.46 **DEA schedule :** 1

Notes : *Urinary metabolite of THC (tritium-labeled).*



Catalog number : 7370-026**Name :** [1',2'-³H₂]-8 α ,11-Dihydroxy- Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₄**FW :** 350.48 **DEA schedule :** 1**Catalog number :** 7370-024**Name :** [4',5'-³H₂]-11-Hydroxy- Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.48 **DEA schedule :** 1**Catalog number :** 7370-035**Name :** [4',5'-³H₂]-11-Hydroxy- Δ^8 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.48 **DEA schedule :** 1**Catalog number :** 7370-021**Name :** [2,4,8,10a-³H₄]-8 β -Hydroxy- Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₃**FW :** 338.49 **DEA schedule :** 1**Catalog number :** 7370-004**Name :** [1',2'-³H₂] Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.48 **DEA schedule :** 1**Notes :** Hallucinogen; psychotropic; analgesic (tritium-labeled).**Catalog number :** 7370-009**Name :** [4',5'-³H₂] Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.48 **DEA schedule :** 1**Catalog number :** 7370-027**Name :** [4',5'-³H₂] Δ^8 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.48 **DEA schedule :** 1**Notes :** Hallucinogen; psychotropic; analgesic (tritium-labeled).

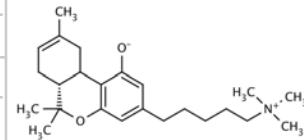
11 – Tritium Labeled

Catalog number : 7370-052

Name : [$^3\text{H}_3$]-5'-Trimethylammonium- Δ^8 -THC phenolate

Mol. formula : $\text{C}_{24}\text{H}_{37}\text{NO}_2$

FW : 371.56 DEA schedule : 1



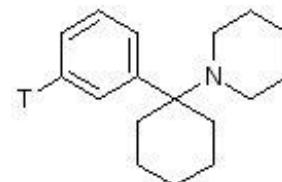
Dissociatives: Phencyclidine Class

Catalog number : 7471-001

Name : [Phenyl-3- $^3\text{H}(n)$]Phencyclidine; [Phenyl-3- $^3\text{H}(n)$]PCP

Mol. formula : $\text{C}_{17}\text{H}_{25}\text{N}$

FW : 245.40 DEA schedule : 2

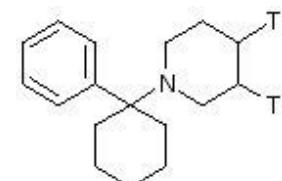


Catalog number : 7471-043

Name : [3',4'- ^3H]Phencyclidine; [3,4- ^3H]PCP

Mol. formula : $\text{C}_{17}\text{H}_{25}\text{N}$

FW : 247.4 DEA schedule : 2



Hallucinogens: Amphetamine Class

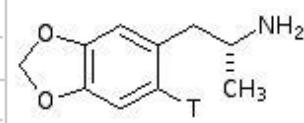
Catalog number : 7400-006

CASRN : 6292-91-7

Name : (-)-[6'- $^3\text{H}_2(n)$]-3',4'-Methylenedioxyamphetamine hydrochloride;
(-)-[6'- $^3\text{H}_2(n)$]MDA

Mol. formula : $\text{C}_{10}\text{H}_{14}\text{ClNO}_2$

FW : 217.68 DEA schedule : 1



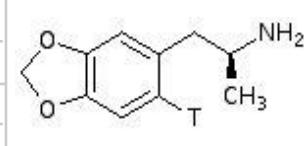
Catalog number : 7400-004

Name : (+)-[6'- $^3\text{H}_2(n)$]-3',4'-Methylenedioxyamphetamine hydrochloride;
(+)-[6'- $^3\text{H}_2(n)$]MDA

Mol. formula : $\text{C}_{10}\text{H}_{14}\text{ClNO}_2$

FW : 215.68 DEA schedule : 1

Notes : Hallucinogen (tritium-labeled).

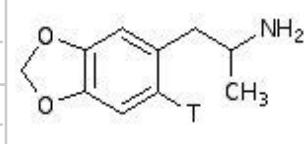


Catalog number : 7400-005

Name : [6- $^3\text{H}_2(n)$]-3,4-Methylenedioxyamphetamine hydrochloride;
[6- $^3\text{H}_2(n)$]MDA

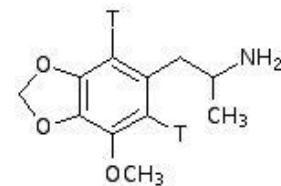
Mol. formula : $\text{C}_{10}\text{H}_{14}\text{ClNO}_2$

FW : 217.68 DEA schedule : 1

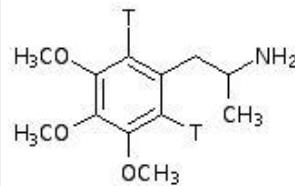


Catalog number : 7401-002

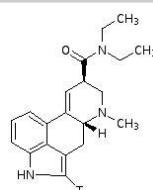
Name : [2',6'-³H(n)]-3-Methoxy-4,5-methylenedioxymethamphetamine hydrochloride; [2,6-³H(n)]MMDA

Mol. formula : C₁₁H₁₆ClNO₃**FW :** 249.72 **DEA schedule :** 1**Catalog number :** 7390-002

Name : (±)-[2,6-³H₂(n)]-3,4,5-Trimethoxyamphetamine hydrochloride

Mol. formula : C₁₂H₂₀ClNO₃**FW :** 265.76 **DEA schedule :** 1**Hallucinogens: Ergot alkaloids****Catalog number :** 7315-007**CASRN :** 377756-22-4

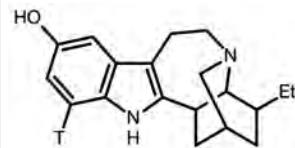
Name : [2-³H]-(+)-Lysergic acid diethylamide; (+)-[3H]-LSD

Mol. formula : C₂₀H₂₅N₃O**FW :** 323.42 **DEA schedule :** 1**Hallucinogens: Ibogaine Class****Catalog number :** 7260-005

Name : [12-³H]-Noribogaine

Mol. formula : C₁₉H₂₄N₂O**FW :** 296.41 **DEA schedule :** 0

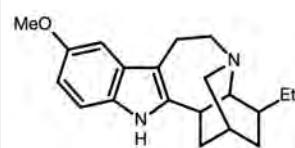
Notes : *Ibogaine-like effect without tremors (tritium-labeled).*

**Catalog number :** 7260-002**CASRN :** 146560-35-2

Name : Tritium-labeled Ibogaine; [12-³H]Ibogaine

Mol. formula : C₂₀H₂₆N₂O**FW :** 312.44 **DEA schedule :** 1

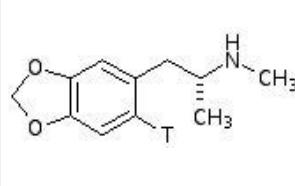
Notes : *Hallucinogen (tritium-labeled).*

**Hallucinogens: Methamphetamine Class****Catalog number :** 7405-005**CASRN :** 4764-17-4

Name : (–)[6'-³H(n)]-3',4'-Methylenedioxymethamphetamine hydrochloride; (–)[³H]MDMA

Mol. formula : C₁₁H₁₆ClNO₂**FW :** 231.71 **DEA schedule :** 1

Notes : *CNS stimulant; hallucinogen (tritium-labeled).*



11 – Tritium Labeled

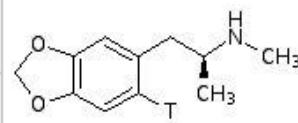
Catalog number : 7405-006

Name : (+)-[6'-³H(n)]-3',4'-Methylenedioxymethamphetamine hydrochloride;
(+)-[³H]MDMA

Mol. formula : C₁₁H₁₆ClNO₂

FW : 231.71 DEA schedule : 1

Notes : CNS stimulant; hallucinogen (tritium-labeled).



Miscellaneous: Glutamate Receptor Related

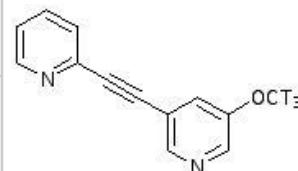
Catalog number : NOCD-058

Name : 3-[³H]Methoxy-5-(pyridin-2-yl-ethynyl)pyridine

Mol. formula : C₁₃H₁₀N₂O

FW : 216.26 DEA schedule : 0

Notes : Tritium-labeled MPEP analog



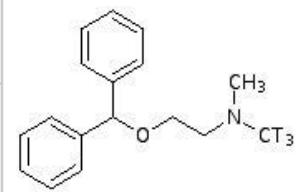
Miscellaneous: Histamine Receptor Related

Catalog number : NOCD-063

Name : [N-C³H₃]Diphenhydramine

Mol. formula : C₁₇H₂₁NO

FW : 261.38 DEA schedule : 0



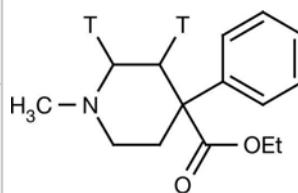
Opioids: Meperidine Class

Catalog number : 9230-003

Name : [2,3-³H₂]Meperidine hydrochloride

Mol. formula : C₁₅H₂₂ClNO₂

FW : 287.81 DEA schedule : 2



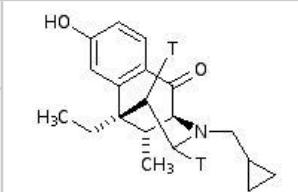
Opioids: Metazocine Class

Catalog number : 9240-003

Name : [11,12-³H₂](-)-Ethylketazocine

Mol. formula : C₁₉H₂₅NO₂

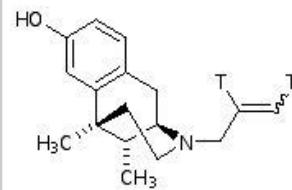
FW : 303.42 DEA schedule : 1



Catalog number : 9240-019

Name : (-)-[17,18-³H]N-AllylnormetazocineMol. formula : C₁₇H₂₃NO

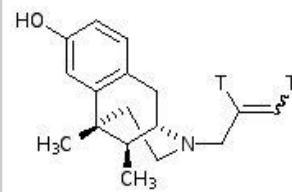
FW : 261.39 DEA schedule : 0



Catalog number : 9240-040

Name : (+)-[17,18-³H]N-AllylnormetazocineMol. formula : C₁₇H₂₃NO

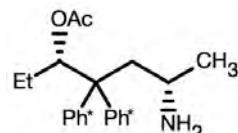
FW : 261.39 DEA schedule : 0

**Opioids: Methadone Class**

Catalog number : 9633-011

Name : (-)-[o,o'-³H₂(n)]-α-Acetyl-N,N-dinormethadol hydrochlorideMol. formula : C₂₁H₂₈CINO₂

FW : 389.97 DEA schedule : 0

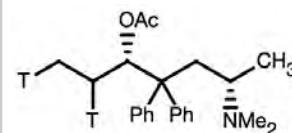


Ph* = o-T-phenyl

Catalog number : 9648-012

Name : (-)-[1,2-³H₂]α-AcetylmethadolMol. formula : C₂₃H₃₂CINO₂

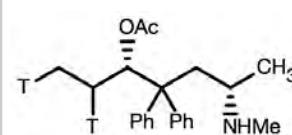
FW : 353.50 DEA schedule : 2



Catalog number : 9633-006

Name : (-)-[1,2-³H]α-Acetylnormethadol hydrochlorideMol. formula : C₂₃H₃₂CINO₂

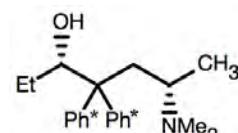
FW : 375.94 DEA schedule : 1



Catalog number : 9605-003

Name : (-)-[o,o'-³H₂(n)]α-Methadol hydrochlorideMol. formula : C₂₁H₂₉NO

FW : 347.93 DEA schedule : 1



Ph* = o-T-Phenyl

References : Portoghese, PS; Williams DA *J Med Chem* 1969, 12, 839-44.

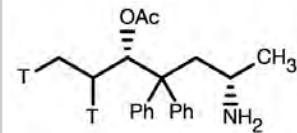
11 – Tritium Labeled

Catalog number : 9633-015

Name : (-)-[1,2-³H] α -Acetyl-N,N-dinormethadol hydrochloride

Mol. formula : C₂₁H₂₈ClNO₂

FW : 361.9-2 DEA schedule : 0

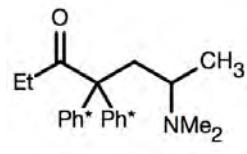


Catalog number : 9250-006

Name : (\pm)-[o,o'-³H₂(n)]Methadone

Mol. formula : C₂₁H₂₇NO

FW : 313.46 DEA schedule : 2



Ph* = o-T-phenyl

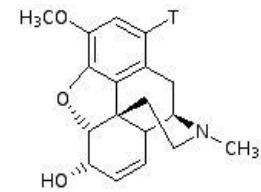
Opioids: Morphine Class

Catalog number : 9050-011

Name : [1-³H]Codeine

Mol. formula : C₁₈H₂₁NO₃

FW : 301.37 DEA schedule : 2

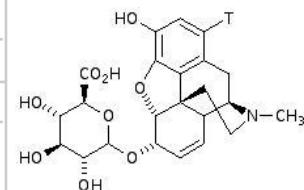


Catalog number : 9300-014

Name : [1-³H]Morphine-6-glucuronide

Mol. formula : C₂₃H₂₇NO₉

FW : 463.47 DEA schedule : 2

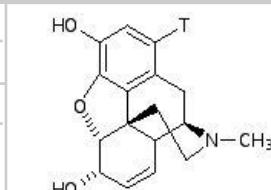


Catalog number : 9300-002

Name : Tritium-labeled Morphine sulfate

Mol. formula : C₁₇H₁₉NO₃

FW : 475.50 DEA schedule : 2



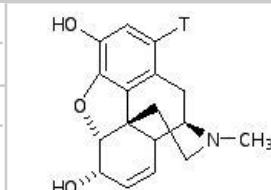
Catalog number : 9300-006

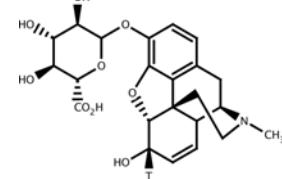
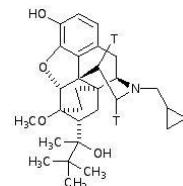
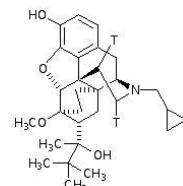
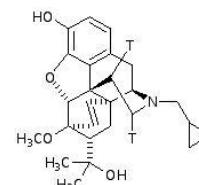
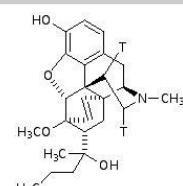
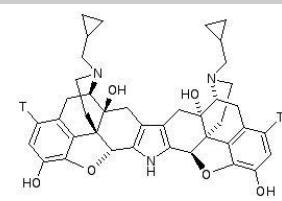
CASRN : 80573-75-7

Name : [1-³H(n)]Morphine

Mol. formula : C₁₇H₁₉NO₃

FW : 287.35 DEA schedule : 2



Catalog number : 9300-011**Name :** Morphine-(6-³H)-3-glucuronide**Mol. formula :** C₂₃H₂₇NO₉**FW :** 463.47**DEA schedule :** 2**Opioids: Orvinol Class****Catalog number :** 9064-002**CASRN :** 161772-95-8**Name :** [15,16-³H₂]Buprenorphine hydrochloride**Mol. formula :** C₂₆H₃₈ClNO₄**FW :** 508.11**DEA schedule :** 5**Notes :** Narcotic analgesic (tritium-labeled).**References :** Robinson, SE *CNS Drug Rev* 2002, 8, 377-90.**Catalog number :** 9064-003**CASRN :** 161772-95-8**Name :** [15,16-³H₂]Buprenorphine**Mol. formula :** C₂₉H₄₁NO₄**FW :** 508.11**DEA schedule :** 3**Notes :** Narcotic analgesic (tritium-labeled).**References :** Robinson, SE *CNS Drug Rev* 2002, 8, 377-90.**Catalog number :** 9058-002**Name :** [15,16-³H₂]Diprenorphine**Mol. formula :** C₂₆H₃₆ClNO₄**FW :** 429.58**DEA schedule :** 2**Notes :** (see footnotes 1 & 2 in Section C)**Catalog number :** 9056-001**Name :** [15,16-³H₂]Etorphine**Mol. formula :** C₂₅H₃₃NO₄**FW :** 429.58**DEA schedule :** 1**Notes :** (see footnotes 1 & 2 in Section C)**Opioids: Oxymorphone Class****Catalog number :** NOCD-084**Name :** [1,1'-³H(n)]Norbinaltorphimine; [³H]norBNI**Mol. formula :** C₄₀H₄₃N₃O₆**FW :** 665.80**DEA schedule :** 0**References :** Birch, PJ; et al. *Eur J Pharmacol* 1987, 144, 405-8.
Portoghese, PS; Lipkowski, AW; Takemori, AE *Life Sci* 1987, 40, 1287-92.

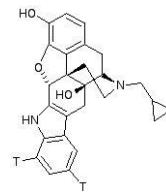
11 – Tritium Labeled

Catalog number : 9652-061

Name : [5',7'-³H₂]Naltrindole

Mol. formula : C₂₆H₂₆N₂O₃

FW : 418.51 DEA schedule : 0

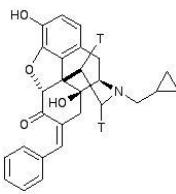


Catalog number : 9652-068

Name : [15,16-³H]-7-Benzylidine-7-dehydronaltrexone; [³H]BNTX

Mol. formula : C₂₇H₂₈ClNO₄

FW : 433.52 DEA schedule : 0

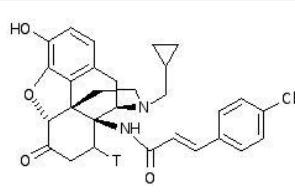


Catalog number : NOCD-079

Name : Tritium-labeled Clo cinnamonox

Mol. formula : C₂₉H₂₉ClN₂O₄

FW : 507.01 DEA schedule : 0



References : Comer, SD; Burke, TF; Lewis, JW; Woods, JH *J Pharmacol Exp Ther* **1992**, 262, 1051-6.

Catalog number : 9333-012

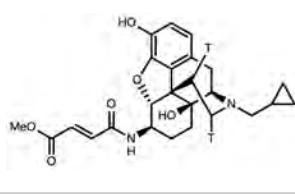
Name : [15,16-³H₂]-6β-Funaltrexamine

Mol. formula : C₂₆H₃₃ClN₂O₆

FW : 458.53 DEA schedule : 2

Notes : Irreversible μ -opioid receptor antagonist (tritium-labeled).

References : Portoghese, PS; el Kouhen, R; Law, PY; Loh, HH; Le Bourdonnec, B *Farmaco* **2001**, 56, 191-6.

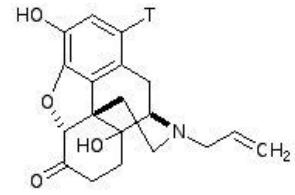


Catalog number : 9652-031

Name : (-)-[1-³H(n)]Naloxone

Mol. formula : C₁₉H₂₁NO₄

FW : 329.38 DEA schedule : 0

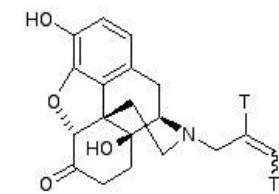


Catalog number : 9652-035

Name : (-)-[19,20-³H₂]Naloxone

Mol. formula : C₁₉H₂₁NO₄

FW : 331.39 DEA schedule : 0

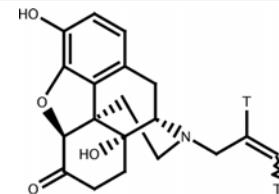


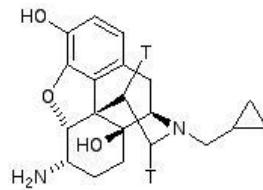
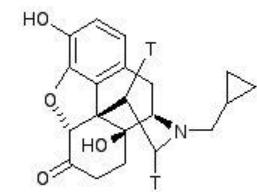
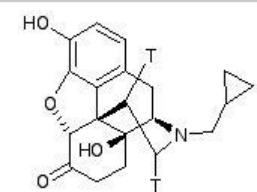
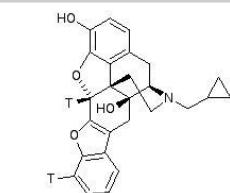
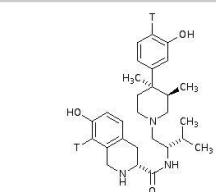
Catalog number : 9652-039

Name : (+)-[19,20-³H₂]Naloxone

Mol. formula : C₁₉H₂₁NO₄

FW : 331.39 DEA schedule : 0



Catalog number : 9333-009**Name :** [15,16-³H₂]-6β-Naltrexamine**Mol. formula :** C₂₀H₂₆N₂O₃**FW :** 346.45 **DEA schedule :** 2**Catalog number :** 9652-012**Name :** [15,16-³H₂]Naltrexone**Mol. formula :** C₂₀H₂₃NO₄**FW :** 345.42 **DEA schedule :** 0**Catalog number :** 9652-013**Name :** [15,16-³H₂]Naltrexone hydrochloride**Mol. formula :** C₂₀H₂₄ClNO₄**FW :** 381.88 **DEA schedule :** 0**Catalog number :** 9652-065**Name :** Tritium-labeled Naltriben**Mol. formula :** C₂₆H₂₅NO₄**FW :** 419.50 **DEA schedule :** 0**Notes :** Selective δ-opioid receptor antagonist (tritium-labeled).**References :** Sofuooglu, M; Portoghese, PS; Takemori, AE *J Pharmacol Exp Ther* 1991, 257, 676–80.**Opioids: Phenylpiperidine Class****Catalog number :** NOCD-059**Name :** [³H]DTic**Mol. formula :** C₂₈H₃₉N₃O₃**FW :** 469.64 **DEA schedule :** 0**Notes :** Selective kappa-opioid antagonist (tritium-labeled).**References :** Thomas, JB; et al. *J Med Chem* 2001, 44, 2687–90.**Peptides: Cannabinoid-related****Catalog number :** MPSP-096**Name :** [³H₂]Hemopressin**Mol. formula :** C₅₃H₇₇N₁₃O₁₂**FW :** 1092.3**Notes :** Radiolabeled hemopressin (MPSP-90; PEPT-053).

H-Pro-Val-Asn-[³H]Phe-Lys-[³H]Phe-Leu-Ser-His-OH

Catalog number : MPSP-091**Name :** [³H]Hemopressin**Mol. formula :** C₅₃H₇₇N₁₃O₁₂**FW :** 1096.3**Notes :** *Radiolabeled hemopressin (MPSP-90; PEPT-053).*H-[3H]Pro-Val-Asn-[3H]
Phe-Lys-[3H]Phe-Leu-Ser-
His-OH**Catalog number :** MPSP-097**Name :** [³H]Leu⁶-Hemopressin**Mol. formula :** C₅₀H₇₉N₁₃O₁₂**FW :** 1056.3**Notes :** *Radiolabeled Leu⁶-Hemopressin (PEPT-054).*H-Pro-Val-Asn-[3H]Phe-
Lys-Leu-Leu-Ser-His-OH**Catalog number :** MPSP-093**Name :** [³H]Tyr⁷-Hemopressin (1-6)**Mol. formula :** C₄₇H₆₁N₉O₁₀₃H₂**FW :** 918.1**Notes :** *Radioligand for MPSP-092.*H-Pro-Val-Asn-Phe-Lys-
Phe-[3H2]Tyr-OH**Peptides: Opioid****Catalog number :** MPSP-014**Name :** [³H]CTAP**Mol. formula :** C₅₁H₆₈N₁₃O₁₁S₂₃H**FW :** 1105.0**Notes :** *Radioactive ligand for MPSP-013***References :** Abbruscato, TJ; et al. *J Pharmacol Exp Ther* 1997, 280, 402-9.H-d-Phe[3H]-c[Cys-Tyr-d-
Trp-Arg-Thr-Pen]-Thr-
NH2**Catalog number :** MPSP-007**Name :** [³H₂]DADLE**Mol. formula :** C₂₉H₃₇N₅O₇₃H₂**FW :** 573.7**Notes :** *Tritiated DADLE***References :** Luciano, MG; et al. *Brain Res Bull* 1981, 7, 677-82.H-Tyr[3H2]-d-Ala-Gly-
Phe-d-Leu-OH**Catalog number :** MPSP-035**Name :** [³H₂]DALCE**Mol. formula :** C₃₃H₄₄N₆O₁₀S₃H₂**FW :** 677.0**Notes :** *Radioactive ligand for MPSP-034*H-Tyr[3H2]-d-Ala-Gly-
Phe-Leu-Cys-OH

Catalog number : MPSP-028**Name :** [³H₂]DALDA**Mol. formula :** C₃₀H₄₃N₉O₅₃H₂**FW :** 615.8**Notes :** Radioactive ligand for MPSP-027

H-Tyr[3H2]-d-Arg-Phe-Lys-NH2

Catalog number : MPSP-012**Name :** [³H₂]DAMGO**Mol. formula :** C₂₆H₃₃N₅O₆₃H₂**FW :** 517.7**Notes :** Radioactive ligand for MPSP-011**References :** Zajac, JM; Roques BP *Life Sci* 1983, 33 Suppl 1, 155–8.H-Tyr[3H2]-d-Ala-Gly-N α -Me-Phe-Gly-ol**Catalog number :** MPSP-031**Name :** [³H₂]Deltorphin I**Mol. formula :** C₃₇H₅₀N₈O₁₀₃H₂**FW :** 772.9**Notes :** Radioactive ligand for MPSP-030**References :** Erspamer, V; et al. *Proc Natl Acad Sci USA* 1989, 86, 5188.

H-Tyr[3H2]-d-Ala-Phe-Asp-Val-Val-Gly-NH2

Catalog number : MPSP-037**Name :** [³H₂]Deltorphin II**Mol. formula :** C₃₈H₅₂N₈O₁₀₃H₂**FW :** 786.9**Notes :** Radioactive ligand for MPSP-036**References :** Erspamer, V; et al. *Proc Natl Acad Sci USA* 1989, 86, 5188.

H-Tyr[3H2]-d-Ala-Phe-Glu-Val-Val-Gly-NH2

Catalog number : MPSP-002**Name :** [³H₂]DPDPE**Mol. formula :** C₃₀H₃₇N₇O₅S₂₃H₂**FW :** 649.8**Notes :** Radioactive ligand for MPSP-001**References :** Cotton, R; et al. *Eur J Pharmacol* 1984, 97, 331–2.

H-Tyr[3H2]-c[d-Pen-Gly-Phe-d-Pen]-OH

Catalog number : MPSP-045**Name :** [³H₂]p-Cl-Phe-DPDPE**Mol. formula :** C₃₀H₃₆ClN₅O₇S₂₃H₂**FW :** 684.4**Notes :** Radioactive ligand for MPSP-044

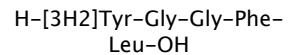
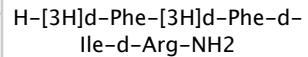
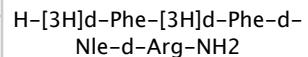
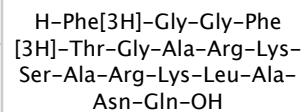
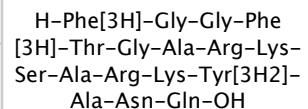
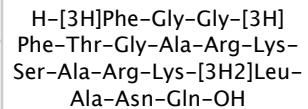
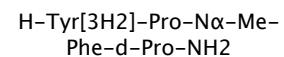
H-Gly-Gly-Phe[3H]-Leu-Arg-Arg-Ile-Arg-Pro-Lys-Leu-Lys-Trp-Asp-Asn-Gln-OH

Catalog number : MPSP-004**Name :** [³H₂]DSLET**Mol. formula :** C₃₃H₄₄N₆O₁₀₃H₂**FW :** 690.9**Notes :** Radioactive ligand for MPSP-003**References :** David, M; et al. *Eur J Pharmacol* 1982, 78, 385–7.

H-Tyr[3H2]-d-Ser-Gly-Phe-Leu-Thr-OH

11 – Tritium Labeled

Catalog number : MPSP-033	Name : [³ H ₂]DSTBULET	Mol. formula : C ₃₃ H ₄₄ N ₆ O ₁₀₃ H ₂	FW : 746.9	Notes : Radioactive ligand for MPSP-032	References : Delay-Goyet, P <i>NIDA Res Mono</i> 1986 , 75, 197-200.	H-Tyr[3H2]-d-Ser(But)-Gly-Phe-Leu-Thr-OH
Catalog number : MPSP-020	Name : [³ H ₂]Dynorphin (1-11)	Mol. formula : C ₆₃ H ₁₀₁ N ₂₁ O ₁₃₃ H ₂	FW : 1366.7	Notes : Radioactive ligand for MPSP-019	References : Chavkin, C; Goldstein, A <i>Proc Natl Acad Sci USA</i> 1981 , 78, 6543-7.	H-Tyr[3H2]-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-OH
Catalog number : MPSP-018	Name : [³ H ₂]Dynorphin (1-13) amide	Mol. formula : C ₇₅ H ₁₂₅ N ₂₅ O ₁₄₃ H ₂	FW : 1607.2	Notes : Radioactive ligand for MPSP-017		H-Tyr[3H2]-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-Pro-Lys-Leu-Lys-NH2
Catalog number : MPSP-022	Name : [³ H ₂]Dynorphin (1-9)	Mol. formula : C ₅₂ H ₈₂ N ₁₈ O ₁₁₃ H ₂	FW : 1141.5	Notes : Radioactive ligand for MPSP-021	References : Robson, LE; et al. <i>Life Sci</i> 1983 , 33 Suppl 1, 283-6. Wood, MS; Rodriguez FD; Traynor JR <i>Neuropharmacology</i> 1989 , 28, 1041-6.	H-Tyr[3H2]-Gly-Gly-Phe-Leu-Arg-Arg-Ile-Arg-OH
Catalog number : MPSP-041	Name : [³ H]Dynorphin (2-17)	Mol. formula : C ₉₀ H ₁₅₅ N ₃₀ O ₂₁₃ H ₂	FW : 1986.8	Notes : Radioactive ligand for MPSP-040		H-Gly-Gly-Phe[3H]-Leu-Arg-Arg-Ile-Arg-Pro-Lys-Leu-Lys-Trp-Asp-Asn-Gln-OH
Catalog number : MPSP-074	Name : [³ H ₂]Endomorphin-2	Mol. formula : C ₃₂ H ₃₅ N ₅ O ₅₃ H ₂	FW : 575.7	Notes : Radioactive ligand for MPSP-073	References : Spetea, M; et al. <i>Biochem Biophys Res Commun</i> 1998 , 250, 720-5.	H-Tyr[3H2]-Pro-Phe-Phe-NH2
Catalog number : MPSP-025	Name : [³ H ₂]β-Endorphin	Mol. formula : C ₁₅₈ H ₂₄₉ N ₃₉ O ₄₆ S ₃ H ₂	FW : 3469.6	Notes : Radioactive ligand for MPSP-024	References : Li, CH; et al. <i>Proc Natl Acad Sci USA</i> 1980 , 77, 2303-4.	H-Tyr-Gly-Gly-Phe-Met-Thr-Ser-Glu-Lys-Ser-Gln-Thr-Pro-Leu-Val-Thr-Leu-Phe-Lys-Asn-Ala-Ile-Ile-Lys-Asn-Ala-[3H2]Tyr-Lys-Lys-Gly-Glu-OH

Catalog number : MPSP-081**Name :** [³H₂]Leu-Enkephalin-acid**Mol. formula :** C₂₈H₃₇N₅O₇**FW :** 559.3**Notes :** Radioactive ligand for MPSP-080.**Catalog number :** MPSP-079**Name :** [³H₂]Kaffiralin-1**Mol. formula :** C₃₀H₄₂N₈O₄₃H₂**FW :** 584.7**Notes :** Radioactive ligand for MPSP-076**References :** Dooley, CT; et al. *J Biol Chem* 1998, 273, 18848–56.**Catalog number :** MPSP-077**Name :** [³H₂]Kaffiralin-2**Mol. formula :** C₃₀H₄₂N₈O₄₃H₂**FW :** 584.7**Notes :** Radioactive ligand for MPSP-076.**Catalog number :** MPSP-059**Name :** [³H₂]Orphanin FQ; [³H₂]Nociceptin**Mol. formula :** C₇₉H₁₂₇N₂₇O₂₂₃H₂**FW :** 1813.4**Notes :** Radioactive ligand for MPSP-058**References :** Dooley, CT; Houghten RA *Life Sci* 1996, 59, PL23–9.**Catalog number :** MPSP-082**Name :** [³H₄]Tyr¹⁴Orphanin FQ**Mol. formula :** C₈₂H₁₂₃N₂₇O₂₃₃H₄**FW :** 1867.1**Notes :** Radioactive ligand for MPSP-058**References :** Dooley, CT; et al. *Life Science* 1996, 59, PL23–PL29.**Catalog number :** MPSP-095**Name :** [³H₄]Orphanin FQ**Mol. formula :** C₇₉H₁₂₅N₂₇O₂₂₃H₄**FW :** 1817.4**Notes :** Radioactive ligand for MPSP-058.**Catalog number :** MPSP-010**Name :** [³H₂]PLO17**Mol. formula :** C₂₉H₃₅N₅O₅₃H₂**FW :** 539.6**Notes :** Radioactive ligand for MPSP-009.**References :** Hawkins, KN; et al. *Eur J Pharmacol* 1987, 133, 351–2.

11 – Tritium Labeled

Catalog number : MPSP-039

Name : [$^3\text{H}_2$]Tipp-Enkephalin

Mol. formula : $\text{C}_{37}\text{H}_{36}\text{N}_4\text{O}_{63}\text{H}_2$

FW : 638.8

Notes : Radioactive ligand for MPSP-038

References : Nevin, ST; et al. *Life Sci* 1993, 53, PL57-62.

H-Tyr[$^3\text{H}_2$]-Tic-Phe-Phe-OH

Stimulants: Amphetamine Class

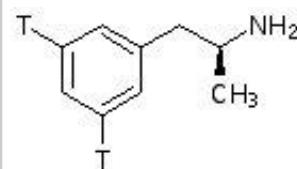
Catalog number : 1100-009

Name : (+)-(S)-[3,5- $^3\text{H}(n)$]Amphetamine

Mol. formula : $\text{C}_9\text{H}_{13}\text{N}$

FW : 139.22 DEA schedule : 2

Notes : CNS stimulant (tritium-labeled).

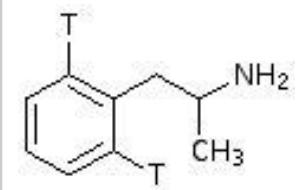


Catalog number : 1100-005

Name : (\pm)-[2',6'- $^3\text{H}_2$]Amphetamine; 2,6-Tritoamphetamine

Mol. formula : $\text{C}_9\text{H}_{13}\text{N}$

FW : 135.21 DEA schedule : 2



Stimulants: Methamphetamine Class

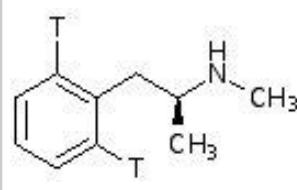
Catalog number : 1105-004

Name : (+)-[2',6'- $^3\text{H}(n)$]Methamphetamine hydrochloride

Mol. formula : $\text{C}_{10}\text{H}_{16}\text{ClN}$

FW : 185.69 DEA schedule : 2

Notes : CNS stimulant (tritium-labeled).



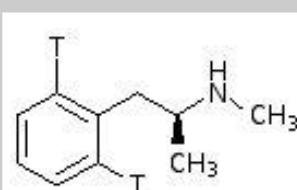
Catalog number : 1105-008

Name : (+)-[2,6- $^3\text{H}(n)$]Methamphetamine

Mol. formula : $\text{C}_{10}\text{H}_{15}\text{N}$

FW : 153.25 DEA schedule : 2

Notes : CNS stimulant (tritium-labeled).



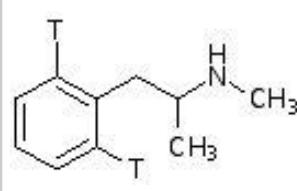
Catalog number : 1105-006

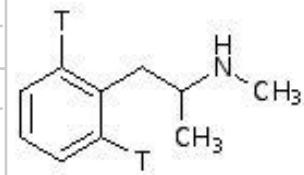
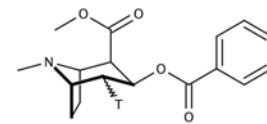
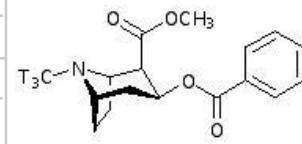
Name : (\pm)-[2',6'- $^3\text{H}(n)$]Methamphetamine hydrochloride

Mol. formula : $\text{C}_{10}\text{H}_{16}\text{ClN}$

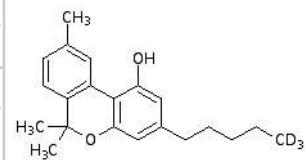
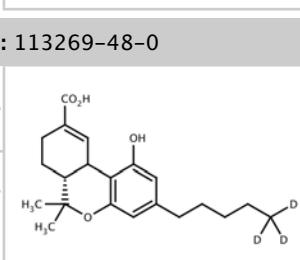
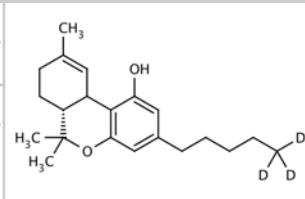
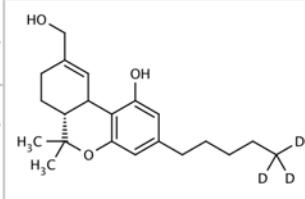
FW : 189.71 DEA schedule : 2

Notes : CNS stimulant (tritium-labeled).



Catalog number : 1105-007**Name :** (\pm) -[2,6- 3 H(n)]Methamphetamine**Mol. formula :** C₁₀H₁₅N**FW :** 153.25 **DEA schedule :** 2**Notes :** CNS stimulant (tritium-labeled).**Stimulants: Tropane Class****Catalog number :** 9041-004**CASRN :** 85438-94-4**Name :** $(-)$ -[4- 3 H]Cocaine**Mol. formula :** C₁₇H₂₁NO₄**FW :** 303.35 **DEA schedule :** 2**Catalog number :** 9041-003**Name :** $(-)$ -[N-C³H₃]Cocaine**Mol. formula :** C₁₇H₂₁NO₄**FW :** 309.38 **DEA schedule :** 2



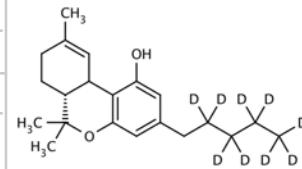
Cannabinoids: Tetrahydrocannabinol Class**Catalog number :** 7360-014**Name :** [5'-²H₃]Cannabinol**Mol. formula :** C₂₁H₂₆O₂**FW :** 310.43 **DEA schedule :** 1**Notes :** *Inactive constituent of cannabis (deuterium-labeled).***Catalog number :** 7370-038**Name :** [5'-²H₃]-11-nor-Δ⁸-THC-9-carboxylic acid**Mol. formula :** C₂₁H₂₈O₄**FW :** 347.46 **DEA schedule :** 1**Catalog number :** 7370-003**CASRN :** 113269-48-0**Name :** [5'-²H₃]9-Carboxy-11-nor-Δ⁹-THC**Mol. formula :** C₂₁H₂₈O₄**FW :** 347.46 **DEA schedule :** 1**Notes :** *Urinary metabolite of THC (deuterium-labeled).***Catalog number :** 7370-005**CASRN :** 81586-39-2**Name :** Deuterium-labeled Δ⁹-THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 317 **DEA schedule :** 1**Notes :** *Hallucinogen; psychotropic; analgesic (deuterium-labeled).***Catalog number :** 7370-010**CASRN :** 130410-26-3**Name :** [5'-²H₃]-11-Hydroxy-Δ⁹-THC**Mol. formula :** C₂₁H₃₀O₃**FW :** 333 **DEA schedule :** 1

Catalog number : 7370-025

Name : [2',2',3',3',4',4',5',5',5'-²H₉]Δ⁹-THCMol. formula : C₂₁H₃₀O₂

FW : 323.52 DEA schedule : 1

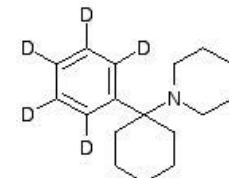
Notes : Hallucinogen; psychotropic; analgesic (deuterium-labeled).

**Dissociatives: Phencyclidine Class**

Catalog number : 7471-003

Name : [Phenyl-2,3,4,5,6-²H₅]Phencyclidine; [Phenyl-2,3,4,5,6-²H₅]PCPMol. formula : C₁₇H₂₅N

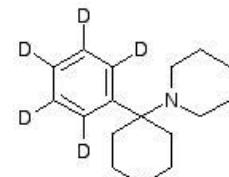
FW : 248.42 DEA schedule : 2



Catalog number : 7471-006

Name : [Phenyl-²H₅]Phencyclidine hydrochloride; [Phenyl-²H₅]PCP HClMol. formula : C₁₇H₂₆CIN

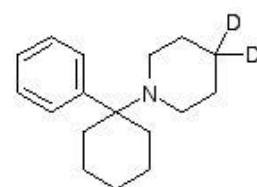
FW : 284.89 DEA schedule : 2



Catalog number : 7471-008

Name : [Piperidino-4,4-²H₂]Phencyclidine hydrochloride;
[Piperidino-4,4-²H₂]PCPMol. formula : C₁₇H₂₆CIN

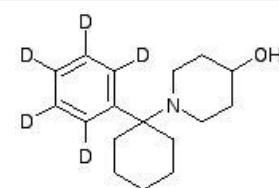
FW : 281.86 DEA schedule : 2



Catalog number : 7471-082

Name : [Phenyl-²H₅]-1-(1-Phenylcyclohexyl)-4-hydroxypiperidineMol. formula : C₁₇H₂₅NO

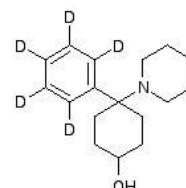
FW : 264.43 DEA schedule : 1

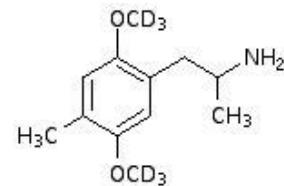
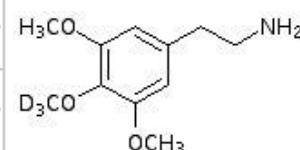
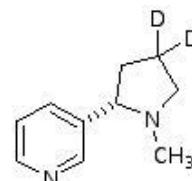
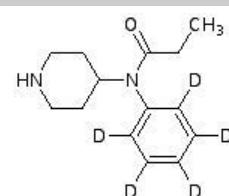
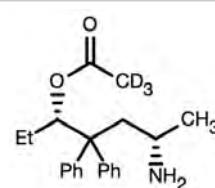


Catalog number : 7471-084

Name : [Phenyl-²H₅]-1-(1-Phenyl-4-hydroxycyclohexyl)piperidine (*cis/trans*)Mol. formula : C₁₇H₂₅NO

FW : 264.43 DEA schedule : 1

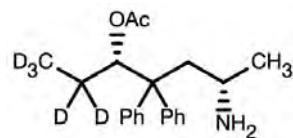


Hallucinogens: Amphetamine Class**Catalog number :** 7395-004**Name :** [OC²H₃]-2,5-Dimethoxy-4-methylamphetamine HCl; [OC²H₃]DOM**Mol. formula :** C₁₂H₂₀ClNO₂**FW :** 251.78 **DEA schedule :** 1**Hallucinogens: Phenethylamine Class****Catalog number :** 7381-002**Name :** [4'-OC²H₃]Mescaline hydrochloride**Mol. formula :** C₁₁H₁₈ClNO₃**FW :** 247.72 **DEA schedule :** 1**Nicotinics: Nicotine Class****Catalog number :** NICT-015**CASRN :** 121949-85-7**Name :** (–)-Nicotine-4,4-d₂ (–)-di-p-toluoyle-L-tartrate**Mol. formula :** C₃₀H₃₂N₂O₈**FW :** 550.60 **DEA schedule :** 0**References :** Jacob, PJ *Labelled Comp Radiopharm* **1988**, 25, 1117-28.**Opioids: Fentanyl Class****Catalog number :** 9801-008**CASRN :** 1211527-23-9**Name :** [Phenyl-²H₃]N-4-Piperidyl-N-phenylpropanamide**Mol. formula :** C₁₄H₂₀N₂O**FW :** 237.36 **DEA schedule :** 2**Notes :** Precursor for the synthesis of mass-labeled fentanyl analogs.**Opioids: Methadone Class****Catalog number :** 9633-012**Name :** (–)-[Acetyl-²H₃] α -Acetyl-N,N-dinormethadol hydrochloride**Mol. formula :** C₂₁H₂₈ClNO₂**FW :** 361.92 **DEA schedule :** 0

Catalog number : 9633-014

Name : (-)-[1,1,1,2,2,3-²H₆] α -Acetyl-N,N-dinormethadol hydrochlorideMol. formula : C₂₁H₂₈CINO₂

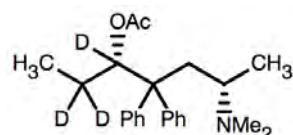
FW : 361.92 DEA schedule : 0



Catalog number : 9648-016

Name : (-)-[2,2,3-²H₃] α -Acetylmethadol hydrochlorideMol. formula : C₂₃H₃₂CINO₂

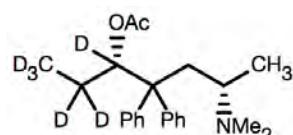
FW : 353.50 DEA schedule : 2



Catalog number : 9648-010

Name : (-)-[1,1,1,2,2,3-²H₆] α -Acetylmethadol hydrochlorideMol. formula : C₂₄H₃₄CINO

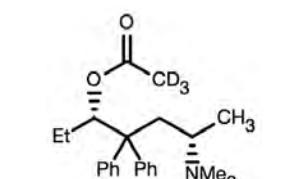
FW : 389.97 DEA schedule : 2



Catalog number : 9648-011

Name : (-)-[Acetyl-²H₃] α -acetylmethadol hydrochlorideMol. formula : C₂₁H₂₈CINO₂

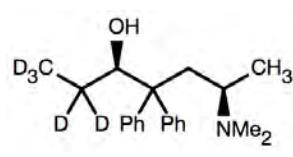
FW : 389.97 DEA schedule : 2



Catalog number : 9605-010

Name : [1,1,1,2,2-²H₅] α -Methadol hydrochlorideMol. formula : C₂₁H₃₀CINO

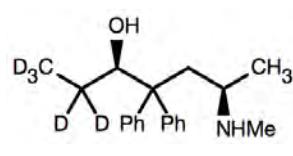
FW : 347.93 DEA schedule : 1

References : Portoghesi, PS; Williams DA *J Med Chem* 1969, 12, 839–44.

Catalog number : 9605-021

Name : [1,1,1,2,2-²H₅] α -Normethadol perchlorateMol. formula : C₂₀H₂₈CINO₅

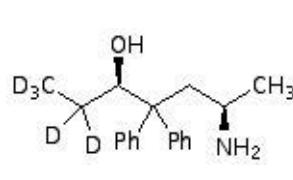
FW : 397.90 DEA schedule : 1

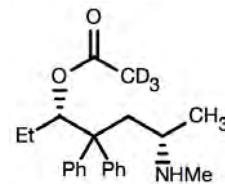
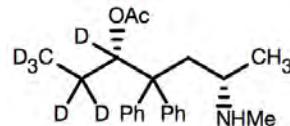
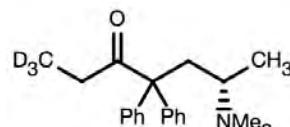
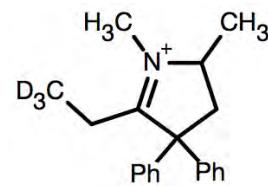
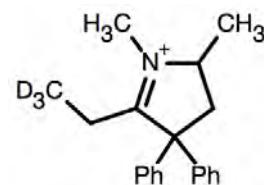
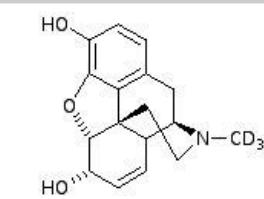


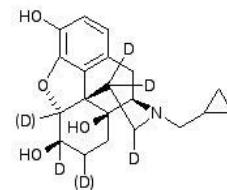
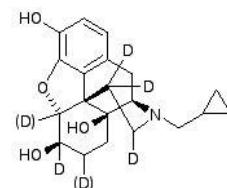
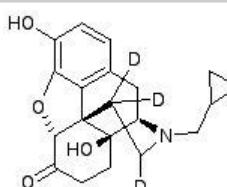
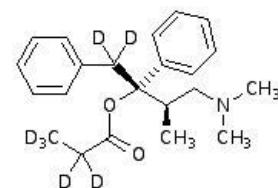
Catalog number : 9605-031

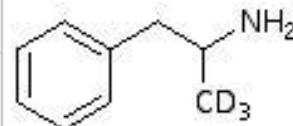
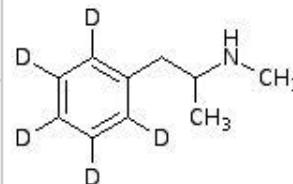
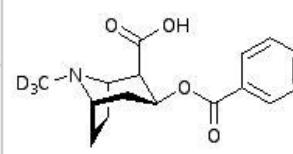
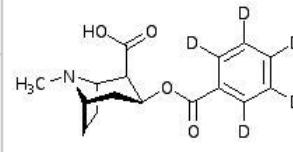
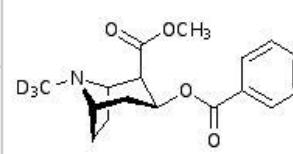
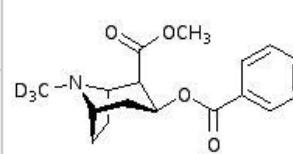
Name : [1,1,1,2,2-²H₅] α -N,N-Dinormethadol maleateMol. formula : C₁₉H₂₅NO

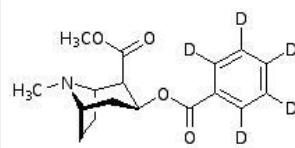
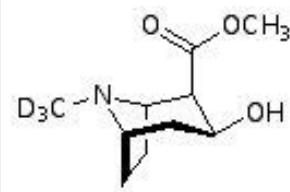
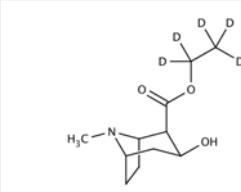
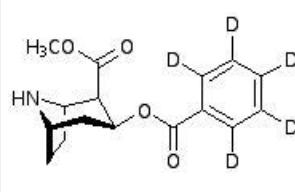
FW : 399.49 DEA schedule : 1

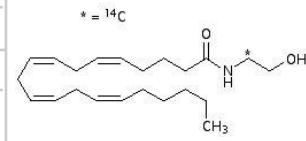
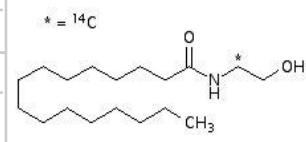
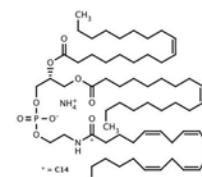
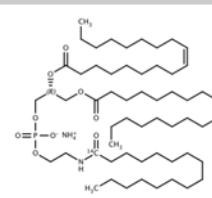


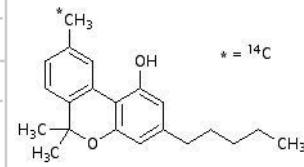
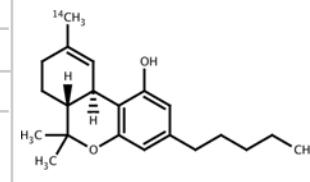
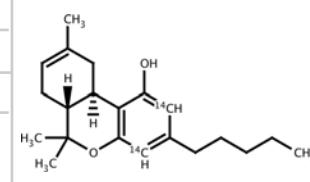
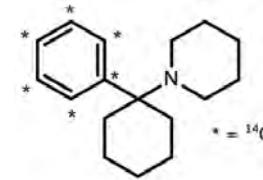
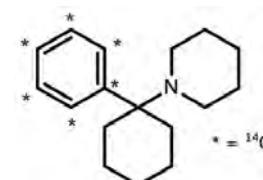
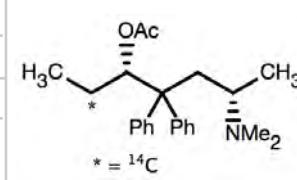
Catalog number : 9633-003**Name :** $(-)[\text{Acetyl}-^2\text{H}_3]\alpha\text{-Acetylnormethadol hydrochloride}$ **Mol. formula :** $\text{C}_{22}\text{H}_{30}\text{ClNO}_2$ **FW :** 375.94 **DEA schedule :** 1**Catalog number :** 9633-004**Name :** $(-)[1,1,1,2,2,3-^2\text{H}_6]\alpha\text{-Acetylnormethadol hydrochloride}$ **Mol. formula :** $\text{C}_{23}\text{H}_{22}\text{ClNO}_2$ **FW :** 375.94 **DEA schedule :** 1**Catalog number :** 9250-011**Name :** $(+)-[1,1,1-^2\text{H}_3]\text{Methadone hydrochloride}$ **Mol. formula :** $\text{C}_{21}\text{H}_{28}\text{ClNO}$ **FW :** 348.92 **DEA schedule :** 2**Catalog number :** 9250-027**Name :** $[\text{Ethyl}-2',2',2'-^2\text{H}_3]-1,5\text{-Dimethyl-3,3-diphenyl-2-ethylpyrrolinium perchlorate}$ **Mol. formula :** $\text{C}_{20}\text{H}_{25}\text{ClNO}_4$ **FW :** 384.91 **DEA schedule :** 0**Catalog number :** 9250-032**Name :** $[\text{Ethyl}-2',2',2'-^2\text{H}_3]-3,3\text{-Diphenyl-2-ethyl-5-methyl-1-pyrroline hydrochloride}$ **Mol. formula :** $\text{C}_{20}\text{H}_{25}\text{ClN}$ **FW :** 302.86 **DEA schedule :** 0**Opioids: Morphine Class****Catalog number :** 9300-005**CASRN :** 67293-88-3**Name :** $[\text{N-C}^2\text{H}_3]\text{Morphine}$ **Mol. formula :** $\text{C}_{17}\text{H}_{19}\text{NO}_3$ **FW :** 288.36 **DEA schedule :** 2

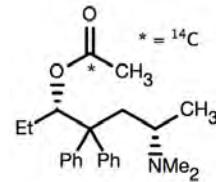
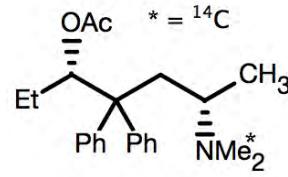
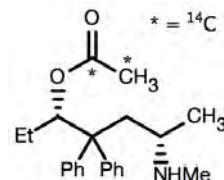
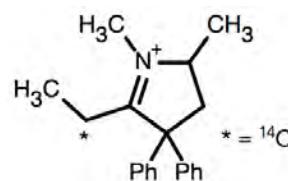
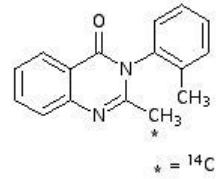
Opioids: Oxymorphone Class**Catalog number :** 9652-057**Name :** [5,6,7,15,15,16-²H₆]-6α-Naltrexol**Mol. formula :** C₂₀H₂₅NO₄**FW :** 349.45 **DEA schedule :** 0**Catalog number :** 9652-052**Name :** [5,6,7,15,15,16-²H₆]-6β-Naltrexol hydrochloride**Mol. formula :** C₂₀H₂₆ClNO₄**FW :** 385.91 **DEA schedule :** 0**Catalog number :** 9652-014**Name :** [15,15,16-²H₃]Naltrexone**Mol. formula :** C₂₀H₂₃NO₄**FW :** 344.42 **DEA schedule :** 0**Opioids: Propoxyphene Class****Catalog number :** 9273-004**Name :** [²H₇]Propoxyphene hydrochloride**Mol. formula :** C₂₂H₂₉NO₂**FW :** 346.51 **DEA schedule :** 2**Sedatives: Methaqualone Class****Catalog number :** 2565-001**CASRN :** 72-44-6**Name :** [²H₄]Methaqualone**Mol. formula :** C₁₆H₁₄N₂O**FW :** 254.32 **DEA schedule :** 1

Stimulants: Amphetamine Class**Catalog number :** 1100-004**CASRN :** 38875-35-3**Name :** (\pm)-[1,1,1- 2 H₃]Amphetamine sulfate**Mol. formula :** C₉H₁₃N**FW :** 374.53**DEA schedule :** 2**References :** Cho, AK; et al., *Anal Chem* 1973, 45, 570–4.
Valtier, S; Cody, JT *J Anal Toxicol* 1995, 19, 375–80.**Stimulants: Methamphetamine Class****Catalog number :** 1105-002**Name :** (\pm)-[2',3',4',5',6'- 2 H₅]Methamphetamine hydrochloride**Mol. formula :** C₁₀H₁₆ClN**FW :** 190.72**DEA schedule :** 2**Stimulants: Tropane Class****Catalog number :** 9180-003**Name :** (-)-[N-C²H₃]Benzoylecgonine**Mol. formula :** C₁₆H₁₉NO₄**FW :** 292.34**DEA schedule :** 2**Catalog number :** 9180-011**Name :** (-)-[Phenyl- 2 H₅]Benzoylecgonine**Mol. formula :** C₁₆H₁₉NO₄**FW :** 294.36**DEA schedule :** 2**Catalog number :** 9041-006**Name :** (-)-[N-C²H₃]Cocaine**Mol. formula :** C₁₇H₂₁NO₄**FW :** 306.37**DEA schedule :** 2**Catalog number :** 9041-007**Name :** (-)-[N-C²H₃]Cocaine hydrochloride**Mol. formula :** C₁₇H₂₂ClNO₄**FW :** 342.83**DEA schedule :** 2

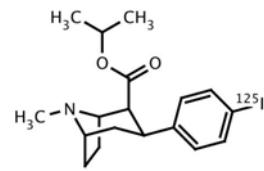
Catalog number : 9041-009**Name :** (-)-[Phenyl-²H₅]Cocaine**Mol. formula :** C₁₇H₂₁NO₄**FW :** 344.84 **DEA schedule :** 2**Catalog number :** 9180-007**Name :** [N-C²H₃]Ecgonine methyl ester hydrochloride**Mol. formula :** C₁₀H₁₈ClNO₃**FW :** 238.73 **DEA schedule :** 2**Catalog number :** 9180-021**CASRN :** 259526-73-3**Name :** Ecgonine (1,1,2,2,2-²H₅)ethyl ester perchlorate**Mol. formula :** C₁₁H₁₄D₅NO₃ • HClO₄**FW :** 318.76 **DEA schedule :** 2**Catalog number :** 9041-017**Name :** (-)-[Phenyl-²H₅]Norcocaine fumarate**Mol. formula :** C₂₀H₂₃NO₈**FW :** 410.43 **DEA schedule :** 2

Cannabinoids: Fatty Acid Derivatives (Anandamides)**Catalog number :** NOCD-007**Name :** Arachidonyl[1-¹⁴C]ethanolamide**Mol. formula :** C₂₃H₃₇NO₂**FW :** 361.56 **DEA schedule :** 0**Cannabinoids: Fatty Acid Derivatives (Palmitoyl amides)****Catalog number :** NOCD-005**Name :** Palmitoyl[1-¹⁴C]ethanolamide**Mol. formula :** C₁₈H₃₇NO₂**FW :** 299.49 **DEA schedule :** 0**Notes :** Cannabinoid CB2 receptor agonist (carbon-labeled).**Cannabinoids: Precursors, Biosynthetic****Catalog number :** NOCD-009**Name :** N-[1-¹⁴C]-Arachidonyl-1,2-dioleoyl-sn-glycero-3-phosphoethanolamine ammonium salt**Mol. formula :** C₆₁H₁₁₁N₂O₉P**FW :** 1047.52 **DEA schedule :** 0**Notes :** Putative biosynthetic precursor of endogenous cannabinoids (carbon-labeled).**References :** Morishita, J; et al. *J Neurochem* 2005, 94, 753–62.**Catalog number :** NOCD-000**Name :** N-[1-¹⁴C]-Palmitoyl-1,2-dioleoyl-sn-glycero-3-phosphoethanolamine ammonium salt**Mol. formula :** C₅₇H₁₁₁N₂O₉P**FW :** 999.47 **DEA schedule :** 0**Notes :** Putative biosynthetic precursor of endogenous cannabinoids (carbon-labeled).**References :** Morishita J; et al. *J Neurochem* 2005, 94, 753–62.

Cannabinoids: Tetrahydrocannabinol Class**Catalog number :** 7360-015**Name :** [11-¹⁴C]Cannabinol**Mol. formula :** C₂₀¹⁴CH₂₆O₂**FW :** 310.43 **DEA schedule :** 1**Notes :** Inactive constituent of cannabis (carbon-labeled).**Catalog number :** 7370-033**Name :** [11-¹⁴C] Δ^9 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 316.45 **DEA schedule :** 1**Notes :** Hallucinogen; psychotropic; analgesic (carbon-labeled).**Catalog number :** 7370-032**Name :** [2,4-¹⁴C] Δ^8 -THC**Mol. formula :** C₂₁H₃₀O₂**FW :** 318.45 **DEA schedule :** 1**Notes :** Hallucinogen; psychotropic; analgesic (carbon-labeled).**Dissociatives: Phencyclidine Class****Catalog number :** 7471-002**Name :** [¹⁴C]Phencyclidine HBr; [¹⁴C]PCP**Mol. formula :** C₁₈H₂₈BrN**FW :** 523.53 **DEA schedule :** 2**Catalog number :** 7471-007**Name :** [Phenyl-U-¹⁴C]-1-(1-Phenylcyclohexyl)Piperidine; [Phenyl-U-¹⁴C]PCP**Mol. formula :** C₁₇H₂₅N**FW :** 245.39 **DEA schedule :** 2**Opioids: Methadone Class****Catalog number :** 9648-013**Name :** (-)-[2-¹⁴C]- α -Acetylmethadol hydrochloride**Mol. formula :** C₂₃H₃₂ClNO₂**FW :** 353.50 **DEA schedule :** 2

Catalog number : 9648-014**Name :** (–)[Acetyl–¹⁴C]–α-Acetylmethadol**Mol. formula :** C₂₃H₃₂CINO₂**FW :** 353.50 **DEA schedule :** 2**Catalog number :** 9648-015**Name :** (–)[N–¹⁴CH₃]–α-Acetylmethadol hydrochloride**Mol. formula :** C₂₃H₃₁NO₂**FW :** 353.50 **DEA schedule :** 2**Catalog number :** 9633-005**Name :** (–)[Acetyl–¹⁴C₂]–α-Acetylnormethadol hydrochloride**Mol. formula :** C₂₂H₃₀CINO₂**FW :** 375.94 **DEA schedule :** 1**Catalog number :** 9250-026**Name :** [1'–¹⁴C]–2–Ethyl–1,5–dimethyl–3,3–diphenylpyrrolinium perchlorate**Mol. formula :** C₂₀H₂₅CINO₄**FW :** 536.01 **DEA schedule :** 0**Sedatives: Methaqualone Class****Catalog number :** 2565-002**CASRN :** 72-44-6**Name :** [2–¹⁴C]Methaqualone**Mol. formula :** C₁₆H₁₄N₂O**FW :** 251.30 **DEA schedule :** 1**Notes :** Sedative; hypnotic (carbon-labeled).



Peptides: Opioid**Catalog number :** MPSP-026**Name :** [¹²⁵I]β-Endorphin**Mol. formula :** C₁₅₈H₂₄₉N₃₉O₄₆S₁₂₅I₂**FW :** 3589.6**Notes :** Radioactive ligand for MPSP-024**References :** Deby-Dupont, G; et al. *C R Seances Soc Biol Fil* **1983**, *177*, 259–68.
Schweigerer, L; et al. *J Biol Chem* **1983**, *258*, 12287–92.**Stimulants: Tropane Class****Catalog number :** NOCD-077**Name :** [¹²⁵I]RTI-121**Mol. formula :** C₁₈H₂₄INO₂**FW :** 413.29 **DEA schedule :** 0**References :** Scheffel, U; et al. *Neuroreport* **1992**, *3*, 969–72.
Carroll, FI; et al. *J Med Chem* **1995**, *38*, 379–88.



Limited quantities of Nicotine Research Cigarettes (NRC)s manufactured under the NIDA DSP are available to research investigators under specified parameters such as nicotine and tar yields ,and type as reduced nicotine (RN), reduced nicotine and menthol (RN-Men), reduced nicotine and high tar (RN-HT), reduced nicotine, high tar and menthol (RN-HT-Men), and conventional nicotine (CN). The NIDA DSP will consider requests for its current inventory of NRCs from researcher investigators funded through the NIH to conduct tobacco science research.

To facilitate planning for either usage of the current inventory of NRCs, or future manufacturing needs, the NIDA DSP requests that research investigators who intend to submit requests for NRCs from the NIDA DSP send an email request with the following information: 1) anticipated total quantity (carton) of each type of NRC needed with their Tobacco Product Master File (TPMF) codes and Type, 2) the requested timeline and amounts for initial receipt, and 3) requested timeline of replenishment of each type of NRC. For those research investigators who have NIDA/NIH grants, they can submit their request according to the NIDA DSP Guidelines for Research Chemicals, Controlled Substances, and Nicotine Research Cigarettes (NRC). For more information, please email NIDANRCSupply@mail.nih.gov

Characteristics of the NRC inventory are described in the following table:

Nicotine Research Cigarettes (RTI)								
TPMF Code	Type	Filter PD (mm H ₂ O)	Filter Type (Mono/Dual)	Filter Type (Regular/Menthол)	Tipping Paper Perforation	Nicotine Content (mg/g tobacco)	Specifications Nicotine Yield	Specifications Tar Yield
NRC100	RN-HV	81	Dual	Regular	800CU	0.47 ± 0.02	0.02 ± 0.01	3.0 ± 1.5
NRC101	RN-HV-Men	81	Dual	Menthол	800CU	0.47 ± 0.01	0.02 ± 0.01	3.0 ± 1.5
NRC102	RN	81	Dual	Regular	100CU	0.42 ± 0.01	0.03 ± 0.01	9.0 ± 1.5
NRC103	RN-Men	81	Dual	Menthол	100CU	0.44 ± 0.01	0.03 ± 0.01	9.0 ± 1.5
NRC104	RN-HT	56	Mono	Regular	0CU	0.51 ± 0.00	0.04 ± 0.02	13.0 ± 2.0
NRC105	RN-HT-Men	56	Mono	Menthол	0CU	0.46 ± 0.01	0.04 ± 0.02	13.0 ± 2.0
NRC200	RN	81	Dual	Regular	100CU	1.40 ± 0.03	0.07 ± 0.02	9.0 ± 1.5
NRC201	RN-Men	81	Dual	Menthол	100CU	1.34 ± 0.07	0.07 ± 0.02	9.0 ± 1.5
NRC300	RN	81	Dual	Regular	100CU	2.53 ± 0.04	0.12 ± 0.03	9.0 ± 1.5
NRC301	RN-Men	81	Dual	Menthол	100CU	2.54 ± 0.05	0.12 ± 0.03	9.0 ± 1.5
NRC302	RN-HT-Men	56	Mono	Menthол	0CU	2.50 ± 0.04	0.16 ± 0.03	13.0 ± 2.0
NRC400	RN	81	Dual	Regular	100CU	5.62 ± 0.18	0.26 ± 0.06	9.0 ± 1.5
NRC401	RN-Men	81	Dual	Menthол	100CU	5.54 ± 0.27	0.26 ± 0.06	9.0 ± 1.5
NRC402	RN-HT	81	Mono	Regular	100CU	6.03 ± 0.18	0.33 ± 0.06	13.0 ± 2.0
NRC405	RN-HT-Men	56	Mono	Menthол	0CU	5.96 ± 0.21	0.40 ± 0.08	13.0 ± 2.0
NRC404	RN-HT	56	Mono	Regular	0CU	8.08 ± 0.15	0.60 ± 0.12	13.0 ± 2.0
NRC500	RN-HT	81	Mono	Regular	100CU	12.06 ± 0.41	0.70 ± 0.15	13.0 ± 2.0
NRC501	RN-Men	81	Dual	Menthол	100CU	11.26 ± 0.11	0.60 ± 0.12	9.0 ± 1.5
NRC600	CN	81	Dual	Regular	100CU	17.36 ± 0.28	0.80 ± 0.15	10.5 ± 1.5
NRC601	CN-Men	81	Dual	Menthол	100CU	16.50 ± 0.17	0.80 ± 0.15	10.5 ± 1.5
NRC602	CN-HT-Men	56	Mono	Menthол	0CU	16.10 ± 0.52	1.10 ± 0.20	16.0 ± 2.0
NRC700	LTNR-HT	81	Mono	Regular	100CU	25.94 ± 0.65	1.60 ± 0.30	16.0 ± 2.0
NRC701	LTNR-Men	81	Dual	Menthол	100CU	25.83 ± 0.48	1.60 ± 0.30	12.0 ± 2.0

Definitions

CN	Conventional Nicotine
CN-HT-Men	Conventional Nicotine-High Tar-Menthol
CN-Men	Conventional Nicotine-Menthol
CU	Coresta Units
	[0CU = Non-perforated tipping paper]
Dual filter	Consists of 10 mm paper and 15 mm cellulose acetate segments
LTNR-HT	Low Tar Nicotine Ratio-High Tar
LTNR-Men	Low Tar Nicotine Ratio-Menthol
Mono filter	Consists of 25 mm cellulose acetate segment
PD	Pressure drop
RN	Reduced Nicotine
RN-HT	Reduced Nicotine-High Tar
RN-HT-Men	Reduced Nicotine-High Tar-Menthol
RN-HV	Reduced Nicotine-High Ventilation
RN-HV-Men	Reduced Nicotine-High Ventilation-Menthol
RN-Men	Reduced Nicotine-Menthol
TPMF	Tobacco Product Master File number

Marijuana Plant Material

Research grade marijuana products are available in a variety of cannabinoid content specifications as shown below. For detailed information on specific batches please refer to the NIDA website (<http://www.drugabuse.gov/researchers/research-resources/nida-drug-supply-program-dsp/bulk-marijuana-plant-material-addendum-to-drug-supply-catalog>). NIDA may also be able to develop bulk marijuana of other specific THC and CBD contents by mixing batches to meet researcher needs. For further information please contact the NIDA Drug Supply Program Director, Dr Hari Singh, at hsingh@nida.nih.gov.

Marijuana cigarettes and bulk marijuana plant material produced under the NIDA DSP are available at no cost to research investigators who have an NIH grant. Marijuana is also available to research investigators who are funded through non-Federal sources on a cost-reimbursement basis. Please see the Note at the end of the tables for more information on costs.

Marijuana cigarettes

The following represents the existing stocks of manufactured marijuana cigarettes, which will be made available to the research community until depleted. For additional information on the specific cannabinoid content please see <http://www.-drugabuse.gov/researchers/research-resources/nida-drug-supply-program-dsp/bulk-marijuana-plant-material-addendum-to-drug-supply-catalog>. Please note that NIDA does not plan to manufacture marijuana cigarettes in the near future. Only bulk marijuana will generally be available for research:

Placebo Marijuana Cigarettes (Placebo)	0.001% THC
Marijuana Cigarettes (Medium)	2.0% THC
Marijuana Cigarettes (Medium)	3.6% THC
Marijuana Cigarettes (High)	5.6% THC
Marijuana Cigarettes (High)	6.4% THC

Bulk Marijuana

Bulk marijuana is currently available in the following general categories, and due to recent interest its strength is being provided for both THC and CBD as *Low* (<1%), *Medium* (1-5%), *High* (5-10%), and *Very High* (>10%). Bulk marijuana has small amounts of other cannabinoids (CBC, CBG, CBN, and THCV) which are reported in the batch specific details (<http://www.drugabuse.gov/researchers/research-resources/nida-drug-supply-program-dsp/bulk-marijuana-plant-material-addendum-to-drug-supply-catalog>).

Placebo marijuana (produced by solvent extraction)

1. THC (0%) / CBD (0%)

Low THC varieties

1. Low THC (<1%) / Medium CBD (1-5%)
2. Low THC (<1%) / High CBD (5-10%)
3. Low THC (<1%) / Very High CBD (>10%)

Medium THC varieties

1. Medium THC (1-5%) / Low CBD (<1%)
2. Medium THC (1-5%) / Medium CBD (1-5%)
3. Medium THC (1-5%) / High CBD (5-10%)
4. Medium THC (1-5%) / Very High CBD (>10%)

High THC varieties

1. High THC (5-10%) / Low CBD (<1%)
2. High THC (5-10%) / High CBD (5-10%)
3. High THC (5-10%) / Very High CBD (>10%)

Very high THC varieties

1. Very High THC (>10%) / Low CBD (<1%)

Important note regarding cost-reimbursement

The services and material from the NIDA Drug Supply Program are generally free to NIH-sponsored investigators. Some requests may incur a nominal fee depending on study status, funding source, and type of material being requested. Pursuant to HHS policy published May 21, 1999 marijuana for non-Federally funded research is to be provided on a cost-reimbursement basis. NIDA intends to implement the following pricing schedule for collection of fees for marijuana once HHS and NIH develop updated policies and transactional procedures on the specifics regarding cost-reimbursement of research substances for non-Federally supported projects. These policies and procedures may significantly impact the timing, mechanisms, and recipient of any collected fees but will not impact the shipments. An update will be posted once the final Guidance is received.

Marijuana Pricing Schedule

<u>Item</u>	<u>Price</u>
Non-placebo Cigarette	\$10.96 each
Placebo Cigarette	\$13.94 each
Bulk Marijuana	\$2,497.00 per kilogram

