LDRUG

A training database for learning how to use the DRUG files:

DDFB Derwent Drug File Standard
DRUGB Derwent Drug File for Subscribers

(See database summary sheet DDFU/DRUGU' for details on search and display fields, which are valid in LDRUG).

Subject Coverage

All aspects of drugs:

- Analysis
- Biochemistry
- Galenics
- Metabolism
- Pharmacokinetics
- Pharmacology
- Structure-activity relationships
- Synthesis
- Therapeutics and adverse effects
- Toxicology

File Type

Bibliographic, learning

Features

Alerts (SDI)	Not available				
CAS Registry Number® Identifiers	☑	Page Images	☐	STN® AnaVist™	☐
Keep & Share	☑	SLART	☐	STN Easy®	☐
Learning Database	☑	Structures	☐		

Record Content

- Record contain bibliographic information, controlled term indexing, and structure codes (file DRUGU only)

File Size

- 500 compounds from the registry segment and about 15,000 relating literature records

Coverage

1983-November 1994

Updates

Closed file

Language

English

Database Producer

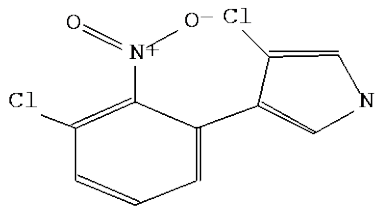
Clarivate Analytics (UK) Limited
Friars House, 160 Black Friars Rd.
London SE1 8EZ
United Kingdom

Sources	• Medical and scientific journals and conference proceedings.
User Aids	• Derwent Drug File Thesaurus * • Journal List and Selection Guidelines * • Product Description * • Work Books * • Online Helps (HELP DIRECTORY lists all help messages available) • STNGUIDE * available from producer
Clusters	• LEARNING STN Database Clusters information (PDF)
Related Databases	• DDFU • DRUGU contain data from 1983 to present
Pricing	Enter HELP COST at an arrow prompt.

Sample Records

DISPLAY ALL STR from REGISTRY SEGMENT

```
AN    20825  LDRUG
DDRN  PYRROLNIT
DDN   PYRROLNITRIN
RN    1018-71-9
CT    ANTIBIOTICS; FUNGICIDES
SS    PYRROLE; CHLORINE; ARYLCHLORIDE; NITROARENE; BH-LINKED-CC
MPC   02& *G; 02- *G; 03& *G; 032 *G; 06& *G; 067 *G; 071 *G; 097 *G; 10- *G;
      104 *G; 105 *G; 131 *G; 163 *G; 17& *G; 17- *G; 171 *G; 174 *G
```



DISPLAY ALL from LITERATURE SEGMENT

```
AN    1986-21902  LDRUG   P A
TI    On the Pharmaceutical and Biopharmaceutical Assessment of Acetylamino
      nitropropoxybenzene (Falimint). Part 6. Quantitive Composition ofthe
      Metabolites.
AU    Gerlach K; Metzner J; Fuerst W
LO    Halle, Germany,East
SO    Pharmazie (40, Number 8, 564-65, 1985) 2 Fig. 5 Tab. 2 Reference
      CODEN: PHARAT       ISSN:  0031-7144
AV    Sekt. Pharmazie der Martin-Luther-University Halle-Wittenberg, DDR-4020
      Halle/Saale, Weinbergweg 15, DDR
LA    German
DT    Journal
AB     The metabolism of Falimint following p.o. administration to human subjects
       was studied.  Polarographic analysis of the urine showed that most of the
       drug was eliminated within 6 hr, the main metabolite being falimintic acid
       (3-(2-acetamido 4-nitrophenoxy) propionic acid).  Other metabolies include
       free and conjugated 2-acetamido-4-nitrophenol (AANP), 2-aminonitrophenol
       (ANP), 3-(2-amino 4-nitrophenoxy) propionic acid (M7) and 2-amino 4-
       nitropropoxybenzene (Ultrasweet).  Variations in the relative amounts of
       the metabolites over the test period were discussed.
SH    P Pharmacology
      A Analysis
CC    5  Analysis
      8  Pharmacokinetics
      33 Respiratory
CT       QUANT. *FT; DET. *FT; ANALYSIS *FT; POLAROGRAPHY *FT; ELIMINATION *FT;
         CONC. *FT; URINE *FT; HUMAN *FT; PHARMACOKINETICS *FT; METABOLITE *FT;
         ELECTROCHEM. *FT
    [01] FALIMINT *OC; FALIMINT *DM; P.O. *FT; ANTITUSSIVES *FT; FALIMINT *RN;
         OC *FT; DM *FT
    [02] FALIMINTATE *OC; FALIMINTATE *DM; PHENOL-ETHER *FT; NITROARENE *FT;
         ARYLAMINE *FT; AMINOACID *FT; C-AMIDE *FT; OC *FT; DM *FT; FALIMINTA
```

LDRUG

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          *RN
     [03] ULTRASWEET *OC; ULTRASWEET *DM; ULTRASWEE *RN; OC *FT; DM *FT
     [04] NITROARENE *FT; ARYLAMINE *FT; PHENOL *FT; C-AMIDE *FT; CONJUGATE *FT;
          PHENOL-ETHER *FT; C-ACID *FT; OC *FT; DM *FT
FA     AB; LA; CT; MPC
FS     Literature
```

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