Indexing - Substance

Food and Drug Administration
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IBOPAMINE HYDROCHLORIDE (3VXW2HU8GS)

Alias Name

Name	Name Type	Assigning Territory	Assigning Organization	Reference Document	Citation	Policy
IBOPAMINE HYDROCHLORIDE	primary name					

Substance Information

Mapping	Definition Hash:	a38c0e4a-2dbf-cf4c-4401-30084a9a8230

Moiety

Name		Quantity	1

Moiety Characteristics

Chemical Structure, MOLFILE

```
  -FDASRS-09061311012D

 23 22  0  0  0  0  0  0  0  0999 V2000
    4.6875   -5.7625    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    1.9875   -1.1875    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    4.6875   -3.4542    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    4.0167   -4.6125    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    2.6667   -2.3292    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    4.0042   -2.3292    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    5.9792   -3.4542    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    2.6667   -0.0250    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    6.0042   -5.7625    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    4.6875   -1.1750    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.6792   -1.1875    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    4.0167   -6.9167    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    6.6625   -2.3292    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    5.9792   -1.1500    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    9.9792   -1.1500    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    8.6417   -1.1500    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   10.6667    0.0000    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    7.9917   -2.3292    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    4.6875   -8.0667    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    2.6917   -6.9167    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.0000   -2.3292    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.0000   -0.0250    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   10.2042   -5.8542    0.0000 Cl  0  0  0  0  0  0  0  0  0  0  0  0
  2  5  1  0  0  0  0
  3  4  1  0  0  0  0
  4  1  1  0  0  0  0
  5  6  1  0  0  0  0
  6  3  2  0  0  0  0
  7  3  1  0  0  0  0
  8  2  2  0  0  0  0
  9  1  2  0  0  0  0
 10  6  1  0  0  0  0
 11  2  1  0  0  0  0
 12  1  1  0  0  0  0
 13  7  2  0  0  0  0
 14 13  1  0  0  0  0
```

	15 16 1 0 0 0 0 16 18 1 0 0 0 0 17 15 1 0 0 0 0 18 13 1 0 0 0 0 19 12 1 0 0 0 0 20 12 1 0 0 0 0 21 11 1 0 0 0 0 22 11 1 0 0 0 0 14 10 2 0 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C17H25NO4.ClH/c1-11(2)16(19)21-14-7-6-13(8-9-18-5)10-15(14)22-17(20)12(3)4;/h6-7,10-12,18H,8-9H2,1-5H3;1H
Chemical Structure, InChI Key	ALIXRWKJZYLBJI-UHFFFAOYSA-N

Molecular Bond Types

Product Information

Product Type	INDEXING - SUBSTANCE	**Item Code (Source)**	3VXW2HU8GS

Labeler - Food and Drug Administration (927645523)

Revised: 9/2013 Food and Drug Administration