Indexing - Substance

Food and Drug Administration
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OSELTAMIVIR (20O93L6F9H)

Alias Name

Name	Name Type	Assigning Territory	Assigning Organization	Reference Document	Citation	Policy
oseltamivir	primary name					

Substance Information

Mapping	Definition Hash:	bad994c6-0c44-0117-1db7-657492e2fe5d

Moiety

Name		Quantity	1

Moiety Characteristics

Chemical Structure, MOLFILE

```
  -FDASRS-01291411012D

 22 22  0  0  1  0  0  0  0  0999 V2000
    2.5125    0.3042    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.4958   -0.8500    0.0000 C   0  0  1  0  0  0  0  0  0  0  0  0
    0.4958    0.3042    0.0000 C   0  0  1  0  0  0  0  0  0  0  0  0
    1.5083    0.8833    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.5167    0.8708    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    1.5083   -1.4292    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
   -0.5042   -1.4292    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    2.5125   -0.8500    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -0.5042   -2.5708    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -0.4958    0.8708    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    3.5167    2.0167    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    0.4958   -3.1625    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    1.5083   -2.5583    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    4.4500    0.2625    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
   -0.4833    1.9750    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -1.5167   -3.0250    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    5.4542    0.7417    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -1.4792    2.5417    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.5208    2.5167    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    6.5375    0.1667    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -2.4542    1.9750    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.5292    3.6458    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
  2  6  1  0  0  0  0
  3  4  1  0  0  0  0
  4  1  2  0  0  0  0
  5  1  1  0  0  0  0
  6  8  1  0  0  0  0
  2  7  1  1  0  0  0
  8  1  1  0  0  0  0
  9  7  1  0  0  0  0
  3 10  1  6  0  0  0
 11  5  2  0  0  0  0
 12  9  2  0  0  0  0
  6 13  1  6  0  0  0
 14  5  1  0  0  0  0
 15 10  1  0  0  0  0
```

	16 9 1 0 0 0 0 17 14 1 0 0 0 0 18 15 1 0 0 0 0 19 15 1 0 0 0 0 20 17 1 0 0 0 0 21 18 1 0 0 0 0 22 19 1 0 0 0 0 2 3 1 0 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C16H28N2O4/c1-5-12(6-2)22-14-9-11(16(20)21-7-3)8-13(17)15(14)18-10(4)19/h9,12-15H,5-8,17H2,1-4H3,(H,18,19)/t13-,14+,15+/m0/s1
Chemical Structure, InChI Key	VSZGPKBBMSAYNT-RRFJBIMHSA-N

Molecular Bond Types

Product Information

Product Type	INDEXING - SUBSTANCE	**Item Code (Source)**	20O93L6F9H

Labeler - Food and Drug Administration (927645523)

Revised: 1/2014 Food and Drug Administration