Indexing - Substance

Food and Drug Administration
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SUNITINIB (V99T50803M)

Alias Name

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

Substance Information

Mapping	**Definition Hash:**	6cb40e1f-08bd-c4d8-b711-729c00483596

Moiety

Name		**Quantity**	1

Moiety Characteristics

Chemical Structure, MOLFILE

```
  -FDASRS-10241311012D

 29 31  0  0  0  0  0  0  0  0999 V2000
    2.4333   -1.4856    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    2.4333   -2.4357    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    3.3868   -2.7332    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.9792   -1.9650    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.3935   -1.1962    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.6959   -2.4435    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.6959   -1.4934    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -0.2576   -2.7408    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    1.5625   -1.0124    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    1.5147   -2.9999    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
   -0.8500   -1.9728    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -0.2650   -1.2020    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -0.6718   -0.3394    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -1.6716   -0.2418    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -2.2589   -1.0194    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -1.8441   -1.8781    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.7077   -0.2926    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.6931   -3.6393    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    4.9792   -1.9625    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    5.4833   -2.7874    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    6.4833   -2.7874    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    6.9833   -1.9625    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    7.9875   -1.9625    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    8.4875   -2.7874    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    9.4875   -2.7874    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    5.4850   -1.1411    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    8.4875   -1.1416    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    9.4875   -1.1416    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -2.0864    0.6244    0.0000 F   0  0  0  0  0  0  0  0  0  0  0  0
  3 18  1  0  0  0  0
  8  6  1  0  0  0  0
  4 19  1  0  0  0  0
  5  1  2  0  0  0  0
 19 20  1  0  0  0  0
  7  9  2  0  0  0  0
 20 21  1  0  0  0  0
```

	9 1 1 0 0 0 0 21 22 1 0 0 0 0 6 7 1 0 0 0 0 22 23 1 0 0 0 0 6 10 2 0 0 0 0 23 24 1 0 0 0 0 24 25 1 0 0 0 0 19 26 2 0 0 0 0 11 12 2 0 0 0 0 23 27 1 0 0 0 0 1 2 1 0 0 0 0 27 28 1 0 0 0 0 12 13 1 0 0 0 0 14 29 1 0 0 0 0 2 3 1 0 0 0 0 13 14 2 0 0 0 0 3 4 2 0 0 0 0 14 15 1 0 0 0 0 4 5 1 0 0 0 0 15 16 2 0 0 0 0 16 11 1 0 0 0 0 7 12 1 0 0 0 0 5 17 1 0 0 0 0 11 8 1 0 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C22H27FN4O2/c1-5-27(6-2)10-9-24-22(29)20-13(3)19(25-14(20)4)12-17-16-11-15(23)7-8-18(16)26-21(17)28/h7-8,11-12,25H,5-6,9-10H2,1-4H3,(H,24,29)(H,26,28)/b17-12-
Chemical Structure, InChI Key	WINHZLLDWRZWRT-ATVHPVEESA-N

Molecular Bond Types

Product Information

Product Type	INDEXING - SUBSTANCE	**Item Code (Source)**	V99T50803M

Labeler - Food and Drug Administration (927645523)

Revised: 10/2013 Food and Drug Administration