Indexing - Substance

Food and Drug Administration
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TRENBOLONE (P53R4420TR)

Alias Name

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

Substance Information

Mapping	**Definition Hash:**	158a2255-4e85-06c8-71c9-4bcd2c2ca330

Moiety

Name		**Quantity**	1

Moiety Characteristics

Chemical Structure, MOLFILE

```
  -FDASRS-09061311012D

 22 25  0  0  1  0  0  0  0  0999 V2000
    3.0629   -3.0430    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.0577   -2.3244    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
    2.4381   -1.9583    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    1.8156   -2.3240    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    1.1952   -1.9615    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.5706   -2.3240    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.5706   -3.0449    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   -0.0751   -3.4157    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    1.1952   -3.4032    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    1.8166   -3.0449    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    2.4402   -3.4042    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    2.4389   -1.2408    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.0654   -0.8876    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.6851   -1.2495    0.0000 C   0  0  1  0  0  0  0  0  0  0  0  0
    3.6787   -1.9695    0.0000 C   0  0  1  0  0  0  0  0  0  0  0  0
    4.9152   -1.9835    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    4.9258   -1.2636    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    4.3085   -0.8983    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
    4.3040   -0.1490    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    3.6707   -2.7074    0.0000 H   0  0  0  0  0  0  0  0  0  0  0  0
    3.6831   -0.5074    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.0498   -1.5824    0.0000 H   0  0  0  0  0  0  0  0  0  0  0  0
  1  2  1  0  0  0  0
  3  2  1  0  0  0  0
  4  3  2  0  0  0  0
  4  5  1  0  0  0  0
  6  5  1  0  0  0  0
  6  7  1  0  0  0  0
  7  8  2  0  0  0  0
  7  9  1  0  0  0  0
  9 10  2  0  0  0  0
 10 11  1  0  0  0  0
 11  1  1  0  0  0  0
  4 10  1  0  0  0  0
  3 12  1  0  0  0  0
 13 12  2  0  0  0  0
```

	14 13 1 0 0 0 0 14 15 1 0 0 0 0 2 15 1 0 0 0 0 15 16 1 0 0 0 0 16 17 1 0 0 0 0 17 18 1 0 0 0 0 18 14 1 0 0 0 0 18 19 1 1 0 0 0 15 20 1 6 0 0 0 14 21 1 1 0 0 0 2 22 1 1 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C18H22O2/c1-18-9-8-14-13-5-3-12(19)10-11(13)2-4-15(14)16(18)6-7-17(18)20/h8-10,15-17,20H,2-7H2,1H3/t15-,16+,17+,18+/m1/s1
Chemical Structure, InChI Key	MEHHPFQKXOUFFV-OWSLCNJRSA-N

Molecular Bond Types

Product Information

Product Type	INDEXING - SUBSTANCE	**Item Code (Source)**	P53R4420TR

Labeler - Food and Drug Administration (927645523)

Revised: 9/2013 Food and Drug Administration