Indexing - Substance

Food and Drug Administration
Indexing - Substance

TRIAMCINOLONE (1ZK20VI6TY)

Alias Name

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

Substance Information

Mapping	**Definition Hash:**	a97882f8-deba-3d64-25c5-1e6ac7b68818

Moiety

Name		**Quantity**	1

Moiety Characteristics

Chemical Structure, MOLFILE

```
  -FDASRS-09061311012D

 30 33  0  0  1  0  0  0  0  0999 V2000
    5.6500   -2.3250    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
    3.5667   -3.4917    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
    6.7792   -1.9709    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
    2.5417   -4.0833    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
    5.6250   -3.5167    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
    4.5917   -4.0917    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
    2.5417   -5.2708    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.5834   -2.2959    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
    7.4625   -2.9417    0.0000 C   0  0  1  0  0  0  0  0  0  0  0  0
    4.6250   -1.7209    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    1.5209   -3.4792    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    6.7459   -3.8917    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    1.5209   -5.8500    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    6.7750   -0.7917    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    4.5917   -5.2750    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.5083   -4.0833    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    0.5083   -5.2708    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.5667   -5.8625    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.5667   -4.6708    0.0000 F   0  0  0  0  0  0  0  0  0  0  0  0
    5.7500   -0.2042    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    7.8042   -1.3792    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
   -0.5208   -5.8500    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    5.6458   -1.1458    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    2.5583   -1.7042    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    8.6417   -2.9417    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    2.5417   -2.9042    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    7.8000   -0.2000    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    8.8250   -0.7917    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    4.5917   -2.9125    0.0000 H   0  0  0  0  0  0  0  0  0  0  0  0
    5.6125   -4.6959    0.0000 H   0  0  0  0  0  0  0  0  0  0  0  0
  2  8  1  0  0  0  0
  3  1  1  0  0  0  0
  4  2  1  0  0  0  0
  5  1  1  0  0  0  0
  6  5  1  0  0  0  0
  7  4  1  0  0  0  0
```

	8 10 1 0 0 0 0 9 3 1 0 0 0 0 10 1 1 0 0 0 0 11 4 1 0 0 0 0 12 5 1 0 0 0 0 13 7 2 0 0 0 0 3 14 1 1 0 0 0 15 6 1 0 0 0 0 16 11 2 0 0 0 0 17 16 1 0 0 0 0 18 15 1 0 0 0 0 2 19 1 6 0 0 0 20 14 2 0 0 0 0 3 21 1 6 0 0 0 22 17 2 0 0 0 0 1 23 1 1 0 0 0 8 24 1 1 0 0 0 9 25 1 6 0 0 0 4 26 1 1 0 0 0 27 14 1 0 0 0 0 28 27 1 0 0 0 0 6 29 1 1 0 0 0 5 30 1 6 0 0 0 12 9 1 0 0 0 0 6 2 1 0 0 0 0 7 18 1 0 0 0 0 13 17 1 0 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C21H27FO6/c1-18-6-5-12(24)7-11(18)3-4-13-14-8-15(25)21(28,17(27)10-23)19(14,2)9-16(26)20(13,18)22/h5-7,13-16,23,25-26,28H,3-4,8-10H2,1-2H3/t13-,14-,15+,16-,18-,19-,20-,21-/m0/s1
Chemical Structure, InChI Key	GFNANZIMVAIWHM-OBYCQNJPSA-N

Molecular Bond Types

Product Information

Product Type	INDEXING - SUBSTANCE	**Item Code (Source)**	1ZK20VI6TY

Labeler - Food and Drug Administration (927645523)

Revised: 9/2013 Food and Drug Administration