Indexing - Substance

Food and Drug Administration
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.BETA.-CITRONELLOL, (R)- (P01OUT964K)

Alias Name

Name	Name Type	Assigning Territory	Assigning Organization	Reference Document	Citation	Policy
.BETA.-CITRONELLOL, (R)-	primary name					
.BETA.-CITRONELLOL, R-(+)-	display name					

Substance Information

Mapping	Definition Hash:	cf8d3398-cfb5-50b3-1f10-71a9547a752f

Moiety

Name		Quantity	1

Moiety Characteristics

Chemical Structure, MOLFILE

```
  -FDASRS-09261316272D

 11 10  0  0  1  0  0  0  0  0999 V2000
    3.6125   -2.7084    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    4.3250   -3.1209    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    5.0375   -2.7084    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    8.6000   -3.1084    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    2.9000   -3.1209    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.6083   -1.8834    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    5.7500   -3.1167    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    7.8875   -2.6959    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    6.4625   -2.7042    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
    7.1750   -3.1084    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    6.4583   -1.8792    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
  2  1  2  0  0  0  0
  3  2  1  0  0  0  0
  4  8  1  0  0  0  0
  5  1  1  0  0  0  0
  6  1  1  0  0  0  0
  7  3  1  0  0  0  0
  8 10  1  0  0  0  0
  9  7  1  0  0  0  0
 10  9  1  0  0  0  0
  9 11  1  1  0  0  0
M  END
```

Chemical Structures, InChI	InChI=1S/C10H20O/c1-9(2)5-4-6-10(3)7-8-11/h5,10-11H,4,6-8H2,1-3H3/t10-/m1/s1
Chemical Structure, InChI Key	QMVPMAAFGQKVCJ-SNVBAGLBSA-N

Molecular Bond Types

Product Information			
Product Type	INDEXING - SUBSTANCE	**Item Code (Source)**	P01OUT964K

Labeler - Food and Drug Administration (927645523)

Revised: 9/2013 Food and Drug Administration