Indexing - Substance

Food and Drug Administration
Indexing - Substance

DIAZOLIDINYL UREA (H5RIZ3MPW4)

Alias Name

Name	Name Type	Assigning Territory	Assigning Organization	Reference Document	Citation	Policy
DIAZOLIDINYL UREA	primary name					

Substance Information

Mapping	**Definition Hash:**	9361a305-93be-acab-6b19-5ceea196301c

Moiety (mixture component)

Name		**Quantity**	Undefined not 0

Part1Moiety (mixture component)

Name		**Quantity**	1

Moiety Characteristics

Chemical Structure, MOLFILE

```
C9H16N4O8
  -FDASRS-02211411142D

 27 27  0  0  0  0  0  0  0  0999 V2000
    6.2197   -4.0936    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    5.5113   -3.8686    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    5.0738   -4.4728    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    4.3322   -4.4728    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    3.9530   -3.8270    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    5.5155   -5.0728    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    6.2197   -4.8353    0.0000 C   0  0  1  0  0  0  0  0  0  0  0  0
    6.8613   -5.2061    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    7.5030   -4.8270    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    8.1447   -5.1978    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
    8.7863   -4.8270    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    9.4280   -5.1978    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    7.5030   -4.0853    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    6.8613   -5.9478    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    7.4989   -6.3186    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    6.2197   -5.6621    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    5.8063   -6.3781    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    5.3197   -5.7853    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    5.3155   -3.1478    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    6.8572   -3.7145    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
    6.8572   -2.9728    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
    4.1827   -3.4223    0.0000 H   0  0  0  0  0  0  0  0  0  0  0  0
    8.1447   -5.6632    0.0000 H   0  0  0  0  0  0  0  0  0  0  0  0
    9.4278   -5.6632    0.0000 H   0  0  0  0  0  0  0  0  0  0  0  0
    7.4974   -6.7840    0.0000 H   0  0  0  0  0  0  0  0  0  0  0  0
    5.3409   -6.3781    0.0000 H   0  0  0  0  0  0  0  0  0  0  0  0
    7.2602   -2.7401    0.0000 H   0  0  0  0  0  0  0  0  0  0  0  0
  1  2  1  0  0  0  0
  2  3  1  0  0  0  0
  3  4  1  0  0  0  0
  4  5  1  0  0  0  0
  3  6  1  0  0  0  0
  6  7  1  0  0  0  0
```

	1 7 1 0 0 0 0 7 8 1 6 0 0 0 8 9 1 0 0 0 0 9 10 1 0 0 0 0 10 11 1 0 0 0 0 11 12 1 0 0 0 0 9 13 2 0 0 0 0 8 14 1 0 0 0 0 14 15 1 0 0 0 0 7 16 1 0 0 0 0 16 17 1 0 0 0 0 6 18 2 0 0 0 0 2 19 2 0 0 0 0 1 20 1 0 0 0 0 20 21 1 0 0 0 0 5 22 1 0 0 0 0 10 23 1 0 0 0 0 12 24 1 0 0 0 0 15 25 1 0 0 0 0 17 26 1 0 0 0 0 21 27 1 0 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C9H16N4O8/c14-1-9(12(4-17)7(20)10-2-15)6(19)11(3-16)8(21)13(9)5-18/h14-18H,1-5H2,(H,10,20)/t9-/m0/s1
Chemical Structure, InChIKey	BEEHUIXXNCWWPG-VIFPVBQESA-N

Part2Moiety (mixture component)

Name		**Quantity**	1

Moiety Characteristics

Chemical Structure, MOLFILE	C9H16N4O8 -FDASRS-02211411142D 27 27 0 0 0 0 0 0 0 0999 V2000 6.2197 -4.0936 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 5.5113 -3.8686 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 5.0738 -4.4728 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 4.3322 -4.4728 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 3.9530 -3.8270 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 5.5155 -5.0728 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 6.2197 -4.8353 0.0000 C 0 0 2 0 0 0 0 0 0 0 0 0 6.8613 -5.2061 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 7.5030 -4.8270 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 8.1447 -5.1978 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 8.7863 -4.8270 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 9.4280 -5.1978 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 7.5030 -4.0853 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.8613 -5.9478 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 7.4989 -6.3186 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.2197 -5.6621 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 5.8063 -6.3781 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 5.3197 -5.7853 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 5.3155 -3.1478 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.8572 -3.7145 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 6.8572 -2.9728 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 4.1827 -3.4223 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 8.1447 -5.6632 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 9.4278 -5.6632 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 7.4974 -6.7840 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 5.3409 -6.3781 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 7.2602 -2.7401 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 1 2 1 0 0 0 0 2 3 1 0 0 0 0 3 4 1 0 0 0 0 4 5 1 0 0 0 0 3 6 1 0 0 0 0 6 7 1 0 0 0 0 1 7 1 0 0 0 0 7 8 1 1 0 0 0 8 9 1 0 0 0 0 9 10 1 0 0 0 0 10 11 1 0 0 0 0 11 12 1 0 0 0 0 9 13 2 0 0 0 0 8 14 1 0 0 0 0 14 15 1 0 0 0 0 7 16 1 0 0 0 0 16 17 1 0 0 0 0 6 18 2 0 0 0 0 2 19 2 0 0 0 0 1 20 1 0 0 0 0 20 21 1 0 0 0 0 5 22 1 0 0 0 0

	10 23 1 0 0 0 0 12 24 1 0 0 0 0 15 25 1 0 0 0 0 17 26 1 0 0 0 0 21 27 1 0 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C9H16N4O8/c14-1-9(12(4-17)7(20)10-2-15)6(19)11(3-16)8(21)13(9)5-18/h14-18H,1-5H2,(H,10,20)/t9-/m1/s1
Chemical Structure, InChIKey	BEEHUIXXNCWWPG-SECBINFHSA-N

Moiety (mixture component)

Name		**Quantity**	Undefined not 0

Part1Moiety (mixture component)

Name		**Quantity**	1

Moiety Characteristics

Chemical Structure, MOLFILE	C8H14N4O7 -FDASRS-02211411142D 25 25 0 0 0 0 0 0 0 0999 V2000 6.4042 -0.3625 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 5.1375 -0.7833 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 5.6208 -0.1125 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 6.4125 -1.1875 0.0000 C 0 0 1 0 0 0 0 0 0 0 0 0 5.6292 -1.4500 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 7.1250 -1.6000 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 7.8375 -1.1792 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 8.5500 -1.5917 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 7.1167 0.0583 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 5.4042 0.6875 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 5.4125 -2.2417 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 4.3125 -0.7792 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 7.8292 -0.3542 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 7.1167 -2.4250 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 9.2625 -1.1792 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 7.1167 0.8833 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 3.8917 -0.0667 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 7.8292 -2.8375 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 9.9750 -1.5917 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.3611 -1.6965 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 8.5500 -2.1033 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 7.5597 1.1391 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 4.1431 0.3788 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 7.8286 -3.3491 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 9.9744 -2.1033 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 2 3 1 0 0 0 0 1 3 1 0 0 0 0 1 4 1 0 0 0 0 4 5 1 0 0 0 0 4 6 1 6 0 0 0 6 7 1 0 0 0 0 7 8 1 0 0 0 0 1 9 1 0 0 0 0 3 10 2 0 0 0 0 5 11 2 0 0 0 0 2 12 1 0 0 0 0 7 13 2 0 0 0 0 6 14 1 0 0 0 0 8 15 1 0 0 0 0 9 16 1 0 0 0 0 12 17 1 0 0 0 0 14 18 1 0 0 0 0 15 19 1 0 0 0 0 2 5 1 0 0 0 0 4 20 1 0 0 0 0 8 21 1 0 0 0 0 16 22 1 0 0 0 0 17 23 1 0 0 0 0 18 24 1 0 0 0 0 19 25 1 0 0 0 0 M END

Chemical Structures, InChI	InChI=1S/C8H14N4O7/c13-1-9-7(18)10(2-14)5-6(17)12(4-16)8(19)11(5)3-15/h5,13-16H,1-4H2,(H,9,18)/t5-/m0/s1
Chemical Structure, InChIKey	SOROIESOUPGGFO-YFKPBYRVSA-N

Part2Moiety (mixture component)

Name		**Quantity**	1

Moiety Characteristics

Chemical Structure, MOLFILE	C8H14N4O7 -FDASRS-02211411142D 25 25 0 0 0 0 0 0 0 0999 V2000 6.4042 -0.3625 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 5.1375 -0.7833 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 5.6208 -0.1125 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 6.4125 -1.1875 0.0000 C 0 0 2 0 0 0 0 0 0 0 0 0 5.6292 -1.4500 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 7.1250 -1.6000 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 7.8375 -1.1792 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 8.5500 -1.5917 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 7.1167 0.0583 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 5.4042 0.6875 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 5.4125 -2.2417 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 4.3125 -0.7792 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 7.8292 -0.3542 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 7.1167 -2.4250 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 9.2625 -1.1792 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 7.1167 0.8833 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 3.8917 -0.0667 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 7.8292 -2.8375 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 9.9750 -1.5917 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.3611 -1.6965 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 8.5500 -2.1033 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 7.5597 1.1391 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 4.1431 0.3788 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 7.8286 -3.3491 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 9.9744 -2.1033 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 2 3 1 0 0 0 0 1 3 1 0 0 0 0 1 4 1 0 0 0 0 4 5 1 0 0 0 0 4 6 1 0 0 0 0 6 7 1 0 0 0 0 7 8 1 0 0 0 0 1 9 1 0 0 0 0 3 10 2 0 0 0 0 5 11 2 0 0 0 0 2 12 1 0 0 0 0 7 13 2 0 0 0 0 6 14 1 0 0 0 0 8 15 1 0 0 0 0 9 16 1 0 0 0 0 12 17 1 0 0 0 0 14 18 1 0 0 0 0 15 19 1 0 0 0 0 2 5 1 0 0 0 0 4 20 1 6 0 0 0 8 21 1 0 0 0 0 16 22 1 0 0 0 0 17 23 1 0 0 0 0 18 24 1 0 0 0 0 19 25 1 0 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C8H14N4O7/c13-1-9-7(18)10(2-14)5-6(17)12(4-16)8(19)11(5)3-15/h5,13-16H,1-4H2,(H,9,18)/t5-/m1/s1
Chemical Structure, InChIKey	SOROIESOUPGGFO-RXMQYKEDSA-N

Moiety (mixture component)

Name		**Quantity**	Undefined not 0

Part1Moiety (mixture component)

Name		Quantity	1

Moiety Characteristics

Chemical Structure, MOLFILE	C8H14N4O7 -FDASRS-02211411142D 25 25 0 0 0 0 0 0 0 0999 V2000 6.2825 -8.3442 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 5.5825 -8.1234 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 5.1491 -8.7192 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 5.5867 -9.3151 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 6.2825 -9.0817 0.0000 C 0 0 1 0 0 0 0 0 0 0 0 0 6.9200 -9.4484 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 7.5533 -9.0734 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 8.1909 -9.4401 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 8.8241 -9.0734 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 9.4575 -9.4401 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 7.5533 -8.3359 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.9200 -10.1817 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 7.5492 -10.5484 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.2825 -9.9002 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 5.8733 -10.6079 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 5.3909 -10.0192 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 5.3867 -7.4109 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.9159 -7.9693 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 6.9159 -7.2359 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 4.6881 -8.7202 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 8.1916 -9.9011 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 9.4570 -9.9011 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 7.5473 -11.0094 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 5.4123 -10.6076 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 7.3151 -7.0054 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 1 2 1 0 0 0 0 2 3 1 0 0 0 0 3 4 1 0 0 0 0 4 5 1 0 0 0 0 1 5 1 0 0 0 0 5 6 1 6 0 0 0 6 7 1 0 0 0 0 7 8 1 0 0 0 0 8 9 1 0 0 0 0 9 10 1 0 0 0 0 7 11 2 0 0 0 0 6 12 1 0 0 0 0 12 13 1 0 0 0 0 5 14 1 0 0 0 0 14 15 1 0 0 0 0 4 16 2 0 0 0 0 2 17 2 0 0 0 0 1 18 1 0 0 0 0 18 19 1 0 0 0 0 3 20 1 0 0 0 0 8 21 1 0 0 0 0 10 22 1 0 0 0 0 13 23 1 0 0 0 0 15 24 1 0 0 0 0 19 25 1 0 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C8H14N4O7/c13-1-8(11(3-15)6(18)9-2-14)5(17)10-7(19)12(8)4-16/h13-16H,1-4H2,(H,9,18)(H,10,17,19)/t8-/m0/s1
Chemical Structure, InChIKey	CHUZDWOCSGISNS-QMMMGPOBSA-N

Part2Moiety (mixture component)

Name		Quantity	1

Moiety Characteristics

	C8H14N4O7 -FDASRS-02211411142D 25 25 0 0 0 0 0 0 0 0999 V2000 6.2825 -8.3442 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 5.5825 -8.1234 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 5.1491 -8.7192 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 5.5867 -9.3151 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 6.2825 -9.0817 0.0000 C 0 0 2 0 0 0 0 0 0 0 0 0 6.9200 -9.4484 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0

Chemical Structure, MOLFILE	7.5533 -9.0734 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 8.1909 -9.4401 0.0000 N 0 0 0 0 0 0 0 0 0 0 0 0 8.8241 -9.0734 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 9.4575 -9.4401 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 7.5533 -8.3359 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.9200 -10.1817 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 7.5492 -10.5484 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.2825 -9.9002 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 5.8733 -10.6079 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 5.3909 -10.0192 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 5.3867 -7.4109 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 6.9159 -7.9693 0.0000 C 0 0 0 0 0 0 0 0 0 0 0 0 6.9159 -7.2359 0.0000 O 0 0 0 0 0 0 0 0 0 0 0 0 4.6881 -8.7202 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 8.1916 -9.9011 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 9.4570 -9.9011 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 7.5473 -11.0094 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 5.4123 -10.6076 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 7.3151 -7.0054 0.0000 H 0 0 0 0 0 0 0 0 0 0 0 0 1 2 1 0 0 0 0 2 3 1 0 0 0 0 3 4 1 0 0 0 0 4 5 1 0 0 0 0 1 5 1 0 0 0 0 5 6 1 1 0 0 0 6 7 1 0 0 0 0 7 8 1 0 0 0 0 8 9 1 0 0 0 0 9 10 1 0 0 0 0 7 11 2 0 0 0 0 6 12 1 0 0 0 0 12 13 1 0 0 0 0 5 14 1 0 0 0 0 14 15 1 0 0 0 0 4 16 2 0 0 0 0 2 17 2 0 0 0 0 1 18 1 0 0 0 0 18 19 1 0 0 0 0 3 20 1 0 0 0 0 8 21 1 0 0 0 0 10 22 1 0 0 0 0 13 23 1 0 0 0 0 15 24 1 0 0 0 0 19 25 1 0 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C8H14N4O7/c13-1-8(11(3-15)6(18)9-2-14)5(17)10-7(19)12(8)4-16/h13-16H,1-4H2,(H,9,18)(H,10,17,19)/t8-/m1/s1
Chemical Structure, InChIKey	CHUZDWOCSGISNS-MRVPVSSYSA-N

Molecular Bond Types

Product Information

Product Type	INDEXING - SUBSTANCE	**Item Code (Source)**	H5RIZ3MPW4

Labeler - Food and Drug Administration (927645523)