Indexing - Substance

Food and Drug Administration
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LEVOLEUCOVORIN (990S25980Y)

Alias Name

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

Substance Information

Mapping	**Definition Hash:**	2a36e7eb-2278-ef77-f558-b159284eb839

Moiety

Name		**Quantity**	1

Moiety Characteristics

Chemical Structure, MOLFILE

```
  -FDASRS-09061311012D

34 36  0  0  1  0  0  0  0  0999 V2000
   8.3368   -5.3091    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   9.3517   -4.7093    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
  10.3731   -6.4827    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
  10.3732   -5.3011    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
  13.4349   -4.7125    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   7.3152   -4.7174    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
  14.4535   -5.3175    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
  13.4500   -3.5322    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
  11.3921   -4.7002    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
  12.4161   -5.3018    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   8.3433   -6.4883    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
   0.1125   -3.0709    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   0.1125   -1.8792    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
  -1.9499   -3.0541    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
  -0.9166   -3.6625    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
  -0.9166   -1.2833    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
   1.1292   -3.6625    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
  -1.9499   -1.8834    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   1.1292   -1.2833    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
   2.1667   -3.0541    0.0000 C   0  0  2  0  0  0  0  0  0  0  0  0
   1.1292   -4.8417    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
  -0.9083   -4.8417    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
   2.1667   -1.8834    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   3.1959   -3.6250    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   4.2042   -3.0042    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
  -2.9874   -1.3167    0.0000 N   0  0  0  0  0  0  0  0  0  0  0  0
   0.1083   -5.4376    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
   7.2833   -3.5458    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   6.3000   -5.3375    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   5.2418   -3.5751    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   6.2500   -2.9666    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   5.2667   -4.7709    0.0000 C   0  0  0  0  0  0  0  0  0  0  0  0
   9.3517   -7.0764    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
  11.3942   -7.0807    0.0000 O   0  0  0  0  0  0  0  0  0  0  0  0
 2  1  1  0  0  0  0
 4  3  1  1  0  0  0
```

	4 2 1 0 0 0 0 5 10 1 0 0 0 0 6 1 1 0 0 0 0 7 5 1 0 0 0 0 8 5 2 0 0 0 0 9 4 1 0 0 0 0 10 9 1 0 0 0 0 1 11 2 0 0 0 0 13 12 2 0 0 0 0 14 15 1 0 0 0 0 15 12 1 0 0 0 0 16 13 1 0 0 0 0 17 12 1 0 0 0 0 18 14 2 0 0 0 0 19 13 1 0 0 0 0 20 17 1 0 0 0 0 21 17 1 0 0 0 0 22 15 2 0 0 0 0 6 29 2 0 0 0 0 23 20 1 0 0 0 0 20 24 1 6 0 0 0 25 24 1 0 0 0 0 26 18 1 0 0 0 0 27 21 2 0 0 0 0 28 31 2 0 0 0 0 29 32 1 0 0 0 0 30 25 1 0 0 0 0 31 30 1 0 0 0 0 32 30 2 0 0 0 0 23 19 1 0 0 0 0 18 16 1 0 0 0 0 28 6 1 0 0 0 0 3 33 2 0 0 0 0 3 34 1 0 0 0 0 M END
Chemical Structures, InChI	InChI=1S/C20H23N7O7/c21-20-25-16-15(18(32)26-20)27(9-28)12(8-23-16)7-22-11-3-1-10(2-4-11)17(31)24-13(19(33)34)5-6-14(29)30/h1-4,9,12-13,22H,5-8H2,(H,24,31)(H,29,30)(H,33,34)(H4,21,23,25,26,32)/t12-,13-/m0/s1
Chemical Structure, InChI Key	VVIAGPKUTFNRDU-STQMWFEESA-N

Molecular Bond Types

Product Information

Product Type	INDEXING - SUBSTANCE	**Item Code (Source)**	990S25980Y

Labeler - Food and Drug Administration (927645523)