
Chemical Structure, MOLFILE	- FDASRS - 11061311022D																			
	16	16	0	0	0	0	0	0	0	0999	V2000									
	4.9834	-1.4375	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	3.8376	-2.0458	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	6.1376	-2.1083	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	8.4542	-0.6334	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	2.1709	-0.9750	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	6.1376	-3.4375	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	3.8376	-3.4375	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	4.9834	-4.1083	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	4.9834	-0.1083	0.0000	O	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	2.9334	0.1042	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	1.1667	-0.1250	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	1.2292	-1.9250	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	9.2666	-1.7167	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	9.5750	0.0958	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	7.6792	0.4458	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	4.9834	-5.4375	0.0000	C	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	2	1	2	0	0	0	0													
	3	1	1	0	0	0	0													
	4	3	1	0	0	0	0													
	5	2	1	0	0	0	0													
	6	3	2	0	0	0	0													
	7	2	1	0	0	0	0													
	8	6	1	0	0	0	0													
	9	1	1	0	0	0	0													
	10	5	1	0	0	0	0													
	11	5	1	0	0	0	0													
	12	5	1	0	0	0	0													
	13	4	1	0	0	0	0													
	14	4	1	0	0	0	0													
	15	4	1	0	0	0	0													
	16	8	1	0	0	0	0													
7	8	2	0	0	0	0														
M	END																			
Chemical Structures, InChI	InChI=1S/C15H24O/c1-10-8-11(14(2,3)4)13(16)12(9-10)15(5,6)7/h8-9,16H,1-7H3																			

Chemical Structure, InChI Key		NLZUEZXRPGMBCV-UHFFFAOYSA-N	
Molecular Bond Types			
Product Information			
Product Type	INDEXING - SUBSTANCE	Item Code (Source)	1P9D0Z171K

Labeler - Food and Drug Administration (927645523)