

Aflatoxin G2

Other names:

1H,12H-Furo[3',2':4,5]furo[2,3-h]pyrano[3,4-c][1]benzopyran-1,12-dione,
3,4,7a,9,10,10a-hexahydro-5-methoxy-, (7aR,cis)-
1H,12H-Furo[3',2':4,5]furo[2,3-h]pyrano[3,4-c][1]benzopyran-1,12-dione,
3,4,7a«alpha»,9,10,10a«alpha»-hexahydro-5-methoxy-
Dihydroaflatoxin G1

(7aR,cis)3,4,7a,9,10,10a-hexahydro-5-methoxy-1H,12H-furo[3',2':4,5]furo[2,3-h]pyrano[3,

Inchi:

InChI=1S/C17H14O7/c1-20-9-6-10-12(8-3-5-22-17(8)23-10)14-11(9)7-2-4-21-15(18)13(7

InchiKey:

WPCVRWVBBXIRMA-UHFFFAOYSA-N

Formula:

C17H14O7

SMILES:

COc1cc2c(c3oc(=O)c4c(c13)CCOC4=O)C1CCOC1O2

Mol. weight [g/mol]:

330.29

CAS:

7241-98-7

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.86		Crippen Method
logp	1.737		Crippen Method
mcvol	211.380	ml/mol	McGowan Method

Sources

Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Crippen Method:

https://www.chemeo.com/doc/models/crippen_log10ws

McGowan Method:

<http://link.springer.com/article/10.1007/BF02311772>

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C7241987&Units=SI>

Legend

log10ws: Log10 of Water solubility in mol/l

logp: Octanol/Water partition coefficient

mcvol: McGowan's characteristic volume

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