

4H-1-BENZOPYRANE-4-ONE, 2-(3,4-DIHYDROXYPHENYL)-5,7-DIHYDROXY-

Other names:	2-(3,4-dihydroxyphenyl)-5,7-dihydroxy-4-benzopyrone 4H-1-Benzopyran-4-one, 2-(3,4-dihydroxyphenyl)-5,7-dihydroxy- 5,7,3',4'-tetrahydroxyflavone Digitoflavone Flacitran Flavone, 3',4',5,7-tetrahydroxy- Luteolin Weld Lake luteoline
Inchi:	InChI=1S/C15H10O6/c16-8-4-11(19)15-12(20)6-13(21-14(15)5-8)7-1-2-9(17)10(18)3-7/h
InchiKey:	IQPNAANSBPBGFQ-UHFFFAOYSA-N
Formula:	C15H10O6
SMILES:	O=c1cc(-c2ccc(O)c(O)c2)oc2cc(O)cc(O)c12
Mol. weight [g/mol]:	286.24

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.67		Cheméo Relay (1.0)
logp	2.282		Crippen Method
mcvol	190.450	ml/mol	McGowan Method
tf	610.53	K	Measurement and correlation of the solubilities of luteolin and rutin in five imidazole-based ionic liquids
tf	601.00	K	Measurement and Correlation of Solubilities of Luteolin in Organic Solvents at Different Temperatures
tf	610.58	K	Solubility of Luteolin in Ethanol + Water Mixed Solvents at Different Temperatures

Sources

Cheméo Relay (1.0, beta):

Solubility of Luteolin in Ethanol + Water <https://www.doi.org/10.1021/je900381r>

Mixed Solvents at Different
Temperatures:

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

Measurement and correlation of the
solubilities of luteolin and rutin in five
imidazole based ionic liquids:

<https://www.doi.org/10.1016/j.fluid.2013.01.026>

McGowan Method:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C491703&Units=SI>

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

Measurement and Correlation of
Solubilities of Luteolin in Organic
Solvents at Different Temperatures:

<https://www.doi.org/10.1021/je060133l>

<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

Cheméo Relay (1.0):

Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
tf:	Normal melting (fusion) point

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