

# cis-Captafol

<b>Other names:</b>	1H-Isoindole-1,3(2H)-dione, 3a,4,7,7a-tetrahydro-2-[(1,1,2,2-tetrachloroethyl)thio]-, cis- 4-Cyclohexene-1,2-dicarboximide, N-[(1,1,2,2-tetrachloroethyl)thio]-, cis- N-(1,1,2,2-Tetrachloro-ethylsulfenyl)cis-4-cyclohexene-1,2-dicarboximide TN 80 Crisfolatan Mycodifol Pillartan cis-N-((1,1,2,2-Tetrachloroethyl)thio)-4-cyclohexene-1,2-dicarboximide cis-3a,4,7,7a-Tetrahydro-2-((1,1,2,2-tetrachloroethyl)thio)-1H-isoindole-1,3(2H)-dione N-(Tetrachloroethylthio)-«DELTA»4-tetrahydrophthalimide Captafol, cis- 1H-Isoindole-1,3(2H)-dione, 3a,4,7,7a-tetrahydro-2-((1,1,2,2-tetrachloroethyl)thio)-, (3aR,7aS)-rel- TN 80 (pesticide) cis-N-((1,1,2,2-Tetrachloroethyl)thio)-4-cyclohexene-1,2-dicarboximide
<b>Inchi:</b>	InChI=1S/C10H9Cl4NO2S/c11-9(12)10(13,14)18-15-7(16)5-3-1-2-4-6(5)8(15)17/h1-2,5-6
<b>InchiKey:</b>	JHRWWRDRBPCWTF-OLQVQODUSA-N
<b>Formula:</b>	C10H9Cl4NO2S
<b>SMILES:</b>	O=C1C2CC=CCC2C(=O)N1SC(Cl)(Cl)C(Cl)Cl
<b>Mol. weight [g/mol]:</b>	349.06
<b>CAS:</b>	2939-80-2

## Physical Properties

Property code	Value	Unit	Source
log10ws	-4.48		Crippen Method
logp	3.521		Crippen Method
mvol	204.170	ml/mol	McGowan Method

## Sources

<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C2939802&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C2939802&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>

# Legend

<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume

Latest version available from:

<https://www.cheméo.com/cid/12-662-3/cis-Captafol.pdf>

Generated by Cheméo on 2024-05-16 10:20:46.220577366 +0000 UTC m=+18144095.141154677.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.