

Iprodione

Other names:	1-Imidazolidinecarboxamide, 3-(3,5-dichlorophenyl)-N-(1-methylethyl)-2,4-dioxo-1-Isopropyl carbamoyl-3-(3,5-dichlorophenyl)-hydantoin 3-(3,5-Dichlorophenyl)-1-(1-methylethyl)carbamoylhydantion 3-(3,5-Dichlorophenyl)-N-(1-methylethyl)-2,4-dioxo-1-imidazolidinecarboxamide 3-(3,5-dichlorophenyl)-2,4-dioxo-N-isopropylimidazolidine-1-carboxamide 3-(3,5-dichlorophenyl)-2,4-dioxo-N-propan-2-ylimidazolidine-1-carboxamide Chipco 26019 FA 2071 Glycophen Glycophene Iprodial Kidan LFA 2043 MRC 910 NRC 910 Promidione ROP 500 F RP 26019 Rovral Rovral 50WP Rovral Flo Rovral PM Verisan
Inchi:	InChI=1S/C13H13Cl2N3O3/c1-7(2)16-12(20)17-6-11(19)18(13(17)21)10-4-8(14)3-9(15)5
InchiKey:	ONUFESLQCSAYKA-UHFFFAOYSA-N
Formula:	C13H13Cl2N3O3
SMILES:	CC(C)NC(=O)N1CC(=O)N(c2cc(Cl)cc(Cl)c2)C1=O
Mol. weight [g/mol]:	330.17
CAS:	36734-19-7

Physical Properties

Property code	Value	Unit	Source
log10ws	-4.38		Aqueous Solubility Prediction Method
log10ws	-4.38		Estimated Solubility Method
logp	2.880		Crippen Method

mcvol	218.540	ml/mol	McGowan Method
rinpol	2439.00		NIST Webbook
rinpol	2452.00		NIST Webbook
rinpol	2439.00		NIST Webbook
rinpol	2452.00		NIST Webbook

Sources

Estimated Solubility Method: http://pubs.acs.org/doi/suppl/10.1021/ci034243x/suppl_file/ci034243xsi20040112_053635.txt

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

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C36734197&Units=SI>

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

Aqueous Solubility Prediction Method: <http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa>

Legend

log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
m_{cvol}:	McGowan's characteristic volume
rin_{pol}:	Non-polar retention indices

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