

Glutaric acid, hex-4-yn-3-yl 3-ethylphenyl ester

Inchi:	InChI=1S/C19H24O4/c1-4-9-16(6-3)22-18(20)12-8-13-19(21)23-17-11-7-10-15(5-2)14-17
InchiKey:	OXURDZWSFVNBEH-UHFFFAOYSA-N
Formula:	C19H24O4
SMILES:	CC#CC(CC)OC(=O)CCCC(=O)Oc1cccc(CC)c1
Mol. weight [g/mol]:	316.39

Physical Properties

Property code	Value	Unit	Source
gf	-55.60	kJ/mol	Joback Method
hf	-433.01	kJ/mol	Joback Method
hfus	43.79	kJ/mol	Joback Method
hvap	80.90	kJ/mol	Joback Method
log10ws	-5.12		Crippen Method
logp	3.670		Crippen Method
mcvol	261.090	ml/mol	McGowan Method
pc	1644.42	kPa	Joback Method
rinpola	2285.00		NIST Webbook
tb	826.92	K	Joback Method
tc	1042.78	K	Joback Method
tf	578.25	K	Joback Method
vc	0.996	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	768.81	J/molxK	826.92	Joback Method
cpg	784.24	J/molxK	862.90	Joback Method
cpg	798.50	J/molxK	898.87	Joback Method
cpg	811.60	J/molxK	934.85	Joback Method
cpg	823.57	J/molxK	970.82	Joback Method
cpg	834.41	J/molxK	1006.80	Joback Method
cpg	844.15	J/molxK	1042.78	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U390891&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpola:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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