

tritium

Inchi: InChI=1S/H2/h1H/i1+2T
InchiKey: YZCKVEUIGOORGS-NJFSPNSNSA-N
Formula: H2
SMILES: [3H][3H]
Mol. weight [g/mol]: 6.03

Physical Properties

Property code	Value	Unit	Source
hfus	-0.88	kJ/mol	Joback Method
logp	0.246		Crippen Method
mcvol	10.860	ml/mol	McGowan Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	-10.08	J/mol×K	198.00	Joback Method
cpg	-8.36	J/mol×K	221.51	Joback Method
cpg	-6.92	J/mol×K	245.01	Joback Method
cpg	-5.74	J/mol×K	268.52	Joback Method
cpg	-4.82	J/mol×K	292.03	Joback Method
cpg	-4.12	J/mol×K	315.53	Joback Method
cpg	-3.65	J/mol×K	339.04	Joback Method
dvisc	0.0000006	Pa×s	122.50	Joback Method
dvisc	0.0000010	Pa×s	135.08	Joback Method
dvisc	0.0000014	Pa×s	147.67	Joback Method
dvisc	0.0000020	Pa×s	160.25	Joback Method
dvisc	0.0000026	Pa×s	172.83	Joback Method
dvisc	0.0000033	Pa×s	185.42	Joback Method
dvisc	0.0000040	Pa×s	198.00	Joback Method

Sources

Cheméo Relay (1.0):

Cheméo Relay (1.0, beta):

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C10028178&Units=SI>

Joback Method:

https://en.wikipedia.org/wiki/Joback_method

McGowan Method:

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

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
dvisc:	Dynamic viscosity
hfus:	Enthalpy of fusion at standard conditions
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume

Latest version available from:

<https://www.chemeo.com/cid/34-545-9/tritium.pdf>

Generated by Cheméo on 2026-07-22 05:03:09.319264239 +0000 UTC m=+207685.161149617.

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