

NAME

iapws - Compute light and heavy water properties.

SYNOPSIS

iapws *SUBCOMMAND* [*OPTION* . . .]

DESCRIPTION

iapws is a command line interface which computes the properties of light and heavy water according to IAPWS.

SUBCOMMANDS

Valid subcommands are:

- +kh** Compute the Henry's constant for gases in H₂O or D₂O. The default behavior is to compute the constant kH for O₂ at 25°C. See options.
- +kd** Compute the vapor-liquid distribution constant for gases in H₂O or D₂O. The default behavior is to compute the constant kD for H₂ at 25°C. See options.

Their syntax is:

- +kh** [OPTION...]
- +kd** [OPTION...]

OPTIONS

kh:

- temperature, -T TEMPERATURE...**
Temperature in °C. Default to 25°C.
- fugacity, -f FUGACITY...**
Liquid-phase fugacity in MPa. Default to 0.1
- gases, -g GAS...**
Gases for which to compute kH. Default to O₂
- D2O** Set heavywater as the solvent.
- listgases**
Display available gases for computing kH.

kd:

- temperature, -T TEMPERATURE...**
Temperature in °C. Default to 25°C.
- x2, -x x2...**
Molar fraction of gas in water. Default to 1
- gases, -g GAS...**
Gases for which to compute kD. Default to H₂
- D2O,** Set heavywater as the solvent.
- listgases**
Display available gases for computing kD.

all:

- usage, -u**
Show usage text and exit.
- help, -h**
Show help text and exit.

--verbose, -V

Display additional information when availabl

--version, -v

Show version information and exit.

EXAMPLE

Minimal example

```
iapws
```

SEE ALSO

***ciaaw*(3), *codata*(3)**