

hp02ver.dat thermodynamic data file

no_print | print generates print output

plot | obsolete 6.8.4+

solution_model.dat | solution model file, blank = none

Title Text

perplex_option.dat | Perple_X option file

5 calculation type: 0- composition, 1- Schreinemakers, 3- Mixed, 4- swash, 5- gridded min, 7- 1d fract, 8- gwash, 9- 2d fract, 10- 7 w/file input, 11- 9 w/file input, 12- 0d infiltration

0 unused place holder, post 06

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0 number component transformations

15 number of components in the data base

1 component amounts, 0 - mole, 1 mass

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0 ifug EoS for saturated phase

2 gridded minimization dimension (1 or 2)

0 special dependencies: 0 - P and T independent, 1 - P(T), 2 - T(P)

0.00000 0.00000 0.00000 0.00000 0.00000 Geothermal gradient
polynomial coeffs.

begin thermodynamic component list

SiO2	1	55.3200	0.00000	0.00000	mass	amount
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Al2O3	1	1.09000	0.00000	0.00000	mass	amount
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FeO	1	37.4300	0.00000	0.00000	mass	amount
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MgO	1	3.03000	0.00000	0.00000	mass	amount
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CaO	1	1.09000	0.00000	0.00000	mass	amount
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O2	1	2.00000	0.00000	0.00000	mass	amount
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end thermodynamic component list

```
begin saturated component list
end saturated component list
```

```
begin saturated phase component list
end saturated phase component list
```

```
begin independent potential/fugacity/activity list
end independent potential list
```

```
begin excluded phase list
end excluded phase list
```

```
begin solution phase list
Cpx(h)
Opx(JH)
MtUl(A)
end solution phase list
```

15000.0	1200.00	0.00000	0.00000	0.00000	max p, t,
xco2, mu_1, mu_2					
3000.00	700.000	0.00000	0.00000	0.00000	min p, t,
xco2, mu_1, mu_2					
0.00000	0.00000	0.00000	0.00000	0.00000	unused
place holder post 06					

2 1 4 5 3 indices of 1st & 2nd independent & sectioning variables