

# P-Ti (Phosphorus-Titanium)

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The Ti-P phase diagram in [Massalski2] was redrawn from [1987Mur]. The phosphorus-rich part (>55 at.% P) unknown in [1987Mur] was estimated by [2006Oht] by thermodynamic modeling. The result is shown in Fig. 1.

Ti-P crystal structure data are shown in Table 1.

Murray, Ed., ASM International, Metals Park, OH, 1987, p 234-236

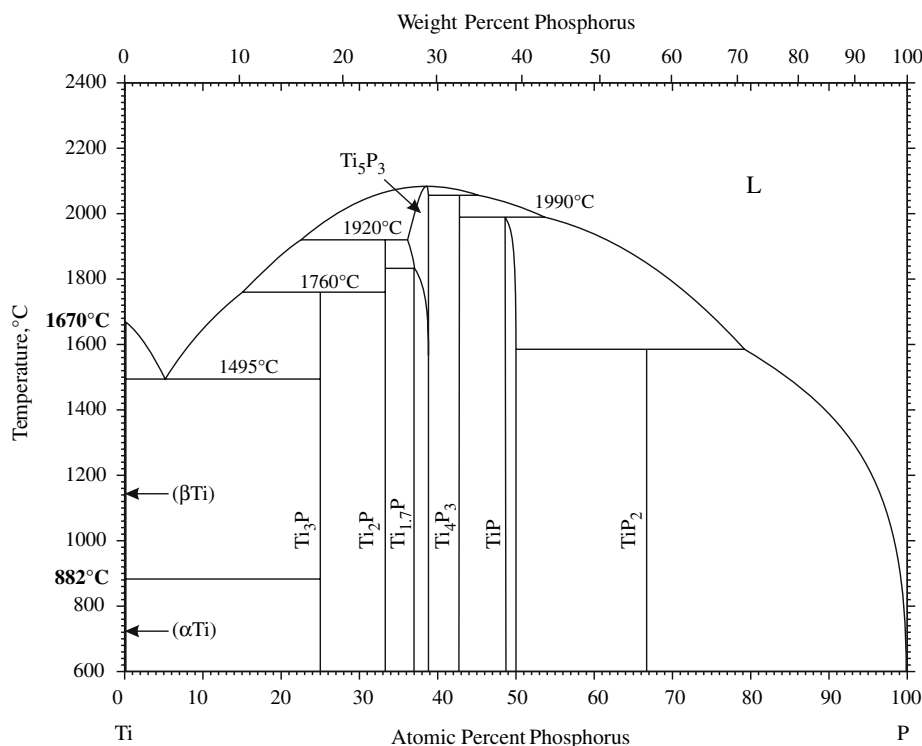
**2006Oht:** H. Ohtani, N. Hanaya, M. Hasebe, S. Teraoka, and M. Abe, Thermodynamic Analysis of the Fe-Ti-P Ternary System by Incorporating First-Principles Calculations into the CALPHAD Approach, *CALPHAD*, 2006, **30**(2), p 147-158

## References

**1987Mur:** J.L. Murray, The P-Ti (Phosphorus-Titanium) System, in *Phase Diagrams of Binary Titanium Alloys*, J.L.

**Table 1** Ti-P crystal structure data

Phase	Composition, at.% P	Pearson symbol	Space group	Struktur bericht designation	Prototype
(βTi)	0	<i>cI2</i>	<i>Im<math>\bar{3}m</math></i>	<i>A2</i>	W
(αTi)	0	<i>hP2</i>	<i>P6<sub>3</sub>/mmc</i>	<i>A3</i>	Mg
Ti <sub>3</sub> P	25	<i>tP32</i>	<i>P4<sub>2</sub>/n</i>	...	Ti <sub>3</sub> P
Ti <sub>2</sub> P	33.3	<i>hP36</i>	<i>P6/mmm</i>	...	...
Ti <sub>1.7</sub> P	37.0	<i>oP*</i>	<i>P2<sub>1</sub>2<sub>1</sub>2<sub>1</sub></i>	...	...
Ti <sub>5</sub> P <sub>3</sub>	36 to 39	<i>hP16</i>	<i>P6<sub>3</sub>/mcm</i>	<i>D8<sub>8</sub></i>	Mn <sub>5</sub> Si <sub>3</sub>
Ti <sub>4</sub> P <sub>3</sub>	42.9	<i>c**</i>	...	...	...
TiP	48 to 50	<i>hP8</i>	<i>P6<sub>3</sub>/mmc</i>	<i>B<sub>i</sub></i>	TiAs
TiP <sub>2</sub>	66.7	<i>tI12</i>	<i>I4/mcm</i>	<i>C16</i>	Al <sub>2</sub> Cu



**Fig. 1** P-Ti phase diagram