substance: OsP_2
property: physical properties

Energy level scheme is similar to that given for $FeAs_2$ (Fig. 1).

energy gap

E_g	≈ 1.2 eV	RT	from diffuse reflectance of sintered samples	63H

magnetic susceptibility
(in 10^{-6} cm^3 mol^{-1}, χ in CGS units)

χ_m	– 74	T = 295 K	powder sample	63H
	– 50	T = 85 K		
	– 97	T = 90...760 K, B < 0.8 T	polycrystalline sample	68H

far infrared absorption: for spectrum in the range 150...500 cm^{-1}, see [77L].

Comparative tables on structural data of transition metal dipnictides:

structure, chemical bond: see document ,

crystallographical data of compounds with octahedrally coordinated cations, see document ,

interatomic distances in marcasite- and loellingite-type compounds, see document .

References:

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Fig. 1.

$FeAs_2$. One-electron energy levels for the valence electrons in loellingite [72F, 72G].