

In the present work, the Rotation method<sup>14</sup> is adopted for  $L_0$ . The single parametric angle  $\phi$  in ( $2 \times 2$ ) dimension was evaluated for the 10 ions (table 1). It is interesting to note in the table 1 that the parametric angle  $\phi$  increases with the increase in the atomic weight of the central atom X. The valence force constants thus evaluated are in good agreement with the values reported by others<sup>3, 4</sup>.

Mean vibrational amplitudes (table 2) and mean square perpendicular amplitudes (table 3) for these ions for various pairs of atomic nuclei were evaluated to check the reliability of the force field model.

**Table 3** Mean Square Perpendicular Amplitudes of Vibration (in Å) of some  $XY_6$  type ions (X = Pb, Pt, Ir, Os, Re, Hf, Ce, Te, Sn, Pd; Y = Cl)

	$\langle X-Y \rangle$	$\langle X-Y \rangle$		$\langle X-Y \rangle$ long
		short	long	
	$\langle X-Y \rangle$	$\Delta x^2$	$\Delta y^2$	$\Delta x^2 = \Delta y^2$
$PdCl_6^{2-}$	0.009882	0.006374	0.029344	0.0071073
$PtCl_6^{2-}$	0.010526	0.010503	0.026030	0.010556
$IrCl_6^{2-}$	0.005946	0.007582	0.014195	0.005827
$OsCl_6^{2-}$	0.007424	0.007538	0.018638	0.007605
$ReCl_6^{2-}$	0.007907	0.007688	0.020028	0.008162
$HfCl_6^{2-}$	0.012404	0.009774	0.035359	0.008464
$CeCl_6^{2-}$	0.013313	0.011740	0.034008	0.014044
$TeCl_6^{2-}$	0.009521	0.013930	0.022432	0.009124
$SnCl_6^{2-}$	0.008008	0.010801	0.020276	0.008260
$PdCl_6^{2-}$	0.006301	0.008852	0.016097	0.006588

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