Erratum: Experimental Vibrational Zero-Point Energies: Diatomic Molecules [J. Phys. Chem. Ref. Data 36, 389-397 (2007)]

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Key words: molecular energetics; uncertainty; vibrational spectroscopy; viscosity; zero-point energy.

There is a typographical error in a state designation for Cl_2^+ on page 390 in the last sentence of Sec. 3. This sentence should read: "For Cl_2^+, only separate constants for $^2\Pi_{3/2g}$ and $^2\Pi_{1/2g}$ were reported.[7]"

There is a transcription error for the harmonic frequency (ω_e) of SiF in Table on page 392. The correct entry for SiF in this table should be:

Molecule	ω_e	$\omega_e x_e$	$\omega_e y_e$	B_e	α_e	Reference
SiF	857.32507(22)	4.83419(9)	0.019807(16)	0.58125735(21)	0.00503859(39)	84

The transcription error for the harmonic frequency (ω_e) of SiF resulted in an error in the calculated zero point energy (ZPE) for SiF in Table 5 on page 395. The correct entry for SiF in this table should be:

Molecule	State designation	ZPE	u_{stat}	u_{trunc}
$^{28}Si^{19}F$	$^2\Pi$	427.6724(2)	0.00019	0.000021

I thank Dr. Russell D. Johnson III for pointing out these errors.