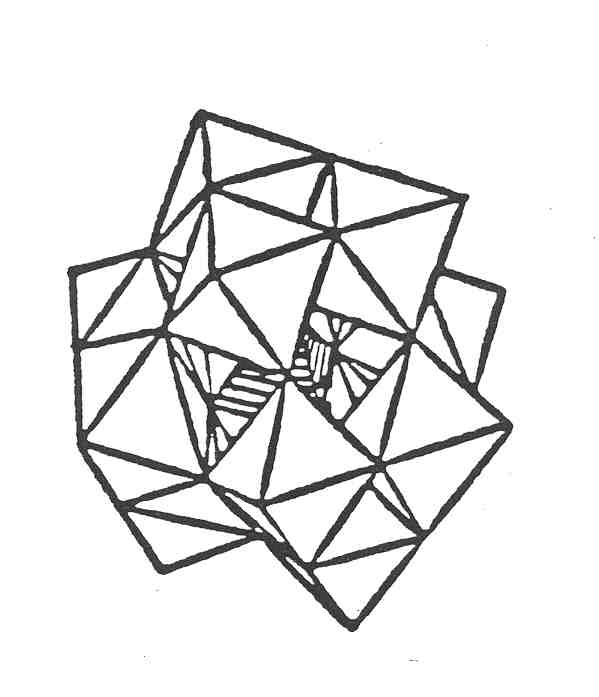
**Application of binomial distribution to interpret 31P NMR for aqueous solution of α-dodecatungstophosphoric acid, α-H3[PW12O40]**

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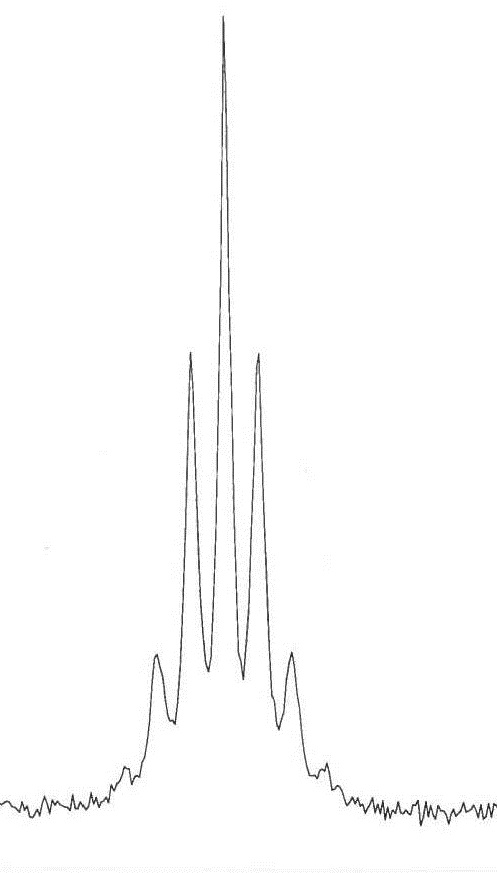
α-Dodecatungstophosphoric acid, H3[PW12O40], is a strong acid whose anion has so-called "Keggin" structure. The anion structure has overall Td symmetry and is based on a central PO4 tetrahedron surrounded by twelve WO6 octahedra, arranged in four groups of three edge-shared octahedra, W3O13. These groups are linked together and to the central PO4 tetrahedron by sharing corners.

The structure is shown below in the polyhedral form in which the tungstens reside in the centers and oxygen atoms are at vertices of the white octahedra. The central tetrahedral PO4 unit is shaded and is surrounded by twelve WO6 octahedra. Note that all tungsten atoms in this structure are identical.



**Figure 1.** Polyhedral representation of α-[PW12O40]3- structure.

31P NMR of aqueous solution of H3[PW12O40] consists of a symmetrical heptet with approximate intensities 1 : 4 : 11 : 16 : 11 : 4 : 1. See Figure 2.



**Figure 2.** 31P NMR of 0.1M H3[PW12O40]*(aq)*

Interpret the spectrum knowing that the only NMR active isotope of tungsten, 183W, with spin ½, is 14.3% abundant. The other isotopes, 180W, 182W, 184W, and 186W, all have a nuclear spin equal to zero. Assume that the only observed coupling is two-bond P-W coupling.

To do your interpretation answer the following questions 1-7.

**1.**  Group all isotopes with zero nuclear spin into one group and mark them all as 184W. Then any [PW12O40]3- anion (PW12) can be written as P(183W)k(184W)12-k. List all PW12 anions which exist in aqueous solution, but differ in isotopic distribution of W. Which anions contribute to the spectrum?

**2.** Write the general expression for the multiplet observed for each individual anion with the formula P(183W)k(184W)12-k.

**3.** Consider the probability mass function used to predict the probability of exactly *k* successes in *n* trials, where success probability in individual trial is *p*.

where

Apply this equation to calculate mole fraction of the PW12 anions with *k* 183W isotopes.

**4.** Recall that the intensities within the multiplets, due to coupling with spin ½ nuclei, follow the Pascal’s triangle.

1

1 1

1 2 1

1 3 3 1

1 4 6 4 1

etc…

For example the quartet has intensities 1 : 3 : 3 : 1.

In order to use the triangle to predict relative intensities of all signals in the observed spectrum, these values need to be normalized, so that the sum of all signals for a given multiplet equals to 1. For example, in the triplet, 1:2:1 needs to be replaced by 0.25:0.5:0.25. Calculate the normalized intensities for the Pascal triangle. Limit your calculations to the multiplets originating from the anions with mole fractions greater than 0.001.

**5.** Combine your answers in 3 and 4 to calculate the relative intensities of all signals within each multiplet.

**6.** All multiplets are centered around the same chemical shift. As a result some signals from different multiplets superimpose. Draw the pattern demonstrating all spectra, showing which signals superimpose. Calculate the expected intensities for all signals in the final spectrum.

**7.** Fully interpret the observed spectrum and the intensities of all peaks.