

INDIAN JOURNAL OF APPLIED RESEARCH 🗴 67

INTRODUCTION

a very simple way.

The method which is generally used<sup>1,2,3,4,5</sup> for the prediction of spin multiplicity value [(2S+1), where  $S = \Sigma s$  = total spin quantum no] is time consuming. To keep the matter in mind a new innovative method has to be introduced for calculation of spin-multiplicity value in the easiest way by ignoring the calculation of total spin quantum number ( $S = \Sigma s$ ). Another method has also to be introduced for resolving magnetic behavior of diatomic hetero nuclear molecules or ions like CO, NO, NO<sup>+</sup>, NO<sup>-</sup>, CN, CN<sup>-</sup> etc. in a very simple way. Another three innovative methods earlier introduced on the easy prediction of 'Hybridization', 'Bond-Order' and 'IUPAC nomenclature of spiro and bicyclo compounds<sup>6,7,8</sup> for the benefit of students.

These new innovative methods would go a long way to help to the students of chemistry who would choose the subject as their career. Experiment in vitro on 100 number of students show that for determination of spin multiplicity value using old (2S + 1) rule, strike rate is 1Q/3min and by using these new innovative methods strike rate is 1Q/5secs. On the basis of this experiment I can strongly recommend that these new methods will be the very rapid one for the determination of spin-multiplicity value and its corresponding spin-state (Table-1) by ignoring the calculation of total spin quantum number (S =  $\Sigma$ s) in (2S + 1) rule.

### New innovative methods for determination of spin multiplicity:

First of all we should classify the species (atoms, molecules, ions or complexes) for which spin multiplicity should be evaluated into three types based on the nature of alignment of unpaired electrons present in them.

a) Species having unpaired electrons alignment upward arrow (1):

In this case, spin multiplicity = (n+1); where n = number of unpaired electrons.



Spin multiplicity = (n + 1) = (1+1) = 2 (spin state = doublet)



Spin multiplicity = (n + 1) = (2+1) = 3 (spin state = triplet)

Spin multiplicity = (n + 1) = (3 + 1) = 4 (spin state = quartet)



Spin multiplicity = (n + 1) = (2 + 1) = 3 (in this case ignore paired electrons) (spin state = triplet)



Spin multiplicity = (n + 1) = (1 + 1) = 2 (spin state = doublet)



Spin multiplicity = (n + 1) = (0 + 1) = 1 (spin state = singlet)

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In this case spin multiplicity = (-n+1); where n = number of unpaired electrons. Here (-ve) sign indicate downward arrow.



Spin multiplicity = (-n + 1) = (-1 + 1) = 0(where n = no of unpaired es)



Spin multiplicity = (-n + 1) = (-2 + 1) = -1



Spin multiplicity = (-n + 1) = (-3 + 1) = -2



Spin multiplicity = (-n + 1) = (-2 + 1) = -1(in this case ignore paired electrons)



Spin multiplicity = (-n + 1) = (-1 + 1) = 0

# c) Species having unpaired electrons alignment both upward and downward arrow

In this case spin multiplicity = [(+n) + (-n) + 1]; where n = number of unpaired electrons in each mode. Here, (+ve) sign indicate upward mode of arrow and (-ve) sign indicate downward mode of arrow.

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## RESEARCH PAPER

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Here total no of unpaired electrons = 2 in which one having upward direction (+1) and other having downward mode (-1).

Hence Spin multiplicity = [(+n) + (-n) + 1] = [(+1) + (-1) + 1] = 1 (spin state = singlet)



Here the total no of unpaired electrons = 3 in which two unpaired electrons lie in upward (+2) and one unpaired electrons lie in downward (-1).

Hence Spin multiplicity = [(+n) + (-n) + 1] = [(+2) + (-1) + 1] = 2 (spin state = doublet)

| $\uparrow \downarrow \downarrow \uparrow \downarrow \downarrow \uparrow$ |
|--------------------------------------------------------------------------|
|--------------------------------------------------------------------------|

Here the total no of unpaired electrons = 5 in which three unpaired electrons lie upward (+3) and two unpaired electrons lie downward (-2) .

Hence Spin multiplicity = [(+n) + (-n) + 1] = [(+3) + (-2) + 1] = 2 (spin state = doublet)

# Table-1: spin multiplicity value and its corresponding spin state

| Spin multiplicity value | Spin state |
|-------------------------|------------|
| 1                       | Singlet    |
| 2                       | Doublet    |
| 3                       | Triplet    |
| 4                       | Quartet    |
| 5                       | Quintet    |

New innovative method for determination of magnetic properties of diatomic heteronuclei molecules or ions: In this case, first of all we have to count the total number of electrons with in the given species, then we isolate the species into two groups on the basis of even and odd no of total electrons present into them. It is generally observed that the species having even number of electrons (total electrons = even no) show diamagnetic character and the species having odd number of electrons (total electrons = odd no) show paramagnetic character (Table-2).

| Table | 2:  | Magnetic  | behavior | of | diatomic | hetero | nuclear |
|-------|-----|-----------|----------|----|----------|--------|---------|
| molec | ule | s or ions |          |    |          |        |         |

| Molecules / ions | Total no of electrons | Magnetic behavior |
|------------------|-----------------------|-------------------|
|                  | (odd/even)            |                   |
| CO               | 14 (even)             | Diamagnetic       |
| NO               | 15 (odd)              | Paramagnetic      |
| NO <sup>+</sup>  | 14 (even)             | Diamagnetic       |
| NO <sup>-</sup>  | 16 (even)             | Diamagnetic       |
| CN               | 13 (odd)              | Paramagnetic      |
| CN-              | 14 (even)             | Diamagnetic       |
| CN <sup>+</sup>  | 12 (even)             | Diamagnetic       |
| HCI              | 18 (even)             | Diamagnetic       |
| HF               | 10(even)              | Diamagnetic       |

**CONCLUSIONS:** This article is very helpful to students in chemistry of undergraduate,

Graduate and also in Postgraduate level. These are the very time savings method. By using these methods student can predict spin state in a very simple way.

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