

---

Subject: V3000 input

Posted by [bruno.b](#) on Fri, 24 Oct 2014 07:54:15 GMT

[View Forum Message](#) <> [Reply to Message](#)

---

FYI, Data warrior cannot read the attached file. I think it is a valid V3000 file. Pubchem can read it.

Best regards,

Bruno

### File Attachments

1) [abs\\_stereo\\_index.v3000.sdf](#), downloaded 1305 times

---

---

Subject: Re: V3000 input

Posted by [thomas](#) on Sat, 25 Oct 2014 23:10:12 GMT

[View Forum Message](#) <> [Reply to Message](#)

---

After a quick check it seems that in a V3 molfile atoms have to be in order, which they are not. In any case DataWarrior should throw a reasonable error message, or should be kind enough to accept atoms in any order. I will check and let you know.

Thanks for the question and kind regards, Thomas

---

---

Subject: Re: V3000 input

Posted by [bruno.b](#) on Mon, 27 Oct 2014 15:00:24 GMT

[View Forum Message](#) <> [Reply to Message](#)

---

Hi Thomas,

according to the Symyx June 2010 documentation, the atom index can be any strictly positive number.

"Identifies atoms. The actual value of the index does not matter as long as each index is unique to each atom. "

For the bonds:

"The actual value of the index does not matter as long as all are unique."

I did not find comment about the order of the atom indices in the Symyx document.

Regards,

Bruno

---

---

Subject: Re: V3000 input

Posted by [thomas](#) on Wed, 05 Nov 2014 15:46:13 GMT

[View Forum Message](#) <> [Reply to Message](#)

---

Hi Bruno,

I checked myself and found the same thing. I will make sure, that this issue will be fixed in the next public DataWarrior release.

Kind regards,

Thomas

---