### NAME

MolecularVolumeDescriptors

### **SYNOPSIS**

use MolecularDescriptors::MolecularVolumeDescriptors;
use MolecularDescriptors::MolecularVolumeDescriptors qw(:all);

#### DESCRIPTION

MolecularVolumeDescriptors class provides the following methods:

 $new,\ Generate Descriptors,\ Get Descriptor Names,\ Get VDWA tom Radii And Volumes Data, Stringify Molecular Volume Descriptors$ 

MolecularVolumeDescriptors is derived from MolecularDescriptors class which in turn is derived from ObjectProperty base class that provides methods not explicitly defined in MolecularVolumeDescriptors, MolecularDescriptors or ObjectProperty classes using Perl's AUTOLOAD functionality. These methods are generated on-the-fly for a specified object property:

```
Set<PropertyName>(<PropertyValue>);
$PropertyValue = Get<PropertyName>();
Delete<PropertyName>();
```

van der Waals molecular volume [ Ref 93 ] (A\*\*3/molecule) of a molecule is calculated using atomic and bonds contributions along with adjustments for aromatic and non-aromatic rings using the following equation:

van der Waals atomic volume for atoms is taken from data file VDWAtomRadiiAndVolumes.csv distributed with MayaChemTools. It contains van der Waals atom radii and atom and volumes data for 38 elements; Table 2 [ Ref 93 ] contains data for only 15 elements. After converting valid van der Waals atom radius data from pm (picometer) to A (Angstrom) available under column name VanderWaalsRadius in PeriodicTableElementsData.csv data file, van der Waals atom volume is calculated using: 4/3 \* PI \* (Radius \*\* 3). For elements specified in Table 2 [ Ref 93 ] - H, B, C, N, O, F, Si, P, S, CI, As, Se, Br, Te, I - the van der Waals atom radii and calculated atom volumes match the values in the table.

### **METHODS**

new

Using specified *MolecularVolumeDescriptors* property names and values hash, new method creates a new object and returns a reference to newly created MolecularVolumeDescriptors object. By default, the following properties are initialized:

## GenerateDescriptors

\$MolecularVolumeDescriptors->GenerateDescriptors();

Calculate van der Waals molecular volume descriptor for a molecule and returns

## Molecular Volume Descriptors.

### GetDescriptorNames

Returns all available descriptor names as an array.

### GetVDWAtomRadiiAndVolumesData

Returns a hash reference to van der Waals atom symbols corresponding to atom types and associated data loaded from VDWAtomRadiiAndVolumes.csv data file as a reference to hash with the following hash data format:

## StringifyMolecularVolumeDescriptors

Returns a string containing information about MolecularVolumeDescriptors object.

### **AUTHOR**

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# SEE ALSO

 $Molecular Descriptors.pm,\ Molecular Descriptors Generator.pm$ 

## COPYRIGHT

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