### NAME

**MACCSKeys** 

### SYNOPSIS

```
use Fingerprints::MACCSKeys;
use Fingerprints::MACCSKeys qw(:all);
```

### **DESCRIPTION**

MACCSKeys [ Ref 45-47 ] class provides the following methods:

new, GenerateFingerprints, GenerateMACCSKeys, GetDescription, SetSize, SetType, StringifyMACCSKeys

MACCSKeys is derived from Fingerprints class which in turn is derived from ObjectProperty base class that provides methods not explicitly defined in MACCSKeys, Fingerprints 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>();
```

For each MACCS (Molecular ACCess System) keys definition, atoms are processed to determine their membership to the key and the appropriate molecular fingerprints strings are generated. An atom can belong to multiple MACCS keys.

For *MACCSKeyBits* value of Type option, a fingerprint bit-vector string containing zeros and ones is generated and for *MACCSKeyCount* value, a fingerprint vector string corresponding to number of MACCS keys [ Ref 45-47 ] is generated.

MACCSKeyBits or MACCSKeyCount values for Type along with two possible 166 | 322 values of Size supports generation of four different types of MACCS keys fingerprint: MACCS166KeyBits, MACCS166KeyCount, MACCS322KeyBits, MACCS322KeyCount.

The current release of MayaChemTools generates the following types of MACCS keys fingerprints bit-vector and vector strings:

FingerprintsBitVector;MACCSKeyBits;166;HexadecimalString;Ascending;000 000000021210210e845f8d8c60b79dffbfffd1

# **METHODS**

new

```
$NewMACCSKeys = new MACCSKeys(%NamesAndValues);
```

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

```
Molecule = '';
    Type = ''
    Size = ''
Examples:
    $MACCSKeys = new MACCSKeys('Molecule' => $Molecule,
                                'Type' => 'MACCSKeyBits',
                                'Size' => 166);
    $MACCSKeys = new MACCSKeys('Molecule' => $Molecule,
                                'Type' => 'MACCSKeyCount',
                                'Size' => 166);
    $MACCSKeys = new MACCSKeys('Molecule' => $Molecule,
                               'Type' => 'MACCSKeyBit',
                                'Size' => 322);
    $MACCSKeys = new MACCSKeys('Molecule' => $Molecule,
                                'Type' => 'MACCSKeyCount',
                                'Size' => 322);
```

## GetDescription

```
$Description = $MACCSKeys->GetDescription();
```

\$MACCSKeys->GenerateMACCSKeys();

Returns a string containing description of MACCS keys fingerprints.

# GenerateMACCSKeys or GenerateFingerprints

print "\$MACCSKeys\n";

```
$MACCSKeys = $MACCSKeys->GenerateMACCSKeys();
```

Generates MACCS keys fingerprints and returns MACCSKeys.

For *MACCSKeyBits* value of Type, a fingerprint bit-vector string containing zeros and ones is generated and for *MACCSKeyCount* value, a fingerprint vector string corresponding to number of MACCS keys is generated.

MACCSKeyBits or MACCSKeyCount values for Type option along with two possible 166 | 322 values of Size supports generation of four different types of MACCS keys fingerprint: MACCS166KeyBits, MACCS166KeyCount, MACCS322KeyBits, MACCS322KeyCount.

Definition of MACCS keys uses the following atom and bond symbols to define atom and bond environments:

```
Atom symbols for 166 keys [ Ref 47 ]:

A : Any valid periodic table element symbol

Q : Hetro atoms; any non-C or non-H atom

X : Halogens; F, Cl, Br, I
```

```
Z : Others; other than H, C, N, O, Si, P, S, F, Cl, Br, I
   Atom symbols for 322 keys [ Ref 46 ]:
   A : Any valid periodic table element symbol
   {\tt Q} : Hetro atoms; any non-C or non-H atom
   X : Others; other than H, C, N, O, Si, P, S, F, Cl, Br, I
   Z is neither defined nor used
   Bond types:
    - : Single
      : Double
   T : Triple
   # : Triple
   ~ : Single or double query bond
   % : An aromatic query bond
   None : Any bond type; no explicit bond specified
   ! : Chain or non-ring bond; ! before a bond type specifies chain bond
   @ : A ring linkage and the number following it specifies the
        atoms position in the line, thus @1 means linked back to the first
        atom in the list.
   Aromatic: Kekule or Arom5
   Kekule: Bonds in 6-membered rings with alternate single/double bonds
           or perimeter bonds
   Arom5: Bonds in 5-membered rings with two double bonds and a hetro
           atom at the apex of the ring.
MACCS 166 keys [ Ref 45-47 ] are defined as follows:
   Key Description
   1 ISOTOPE
   2 103 < ATOMIC NO. < 256
   3 GROUP IVA, VA, VIA PERIODS 4-6 (Ge...)
   4 ACTINIDE
   5 GROUP IIIB, IVB (Sc...)
   6 LANTHANIDE
   7 GROUP VB, VIB, VIIB (V...)
   8 QAAA@1
   9 GROUP VIII (Fe...)
   10 GROUP IIA (ALKALINE EARTH)
   11 4M RING
   12 GROUP IB, IIB (Cu...)
   13 ON(C)C
   14 S-S
   15 OC(0)0
   16 QAA@1
   17 CTC
   18 GROUP IIIA (B...)
   19 7M RING
   20 SI
   21 C=C(Q)Q
   22 3M RING
   23 NC(0)0
   24 N-O
   25 NC(N)N
   26 C$=C($A)$A
```

27 I 28 QCH2Q 29 P 30 CQ(C)(C)A 31 QX 32 CSN 33 NS 34 CH2=A 35 GROUP IA (ALKALI METAL) 36 S HETEROCYCLE 37 NC(O)N 38 NC(C)N 39 OS(0)0 40 S-O 41 CTN 42 F 43 QHAQH 44 OTHER 45 C=CN 46 BR 47 SAN 48 OQ(O)O 49 CHARGE 50 C=C(C)C 51 CSO 52 NN 53 QHAAAQH 54 QHAAQH 55 OSO 56 ON(O)C 57 O HETEROCYCLE 58 QSQ 59 Snot%A%A 60 S=0 61 AS(A)A 62 A\$A!A\$A 63 N=O 64 A\$A!S 65 C%N 66 CC(C)(C)A 67 QS 68 QHQH (&...) 69 QQH 70 QNQ 71 NO 72 OAAO 73 S=A 74 CH3ACH3 75 A!N\$A 76 C=C(A)A 77 NAN 78 C=N 79 NAAN 80 NAAAN 81 SA(A)A 82 ACH2QH 83 QAAAA@1 84 NH2 85 CN(C)C 86 CH2QCH2 87 X!A\$A 88 S

89 OAAAO 90 QHAACH2A

```
91 QHAAACH2A
92 OC(N)C
93 QCH3
94 QN
95 NAAO
96 5M RING
97 NAAAO
98 QAAAAA@1
99 C=C
100 ACH2N
101 8M RING
102 QO
103 CL
104 QHACH2A
105 A$A($A)$A
106 QA(Q)Q
107 XA(A)A
108 CH3AAACH2A
109 ACH20
110 NCO
111 NACH2A
112 AA(A)(A)A
113 Onot%A%A
114 CH3CH2A
115 CH3ACH2A
116 CH3AACH2A
117 NAO
118 ACH2CH2A > 1
119 N=A
120 HETEROCYCLIC ATOM > 1 (&...)
121 N HETEROCYCLE
122 AN(A)A
123 OCO
124 QQ
125 AROMATIC RING > 1
126 A!O!A
127 A$A!O > 1 (&...)
128 ACH2AAACH2A
129 ACH2AACH2A
130 QQ > 1 (&...)
131 QH > 1
132 OACH2A
133 A$A!N
134 X (HALOGEN)
135 Nnot%A%A
136 O=A > 1
137 HETEROCYCLE
138 QCH2A > 1 (&...)
139 OH
140 0 > 3 (&...)
141 CH3 > 2 (&...)
142 N > 1
143 A$A!O
144 Anot%A%Anot%A
145 6M RING > 1
146 0 > 2
147 ACH2CH2A
148 AQ(A)A
149 CH3 > 1
150 A!A$A!A
151 NH
152 OC(C)C
153 OCH2A
```

154 C=0

155 A!CH2!A 156 NA(A)A 157 C-O 158 C-N 159 O > 1 160 CH3 161 N 162 AROMATIC 163 6M RING 164 O 165 RING 166 FRAGMENTS

MACCS 322 keys set as defined in tables 1, 2 and 3 [ Ref 46 ] include:

- o 26 atom properties of type P, as listed in Table 1
- o 32 one-atom environments, as listed in Table 3
- o 264 atom-bond-atom combinations listed in Table 4

Total number of keys in three tables is: 322

Atom symbol, X, used for 322 keys [ Ref 46 ] doesn't refer to Halogens as it does for 166 keys. In order to keep the definition of 322 keys consistent with the published definitions, the symbol X is used to imply "others" atoms, but it's internally mapped to symbol X as defined for 166 keys during the generation of key values.

Atom properties-based keys (26):

```
Key
    Description
      A(AAA) or AA(A)A - atom with at least three neighbors
1
      Q - heteroatom
3
      Anot%not-A - atom involved in one or more multiple bonds, not aromatic
4
     A(AAAA) or AA(A)(A)A - atom with at least four neighbors
5
      A(QQ) or QA(Q) - atom with at least two heteroatom neighbors
6
      A(QQQ) or QA(Q)Q - atom with at least three heteroatom neighbors
      QH - heteroatom with at least one hydrogen attached
8
      CH2(AA) or ACH2A - carbon with at least two single bonds and at least
      two hydrogens attached
      {
m CH3}({
m A}) or {
m ACH3} - carbon with at least one single bond and at least three
9
     hydrogens attached
10
     Halogen
11
     A(-A-A-A) or A-A(-A)-A - atom has at least three single bonds
      AAAAAA@1 > 2 - atom is in at least two different six-membered rings
12
      A(\$A\$A\$A) or A\$A(\$A)\$A - atom has more than two ring bonds
13
14
      A$A!A$A - atom is at a ring/chain boundary. When a comparison is done
      with another atom the path passes through the chain bond.
15
      Anot%A%Anot%A - atom is at an aromatic/nonaromatic boundary. When a
      comparison is done with another atom the path
      passes through the aromatic bond.
16
      A!A!A - atom with more than one chain bond
      A!A$A!A - atom is at a ring/chain boundary. When a comparison is done
17
      with another atom the path passes through the ring bond.
18
      A%Anot%A%A - atom is at an aromatic/nonaromatic boundary. When a
      comparison is done with another atom the
      path passes through the nonaromatic bond.
19
      HETEROCYCLE - atom is a heteroatom in a ring.
20
      rare properties: atom with five or more neighbors, atom in
      four or more rings, or atom types other than
     H, C, N, O, S, F, Cl, Br, or I
     rare properties: atom has a charge, is an isotope, has two or
     more multiple bonds, or has a triple bond.
2.2
     N - nitrogen
23
     S - sulfur
      0 - oxygen
2.5
      A(AA)A(A)A(AA) - atom has two neighbors, each with three or
```

more neighbors (including the central atom).

26 CHACH2 - atom has two hydrocarbon (CH2) neighbors

Atomic environments properties-based keys (32):

```
Description
Key
      C(CC)
27
28
      C(CCC)
29
      C(CN)
30
      C(CCN)
31
      C(NN)
32
      C(NNC)
33
      C(NNN)
34
      C(CO)
35
      C(CCO)
36
      C(NO)
37
      C(NCO)
      C(NNO)
38
39
      C(00)
40
      C(C00)
41
      C(NOO)
42
      C(000)
43
      Q(CC)
44
      Q(CCC)
45
      Q(CN)
46
      Q(CCN)
47
      Q(NN)
48
      Q(CNN)
49
      Q(NNN)
50
      Q(CO)
51
      Q(CCO)
52
      Q(NO)
      Q(CNO)
53
54
      Q(NNO)
55
      Q(00)
56
      Q(C00)
57
      Q(NOO)
58
      Q(000)
```

Note: The first symbol is the central atom, with atoms bonded to the central atom listed in parentheses. Q is any non-C, non-H atom. If only two atoms are in parentheses, there is no implication concerning the other atoms bonded to the central atom.

Atom-Bond-Atom properties-based keys: (264)

```
Кеу
      Description
59
      C-C
60
      C-N
      C-0
61
62
      C-S
63
      C-Cl
64
      C-P
65
      C-F
66
      C-Br
67
      C-Si
68
      C-I
69
      C-X
70
      N-N
71
      N-O
72
      N-S
73
      N-Cl
74
      N-P
75
      N-F
76
      N-Br
77
      N-Si
78
      N-I
```

79	N-Y
	N-X
80	0-0 0-S
81	
82	O-Cl O-P
83	O-P
84	O-F
85	O-Br O-Si
86	O-Si
87	O-I
88	O-X
89	O-X S-S
90	S-Cl
91	
92	S-P S-F
93	S-Br
94	S-Si
95	S-Si S-I
	S-X
96	S-X
97	Cl-Cl Cl-P
98	CI-P
99	Cl-F
100	Cl-Br Cl-Si
101	Cl-Si
102	Cl-I
103	Cl-X
104	P-P
105	P-F
106	P-Br
107	P-Br P-Si
108	P-I
109	
110	P-X F-F
111	F-Br
112	F-Si
113	F-Si F-I
114	F-X
115	Br-Br
116	Br-Br Br-Si
117	Br-I
118	
119	Br-X Si-Si
120	Si-Ji
121	Si-X
122	
123	I-I
124	
125	X-X C=C
126	C=C
127	
	C=0 C=S
128	C=S C=Cl
129	
130	C=P
131	C=F
132	C=Br
133	C=Si
134	C=I
135	C=X
136	N=N
137	N=O
138	N=S
139	N=Cl
140	N=P
141	N=F
142	N=Br

N=Br

142

143	N=Si
144	N=I
145	N=X
146	0=0
147	O=S
148	O=Cl
149	O=P
150	O=F
151	O=Br
152	O=Si
153	O=I
154	O=X
155	S=S
156	S=Cl
157	S=P
158	S=F
159	S=Br
160	S=Si
161	S=I
162	S=X
163	Cl=Cl
164	Cl=P
165	Cl=F
166	Cl=Br Cl=Si
167	
168	Cl=I
169	Cl=X P=P
170	
171 172	P=F P=Br
173	P=Bi P=Si
174	P=I
175	P=I P=X
176	F=F
177	F=Br
178	F=Si
179	F=I
180	F=X
181	Br=Br
182	Br=Si
183	Br=I
184	Br=X
185	Si=Si
186	Si=I
187	Si=X
188	I = I
189	I = X
190	X=X
191	C#C
192	C#N
193	C#0
194	C#S
195	C#Cl
196	C#P
197	C#F
198	C#Br
199	C#Si
200	C#I
201 202	C#X
202	N#N
	N#0
204 205	N#S N#Cl
205	N#CT

206

N#P

207	N#F
208	N#Br
209	N#Si
210	N#I
211	N#X
212	0#0
213	O#S
214	O#Cl
215	O#P
216	O#F
217	O#Br
218	O#Si
219	O#I
220	O#X
221	
	S#S
222	S#Cl
223	S#P
224	S#F
225	S#Br
226	S#Si
227	S#I
228	S#X
229	Cl#Cl
230	Cl#P
231	Cl#F
232	Cl#Br
233	Cl#Si
234	Cl#I
235	Cl#X
236	P#P
237	P#F
238	P#Br
239	P#Si
240	P#I
241	P#X
242	F#F
243	F#Br
244	F#Si
245	F#I
246	F#X
247	Br#Br
248	Br#Si
249	Br#I
250	Br#X
251	Si#Si
252	Si#I
253	Si#X
254	I#I
255	I#X
256	X#X
257	C\$C
258	C\$N
259	C\$0
260	C\$S
261	C\$Cl
262	C\$P
263	C\$F
264	C\$Br
265	C\$Si
266	C\$I
267	C\$X
268	N\$N
269	N\$O
270	N\$S

```
271
          N$Cl
    272
         NŠP
    273
         N$F
    274
         N$Br
    275
         N$Si
    276
         N$I
    277
         N$X
    278
         0$0
    279
         0$S
    280
         O$Cl
    281
         O$P
    282
         O$F
    283
          O$Br
    284
         O$Si
    285
         O$I
    286
         O$X
    287
         S$S
    288
         S$Cl
    289
         S$P
    290
         S$F
    291
         S$Br
    292
         S$Si
    293
         S$I
    294
         S$X
    295
         Cl$Cl
    296
         Cl$P
    297
         Cl$F
    298
         Cl$Br
    299
         Cl$Si
    300
         Cl$I
    301
         Cl$X
    302
         P$P
    303
         P$F
    304
         P$Br
    305
         P$Si
    306
         P$I
    307
         P$X
    308
         F$F
    309
         F$Br
    310
         F$Si
    311
         F$I
    312
         F$X
    313
         Br$Br
    314
        Br$Si
    315
        Br$I
    316
        Br$X
    317
         Si$Si
    318
         Si$I
    319
         Si$X
    320
         I$I
    321
          I$X
    322
         X$X
    $MACCSKeys->SetSize($Size);
Sets size of MACCS keys and returns MACCSKeys. Possible values: 166 or 322.
```

\$MACCSKeys->SetType(\$Type);

Sets type of MACCS keys and returns MACCSKeys. Possible values: MACCSKeysBits or MACCSKeysCount.

SetSize

SetType

# StringifyMACCSKeys

\$String = \$MACCSKeys->StringifyMACCSKeys();

Returns a string containing information about MACCSKeys object.

# **AUTHOR**

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# **SEE ALSO**

Fingerprints.pm, FingerprintsStringUtil.pm, AtomNeighborhoodsFingerprints.pm, AtomTypesFingerprints.pm, EStateIndiciesFingerprints.pm, ExtendedConnectivityFingerprints.pm, PathLengthFingerprints.pm, TopologicalAtomPairsFingerprints.pm, TopologicalAtomTripletsFingerprints.pm, TopologicalAtomTorsionsFingerprints.pm, TopologicalPharmacophoreAtomPairsFingerprints.pm, TopologicalPharmacophoreAtomTripletsFingerprints.pm

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