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 000000021210210e845f8d8c60b79dffbffffd1

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
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
$ : Ring bond; $ before a bond type specifies ring bond
! : Chain or non-ring bond; ! before a bond type specifies chain bond
```

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```
@ : 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(O)O
    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 > 1137 HETEROCYCLE 138 QCH2A > 1 (&...) 139 OH 140 0 > 3 (&...) 141 CH3 > 2 (&...) 142 N > 1143 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 QCH2A 154 C=0 155 A!CH2!A 156 NA(A)A 157 C-0 158 C-N 159 0 > 1 160 CH3 161 N 162 AROMATIC 163 6M RING 164 0 165 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):

17

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```
Key
      Description
      A(AAA) or AA(A)A - atom with at least three neighbors
1
2
      0 - heteroatom
3
      Anot%not-A - atom involved in one or more multiple bonds, not aromatic
      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
7
      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
9
      CH3(A) or ACH3 - carbon with at least one single bond and at least three
      hydrogens attached
10
      Halogen
11
      A(-A-A-A) or A-A(-A)-A - atom has at least three single bonds
12
      AAAAAA@1 > 2 - atom is in at least two different six-membered rings
13
      A(\$A\$A\$A) or A\$A(\$A)\$A - atom has more than two ring bonds
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

A%Anot%A%A - atom is at an aromatic/nonaromatic boundary. When a

with another atom the path passes through the ring bond.

```
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
21
      rare properties: atom has a charge, is an isotope, has two or
      more multiple bonds, or has a triple bond.
22
      N - nitrogen
      S - sulfur
23
      0 - oxygen
2.4
25
      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
27
      C(CC)
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)
38
      C(NNO)
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
      O(NN)
48
      Q(CNN)
49
      Q(NNN)
50
      Q(CO)
51
      Q(CCO)
52
      Q(NO)
53
      Q(CNO)
54
      Q(NNO)
55
      Q(00)
56
      Q(COO)
57
      Q(NOO)
      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)

Key	Description
59	C-C
60	C-N
61	C-0
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-F N-Br
77	N-Si
78	N-I
79	N-X
80	0-0
81	
82	0-S 0-Cl
83	0-C1 0-P
84	
85	O-F O-Br
	O-Bi O-Si
86 87	0-21
	0-I
88	0-X S-S
89	
90	S-Cl
91	S-P S-F
92	S-F
93	S-Br S-Si
94	S-SI S-I
95	
96	S-X
97	Cl-Cl Cl-P
98	
99	Cl-F
100	Cl-Br Cl-Si
101	CI-Si
102	Cl-I
103	Cl-X
104	P-P
105	P-F
106	P-Br
107	P-Si
108	P-I
109	P-X
110	F-F
111	F-Br
112	F-Si
113	F-I
114	F-X
115	Br-Br
116	Br-Si
117	Br-I
118	Br-X
119	Si-Si
120	Si-I
121	Si-X
122	I-I
123	I-X
124	X-X
125	C=C
126	C=N
127	C=O
128	C=S
129	C=Cl
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
143	N=Si
144	N=I
145	N=X
146	0=0
147	O=S

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```
148
      O=Cl
149
      O=P
150
      O=F
      O=Br
151
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
167
      Cl=Si
168
      Cl=I
169
      Cl=X
170
      P=P
171
      P=F
172
      P=Br
173
      P=Si
174
      P=I
175
      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
      C#X
202
      N#N
203
      N#O
204
      N#S
205
      N#Cl
206
      N#P
207
      N#F
208
      N#Br
209
      N#Si
210
      N#I
211
      N#X
212
      0#0
213
      0#S
214
      O#Cl
215
      O#P
216
      O#F
217
      O#Br
218
      O#Si
```

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220

O#I

О#Х

- 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
- Cl#Si 233
- 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 N\$P
- 273 N\$F
- 274 N\$Br
- 275 N\$Si 276 N\$I
- 277 N\$X
- 278 0\$0
- 279 0\$8
- 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
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```
294
      SSX
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$F
304
      PSBr
305
      P$Si
306
      P$I
307
      P$X
308
      F$F
309
      F$Br
310
      F$Si
311
      F$I
      F$X
312
313
      Br$Br
314
      Br$Si
315
      Br$I
316
      Br$X
317
      Si$Si
318
      SiŠI
319
      Si$X
320
      I$I
321
      I$X
322
      XSX
```

SetSize

```
$MACCSKeys->SetSize($Size);
```

Sets size of MACCS keys and returns MACCSKeys. Possible values: 166 or 322.

SetType

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

 $Sets\ type\ of\ MACCS\ keys\ and\ returns\ \textit{MACCSKeys}.\ Possible\ values:\ \textit{MACCSKeysBits}\ or\ \textit{MACCSKeysCount}.$

StringifyMACCSKeys

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

Returns a string containing information about MACCSKeys object.

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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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