

Molecular Coding Format manual

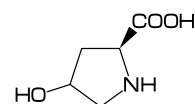
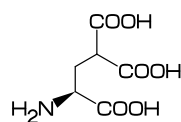
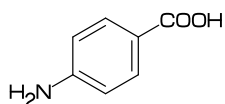
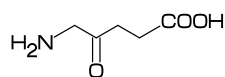
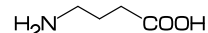
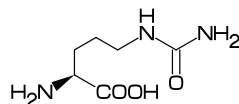
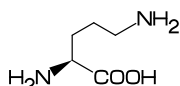
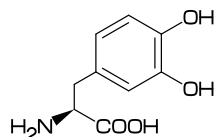
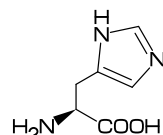
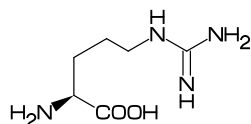
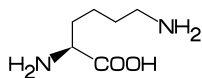
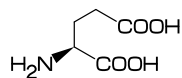
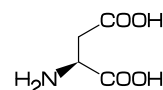
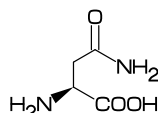
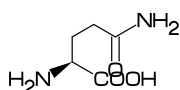
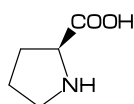
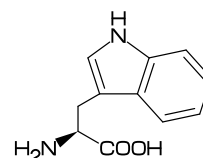
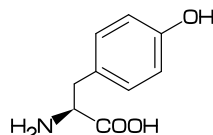
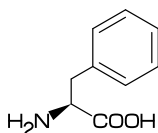
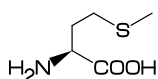
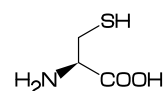
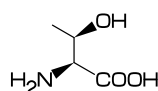
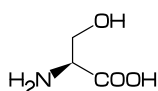
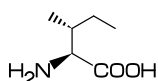
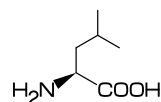
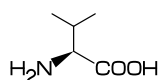
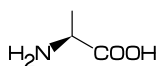
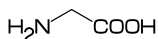
Akira Yamaji

June 8, 2025

mcf2graph version 5.21

Located at <http://www.ctan.org/pkg/mcf2graph>

Suggestion or request mail to: mcf2graph@gmail.com



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1 Introduction

Molecular Coding Format(MCF) is new linear notation represent chemical structure diagrams. This Coding is named from programing technique such as operator, array, scope, macro, adresssing, etc. mcf2graph convert from MCF to PNG, SVG, EPS, MOL file. It is also able to calculate molecular weight, exact mass, molecular formula.

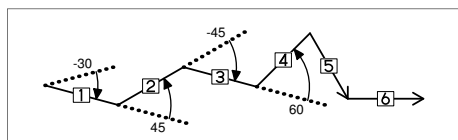
2 MCF syntax

2.1 Make bond

2.1.1 Chain

real number plus (+): counterclockwise
real number minus(-): clockwise
\$n (0<=n<360): absolute angle

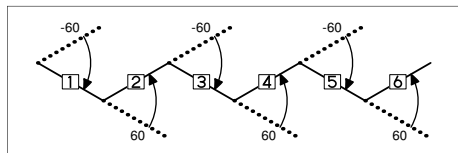
<10,-30,45,-45,60,\$300,\$0



2.1.2 Chain with !,!n

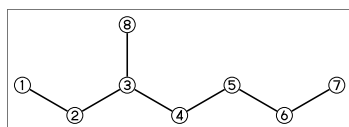
! : take value 60 or -60 depend on current angle
!6 : !,!,!,!,!,!

<-30,!6



take value 0 just after jump to atom
@n,! : !=0

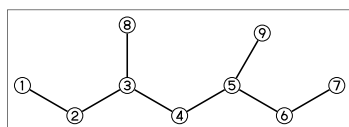
<-30,!5,@3,!



2.1.3 Jump to atom

@n : Jump to An
** An: atom number(-999<=n<=4095)

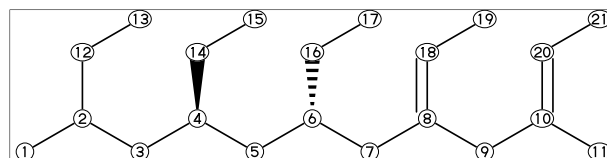
<-30,!6,@3,0,@5,-30



2.1.4 Branch bond

@n,! : @n,!
@n,!w : @n,!~wf
@n,!z : @n,!~zf
@n,!d : @n,!~db
@n,!dr : @n,!~dr

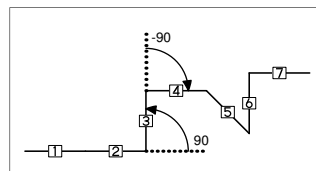
<30,!10,
@2,!,,@4,!w,!,@6,!z,!,@8,!d,!,@10,!dr,!



2.1.5 Rotate current angle

<angle : rotate current angle

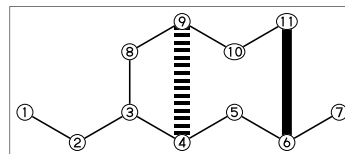
0,0,<90,0,<-90,0,<\$315,0,<\$90,0,<\$0,0



2.1.6 Connect atom

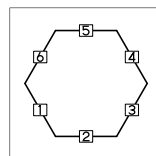
&n : Connect to An

<-30,!6,@3,!,,!3,&6~bd,@9,&4~bz



2.1.7 Ring

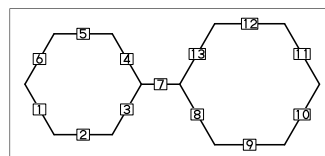
?n : n membered ring(3<=n<=20)
?6



2.1.8 Ring length

?n`length : change ring length

?6,@4,!,,?6`1.2



2.2 Change bond type

2.2.1 Double, triple, wedge, vector

(Double, triple)

a~type : ~type, a

dm : double middle

dl : double left side

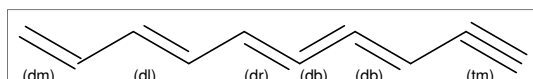
dr : double right side

db : double left or right side

tm : triple

!db, !d : !~db / !tm, !t : !~tm

<-30, !~dm, !, !~dl, !, !~dr, !~db, !, !~tm
<-30, !dm, !, !dl, !, !dr, !d, !d, !, !t



(Wedge, Vector)

wf: wedge forward

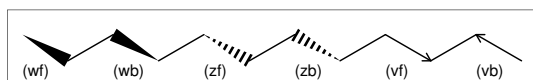
wb: wedge backward

zf: hashed(zebra stripe) wedge forward

zb: hashed(zebra stripe) wedge backward

vf: vector forward / vb: vector backward

<-30,
!~wf, !, !~wb, !, !~zf, !, !~zb, !, !~vf, !~vb



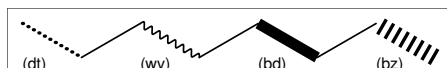
(Dotted, wave)

Bn=bond type : change bond type at Bn

dt : dotted / wv : wave

bd : broad / bz : broad dotted

<-30, !7, 1=dt, 3=wv, 5=bd, 7=bz



2.2.2 Over line

si_ : single over line

wf_ : wedge forward over line

wb_ : wedge backward over line

zf_ : hashed wedge forward over line

zb_ : hashed wedge backward over line

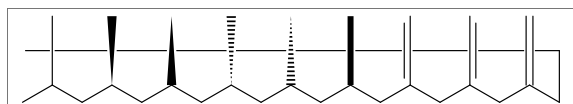
bd_ : broad over line

dl_ : double left over line

dr_ : double right over line

dm_ : double over line

<30, !8, !, 60, 90`18,
{2~si_'4~wf_'6~wb_'8~zf_'10~zb_'
12~bd_'14~dl_'16~dr_'18~dm_}:?`2



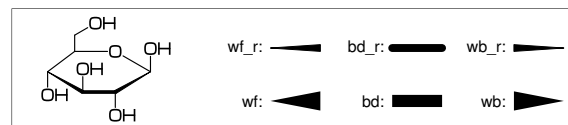
2.2.3 Steric ring

wf_r : wedge forward (half width)

bd_r : broad (half width, rounded)

wb_r : wedge backward (half width)

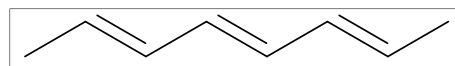
#1.25, -30~wf_r, 30~bd_r`1, 30~wb_r,
120, 0, 30, &1, ##, #.5, 6~\$90: /!OH,
{1~\$270'2~\$90'3~\$270'4~\$90}: /OH,



2.2.4 Change multiple bond type

{2'4'6}=dl or 2'4'6=dl : 2=dl, 4=dl, 6=dl

<30, !7, {2'4'6}=dl or <30, !7, 2'4'6=dl

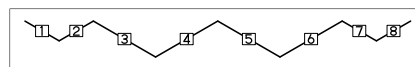


2.3 Change bond length

2.3.1 Change chain length

(!, !n)`length : change length of !, !n

<-30, !2, !4`1.2, !2

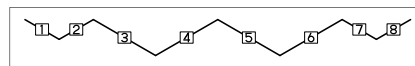


2.3.2 Change multiple bond length

#n : bond length=n

: reset bond length

<-30, !2, #1.2, !4, ##, !2

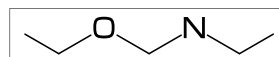


2.4 Change atom

2.4.1 Insert atom

Insert hetero atom

<-30, !2, 0, !2, N, !2

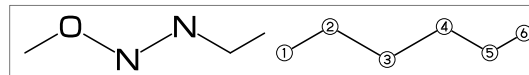


2.4.2 Change atom

2:0 : change A2 C to O

{3'4}:N or 3'4:N : change A3, A4 C to N

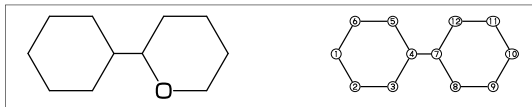
<30, !4, 2:0, {3'4}:N or <30, !4, 2:0, 3'4:N



2.4.3 Brock address

| : divide brock

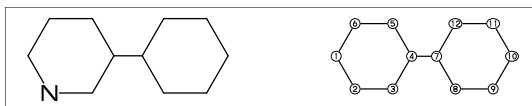
?6,@4,! ,| ,?6,2:0



2.4.4 Reset brock address

|| : reset brock address

?6,@4,! ,| ,?6,||,2:N

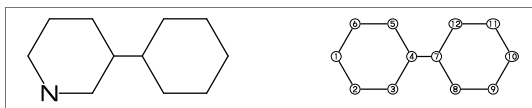


2.4.5 Absolute address

\$2:n : change A\$2 C to N

**1<=n<=3095

?6,@4,! ,| ,?6,\$2:N

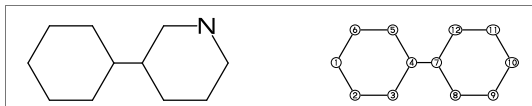


2.4.6 Relative address

-2:n : change A(-2) C to N

** -999<=n<=-1

?6,@4,! ,?6,-2:N



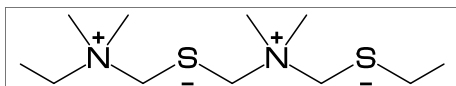
2.4.7 Charged atom

p_ : positive

n_ : negative

<-30,!2,N,??,p_,!2,S,n_~180,

!6,7:N,7:??,9:S,7:n_,9:n_~180



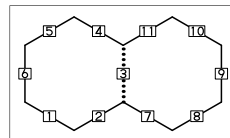
2.5 Fuse ring

(Attached 1 bond)

?6,3=?6 : fuse ?6 at B3

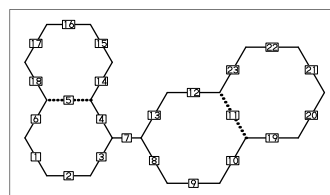
** Bn(n:-999<=n<=4095): bond number

?6,3=?6



** fused ring size depend on attached bond length

?6,@4,! ,?6`1.2,5=?6,11=?6

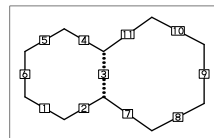


?6,3=#1.3'?6 : fuse #1.3'?6 at B3

#1.3'?6 : 6 membered ring scaled 1.3

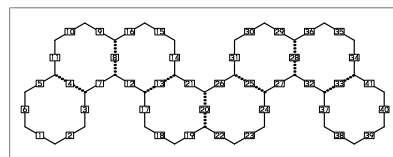
** #n'?m (5<=m<=6,1.1<=n<=1.5)

?6,3=#1.3'?6



?6,-3'-4'-4'-2'-2'-4'-4=?6

?6,4'8'13'20'25'28'33=?6



(Attached 2 bond)

4--11=?6 : fuse 4/6 ring to B11..B4

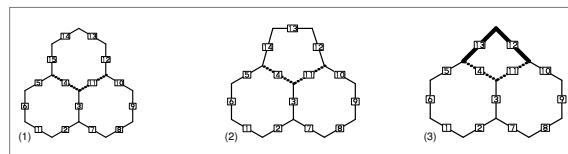
4--11=?5 : fuse 3/5 ring to B11..B4

4--11=?4 : fuse 2/4 ring to B11..B4

1:<30,?6,3=?6,11--4=?6

2:<30,?6,3=?6,11--4=?5

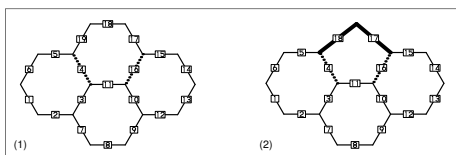
3:<30,?6,3=?6,11--4=?4



(Attached 3 bond)

16---4=?6 : fuse 3/6 ring to B16..B4
16---4=?5 : fuse 2/5 ring to B16..B4

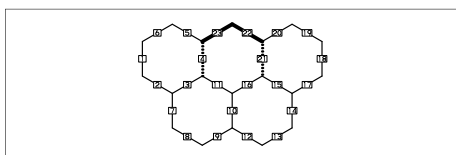
1:?6,3'10'(16---4)=?6
2:?6,3'10=?6,16---4=?5



(Attached 4 bond)

21----4=?6 : fuse 2/6 ring to B21..B4

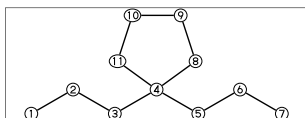
<-30,?6,3'10'15'(21----4)=?6



2.6 Spiro ring

@4,?5 : add ?5 at A4

<30,!6,@4,?5

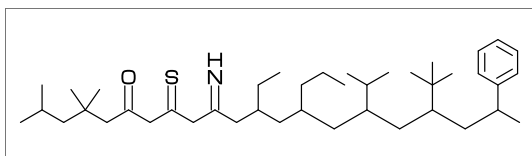


2.7 Group

2.7.1 Insert group

? : methyl
?? : dimethyl
?0 : carbonyl
?S : thioketone
?NH : imino
/! : ethyl
/!2 : propyl
/?! : isopropyl
/??! : tert-butyl
/Ph : phenyl

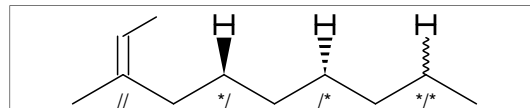
<30,!?,!2,??,!2,?0,!2,?S,!2,
?NH,!2,/!,!2,/!2,!2`1,/?!,!2`1,
/??!,!2`1,/#.6'Ph,!



2.7.2 Insert modified group

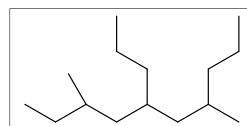
// : double (double middle)
*/ : wedge forward
/* : hashed wedge forward
** : wave

<30,! ,//!,!2,*/H,!2,/*H,!2,*/H,!



~ : change type
^ : change angle
` : change length
> : change environment

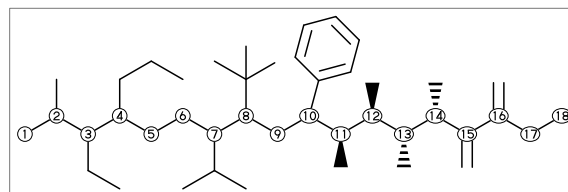
<-30,#1,! ,
?`2^30,!2,/!2>lr,!2,/!2>r1,!)



2.7.3 Add group

?w : ?~wf
?z : ?~zf
?d : ?~dm

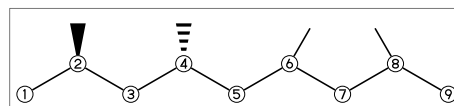
<30,!17,2:?,3:/!,4:/!2,7:/?! ,
8:/??!,10~15:/#.6'Ph,
11'12':?w,13'14:?z,15'16:?d



2.7.4 Add modified group

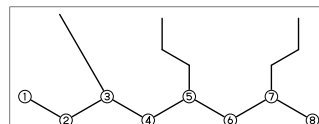
~,^,` : change type,angle,length

<30,!6,{2~wf'4~zf'6~30'8~\$120}:?



^,^,> : change angle,length,environment

<-30,!7`1,3:?`2^30,5:/!2>lr,7:/!2>r1

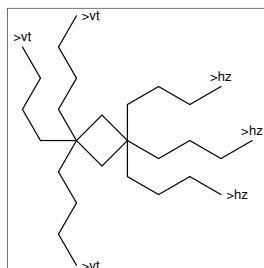


2.8 Chain environment

2.8.1 Horizontal,vertical

>hz : horizontal environment (default)
>vt : vertical environment

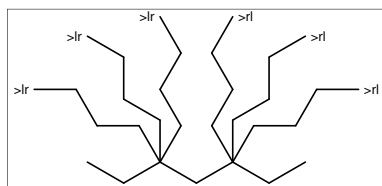
```
?4,  
{3~-90'3~-30'3^90}:/!3>hz,  
{1~-60'1'1^60}:/!3>vt
```



2.8.2 Left-right,right-left

>lr : left-right environment
>rl : right-left environment

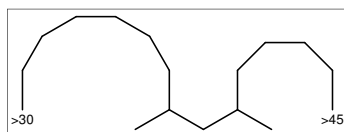
```
<-30,!6,  
{3~-30'3'3^30}:/!3>lr,
```



2.8.3 Rotate fixed angle

>n : rotate n

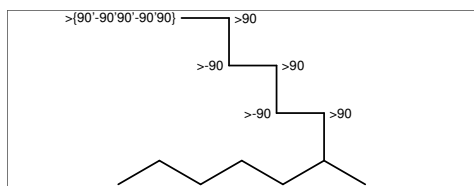
```
<30,!4,  
2:/!6>30, % 2:\,30,30,30,30,30,30  
4:/!4>-45 % 4:\,-45,-45,-45,-45
```



2.8.4 Rotate multiple angle

>{90'-90,...} : rotate 90,-90,...

```
<30,!6,6>{90'-90'90'-90'90}:/!5
```

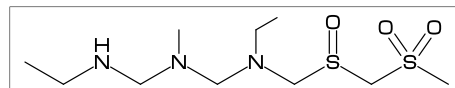


2.9 Miscellaneous

2.9.1 Abbreviated parts

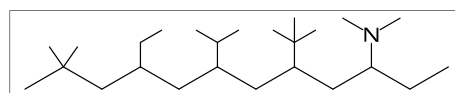
```
NH : N,/H~n1 N? : N,?  
N?2 : N,/! S?0 : S,?0  
S?0?0 : S,?0^35,/^~35
```

```
<-30,!2,NH,!2,N?,!2,N?2,S?0,!2,S?0?0,!
```



```
?? : ?^35,?-35 /?! : isopropyl  
/?! : tert-butyl /N?! : dimethylamino
```

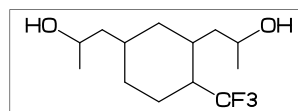
```
<30,!11^1,2:??,4:/!,6:/?!,8:/?!,10:/N?!
```



2.9.2 Define group,parts

```
`(..) : define group  
``(..)(..) : define group with atoms  
'(..) : define parts
```

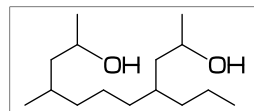
```
CF3:=`("{CF3}"); % ** group weight =0  
CF3:=``("{CF3}")(C,F,F,F);  
** group weight =12(C)+19(F)*3=69  
iBuOH:=''(!?,OH);  
drawm(<30,?6,4'6:/iBuOH,3:/CF3)
```



2.9.3 Concatenate group,parts

a'b : '(a,b)

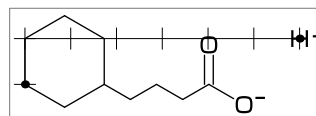
```
<30,!8,2'6:/!?'!OH
```



2.9.4 Move position [@]

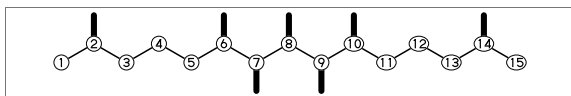
```
@(x'y) : Move l*(x,y) from current position  
@$(x'y): Move l*(x,y) from origin(@1)  
** l=bond length of ring
```

```
<30,?6,@3,!4,?0,!0,n^60,@$(6'1),H,p^15
```



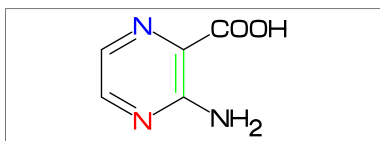
2.9.5 Serial number

```
6_10 : 6,7,8,9,10  
<30,'14',{2'6_10'14}:?~bd_r`0.5
```



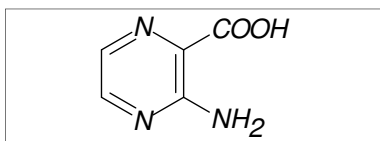
2.9.6 Change color

```
beginfigm  
  drawm(<30,Ph,2'5:N,3:/NH2,4:/COOH,  
        2:red, 5:blue, 3=green)  
endfigm
```



2.9.7 Change font

```
beginfigm  
  %-----  
  atomfont:="phvro8g";  
  %-----  
  drawm(<30,Ph,2'5:N,3:/NH2,4:/COOH)  
endfigm
```

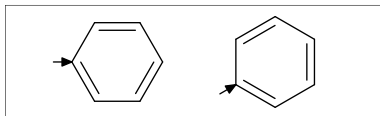


3 Option parameter

3.1 Angle parameter

mangle=0 ** default

```
@(0.2,0.5) drawm(Ph)  
mangle:=30;  
@(0.8,0.5) drawm(Ph)
```

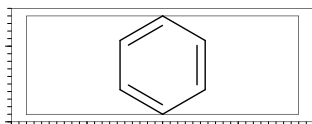


3.2 Size/Ratio parameter

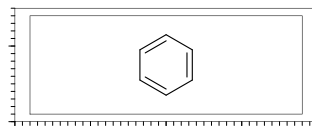
3.2.1 Bond length [|=]

|=(n) : abbreviated form of blength:=n;

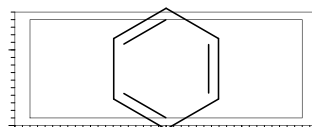
(fit to figure size)
blength=0 ** default



(ratio bond/figure width)
blength=0.1 ** (0<blength<=1)
blength=60mm(width)*0.1=6mm

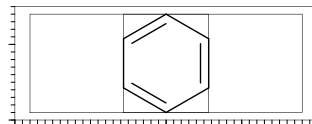


(bond length)
blength=9mm
** (blength>1) ignore msize(w,h)

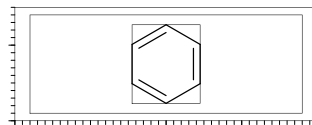


3.2.2 Molecular size

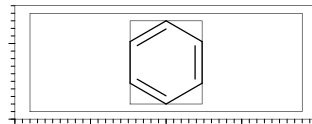
##(1,1) : msize=(1,1) ** default
p : abbreviated form of msize:=p;



##(0.25,1) : msize=(0.25,1)
** msize=(40mm-4mm)*0.25=9mm

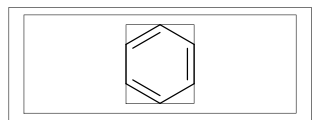


##(11mm,11mm) : msize=(11mm,11mm)



3.2.3 Molecular position

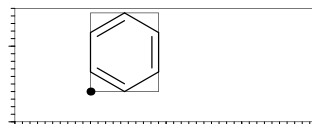
@(0.5,0.5) : mposition=(0.5,0.5) **default



@(1,0) : mposition=(1,0)



@(10mm,4mm) : mposition=(10mm,4mm)

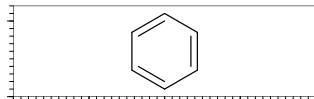


3.3 Size parameter

3.3.1 Figure size [#]

fsize=(figure width,figure height)
** default: (30mm,20mm)
p : abbreviated form of fsize:=p;

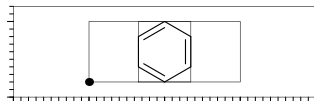
#(40mm,15mm) : fsize=(40mm,15mm)



3.3.2 Figure margin [#@]

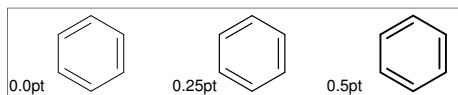
fmargin=(margin left right,top bottom)
** default: (0.4mm,0.4mm)
#@ p : abbreviated form of fmargin:=p;

#@(10mm,2mm) : fmargin=(10mm,2mm)



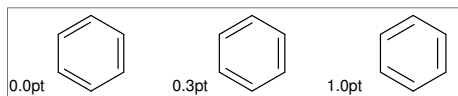
3.3.3 Offset thickness of bond

default: offset_thickness=0.25pt



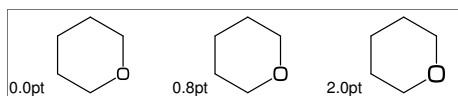
3.3.4 Offset of double bond gap

default: offset_bond_gap=0.3pt



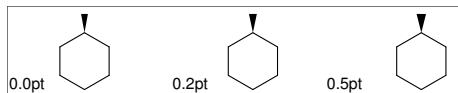
3.3.5 Offset of atom width

default: offset_atom=0.8pt



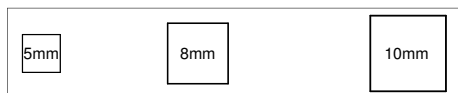
3.3.6 Offset of wedge width

default: offset_wedge=0.2pt



3.3.7 Max bond length [|<]

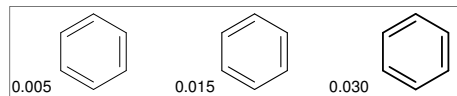
|<(n): abbreviated form of max_blength:=n;
default: max_blength=10mm



3.4 Ratio parameter

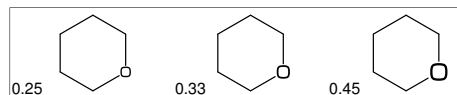
3.4.1 Thickness/bond length

default: ratio_thickness_bond=0.012



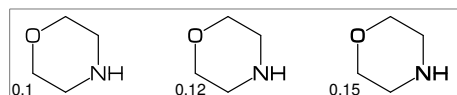
3.4.2 Atom/bond length

default: ratio_atom_bond= 0.36



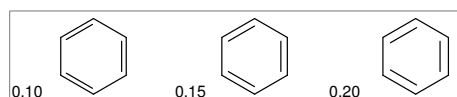
3.4.3 Char thickness/Atom width

default: ratio_char_atom=0.12



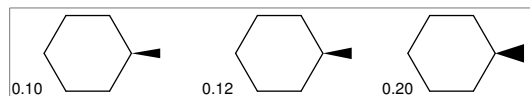
3.4.4 Bond gap/bond length

default: ratio_bondgap_bond= 0.15



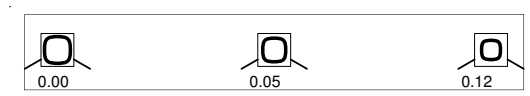
3.4.5 Wedge/bond length

default: ratio_wedge_bond=0.12



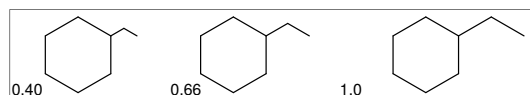
3.4.6 Figure atom gap/atom length

default: ratio_atomgap_atom= 0.050



3.4.7 Chain/ring length

default: ratio_chain_ring= 0.66



3.4.8 Hash gap/bond length

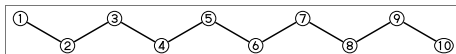
default: ratio_hashgap_bond=0.12



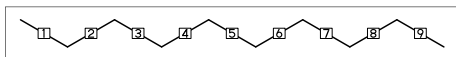
3.5 Drawing mode

3.5.1 Numbering atom,bond

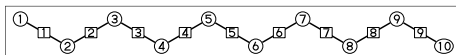
sw_numbering:=Atom;



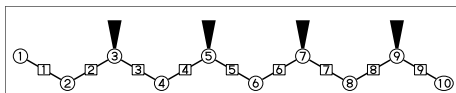
sw_numbering:=Bond;



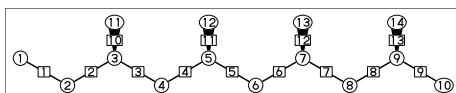
sw_numbering:=Atom+Bond;



sw_numbering:=Atom+Bond;

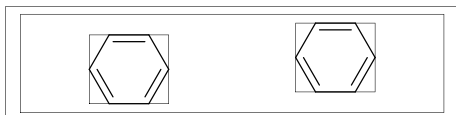


sw_numbering:=Atom+Bond+All;



3.5.2 Trimming mode

```
sw_trimming:=0; ** default
##(1,0.7)
@(0.2,0.3) drawm(Ph) @(0.8,0.7) drawm(Ph)
```



```
sw_trimming:=1;
@(0.2,0.3) drawm(Ph) @(0.8,0.7) drawm(Ph)
```



3.5.3 Omit group

** default: sw_omit=Group



3.5.4 Omit bond type

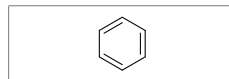
** default: sw_omit=Bond



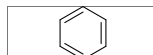
3.6 Frame

3.6.1 Figure frame

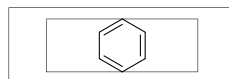
```
** default:sw_frame=0
(Draw figure frame)
fmargin:=(5mm,2mm);
sw_frame=Outside
```



(Frame inside margin)
sw_frame=Inside

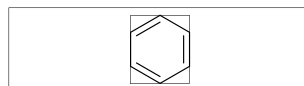


(Draw both frame)
sw_frame=Bothside=Inside+Outside



3.6.2 Molecular frame

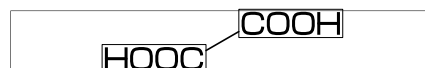
```
sw_frame=Mol
** default:sw_frame=0
```



3.6.3 Atom frame

```
sw_frame=Atom
** default: sw_frame=0
```

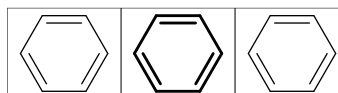
drawm(<30,COOH,!,COOH)



3.7 Parameter setting

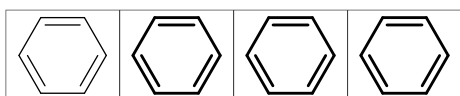
3.7.1 Local parameter setting

```
beginfigm
  drawm(Ph)
endfigm
beginfigm
  %-----
  ratio_thickness_bond:=0.05;
  %-----
  drawm(Ph)
endfigm
beginfigm
  drawm(Ph)
endfigm
```



3.7.2 Global parameter setting

```
beginfigm
  drawm(Ph)
endfigm
%-----
ratio_thickness_bond:=0.05;
%-----
beginfigm
  drawm(Ph)
endfigm
beginfigm
  drawm(Ph)
endfigm
beginfigm
  drawm(Ph)
endfigm
```



4 Command of mcf2graph

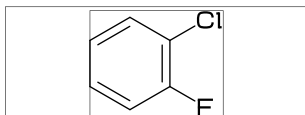
4.1 drawm

(Draw molecule)

```
msize=(a,b)      **default (1,1)
mposition=(c,d)   **default (0.5,0.5)
```

a: ratio molecular width/figure width
b: ratio molecular high/figure high
c: x axis position
d: y axis position

```
drawm(<30,Ph,3:/F,4:/Cl)
```

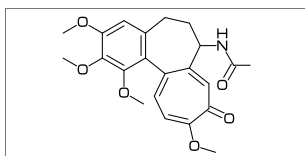


4.2 readm

```
readm(string1,string2, ...);
** string = mcf code
```

(example)

```
readm("<30,Ph,1'2'6:/O!,-4'-5=?7,",
      "-1'-4'-6=d1,-2:?0,-3:/O!,   ",
      "@9,! ,NH,! ,?0,!           ")
putm
```

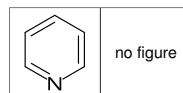


4.3 checkm

```
(immediately compile)
beginfigm drawm(<30,Ph,2:N) endfigm
```

(check mcf and compile)

```
beginfigm
  readm("<30,Ph,2:") % ** '2:' missing arg
  if checkm(mc)=0: drawm(scantokens(mc)) fi
endfigm
** checkm(mc) : error count
```



4.4 getm

```
getm(number)
** number = numeric
** ucount = molecular data unit count
```

```
for i=1 upto ucount:
  beginfigm
    getm(i)      % get data unit no=i
    putm         % put figure
  endfigm
endfor
```

```
getm("name"): "name"=string
```

(example)

```
beginfigm
  getm("Adenine")
  putm
endfigm
```

4.5 putm

putm : put figure

```
if op_row>=1: scantokens(op) fi
if mc_row>=1:
  if checkm(mc)=0:
    drawm(scantokens(mc))
  fi
fi
if ad_row>=1: add(scantokens(ad)) fi
if ex_row>=1: ext(scantokens(ex)) fi
```

4.6 allm

allm : put all loaded molecule

```
for i=1 upto ucount:
  beginfigm getm(i); putm; endfigm
endfor
```

4.7 add [++]

(Add label to molecule)

++(): add()

w: molecular width
h: molecular height
aw: atom font size
em: label font size
p0: origin of molecular structure
l: bond length

An: atom number
A[m]: atom position
A[m]ang: branch angle of A[m]
A[m]up: dir A[m]ang
A[m]left: dir A[m]ang+90
A[m]right: dir A[m]ang-90
A[m]down: dir A[m]ang+180

Bn: bond number
B[m]: bond(path)
B[m]s: bond start position
B[m]m: bond middle position
B[m]e: bond end position
B[m]ang: bond angle
B[m]up: dir B[m]ang
B[m]left: dir B[m]ang+90
B[m]right: dir B[m]ang-90
B[m]down: dir B[m]ang+180

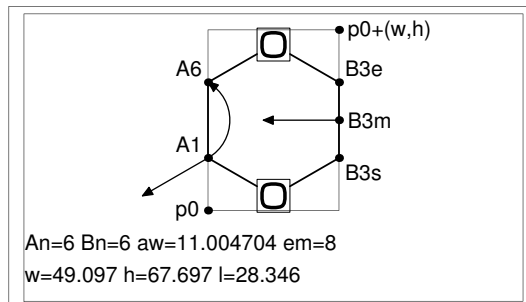
plus : '+' circled
minus : '-' circled
circlediam = 0.6aw (default)
circlepen = 0.2bp (default)

lonpair r: ':' rotated r
lonpairdiam = 0.3aw (default)
lonpairspace = 0.7aw (default)

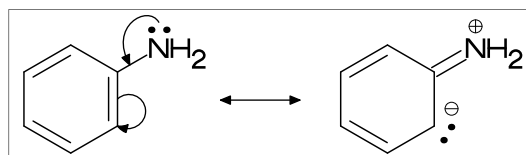
: scaled
<< : rotated
a /* b : point b of a

```
beginfigm
#(70mm,40mm) ##(.91,.9) |<(10mm)
sw_frame:=sw_frame+Atom+Mol;
@(.5,.85) drawm(<30,?6,{2,5}:0)
++(
defaultscale:=.8;
labeloffset:=.3aw;
dotlabel.lft("p0",p0);
dotlabel.rft("p0+(w,h)",p0+(w,h));
dotlabel.ulft("A1",A1);
drawarrow A1..A1+__*l<<A1ang;
dotlabel.lrt("B3s",B3s);
dotlabel.rft("B3m",B3m);
drawarrow B3m..B3m+__*l<<(B3ang+90);
dotlabel.ulft("A6",A6);
drawarrow A1{A1down}..A6;
```

```
dotlabel.rft("B3e",B3e);
label.rft("An"&decimal(An)&
" Bn"&decimal(Bn)&
" aw"&decimal(aw)&
" em"&decimal(em),
p0+(-9em,-1.5em));
label.rft("w"&decimal(w)&
" h"&decimal(h)&
" l"&decimal(l),
p0+(-9em,-3em));
)
endfigm
```



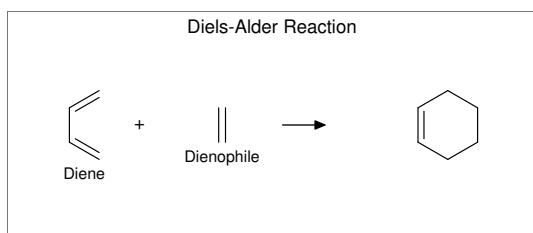
```
beginfigm
#(60mm,20mm) ##(1,0.85)
%-----
@ (0,0) drawm(<30,Ph,3=d1,4://NH2)
%-----
++(
labeloffset:=.7aw;
label.top(lonpair 90,A7);
drawarrow
(A7+up#1.2aw){A7left}
..{B7right}B7/*0.3;
drawarrow
B3m..A3+B2up#1.5aw..{A3down}A3;
)
%-----
@ (1,0) drawm(<30,?6,1'5=d1,4://NH2)
%-----
++(
labeloffset:=.7aw;
label.top(plus,A7);
label.rft(minus,A3);
label(lonpair A3ang,A3+A3up#.7aw);
)
%-----
** (drawdblarrow (.4w,.4h)..(.55w,.4h);)
%-----
endfigm
```



4.8 ext [**]

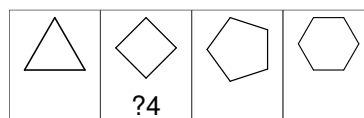
(Extra label to figure)

```
**(): ext()  
w:   figure width  
h:   figure height  
w0:  figure width-2xpart(fmargi  
h0:  figure height-2ypart(fmargi  
aw:  atom font size  
em:  label font size  
p0:  fmargin  
  
n:    molecular number  
p[m]: molecular origin position  
w[m]: molecular width  
h[m]: molecular height  
  
ratio_thickness_char:  
pen thickness / char width  
%-----  
beginfigm  
  #(70mm,30mm)  
  |= (0.065)  
  %-----  
  @(0.1,0.5)  
  drawm(<-210,60`1,60`1,60`1,{1,3}=dl)  
  ++(  
    defaultscale:=0.6;  
    label.bot("Diene",p0+(0.5w,0));  
  )  
  @(0.4,0.5)  
  drawm(<-30,-60`1,1=dm)  
  ++(  
    defaultscale:=0.6;  
    label.bot("Dienophile",p0+(.5w,0));  
  )  
  @(0.9,0.5)drawm(<30,?6,6=dl)  
  %-----  
  **(  
    drawarrow (.52w,.5h)..(.6w,.5h);  
    defaultscale:=0.7;  
    label("+", (0.25w,0.5h));  
    ratio_thickness_char:=0.125;  
    label.bot("Diels-Alder Reaction",  
              (.5w,h)  
            );  
  )  
  %-----  
endfigm
```



4.8.1 Local ext setting

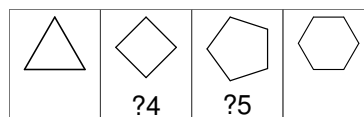
```
beginfigm  
  EN:="?3";@(0.5,1) drawm(<30,?3)  
endfigm  
beginfigm  
  EN:="?4";@(0.5,1) drawm(?4)  
  %-----  
  ** (label.top(EN,(0.5w,0));)  
  %-----  
endfigm  
beginfigm  
  EN:="?5";@(0.5,1) drawm(?5)  
endfigm  
beginfigm  
  EN:="?6";@(0.5,1) drawm(?6)  
endfigm
```



4.8.2 Global ext setting

ext_clear: reset global ext()

```
beginfigm  
  EN:="?3";@(0.5,1) drawm(<30,?3)  
endfigm  
%-----  
ext(label.top(EN,(0.5w,0));)  
%-----  
beginfigm  
  EN:="?4";@(0.5,1) drawm(?4)  
endfigm  
beginfigm  
  EN:="?5";@(0.5,1) drawm(?5)  
endfigm  
%-----  
ext_clear;  
%-----  
beginfigm  
  EN:="?6";@(0.5,1) drawm(?6)  
endfigm
```

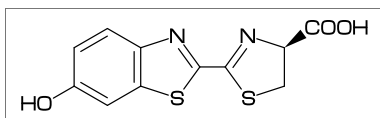


5 Example

5.1 drawm example

(Luciferin)

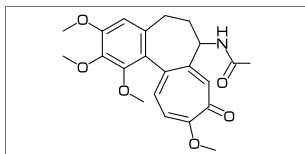
```
beginfigm
  #(50mm,15mm)
  drawm(<30,Ph,3=?5,@8,! ,?5,9'16=d1,9'14:N,7'11:S,1:/OH,-2:*/COOH)
endfigm
```



5.2 readm example

(Colchicine)

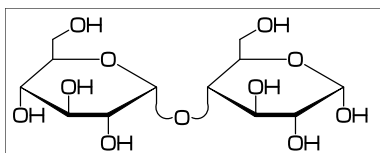
```
beginfigm
  readm(
    "<30,Ph,1'2'6:/O!,-4'-5=?7,      ",
    " -1'-4'-6=d1,-2:?0,-3:/O!,      ",
    " 9:/NH!'?0!                      ")
  #(40mm,20mm) putm
endfigm
```



(Maltose)

(bond type for glycan)
arc_lb : arc left > bottom
arc_br : arc bottom right

```
beginfigm
  %"EN:Maltose","MW:342.3",
  readm(
    %-----
    "hexose_hp,#.5,{1~$270'2~$90'3~$270}:/OH,6~$90:/!OH,##,",
    "@4,$310~arc_lb`1,0,$50~arc_br`1,<$0,      ",
    "|,hexose_hp,#.5,{2~$90'3~$270'4~$270}:/OH,6~$90:/!OH  ")
    %-----
  #(50mm,20mm) putm
endfigm
```



(Erythromycin)

```

beginfigm
  EN:=Erythromycin"; MW="733.93";
  #(60mm,30mm) @(1,0.5)
  readm(
    %-----
    "<30,#1,<-120,60,60,60,-60,60,60,-60,60,60,60,-60,60,60,##,&1,      ",
    " 14:0,13:/*!,1'9:~0,      ",
    " {~2'4'6~-35'8'*10'12~35}:~?z,      ",
    " {6~35'11'12~-35}:*/OH,      ",
    " @$3,!z,0,30~zb,|,~6~.7,6:0,#.5,{5~wf'3~35}:?,4:/*OH,3~-35:/*0! ,##," ,
    " @$5,!z~30~1.7,0,!~zb,|,~6~.7,6:0,#.5,5:~?z,2:*/OH,3:/*N?!      ")
    %-----
  putm
  **(defaultscale:=0.8;
    label.lrt("EN: "&EN,(0,h));
    label.lrt("fm: "&fm,(0,h-5mm));
    label.lrt("mw: "&mw,(0,h-9mm));
    label.lrt("MW: "&MW,(0,h-13mm));
  )
endfigm;

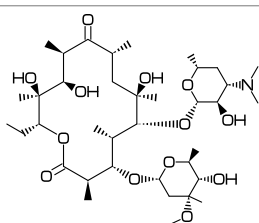
```

EN: Erythromycin

fm: C₃₇H₆₇NO₁₃

mw: 733.9267

MW: 733.93

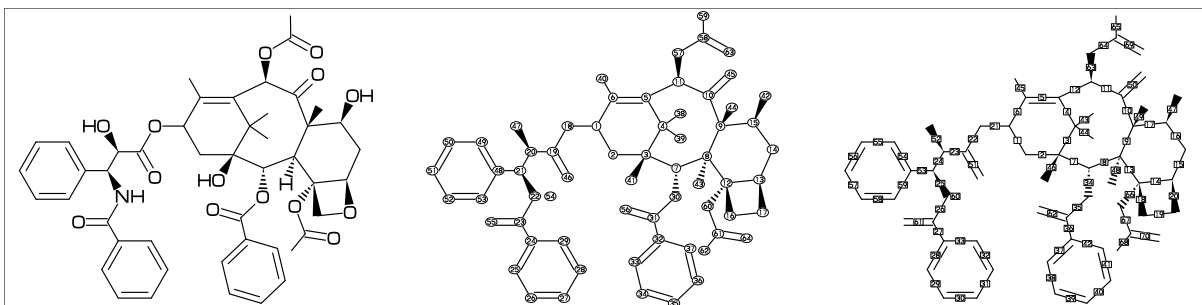


(Paclitaxel)

```

beginfigm
  % "EN:Paclitaxel", "MW:853.918",
  readm(
    %-----
    "?6,5=d1,@3,#1,36,45,45,45,45,##,&5,-4=?6,-4=?4,-1=wb,-3=wf,-1:0, ",
    " 4:??,6:?,{3~-60'15}:*/OH,8:/*H~-60,9:~w~60,10:~0,      ",
    " @1,! ,0,! ,?0,! ,*/OH,! ,/Ph,60~wf,NH,-60,~?0,60,Ph,      ",
    " @7,!z,0,-45,~?0,60,Ph,{11>r1'~12~-15>lr}:*/0!'?0!      ")
    %-----
  #(140mm,30mm)
  if checkm(mc)=0:
    @(0,0.5) drawm(scantokens(mc))
    sw_numbering:=Atom+All;
    @(0.6,0.5) drawm(scantokens(mc))
    sw_numbering:=Bond+All;
    @(1,0.5) drawm(scantokens(mc))
  fi
endfigm

```



5.3 loadm example

(Example)

```
loadm("CAT=biological","MW>=285","MW<=288","a:EN");
```

(output)

```
* jobname=mcf_exa_soc
* numbersystem=double
* output report file
* file name=mcf_exa_soc-report.txt)
* mcf_template 2023.05.07
* Input  : main_lib.mcf [525]
* Output : ucount [4]
* Filter(1): CAT =biological
* Filter(2): MW >= 285
* Filter(3): MW <= 288
* Sort key : EN (ascending)
[1]:Luteolin
[2]:Lycorine
[3]:Morphine
[4]:Piperine )

row[1][1]="CAT:biological;EN:Luteolin;MW:286.24;EXA:-"
row[1][2]=" "
row[1][3]="<30,Ph,3=?6,9=dl,10:0,7:?0,@9,! ,Ph,2'6'14'15:/OH"
row[1][4]=";"
row[2][1]="CAT:biological;EN:Lycorine;MW:287.315;EXA:1"
row[2][2]=" "
row[2][3]="<30,Ph,-4'-2=?6,6'(9--12)=?5,13=dl,8:N,15'17:0,"
row[2][4]="{*9^180'10^60}:*/H,13'*14:*/OH"
row[2][5]=";"
row[3][1]="CAT:biological;EN:Morphine;MW:285.343;EXA:1"
row[3][2]=" "
row[3][3]="<30,Ph,2'-4=?6,1---12=?5,-1:0,-1=zb,"
row[3][4]="@7,60~wf`0.75,70~si_`1.3,45,N!,&9~wb,15=dl,6:/OH,8^180:*/H,12:/*OH"
row[3][5]=";"
row[4][1]="CAT:biological;EN:Piperine;MW:285.343;EXA:1"
row[4][2]=" "
row[4][3]="<30,Ph,-1=?5,-1'-3:0,@4,! ,!d,! ,!d,! ,?0,! ,?6,-6:N"
row[4][4]=";"
%-----
```

(Tag)

CAT : category	USE : the use
EN : english name	JN : japanese name
FM : formula from data	MW : molecular weight from data
MI : monoisotopic mass from data	CAS : CAS number

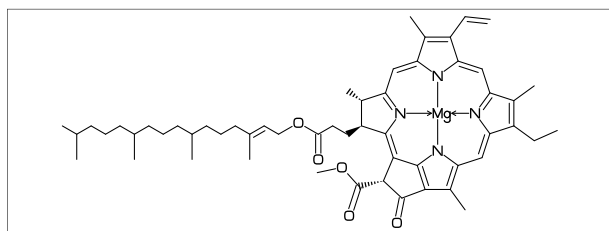
5.4 getm example

(Chlorophyll a)

```

beginfigm
  getm("Chlorophyll a")
  sw_output:=Fig+Mcode;
  #(80mm,30mm)
  drawm(scantokens(mc))
  VerbatimTeX("\gdef\EN{"&EN&"}\gdef\MW{"&MW&"}");
  VerbatimTeX("\gdef\mw{"&mw&"}\gdef\fm{"&fm&"}");
endfigm
\end{mplibcode}
\verbatiminput{temp-mc.aux}          %%% input temp-mc.aux %%%
{\tt ** EN:\EN \quad mw:\MW \quad MW:\mw \quad fm:\fm}%

```



```

<-36,#1,?5,@3,! ,54,?5,@-2,! ,54,?5,@-2,! ,54,?5,@-2,! ,&5,@6,22,70,&8,##,
4'6'8'10'14'16'18'21'23'27=d1,@4,!`1.48~vf,Mg,&17~vb,@11,&27,@27,&23,
4'11'17'23:N,{1~zf'9'15'21}:?:,14:/!,20:/!d,25:/*?0!'0!,26:?0,
@2,-6~wf,!2,?0!,0!2,!d,|,!13,1'5'9'13:?

```

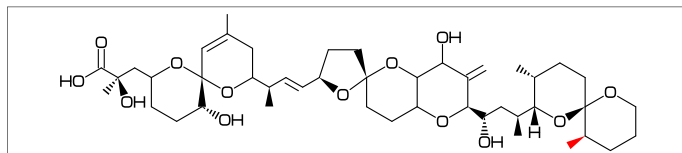
```
** EN:Chlorophyll a mw:893.509 MW:893.4889 fm:C55H72MgN4O5
```

(Dinophysistoxin-1)

```

beginfigm
  getm("Okadaic acid")
  readm(",38:?w,65=red")          %%% add methyl group (color red) %%%
  sw_output:=Fig+Mcode;          %%% output temp-mc.aux %%%
  EN:="Dinophysistoxin-1"; #(90mm,20mm)
  MW:="819";
  if checkm(mc)=0: drawm(scantokens(mc))
    VerbatimTeX("\gdef\EN{"&EN&"}\gdef\MW{"&MW&"}");
    VerbatimTeX("\gdef\mw{"&mw&"}\gdef\fm{"&fm&"}");
  fi
endfigm;
\end{mplibcode}
\verbatiminput{temp-mc.aux}          %%% input temp-mc.aux %%%
{\tt ** EN:\EN \quad mw:\MW \quad MW:\mw \quad fm:\fm}%

```



```

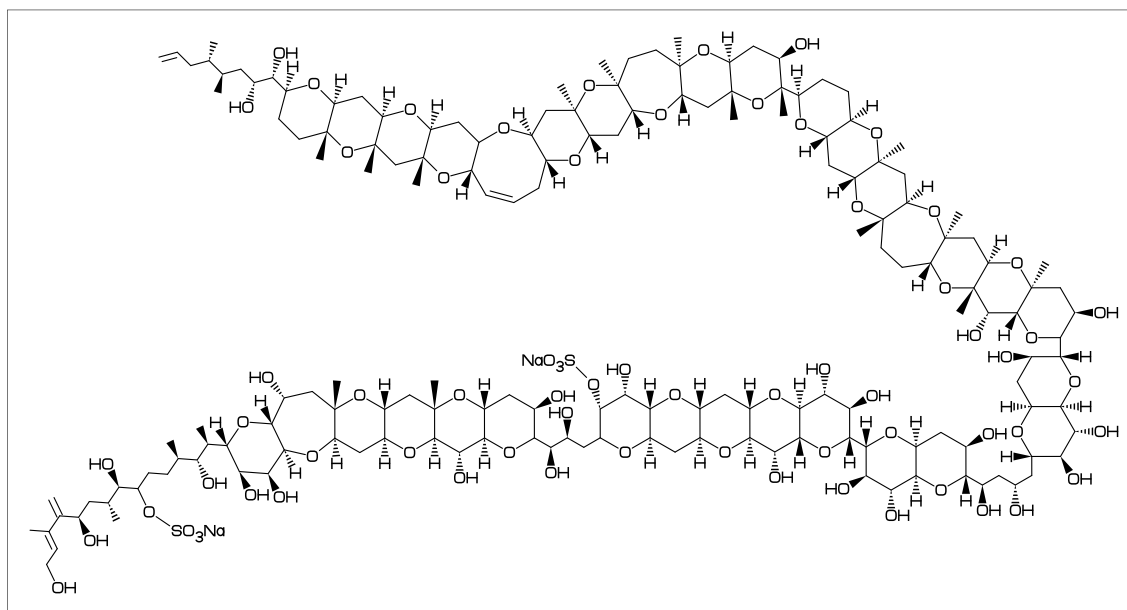
<30,?6,@4,?6,@-4,!4,<-12,?5,@-3,<-12,?6,-3=?6,@-3,!w,!3,
?6,@-4,?6,@6,!2,?z~-40,*/OH^20,!?0!,OH,
3'38=wb,11=d1,15=dr,17'19=wf,5'7'16'24'25'33'42:0,
32:*/H^60,10:?,12'31'*37:?w,27:?d,28:/OH,3'29:/*OH,
38:?w,65=red

```

```
** EN:Dinophysistoxin-1 mw:819 MW:819.0294 fm:C45H70O13
```

(Maitotoxin)

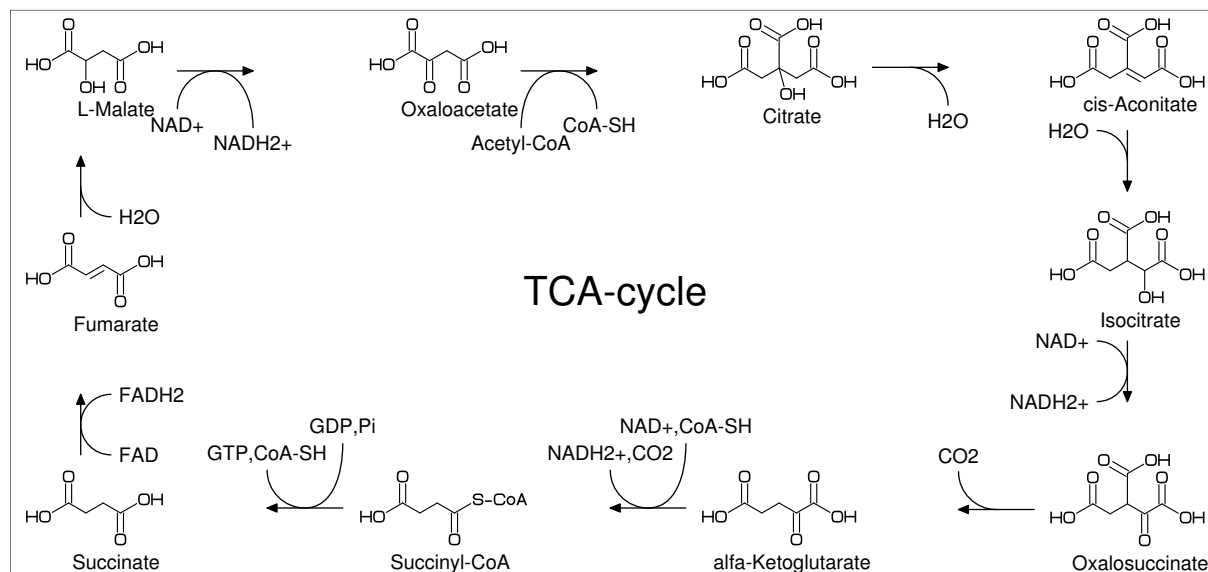
```
%-----
\begin{mplibcode}
  beginfigm
    getm("Maitotoxin") #(150mm,80mm) #@ (3mm,3mm)
    sw_output:=Fig+Mcode;          %%% output temp-mc.aux %%%
    sw_frame:=Outside;
    if checkm(mc)=0: drawm(scantokens(mc))
      VerbatimTeX("\gdef\EN{"&EN&"}\gdef\MW{"&MW&"}");
      VerbatimTeX("\gdef\mw{"&mw&"}\gdef\fm{"&fm&"}");
    fi
  endfigm
\end{mplibcode}
\verbatiminput{temp-mc.aux}          %%% input temp-mc.aux %%%
{\tt ** EN:\EN \quad mw:\MW \quad MW:\mw \quad fm:\fm}%
%-----
```



```
<55.8,?6,-4=?7,-4'-3'-3'-3=?6,@-3,!4,?6,-4'-3'-3'-3=?6,@-3,! ,?6,-3=?6,
@-3,!4,60,<-30,?6,-3=?6,@-3,30,<30,?6,-3'-3=?6,-3=?7,-4'-3'-3=?6,
@-2,! ,?6,-3=?6,-3=?7,-3'-3=?6,-3=?8,-3=d1,-5'-3'-3'-3=?6,
5'7'15'16'23'24'32'40'41'48'49'58'59'72'73'82'83'90'91'99'
100'107'113'114'122'123'130'131'140'141'148'149:0,
{1^60'2'26'28'29'51'54'61'63'68'75^60'78'109}:*/OH,
11'20'35'45'52'55'65'69'86:/*OH,
3'8'13'17'21'33'38'42'56'70'84'92'101'106'111'128'138'142'146'150:/*H^-60,
4'14'22'34'39'43'*47'*57'*71'81'89'98'102'116'121'125'129'133:/*H^60,
6'46'50'53'60'67'74:/*H^-60,
9'18'85'93'112'139'143'147:?w^60^1,80'88'97'*108'115'120'124:?z^-60^1,
@$6,! ,! ,!11,60~dr,-60,60,OH,*2'7'10:*/OH,1'3'*8:?w,11:?d,12:?,@6,!0,30,S03Na,
@$36,-45~zf,0,30,S03Na,
@$150,! ,! ,!7,1'2:/*OH,4:?w,5:?z,7=d1
```

```
** EN:Maitotoxin mw:3425.86 MW:3425.856 fm:C164H256Na2O68S2
```

5.5 User define parts example



```

beginfigm
#(160mm,75mm) |<(4mm)
COOH:=''(?0,! ,OH);      % define COOH
HOCO:=''(OH,! ,?0,);     % define HOCO
S_CoA:=''("{S-CoA}");    % define S_CoA
@(0.33, 1) drawm(<30,HOCO,! ,?0,!2,COOH)                % Oxaloacetate
@(0.66, 1) drawm(<30,HOCO,!4,COOH,@-4`1,! ,COOH,4:/OH~-165) % Citrate
@(1, 1) drawm(<30,HOCO,!2,!~dr,! ,COOH,@-4`1,! ,COOH)    % cis-Aconitate
@(1, 0.58) drawm(<30,HOCO,!4,COOH,@-4,!`1,COOH,5:/OH)    % Isocitrate
@(1, 0.05) drawm(<30,HOCO,!3,?0,! ,COOH,@-4,!`1,COOH)    % Oxalosuccinate
@(0.66,0.05) drawm(<30,HOCO,!3,?0,! ,COOH)                % alpha-Ketoglutarate
@(0.33,0.05) drawm(<30,HOCO,!3,?0,! ,S_CoA)               % Succinyl-CoA
@(0, 0.05) drawm(<30,HOCO,!3,COOH)                         % Succinate
@(0, 0.55) drawm(<30,HOCO,! ,!~dr,! ,COOH)                % Fumarate
@(0, 1) drawm(<30,HOCO,!3,COOH,3:/OH)                      % L-Malate
ext(
  defaultfont:="uhvr8r"; defaultscale:=0.75;
  ext_setup;
  save dx; pair dx; dx:=(12mm,0);
  label.bot("Oxaloacetate",p1+dx); label.bot("Citrate",p2+dx);
  label.bot("cis-Aconitate",p3+dx); label.bot("Isocitrate",p4+dx);
  label.bot("Oxalosuccinate",p5+dx); label.bot("alpha-Ketoglutarate",p6+dx);
  label.bot("Succinyl-CoA",p7+dx); label.bot("Succinate",p8+dx);
  label.bot("Fumarate",p9+dx); label.bot("L-Malate",p10+dx);
  sw_label_emu:=1;
  ext_setup;
  r_arrow(10mm)( 0)(p1+ ( 1.1w1, 0.3h1))("Acetyl-CoA",1.5)(" CoA-SH",1);
  r_arrow(10mm)( 0)(p2+ ( 1.1w2, 0.4h2))("","0")("H2O",1);
  r_arrow( 8mm)(270)(p3+ ( 0.5w3,-0.6h3))("H2O",1)("","0");
  r_arrow( 8mm)(270)(p4+ ( 0.5w4,-0.4h4))("NAD+",1)("NADH2+",1);
  r_arrow(10mm)(180)(p5+ (-0.1w5, 0.4h5))("","0")("CO_2_",1);
  r_arrow(10mm)(180)(p6+ (-0.1w6, 0.5h6))("NAD+,CoA-SH",1.7)("NADH2+,CO2",1);
  r_arrow(10mm)(180)(p7+ (-0.1w7, 0.5h7))("GDP,Pi",1.7)("GTP,CoA-SH",1);
  r_arrow( 8mm)( 90)(p8+ ( 0.4w8, 1.2h8))("FAD",1)("FADH2",1);
  r_arrow( 8mm)( 90)(p9+ ( 0.4w9, 1.2h9))("H2O",1)("","0");
  r_arrow(10mm)( 0)(p10+( 1.1w10,0.3h10))("NAD+",1)("NADH2+",1.5);
  defaultscale:=1.5;
  label("TCA-cycle",(0.5w,0.5h));
)
endfigm

```

6 Example to use mcf2graph

6.1 MetaPost souce file

```
%-----
input mcf2graph;                                > input main macro
%-----
%%% sw_output:=Report;                          > report output
%%% sw_output:=MOL2000;                         > MOL file output
#(60mm,40mm); % (figure width,figure height)    >
outputformat:="png"; hppp:=vppp:=0.1;          > PNG output
outputtemplate:="c%3c-%{EN_}.png";             >
%-----
beginfigm
  % EN:Ampicillin MW:349.405
  drawm(<45,74,-3=?5,2:N,7:S,
        3^45:/*H,1:?0^15,5:/*COOH^-18,6:??,
        @4,!w^15,NH,! ,?0,! ,/*NH2,! ,Ph)      > immediately compile
endfigm                                          >
%-----
beginfigm
  % EN:Cholesterol MW:386.65
  readm(                                         > read Mcode
    "<30,?6,-4'-2=?6,-4=?5,7=dl,      ",      > mc1
    "10:/*H^180,11:/*H^-60,17:/*H^-54,  ",    > mc2
    "4'12:?w^60,                          ",    > mc3
    "@-1,18,?z,-60,!3,?!                  ")    > mc4
  putm                                           > put figure
endfigm                                          >
%-----
loadm("EN<>");                                  > load all unit
beginfigm
  getm("Adenine")                              > get EN=Adenine
  putm                                           > ** put figure
endfigm                                          >
%-----
beginfigm
  getm(4)                                       > select No.4
  putm                                         > put figure
endfigm                                         >
%-----
for i=1 upto ucount:                           > figure count
  beginfigm
    getm(i)                                    > select No.i
    putm                                       > put figure
  endfigm
endfor
%-----
bye
```

6.2 Molecular library file

```
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
% molecular library file    main_lib.mcf    by Akira Yamaji    2022.10.10
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
% tag1:var1;tag2:var2;tag3:var3 ....
% first character of line "%" comment out
% first character of line ":" start MCF
% first character of line ";" stop MCF
% first character of line "=" start parameter setting
% first character of line "*" start ext(...)
% first character of line "+" start add(...)
% CAT = Category, EN = Name, MW = Molecular weight
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
CAT:-;EN:-;MW:0;EXA:-
=
  sw_frame:=Atom;
:
<30,?6,3=?5,1'3'5'9=dl,2'6'9:N,5:/NH2,7:NH
*
  defaultscale:=.5;
  label.bot(decimal(fig_num)&"."&EN,(.5w,0));
+
  defaultscale:=.3;
  label.bot("A2",A2) withcolor red;
  label.top("A6",A6) withcolor red;
  label.top("A9",A9) withcolor red;
;
%=====
CAT:biological;EN:Adenine;MW:135.13;EXA:1
:
<30,?6,3=?5,1'3'5'9=dl,2'6'9:N,5:/NH2,7:NH
;
%-----
CAT:biological;EN:Guanine;MW:151.13;EXA:1
:
<30,?6,3=?5,1'3'9=dl,2'9:N,6'7:NH,5:?0,1:/NH2
;
%-----
CAT:biological;EN:Cytosine;MW:111.10;EXA:1
:
<30,?6,4'6=dl,4:N,3:?0,2:NH,5:/NH2
;
%-----
CAT:biological;EN:Thymine;MW:126.11;EXA:1
:
<30,?6,3=dl,2'6:NH,1'5:?0,4:?
;
%-----
CAT:biological;EN:Uracil;MW:112.09;EXA:1
:
<30,?6,6=dl,3'5:?0,2'4:NH
;
%== Amino acid =====
CAT:biological;EN:Glycine;MW:75.07;EXA:-
:
<30,NH2,!2,COOH
;
%-----
```

6.3 MCF aux file output

(Option parameter setting)

```
sw_output:=Mcode;          %% output 'temp-mc.aux'
```

(Command line)

```
>mpost -s ahlength=3 FILENAME (sw_output=Fig Expand mode)
```

(Output mcf file)

```
sw_output=Mcode            %% file name = 'temp-mc.aux'
```

(result)

```
<30,?6,3=?5,1'3'5'9=d1,2'6'9:N,5:/NH2,7:NH
```

(LuaLaTeX example)

```
%-----  
%% "EN:Vancomycin  
\begin{mplibcode}  
  beginfigm  
    sw_output:=Mcode;      %%% output temp-mc.aux %%%  
  endfigm;  
\end{mplibcode}  
%-----  
\verbatiminput{temp-mc.aux}  
%-----
```

(result)

```
file name = 'temp-mc.aux'
```

```
<30,?6,@4,?6,@-4,!,!3,<-12,?5,@-3,<-12,?6,-3=?6,@-3,!w,!3,  
?6,@-4,?6,@6,!,!,?z^-40,*/OH^20,!,?0,!1,OH,  
3=wb,11=d1,15=dr,17'19=wf,38=wb,5'7'16'24'25'33'42:0,  
32:*/H^60,10:?,12'31:?w,27:?d,37:?z,28:/OH,3'29:/*OH
```

6.4 Report output

(Option parameter setting)

```
sw_output:=Report;          %% file name = 'jobname-report.aux'
```

(Command line)

```
>mpost -s ahlength=7 FILENAME
```

(Output)

```
=====
No[148],Name<Phenol>,Category<synthetic>,File<main_lib.mcf>
-----
Row[1],Length[12],Block[3],BackboneA[6],BackboneB[6],Group[1]
-----
<30,Ph,3:/OH
-----
[1 ] <30
[2 ] Ph
[3 ] 3:/OH
-----
Width[35.80607],Height[24.55503], Shift x[-1.77635],Shift y[-7.54719]
Bond length[11.33855],Atom size[4.881881]
-----
Atom[7],Bond[7],Ring[1],Hide H[5]
< NO. ><atom(s) >( x axis , y axis )<bond><hideH><chg>
A1 C ( 0 , 0 ) 3 1
A2 C ( 1 , -1 ) 3 1
A3 C ( 2 , 0 ) 4
A4 C ( 2 , 1 ) 3 1
A5 C ( 1 , 1 ) 3 1
A6 C ( 0 , 1 ) 3 1
A7 OH ( 3 , 0 ) 1
-----
< NO. >< bond (sdt)><angle +( +- )><length ( pt )>
B1 1 -> 2 ( 2) 330 ( -30) 1 ( 11)
B2 2 -> 3 ( 1) 30 ( 30) 1 ( 11)
B3 3 -> 4 ( 2) 90 ( 90) 1 ( 11)
B4 4 -> 5 ( 1) 150 ( 150) 1 ( 11)
B5 5 -> 6 ( 2) 210 ( -150) 1 ( 11)
B6 6 -> 1 ( 1) 270 ( -90) 1 ( 11)
B7 3 -> 7 ( 1) 330 ( -30)0.660000 ( 7)
-----
<atom>( atom wt ) [ mi wt ] < cnt > < sum wt > [ sum mi wt ]
C ( 12.0107) [ 12 ] * 6 72.0642 [ 72 ]
H (1.0079400) [1.0078250] * 6 6.04764 [6.0469501933]
O ( 15.9994) [15.994914] * 1 15.9994 [15.994914619]
Molecular Weight [Mono Isotopic] = 94.11123 [ 94.0418648]
-----
Weight Calc: 94.11123 - Input: 94.11 = 0.0012399
Formula Calc: C6H6O
=====
```

6.5 MOL file output

(Option parameter setting)

```
sw_output:=MOL2000;    % MOL(V2000)
sw_output:=MOL3000;    % MOL(V3000)
```

(Command line)

```
>mpost -s ahlength=5  FILENAME    % MOL(V2000)
>mpost -s ahlength=6  FILENAME    % MOL(V3000)
```

(Output)

```
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
-MCFtoMOL- EN:Caffeine

14 15  0  0  0  0  0  0  0  0999 V2000
      0      0      0 C  0  0  0  0
  0.86603    -0.5      0 N  0  0  0  0
  1.73206      0      0 C  0  0  0  0
  1.73206      1      0 C  0  0  0  0
  0.86603     1.5      0 C  0  0  0  0
      0      1      0 N  0  0  0  0
  2.6831   -0.30902    0 N  0  0  0  0
  3.27089      0.5      0 C  0  0  0  0
  2.6831    1.30902    0 N  0  0  0  0
  0.86603   -1.36383    0 C  0  0  0  0
 -0.76894    1.44394    0 C  0  0  0  0
 -0.76894   -0.44394    0 D  0  0  0  0
  0.86603    2.36383    0 D  0  0  0  0
  2.95299    2.1396     0 C  0  0  0  0
  1  2  1  0      0  0
  2  3  1  0      0  0
  3  4  2  0      0  0
  4  5  1  0      0  0
  5  6  1  0      0  0
  6  1  1  0      0  0
  3  7  1  0      0  0
  7  8  2  0      0  0
  8  9  1  0      0  0
  9  4  1  0      0  0
  2 10  1  0      0  0
  6 11  1  0      0  0
  1 12  2  0      0  0
  5 13  2  0      0  0
  9 14  1  0      0  0
M  END
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
```


6.6 LuaTeX file example

```

\documentclass{article}
\usepackage{luamplib}%
\usepackage[T1]{fontenc}%
\usepackage{textcomp}%
\mplibcodeinherit{enable}%
\mplibverbatim{enable}%
\mplibnumbersystem{double}%
\begin{mplibcode}
\end{mplibcode}
\begin{document}
\noindent%
%-----
\begin{mplibcode}
  input mcf2graph;
  sw_output:=Fig;
  max_blength:=4.5mm;
  defaultfont:="uhvr8r";
  defaultsize:=8bp;
  defaultscale:=1;
  %-----
  EN:="Limonin";
  MW="470.51";
  beginfigm
    #(50mm,50mm)
    readm(
      %-----
      "<30,?6,-3'-4=?6,          ",
      " -5=?3,-2=wf,-1=wb,6=?5,-4=?6,-5=wf,  ",
      " 13'15'17'20:0,3'12'21:?0,          ",
      " {4~wf^60'8~zf^60'18^35'18^-35}:?,      ",
      " {1^60'5^180'16^60}:/*H,          ",
      " @14,!z,|,?5,1'4=d1,3:0          ")
      %-----
    putm
  endfigm
\end{mplibcode}\\
%-----
\begin{mplibcode}
  EN:="beta-carotene";
  MW:="536.87";
  beginfigm
    #(80mm,50mm)
    readm(
      %-----
      "<30,?6,3=d1,{3'5^35'5^-35}:?,          ",
      " @4,! ,!18,1'3'5'7'9'11'13'15'17=dr,  ",
      " 3'7'12'16:?,          ",
      " |,?6,6=d1,{6'2^35'2^-35}:?          ")
      %-----
    putm
  endfigm
\end{mplibcode}\\
%-----
\end{document}

```

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