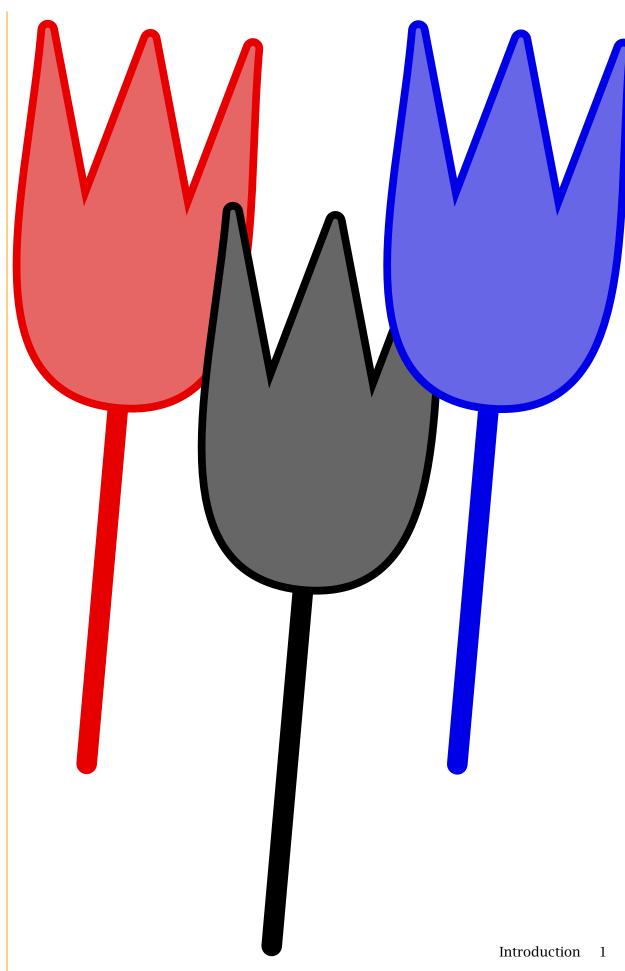
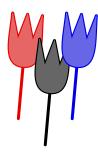
CONT_EXT up-to-date 1998/2

PPCHT_EX examples



PRAGMA ADE Ridderstraat 27 8061GH Hasselt NL



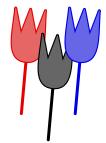
Introduction

This second issue of ConT_EXt up-to-date is dedicated to PPCHT_EX. Most of the chemical examples presented here originate from a range of learning materials that PRAGMA ADE developed and typeset through a number of years. The last few years, examples were donated by Tobias Burnus, Dirk Kuipers et.al. and we offer our thanks to their contribution. Special thanks goes to Richard Müller, who provided some DNA/RNA components. In all we think that the novice user of PPCHT_EX will find this document a useful tool to become more acquainted with PPCHT_EX.

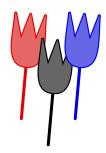
We have tried to categorize and index the structures and equations, but in the end it seemed best to let the reader leave through these pages until a structure or an equation comes up that looks (somewhat) simular to the one he wants to typeset. Keep in mind that many examples come from educational documents and therefore are not always correct due to the purpose they serve.

This file is generated by pdfT_EX, using CONT_EXT and PPCHT_EX in combination with P_fCT_EX (positioning and text) and METAPOST (drawing graphics). The use of METAPOST as graphics engine permits us to think of more advanced features: suggestions are welcome.

If you can find no alternative for your structure or equation, you can contact us via the mailing list of PPCHT_EX. We will come back to you as soon as possible.



```
\chemical{\TL{\oplus}{CH}} \chemical{\L {\oplus}{CH}}
\chemical{\LC{\oplus}{CH}} \chemical{\BL{\oplus}{CH}}
\chemical{TR{\circ}us}{CH}} \chemical{R {\circ}us}{CH}}
\chemical{\RC{\oplus}{CH}} \chemical{\BR{\oplus}{CH}}
\chemical{\LT{\oplus}{CH}} \chemical{\T {\oplus}{CH}}
\chemical{\RT{\oplus}{CH}} \chemical{\LB{\oplus}{CH}}
\chemical{\B {\oplus}{CH}} \chemical{\RB{\oplus}{CH}}
\blank
\chemical{\TL{\delta}{CH}} \chemical{\L {\delta}{CH}}
\chemical{\LC{\delta}{CH}} \chemical{\BL{\delta}{CH}}
\chemical{\TR{\delta}{CH}} \chemical{\R {\delta}{CH}}
\chemical{\RC{\delta}{CH}} \chemical{\BR{\delta}{CH}}
\chemical{\LT{\delta}{CH}} \chemical{\T {\delta}{CH}}
\chemical{\RT{\delta}{CH}} \chemical{\LB{\delta}{CH}}
\chemical{\B {\delta}{CH}} \chemical{\RB{\delta}{CH}}
\blank
\chemical{\X\TL{\oplus}{CH}} \chemical{\X\L {\oplus}{CH}}
\chemical{\X\LC{\oplus}{CH}} \chemical{\X\BL{\oplus}{CH}}
\chemical{X\TR{\circplus}{CH}} \chemical{X\R {\circplus}{CH}}
\chemical{\X\RC{\oplus}{CH}} \chemical{\X\BR{\oplus}{CH}}
\chemical{\X\LT{\oplus}{CH}} \chemical{\X\T {\oplus}{CH}}
\chemical{\X\RT{\oplus}{CH}} \chemical{\X\LB{\oplus}{CH}}
\chemical{\X\B {\oplus}{CH}} \chemical{\X\RB{\oplus}{CH}}
\blank
\chemical{\X\TL{\delta}{CH}} \chemical{\X\L {\delta}{CH}}
\chemical{\X\LC{\delta}{CH}} \chemical{\X\BL{\delta}{CH}}
\chemical{X\TR{\delta}{CH}} \chemical{X\R {\delta}{CH}}
\chemical{\X\RC{\delta}{CH}} \chemical{\X\BR{\delta}{CH}}
\chemical{X\LT{\delta}{CH}} \chemical{X\T {\delta}{CH}}
\chemical{\X\RT{\delta}{CH}} \chemical{\X\LB{\delta}{CH}}
\chemical{\X\B {\delta}{CH}} \chemical{\X\RB{\delta}{CH}}
```









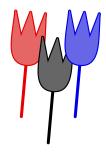
\startchemical \chemical[ONE,SB,Z0,ZTN][C] \stopchemical

\hss

\startchemical \chemical[ONE,SB,Z0,ZBN][C] \stopchemical

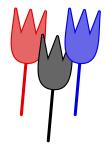
\hss

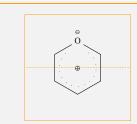
\startchemical \chemical[ONE,SB,Z0,ZTT][C,a,b,c,d,e,f,g,h] \stopchemical

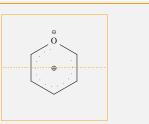




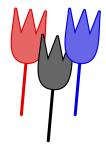
```
\startchemical
  \chemical
    [ONE,ZT5,Z0,SB15,MOV1,Z0,SB15,MOV1,SB15,Z0,ZT1]
    [\[,C,C,C,\]{5}]
\stopchemical
```

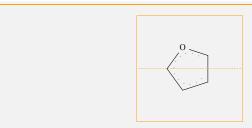


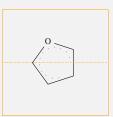




```
\hbox to \hsize \bgroup \hss
\startchemical
  \chemical
    [SIX,B1234,CD1..5,+SB5,-SB6,Z6,Z0]
    [\T{\ominus}{0},\oplus]
\stopchemical
\hss
\startchemical
  \chemical
    [SIX,B1234,CCD1..5,+SB5,-SB6,Z6,Z0]
    [\T{\ominus}{0},\oplus]
\stopchemical
\hss \egroup
```







\startchemical

\chemical[FIVE,B123,CD1235,+SB4,-SB5,Z5][0]

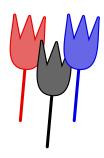
\stopchemical

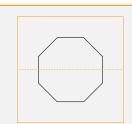
\hss

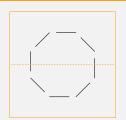
\startchemical

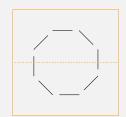
\chemical[FIVE,B123,CCD1235,+SB4,-SB5,Z5][0]

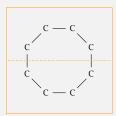
\stopchemical











\startchemical

\chemical[EIGHT,B]

\stopchemical

\hss

\startchemical

\chemical[EIGHT,-SB]

\stopchemical

\hss

\startchemical

\chemical[EIGHT,+SB]

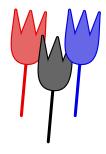
\stopchemical

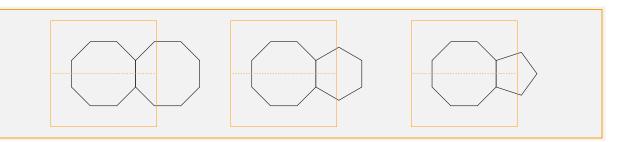
\hss

\startchemical

\chemical[EIGHT,SB,Z][C,C,C,C,C,C,C,C]

\stopchemical





\startchemical

\chemical[EIGHT,B,MOV1,B]

\stopchemical

\hss

\startchemical

\chemical[EIGHT,B,ADJ1,SIX,B]

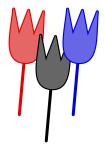
\stopchemical

\hss

\startchemical

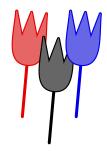
\chemical[EIGHT,B,ADJ1,FIVE,ROT3,B]

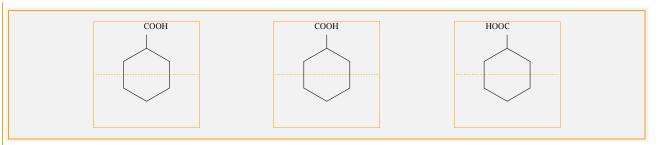
\stopchemical



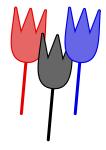
```
COOH
                                                                                                     СООН
CH_3
           CH<sub>3</sub>
                            CH_3
                                                                                               СООН
                                             сн<sub>3</sub> соон
                                                      CH<sub>3</sub> COOH
                                                                                                                         COOH
CH_3
                CH<sub>3</sub>
                                                                                  соон
                                                                                                    СООН
         CH<sub>3</sub>
                                             CH<sub>3</sub>
                                                                           соон
```

```
\setupchemical
 [width=5000]
\hbox to \hsize \bgroup \hss
\startchemical[scale=small]
 \chemical
   [CHAIR, B, +R, -R,
    +RZ1,+RZ2,+RZ3,+RZ4,+RZ5,+RZ6,
    -RZ1,-RZ2,-RZ3,-RZ4,-RZ5,-RZ6]
   [a,b,c,d,e,f,1,2,3,4,5,6]
\stopchemical
\hss
\startchemical[scale=small]
 \chemical
   [CHAIR, B, +R, -R,
    +RZ1,+RZ2,+RZ3,+RZ4,+RZ5,+RZ6,
    -RZ1,-RZ2,-RZ3,-RZ4,-RZ5,-RZ6]
   \stopchemical
\hss
\startchemical[scale=small,size=small]
 \chemical
   [CHAIR, B, +R, -R,
    +RZ1,+RZ2,+RZ3,+RZ4,+RZ5,+RZ6,
    -RZ1,-RZ2,-RZ3,-RZ4,-RZ5,-RZ6]
   [COOH, COOH, COOH, COOH, COOH,
    COOH, COOH, COOH, COOH, COOH]
\stopchemical
\hss \egroup
```



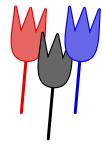


```
\hbox to \hsize \bgroup \hss
\startchemical
  \chemical[SIX,B,R6,RZ6][\SL{COOH}]
\stopchemical
  \hss
\startchemical
  \chemical[SIX,B,R6,RZ6][\SM{COOH}]
\stopchemical
  \hss
\startchemical
  \chemical[SIX,B,R6,RZ6][\SR{HOOC}]
\stopchemical
  \chemical[SIX,B,R6,RZ6][\SR{HOOC}]
\stopchemical
  \hss \egroup
```



```
S + O_2
                                    SO_2
                   \Delta \text{T/H}_2\text{O}
                 + δC + Cδ
```

```
\startformula
  \chemical{S}+\chemical{0_2}
  \chemical{GIVES}{\Delta T / H_20}
  \chemical{+{4}{S}}-{2}{0_2}}
\stopformula
\startformula
  \chemical{T\delta{C}} +
  \chemical{B\delta(C)} +
  \chemical{L\delta{C}} +
  \chemical{\R\delta{C}}}
\stopformula
\startformula
  \chemical{T\circ L} +
  \chemical{B\circ Plus(C)} +
  \chemical{\L\circ plus{C}} +
  \chemical{\R\oplus{C}}}
\stopformula
```

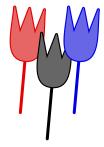


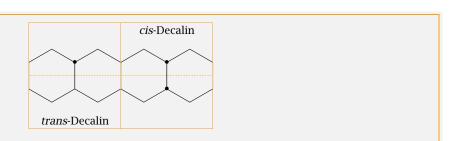


\setupchemical

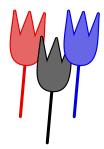
[height=fit,width=1500]

```
\hbox to \hsize \bgroup
                                                                \hss
\startchemical \chemical[ONE,SB]
                                                 \stopchemical \hss
\startchemical \chemical[ONE,30FF1,SB]
                                                 \stopchemical \hss
\startchemical \chemical[ONE,MOV1,SB]
                                                 \stopchemical \hss
\startchemical \chemical[ONE,30FF1,MOV1,SB]
                                                 \stopchemical \hss
\startchemical \chemical[ONE,MOV1,30FF1,SB]
                                                 \stopchemical \hss
\startchemical \chemical[ONE,MOV1,30FF1,0FF0,SB] \stopchemical \hss
\startchemical \chemical[ONE,MOV1,30FF1,MOV0,SB] \stopchemical \hss
\startchemical \chemical[ONE,MOV1,MOV0,SB]
                                                 \stopchemical \hss
                \egroup
```





```
\startformula
  \startchemical[width=fit]
  \chemical[SIX,B,Z1,MOV1,B][\hbox{$\bullet$}]
  \bottext{{\sl trans}-Decalin}
  \stopchemical
  \startchemical[width=fit]
  \chemical[SIX,B,Z12,MOV1,B][\hbox{$\bullet$},\hbox{$\bullet$}]
  \toptext{{\sl cis}-Decalin}
  \stopchemical
  \stopformula
```



In display mode:

$$K_b = \frac{[H_3O^+][OH^-]}{[H_2O]}$$

In text mode: $\frac{K_b = [H_3O^+][OH^-]}{[OH^-]}$

With adjusted brackets:
$$\frac{K_b = \left[H_3 O^+ \right] \left[OH^- \right]}{\left[H_2 O \right]}$$

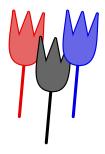
In display mode:

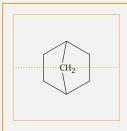
```
\startformula
```

```
K_b = \{\{[\langle H_30^+\}] [\langle H_20^+\}]\}
\stopformula
```

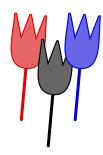
```
In text mode: K_b = {[\chemical {H_30^+}] [\chemical {OH^-}]} \over
{[\chemical {H_20}]}$
```

```
With adjusted brackets: K_b = \{ \bigcup_{k=0}^{H_30^+} \bigcup_{k=0
\{OH^-\} \setminus \{OH^-\} \}  \over \{ \bigcup_{i=1}^{H_20} \}
```





\startchemical
 \chemical[SIX,B,MID,MIDZ][\SL{CH_2}]
\stopchemical



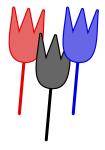
 $\rm H_2O$ or $\rm H_2O$ when typeset in display mode becomes:

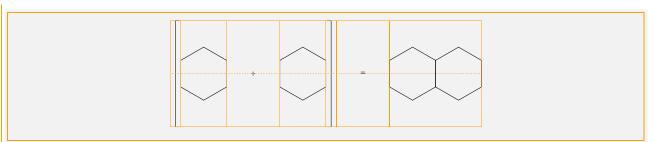
nquia

 $\begin{array}{ccc} \mathrm{H_2O} & \mathrm{or} & \mathrm{H_2O} \\ \mathrm{water} & & \mathrm{water} \end{array}$

 $\chemical{H_20}{liquid}{water} or $\chemical{H_20}{liquid}{water}$ when typeset in display mode becomes:}$

 $\label{lem:cal_H_20} $$ \operatorname{H_20}_{\quad \chemical_H_20}_{\quad \chemical_H_20$



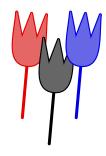


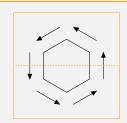
\startformula

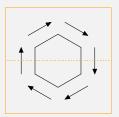
```
\startchemical[width=fit]
                           \chemical[OPENCOMPLEX]
                                                   \stopchemical
\startchemical[width=fit]
                           \chemical[SIX,B]
                                                   \stopchemical
\startchemical[width=2000] \chemical[PLUS]
                                                   \stopchemical
\startchemical[width=fit]
                           \chemical[SIX,B]
                                                   \stopchemical
                           \chemical[CLOSECOMPLEX] \stopchemical
\startchemical[width=fit]
\startchemical[width=2000] \chemical[EQUAL]
                                                   \stopchemical
\startchemical[width=fit] \chemical[SIX,B,MOV1,B] \stopchemical
```

\stopformula

17



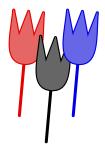




\startchemical \chemical[SIX,B,AU] \stopchemical

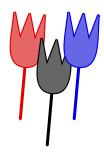
\hss

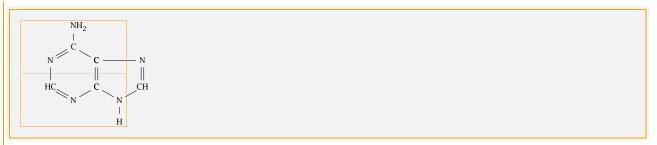
\startchemical \chemical[SIX,B,AD] \stopchemical





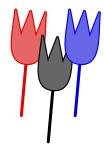
\startchemical
 \chemical[ONE,Z0,ES1,ED3,ET5][C]
\stopchemical

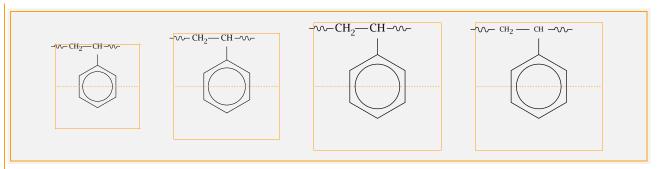




\startchemical

\chemical[SIX,DB135,SB246,Z,SR6,RZ6][C,C,N,HC,N,C,NH_2] \chemical[SIX,MOV1,DB1,SB23,SS6,Z1..5,SR3,RZ3][N,CH,N,C,C,H] \stopchemical





\startchemical[scale=small, size=small]

 $\label{lem:cal_six_b_c_red} $$ \operatorname{SIX}_B, C, R6, PB: RZ6, ONE, CZ0, OE1, SB5, MOV5, CZ0, OFF5, OE5, PE] [CH, CH_2] $$ \operatorname{CPChemical} $$$

\hss

\startchemical[scale=medium, size=medium]

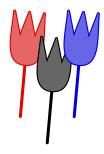
\hss

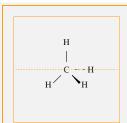
\startchemical[scale=big,size=big]

\hss

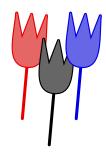
\startchemical[scale=big,size=small]

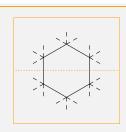
\chemical[SIX,B,C,R6,PB:RZ6,ONE,CZ0,OE1,SB5,MOV5,CZ0,OFF5,OE5,PE][CH,CH_2]\stopchemical

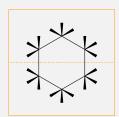


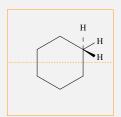


\startchemical \chemical[ONE,SD1,SB4,BB2,SB7,Z01247][C,H,H,H,H] \stopchemical









```
\hbox to \hsize \bgroup \hss
```

\startchemical

\chemical[SIX,B,-RD,RD,+RD]

\stopchemical

\hss

\startchemical

\chemical[SIX,B,-RB,RB,+RB]

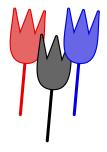
\stopchemical

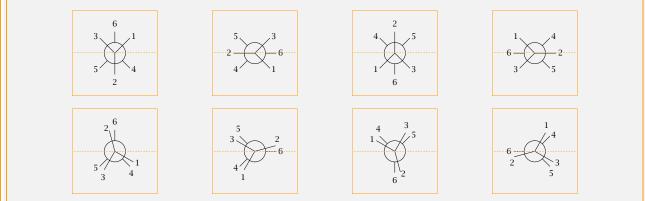
\hss

\startchemical

\chemical[SIX,B,-RD1,R1,+RB1,-RZ1,RZ1,+RZ1][H,H,H]

\stopchemical





\setupchemical[scale=small]

\hbox to \hsize \bgroup \hss

\startchemical\chemical[NEWMAN,STAGGER,ROT1,CB][1,2,3,4,5,6]\stopchemical\hss \startchemical\chemical[NEWMAN,STAGGER,ROT2,CB][1,2,3,4,5,6]\stopchemical\hss \startchemical\chemical[NEWMAN,STAGGER,ROT3,CB][1,2,3,4,5,6]\stopchemical\hss \startchemical\chemical[NEWMAN,STAGGER,ROT4,CB][1,2,3,4,5,6]\stopchemical\hss

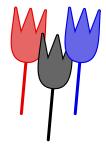
\egroup

\blank

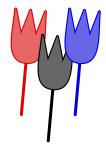
\hbox to \hsize \bgroup \hss

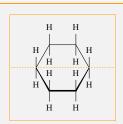
\startchemical\chemical[NEWMAN,ECLIPSE,ROT1,CB][1,2,3,4,5,6]\stopchemical\hss \startchemical\chemical[NEWMAN,ECLIPSE,ROT2,CB][1,2,3,4,5,6]\stopchemical\hss \startchemical\chemical[NEWMAN,ECLIPSE,ROT3,CB][1,2,3,4,5,6]\stopchemical\hss \startchemical\chemical[NEWMAN,ECLIPSE,ROT4,CB][1,2,3,4,5,6]\stopchemical\hss

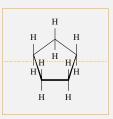
\egroup



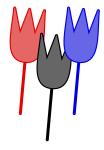
```
In display mode we get:
\startformula
  \chemical{A,GIVES,B,EQUILIBRIUM,C,MESOMERIC,D}
\stopformula
or split over more commands:
\startformula
  \chemical{A} \chemical{GIVES}
  \chemical{B} \chemical{EQUILIBRIUM}
  \chemical{C} \chemical{MESOMERIC}
  \chemical{D}
\stopformula
In text mode we get $\chemical {ABC,->,DEF,<->,GHI,<>,KLM}$ but in
display mode we see:
\startformula
  \chemical{ABC} \chemical{->}
  \chemical{DEF} \chemical{<->}
  \chemical{GHI} \chemical{<>}
  \chemical{KLM}
\stopformula
```

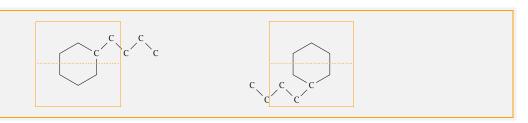






```
\hbox to \hsize \bgroup \hss
\startchemical
  \chemical
    [SIX, FRONT, B, BB612, -R, +R, -RZ, +RZ]
    [H,H,H,H,H,H,H,H,H,H,H]
\stopchemical
\hss
\startchemical
  \chemical
    [FIVE, FRONT, BB, -R, +R, -RZ, +RZ]
    [H,H,H,H,H,H,H,H,H,H,H]
\stopchemical
\hss \egroup
```

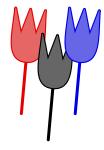


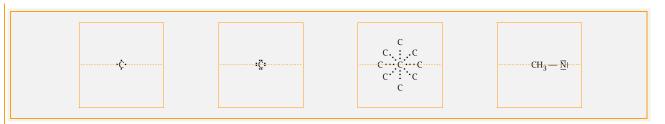


```
\hbox to \hsize \bgroup \hss
\startchemical[scale=small]
\chemical
   [SIX,B2345,+SB6,-SB1]
   \chemical
   [PB:Z1,ONE,CZ0,
    SB8, DIR8, CZ0,
    SB2,DIR2,CZ0,
    SB8, DIR8, CZ0,
    SB2,DIR2,CZ0,PE]
   [C,C,C,C,C,C]
\stopchemical
\hss
\startchemical[scale=small]
  \chemical
    [SIX,B4561,+SB2,-SB3]
    \chemical
    [PB:Z3,ONE,CZ0,
     SB4, DIR4, CZ0,
     SB6, DIR6, CZ0,
     SB4, DIR4, CZ0,
     SB6,DIR6,CZ0,PE]
    [C,C,C,C,C,C]
\stopchemical
```

\hss \egroup

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```
\hbox to \hsize \bgroup \hss
\setupchemical
```

[scale=small]

\startchemical

\chemical[ONE,Z0,ES1357][C]

\stopchemical

\hss

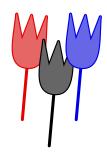
\startchemical \chemical[ONE,Z0,ED1357][C] \stopchemical

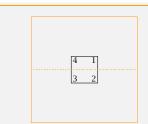
\hss

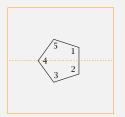
\startchemical \chemical[ONE,Z0,HB,Z][C,C,C,C,C,C,C,C,C] \stopchemical

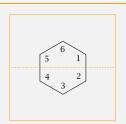
\hss

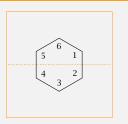
\startchemical \chemical[ONE,Z0,OFF1,SB1,MOV1,Z0,EP137][CH_3,N] \stopchemical





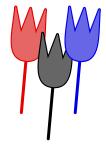




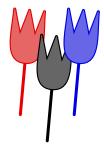


\egroup

```
\label{thm:cal_chemical_four_balance} $$ \operatorname{chemical_FOUR_B_ZN_C,C,C,C,C_S} \hookrightarrow \operatorname{chemical_FIVE_B_ZN_C,C,C,C,C_S} \hookrightarrow \operatorname{chemical_Chemical_SIX_B_ZN_C,C,C,C,C_S} \hookrightarrow \operatorname{chemical_Chemical_SIX_B_ZN_C,C,C,C,C_S} \hookrightarrow \operatorname{chemical_Chemical_SIX_B_ZN_C,C,C,C,C_S} \hookrightarrow \operatorname{chemical_Chemical_SIX_B_ZN_C} = 1,2,3,4,5,6] \hookrightarrow \operatorname{chemical_Chemical_Chemical_C} \hookrightarrow \operatorname{chemical_Chemical_C} \hookrightarrow \operatorname{chemical_C} \hookrightarrow \operatorname{chemical_Chemical_C} \hookrightarrow \operatorname{chemical_C} \hookrightarrow \operatorname{
```

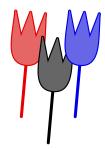


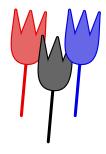
```
\startchemical[scale=small,width=fit,height=fit]
  \chemical
    [SIX,SB2356,DB14,Z2346,SR36,RZ36][C,N,C,C,H,H_2]
 \chemical
    [PB:Z1,ONE,Z0,DIR8,Z0,SB24,DB7,Z27,PE][C,C,CH_3,O]
 \chemical
    [PB:Z5,ONE,Z0,DIR6,Z0,SB24,DB7,Z47,PE][C,C,H_3C,O]
 \chemical
    [SR24, RZ24] [CH_3, H_3C]
\stopchemical
\hss
\startchemical[scale=small,width=fit,height=fit]
  \chemical
    [SIX,SB2356,DB14,Z36,SR36,RZ36][N,C,H,H_2]
 \chemical
    [PB:Z1,ONE,Z0,DIR8,Z0,SB24,DB7,Z27,PE][C,C,CH_3,O]
 \chemical
    [PB:Z5,ONE,Z0,DIR6,Z0,SB24,DB7,Z47,PE][C,C,H_3C,O]
 \chemical
    [PB:Z2,ONE,Z0,DIR2,SB6,CZ0,PE][C,CH_3]
 \chemical
    [PB:Z4,ONE,Z0,DIR4,SB8,CZ0,PE][C,CH_3]
\stopchemical
\hss
\startchemical[scale=small,width=fit,height=fit]
 \chemical
    [SIX,SB2356,DB14,Z,SR36,RZ36,SR1245,RZ24][C,C,N,C,C,C,H,H_2,CH_3,H_3C]
 \chemical
    [PB:RZ1,ONE,Z0,SB2,DB7,Z27,PE][C,CH_3,O]
 \chemical
    [PB:RZ5,ONE,Z0,SB4,DB7,Z47,PE][C,H_3C,O]
\stopchemical
\hss \egroup
```



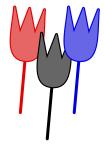
31

```
\hbox to \hsize \bgroup \hss
\startchemical[height=6000,width=fit]
 \bottext{2,2-Dimethyl-3-etylpentan}
 \chemical
    [ONE, Z3570, SB1357]
                                           [H_3C, T_1]\{CH_3\}, H_3C, TL_{2}\{CH\}]
 \chemical
    [MOV1,OFF1,Z0,SB3]
                                           [\T{3}{CH_2}]
 \chemical
    [MOV3, Z0, SB3, MOV3, Z0, MOV7, MOV7]
                                           [CH_2,CH_3]
    [OFF1,SB1,MOV1,OFF1,Z0,20FF1,SB1,Z1] [\T{4}{CH_2},\T{5}{CH_3}]
\stopchemical
\hss
\startchemical[height=6000,width=fit]
 \bottext{2,2-Dimethyl-3-etylpentan}
                                                  [H_3C, XT_{1}_{CH_3}, H_3C, XT_{2}_{C}]
 \chemical[ONE, Z3570, SB1357]
 \chemical[MOV1,OFF1,Z0,SB3]
                                                  [XT{3}{CH_2}]
 \chemical[MOV3,Z0,SB3,MOV3,Z0,MOV7,MOV7]
                                                  [CH_2,CH_3]
 \chemical[OFF1,SB1,MOV1,OFF1,Z0,20FF1,SB1,Z1] [\X\T{4}{CH_2},\X\T{5}{CH_3}]
\stopchemical
\hss \egroup
```



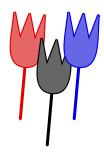


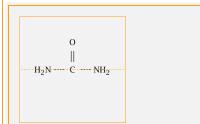
```
HO — CHOH — CH_2NH_2 HO — CHOH — CH_2NH_2
```



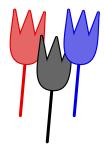
$$\begin{array}{c|c}
C & & & \\
H & & & \\
C & & & \\
O & & & \\
C_{2}H_{5} & & \\
C_{6}H_{5} & & \\
\end{array}$$

```
\startformula
  \startchemical
    \chemical
      [SIX,SB1..6,Z1..6,
       SR15, DR246, -SR3, +SR3,
       RZ12,-RZ3,+RZ3,RZ456]
      [N,C,C,C,N,C,H,O,C_6H_5,C_2H_5,O,H,O]
  \stopchemical
\stopformula
```





\startchemical \chemical[ONE,SB15,DB7,Z0157][C,NH_2,H_2N,O] \stopchemical

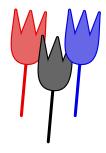


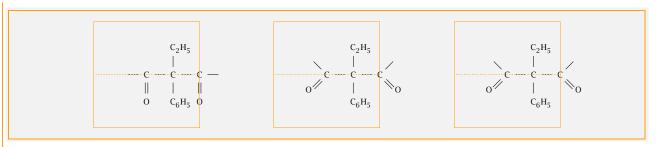
```
C_2H_5
         c - c - o - c_2 H_5
0
         С<sub>6</sub>Н<sub>5</sub> О
```

\startchemical

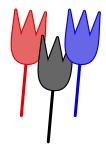
```
\chemical
```

```
[ONE, Z05, SB15,
    MOV1,Z03,SB1,DB3,
     MOV1, Z037, SB137,
     MOV1,Z03,SB1,DB3,
     MOV1,Z01,SB1]
    [0,C_2H_5,C,0,C,C_6H_5,C_2H_5,C,0,0,C_2H_5]
\stopchemical
```





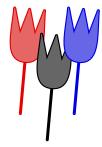
```
\hbox to \hsize \bgroup \hss
\startchemical
  \chemical
    [ONE, Z03, SB15, DB3,
     MOV1, Z037, SB137,
     MOV1,Z03,SB1,DB3]
    [C,0,C,C_6H_5,C_2H_5,C,0]
\stopchemical
\hss
\startchemical
  \chemical
    [ONE, Z04, SB16, DB4,
     MOV1, Z037, SB137,
     MOV1,Z02,SB8,DB2]
    [C,0,C,C_6H_5,C_2H_5,C,0]
\stopchemical
\hss
\startchemical
  \chemical
    [ONE, Z04, SB16, DB4,
     MOV1, Z037, SB137,
     MOV1,Z02,SB8,DB2]
    [C,0,C,C_6H_5,C_2H_5,C,0]
\stopchemical
\hss \egroup
```



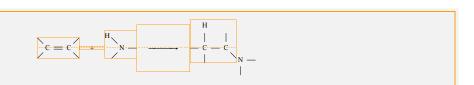
```
C_2H_5
c - c - o - c_2 H_5
C_6H_5
```

\startchemical

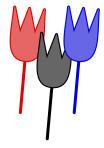
```
\chemical
    [ONE, Z05, SB15,
    MOV1,Z0,SB1,
    MOV1,Z037,SB137,
    MOV1,Z0,SB1,
    MOV1,Z01,SB1]
    [0,C_2H_5,C,C,C_6H_5,C_2H_5,C,0,C_2H_5]
\stopchemical
```



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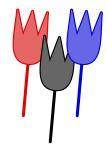
```
\startformula
  \setupchemical
    [width=fit,
     height=fit,
     scale=small,
     size=small]
  \startchemical
    \chemical[ONE, Z0, DB1, SB46, MOV1, Z0, SB28][C, C]
  \stopchemical
  \startchemical
    \chemical[SPACE,PLUS,SPACE]
  \stopchemical
  \startchemical
    \chemical[ONE,Z0,SB146,Z6][N,H]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical[ONE, Z07, SB1357, MOV1, Z0, SB27, PB: Z2, ONE, Z0, SB13, PE][C, H, C, N]
  \stopchemical
\stopformula
```

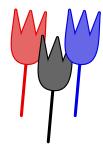


```
\begin{array}{c} \mathsf{CH_3} - \mathsf{CH} = \mathsf{CH_2} \\ \mathsf{H} \\ \\ \mathsf{CH_3} - \mathsf{CH_2} - \mathsf{CH_2} - \mathsf{CH_2} - \mathsf{CH_2} - \mathsf{NH_2} \\ \\ \mathsf{CH_3} - \mathsf{CH_2} - \mathsf{CH_2} - \mathsf{CH_2} - \mathsf{CH_2} - \mathsf{CH_2} - \mathsf{NH_2} \\ \\ \mathsf{N} - \mathsf{CH_2} - \mathsf{CH_2} - \mathsf{CH_2} - \mathsf{CH_2} - \mathsf{NH_2} \\ \\ \mathsf{H} \\ \end{array}
```

```
\setupchemical
  [width=fit,
   height=fit,
   scale=small,
   size=small]
\startformula
  \displaylines
    {\startchemical
       \chemical[ONE, Z015, DB1, SB5][CH, CH_2, CH_3]
     \stopchemical
     \startchemical
       \chemical[SPACE, PLUS, SPACE]
     \stopchemical
     \startchemical
       \chemical
          [ONE, Z046, SB146, OFF1,
           MOV1, Z0, 20FF1, SB1,
           MOV1, OFF1, Z0, 20FF1, SB1,
           MOV1, OFF1, Z0, 20FF1, SB1,
           MOV1, OFF1, Z0, 20FF1, Z1, SB1, PE]
          [N,H,H,CH_2,CH_2,CH_2,CH_2,NH_2]
     \stopchemical
     \startchemical
       \chemical[SPACE,GIVES,SPACE]
     \stopchemical
  \cr
     \startchemical[boven=-2000]
     \chemical
       [ONE, OFF1, Z0, 20FF1, SB1,
        MOV1, OFF1, Z0, 20FF1, SB1,
        MOV1, OFF1, Z0, 30FF1,
        PB: Z2, ONE, Z04, SB146,
        MOV1, OFF1, Z0, 20FF1, SB1,
        MOV1, OFF1, Z0, 20FF1, SB1,
        MOV1,0FF1,Z0,20FF1,SB1,
        MOV1, OFF1, Z0, 20FF1, Z1, SB1, PE]
        [CH_3,CH_2,CH_2,N,H,CH_2,CH_2,CH_2,CH_2,NH_2]
     \stopchemical
  \cr}
```

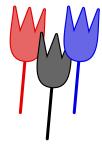
\stopformula





```
\mathbf{N} - \mathbf{C}\mathbf{H}_2 - \mathbf{C}\mathbf{H}_2 - \mathbf{C}\mathbf{H}_2 - \mathbf{C}\mathbf{H}_2 - \mathbf{C}\mathbf{H}_2 - \mathbf{N}\mathbf{H}_2
```

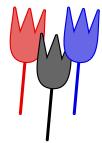
```
\setupchemical
  [width=fit,
   height=fit,
   size=medium]
\startchemical
  \chemical
    [ONE, Z015, SB125,
     DIR2, Z04, SB14,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1,0FF1,Z0,20FF1,SB1,
     MOV1, OFF1, Z0]
    [CH,CH_3,CH_3,N,H,CH_2,CH_2,CH_2,CH_2,NH_2]
\stopchemical
```



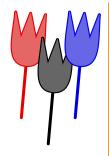
43

```
N = C - CH_2 -
```

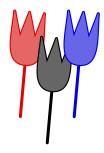
```
\setupchemical
  [width=fit,
   height=fit]
\startchemical
  \chemical
    [ONE,ZO,TB1,
     MOV1, Z0, SB1,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1, OFF1, Z0, 20FF1, SB2,
     DIR2, Z0, SB14,
   SAVE,
     DIR4, OFF5, Z0, 20FF5, SB5,
     MOV5, OFF5, Z0, 20FF5, SB5,
     MOV5, Z0, TB5,
     MOV5,Z0,
   RESTORE,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1,Z0,
   SAVE,
     DIR8, ONE, 20FF1, Z0, OFF1, SB1, SB4,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1, Z0, TB1,
     MOV1,Z0,
   RESTORE,
     DIR2, ONE, 20FF1, Z0, OFF1, SB1, SB6,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1, Z0, TB1,
     MOV1,Z0]
    [N,C,CH_2,CH_2,N,CH_2,CH_2,C,N,CH_2,CH_2,CH_2,CH_2,N,
     CH_2,CH_2,C,N,
     CH_2, CH_2, C, N
\stopchemical
```



```
\def\SomeX
 {\inframed[height=10ex,width=2em]{x}}
\startformula
 \setupchemical
    [width=fit,
     height=fit]
 \startchemical
    \chemical
      [ONE, OE5, Z0,
   SAVE,
      DIR8, Z0, SB14,
      MOV1, Z0, SB1,
     MOV1, Z0, SB1,
      MOV1,Z0,
   RESTORE,
      DIR2, ZO, SB16,
     MOV1, Z0, SB1,
     MOV1, Z0, SB1,
     MOV1,Z0
     [N,CH_2,CH_2,CH_2,NH_2,CH_2,CH_2,CH_2,NH_2]
 \stopchemical
 \startchemical
   \chemical[SPACE,PLUS,SPACE]
 \stopchemical
 \startchemical
   \chemical[Z0][A]
 \stopchemical
 \startchemical
   \chemical[SPACE,GIVES,SPACE,SPACE,SPACE]
 \stopchemical
 \startchemical
    \chemical
      [ONE, OE5, Z0,
     SAVE,
       DIR8, OFF1, Z0, SB14,
       MOV1, OFF1, Z0, OFF1, SB1,
       MOV1,0FF1,Z0,0FF1,SB1,
     RESTORE,
       DIR2, ZO, SB16,
       MOV1, OFF1, Z0, OFF1, SB1,
       MOV1, OFF1, Z0, OFF1, SB1,
       DIR8,30FF1,Z0]
      [N,CH_2,CH_2,CH_2,CH_2,CH_2,SomeX]
```

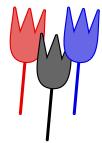


\stopchemical
\startchemical
\chemical[SPACE,PLUS,SPACE][]
\stopchemical
\startchemical
\chemical[Z0][2\,H_20]
\stopchemical
\stopformula

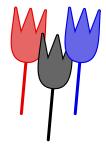


```
- СН — NН<sub>2</sub>
NH — C — CH<sub>2</sub>
         0
```

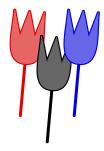
```
\setupchemical
  [width=fit,
   height=fit]
\startchemical
  \chemical
    [ONE,Z0,SB1,
     MOV1,Z07,SB1,DB7,
     MOV1, OFF1, Z0, OFF1, Z1, SB13,
     MOV3,OFF1,Z0,OFF5,SB5,
     MOV5,Z03,SB5,DB3,
     MOV5,Z0]
    [NH,C,O,CH,NH_2,CH_2,C,O,NH]
\stopchemical
```



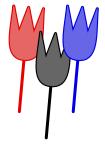
```
\setupchemical
  [width=fit,
   height=fit]
\startchemical
  \chemical
   [ONE,ZO,SB1,
   MOV1,ZO,SB1,
   MOV1,ZO,SB1,
   MOV1,OFF1,ZO,OFF1,Z1,SB13,
   MOV3,OFF1,ZO,OFF5,SB5,
   MOV5,ZO,SB5,
   MOV5,ZO,SB5,
   MOV5,ZO,SB5,
   MOV5,ZO]
  [NH,O,C,CH,NH_2,CH_2,C,O,NH]
\stopchemical
```

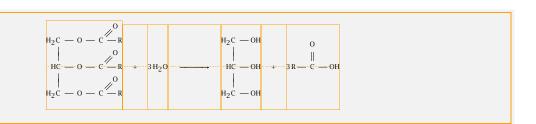


```
\setupchemical
  [width=fit]
\startchemical
  \chemical
    [ONE,ZO,OFF1,SB13,Z3,
        MOV1,OFF1,ZO,OFF1,SB13,Z3,
        MOV1,OFF1,ZO,SB3,Z3]
    [H_2C,Cl,CH,OH,CH_2,Cl]
  \bottext{A}
\stopchemical
```

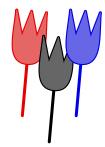


```
\setupchemical
  [axis=on]
\startchemical
  \chemical
    [ONE,
   SAVE,
     OFF1,Z0,20FF1,SB7,SB3,SB1,
     MOV1,Z0,SB1,
     MOV1,Z0,DB8,CZ8,SB1,Z1,
   RESTORE,
   SAVE,
     SUB4, ONE, Z0, 20FF1, SB3, SB1,
     MOV1, Z0, SB1,
     MOV1,Z0,DB8,CZ8,SB1,Z1,
   RESTORE,
     SUB2, ONE, Z0, 20FF1, SB7, SB1,
     MOV1, Z0, SB1,
     MOV1,Z0,DB8,CZ8,SB1,Z1]
    [HC,0,C,0,R,
     H_2C, 0, C, 0, R,
     H_2C, 0, C, 0, R
\stopchemical
```

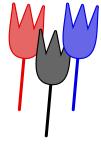




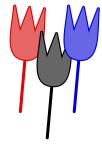
```
\startformula
\setupchemical
  [scale=small,
   size=small,
  width=fit]
\startchemical
  \chemical
    [ONE,
   SAVE,
     OFF1,Z0,20FF1,SB7,SB3,SB1,
     MOV1, Z0, SB1,
     MOV1, Z0, DB8, CZ8, SB1, Z1,
   RESTORE,
   SAVE,
     SUB4, ONE, Z0, 20FF1, SB3, SB1,
     MOV1, Z0, SB1,
     MOV1,Z0,DB8,CZ8,SB1,Z1,
   RESTORE,
     SUB2, ONE, Z0, 20FF1, SB7, SB1,
     MOV1, Z0, SB1,
     MOV1,Z0,DB8,CZ8,SB1,Z1]
    [HC,0,C,0,R,
     H_{2}C,0,C,0,R,
     H_{2}C,0,C,0,R
  \stopchemical
  \startchemical
    \chemical[SPACE,PLUS,SPACE]
  \stopchemical
  \startchemical
    \chemical[Z0][3,H_20]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE,
     SAVE,
       OFF1,Z0,20FF1,SB7,SB3,SB1,Z1,
     RESTORE,
```



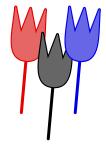
```
SAVE,
       SUB4, ONE, Z0, 20FF1, SB3, SB1, Z1,
     RESTORE,
       SUB2, ONE, Z0, 20FF1, SB7, SB1, Z1]
      [HC,OH,
       H_2C,OH,
       H_2C,OH
  \stopchemical
  \startchemical
    \chemical[SPACE,PLUS,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE,Z0,SB1,MOV1,Z017,SB1,DB7][3\,R,C,OH,O]
  \stopchemical
\stopformula
```



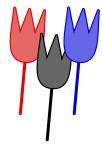
```
\startformula
  \setupchemical
    [width=fit,
     height=fit]
  \startchemical
    \chemical[ONE,Z0157,SB15,DB7][C,OH,H,O]
  \stopchemical
  \startchemical
    \chemical[SPACE, PLUS, SPACE][]
  \stopchemical
  \startchemical
    \chemical
      [ONE, ZO, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1,0FF1,Z0]
      [HO,CH_2,CH_3,]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
  \chemical
    [ONE, Z057, SB15, DB7,
    MOV1, Z0, SB1,
    MOV1, OFF1, Z0, 20FF1, Z1, SB1]
    [C,H,0,0,CH_2,CH_3]
  \stopchemical
  \startchemical
    \chemical[SPACE,PLUS,SPACE]
  \stopchemical
  \startchemical
    \chemical[Z0][H_20]
  \stopchemical
\stopformula
```



```
\startformula
  \setupchemical
    [width=fit,
     height=fit]
  \startchemical
    \chemical
     [ONE, Z057, SB15, DB7, MOV1, Z01, SB1] [C, H, O, O, CH_3]
  \stopchemical
  \startchemical
    \chemical[SPACE, PLUS, SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, ZO, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0]
      [HO,CH_2,CH_3]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, Z057, SB15, DB7,
       MOV1,Z0,SB1,
       MOV1, OFF1, Z0, 20FF1, Z1, SB1]
      [C,H,0,0,CH_2,CH_3]
  \stopchemical
  \startchemical
    \chemical[SPACE,PLUS,SPACE][]
  \stopchemical
  \startchemical
    \chemical
      [ONE, Z0, SB1, MOV1, OFF1, Z0] [H0, CH_3]
  \stopchemical
\stopformula
```

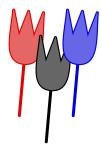


```
\startformula
  \startchemical
    \chemical
      [SIX,SB1..6,Z1..6,
       SR15, DR246, -SR3, +SR3,
       RZ12,-RZ3,+RZ3,RZ456]
      [N,C,C,C,N,C,H,O,C_6H_5,C_2H_5,O,H,O]
  \stopchemical
\stopformula
```



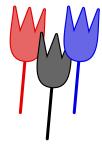
55

```
\setupchemical
  [width=fit]
\definechemical
  [molecule]
  {\chemical
     [ONE, Z0, Z7, SB1, DB7,
      SUB1, SIX, ROT2, B, C, R36,
      SUB1, ONE, Z07, SB15, DB7,
      MOV1,Z0,SB1]
     [C,0,C,0,0]
\startformula
  \startchemical
    \chemical
      [ONE, ZO, SB1, OE5]
    \chemical
      [molecule]
    \chemical
      [SUB1,ONE,Z0,OFF1,SB1,MOV1,OFF1,Z0,20FF1,SB1,MOV1,Z0,SB1][CH_2,CH_2,O]
    \chemical
      [SUB1, molecule]
    \chemical
      [SUB1,ONE,Z0,20FF1,SB1,MOV1,OFF1,Z0,OE1][CH_2,CH_2]
    \bottext{PET}
  \stopchemical
\stopformula
```



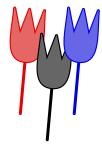
```
\setupchemical
  [width=fit,
   height=fit]
\definechemical
  [molecule]
  {\chemical
     [ONE, Z0, Z7, SB1, DB7,
      SUB1,SIX,ROT2,B,C,R36,
      SUB1, ONE, Z07, SB15, DB7,
      MOV1,Z0,SB1]
     [C,0,C,0,0]
\startchemical
  \chemical
    [ONE,ZO,SB1,
     MOV1,Z0,SB1,
     SUB1, molecule,
     SUB1, ONE, ZO]
    [CH_3,0,CH_3]
\stopchemical
```





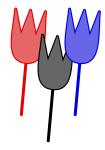
```
\setupchemical
  [width=fit,
   height=fit]
\startchemical
  \chemical
    [ONE,Z0,SB1,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1,Z0]
    [HO,CH_2,CH_2,OH]
\stopchemical
```

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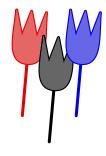


```
\begin{bmatrix} 0 & 0 & 0 \\ \parallel & \parallel & \parallel \\ C & 0 & CH_2 & CH_2 \end{bmatrix}_n
```

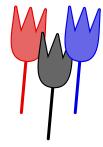
```
\setupchemical
  [width=fit,
   height=fit]
\definechemical
  [molecule]
  {\chemical
     [ONE, Z0, Z7, SB1, DB7,
      SUB1, SIX, ROT2, B, C, R36,
      SUB1, ONE, Z07, SB15, DB7,
      MOV1,Z0,SB1]
     [C,0,C,0,0]}
\startchemical
  \chemical
    [ONE, ZO, SB1, SB5, ZT5,
     SUB1, molecule,
     SUB1, ONE, Z0, 20FF1, SB1,
     MOV1,20FF1,Z0,20FF1,SB1,ZT1]
    [0,\Bigg\lbrack,CH_2,CH_2,\Bigg\rbrack_n]
\stopchemical
```



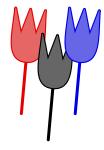
```
\startchemical[width=fit,height=fit,size=small,scale=small]
\chemical
[ONE,SB1,DB7,Z07,OE5,
MOV1,SB137,Z037,
MOV1,SB1,DB7,Z07,
MOV1,SB137,Z037,
MOV1,SB1,DB7,Z07,
MOV1,SB37,Z037,OE1]
[C,0,C,H,H,C,0,C,H,H,C,0,C,H,H]
\stopchemical
```



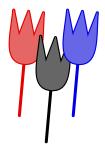
```
\startchemical[width=fit,height=fit,size=small,scale=small]
 \chemical
   [ONE, SB1, DB7, Z07, OE5,
    MOV1,SB17,DB3,Z07,
  SAVE,
    MOV3, ONE, Z0234, SB234,
  RESTORE,
    MOV1, SB1, DB7, Z07,
    MOV1,SB17,DB3,Z07,
  SAVE,
    MOV3, ONE, Z0234, SB234,
  RESTORE,
    MOV1, SB1, DB7, Z07,
    MOV1, SB7, DB3, Z07, OE1,
    MOV3, ONE, Z0234, SB234]
   \stopchemical
```



```
\startchemical[width=fit,height=fit,size=small,scale=small]
\chemical
[ONE,SB1,DB7,Z07,OE5,
MOV1,SB137,Z037,
MOV1,SB137,Z037,
MOV1,SB1,DB7,Z07,
MOV1,SB137,Z037,
MOV1,SB137,Z037,
MOV1,SB1,DB7,Z07,
MOV1,SB1,DB7,Z07,
MOV1,SB37,Z037,
MOV1,SB37,Z037,
SB37,Z037,OE1]
[C,0,C,H,H,C,H,H,C,O,C,H,H,C,H,H,C,O,C,H,H,C,H,H]
\stopchemical
```

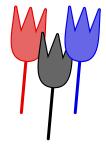


```
\startchemical[width=fit,height=fit,size=small,scale=small]
  \chemical
    [ONE,DB1,Z0,OE5,
     MOV1, ONE, SB1, ZO,
     MOV1, SB137, Z037,
     MOV1, SB137, Z037,
     MOV1, DB1, ZO,
     MOV1, ONE, SB1, ZO,
     MOV1, SB137, Z037,
     MOV1, SB137, Z037,
     MOV1, DB1, ZO,
     MOV1, ONE, SB1, ZO,
     MOV1, SB137, Z037,
     MOV1,SB37,Z037,OE1]
    [C,0,C,H,H,C,H,H,C,0,C,H,H,C,H,H,C,0,C,H,H,C,H,H]
\stopchemical
```



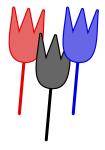


```
\startchemical[width=fit]
  \chemical
    [ONE,Z01,OFF1,SB13,EP57,MOV3,Z0]
    [\TL{\ominus}N,H,H]
  \stopchemical
```



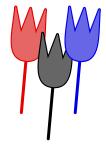
```
Н
                                         Н
                                                     NH_3
                NH_2^-
                                   benzyn
```

```
\startformula
 \setupchemical
    [width=fit,
    height=5000]
 \startchemical
    \chemical[SIX,C,B,R1..6,RZ1..6][Cl,H,H,H,H,H]
 \stopchemical
 \startchemical
    \chemical[SPACE,PLUS,SPACE][]
 \stopchemical
 \startchemical
    \chemical[ONE,Z0][NH_{2}^{-}]
 \stopchemical
 \startchemical
    \chemical[SPACE,GIVES,SPACE][]
 \stopchemical
 \startchemical
    \chemical[SIX,C,B,EB1,R3..6,RZ3..6][H,H,H,H]
    \bottext{benzyn}
 \stopchemical
 \startchemical
    \chemical[SPACE, PLUS, SPACE]
 \stopchemical
 \startchemical
    \chemical[ONE,Z0][NH_{3}]
 \stopchemical
 \startchemical
    \chemical[SPACE,PLUS,SPACE]
 \stopchemical
 \startchemical
    \chemical[ONE,Z0][C1^{-}]
 \stopchemical
\stopformula
```



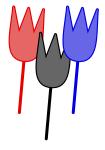
```
Н
Н
            + NH<sub>3</sub>
```

```
\startformula
 \setupchemical
   [width=fit,
    height=fit]
 \startchemical
   \chemical[SIX,C,B,EB1,R3..6,RZ3..6][H,H,H,H]
 \stopchemical
 \startchemical
   \chemical[SPACE,PLUS,SPACE][]
 \stopchemical
 \startchemical
   \chemical[ONE,Z0][NH_3]
 \stopchemical
 \startchemical
   \chemical[SPACE,GIVES,SPACE][]
 \stopchemical
 \startchemical
   \chemical[SIX,C,B,R1..6,RZ1..6][NH_2,H,H,H,H,H]
 \stopchemical
\stopformula
```

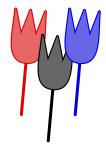


```
Н
```

```
\startformula
 \setupchemical
    [width=fit,
    height=fit]
 \startchemical
    \chemical[SIX,C,B,EB1,R3..6,RZ3..6][H,H,H,H]
 \stopchemical
 \startchemical
    \chemical[SPACE,PLUS,SPACE]
 \stopchemical
 \startchemical
    \chemical[SIX,C,B,EB4,R1236,RZ1236][H,H,H,H]
 \stopchemical
 \startchemical
    \chemical[SPACE,GIVES,SPACE]
 \stopchemical
 \startchemical
    \chemical
      [SIX,C,B,R3456,RZ3456,
       ADJ1, FOUR, B, ADJ1,
       SIX,C,B,R1236,RZ1236]
      [H,H,H,H,H,H,H,H]
 \stopchemical
\stopformula
```

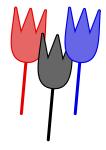


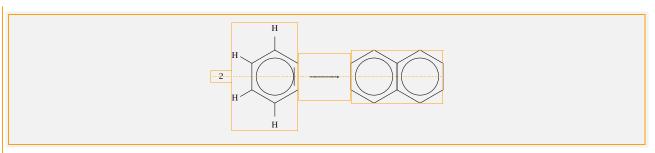
```
\startformula
  \setupchemical
    [width=fit,
     height=fit]
  \startchemical
    \chemical[ONE,Z0,SPACE,SPACE][2]
  \stopchemical
  \startchemical
    \chemical[SIX,C,B,EB1,R3..6,RZ3..6][H,H,H,H]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [SIX,C,B,R3456,RZ3456,ADJ1,
       FOUR, B, ADJ1,
       SIX,C,B,R1236,RZ1236]
      [H,H,H,H,H,H,H,H]
  \stopchemical
  \startchemical
    \chemical[SPACE, PLUS, SPACE][]
  \stopchemical
  \startchemical
    \chemical[ONE,Z0][H_2]
  \stopchemical
\stopformula
```



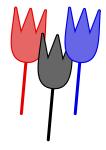
```
Н
                                                    Н
            Н
--2--
```

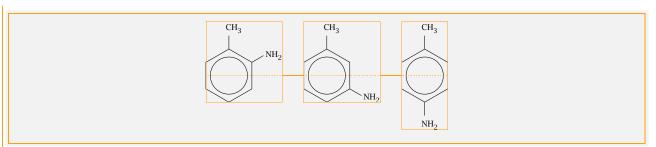
```
\startformula
 \setupchemical
    [width=fit,
    height=fit]
 \startchemical
    \chemical[ONE,Z0,SPACE,SPACE][2]
 \stopchemical
 \startchemical
    \chemical[SIX,C,B,EB1,R3..6,RZ3..6][H,H,H,H]
 \stopchemical
 \startchemical
    \chemical[SPACE,GIVES,SPACE]
 \stopchemical
 \startchemical
    \chemical
      [SIX,B,EB35,R3456,RZ3456,
       ADJ1, FOUR, B, ADJ1,
       SIX,B,EB26,R1236,RZ1236]
      [H,H,H,H,H,H,H,H]
 \stopchemical
\stopformula
```



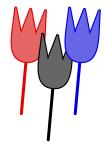


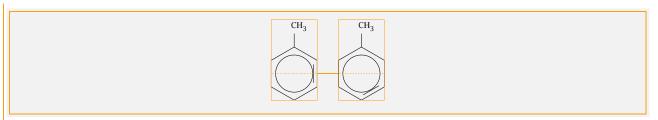
```
\startformula
 \setupchemical
   [width=fit,
    height=fit]
 \startchemical
   \chemical[ONE,Z0,SPACE,SPACE][2]
 \stopchemical
   \startchemical
   \chemical[SIX,C,B,EB1,R3..6,RZ3..6][H,H,H,H]
 \stopchemical
 \startchemical
   \chemical[SPACE,GIVES,SPACE]
 \stopchemical
 \startchemical
   \chemical[SIX,C,B,ADJ1,SIX,C,B]
 \stopchemical
\stopformula
```



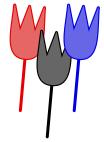


```
\startformula
 \setupchemical
    [width=fit,
    height=fit]
 \startchemical
    \chemical[SIX,C,B,R16,RZ16][NH_2,CH_3]
 \stopchemical
 \startchemical
    \chemical[SPACE,SPACE]
 \stopchemical
 \startchemical
    \chemical[SIX,C,B,R26,RZ26][NH_2,CH_3]
 \stopchemical
 \startchemical
    \chemical[SPACE, SPACE]
 \stopchemical
 \startchemical
    \chemical[SIX,C,B,R36,RZ36][NH_2,CH_3]
 \stopchemical
\stopformula
```

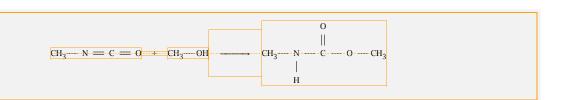




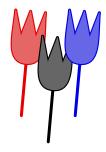
```
\startformula
  \setupchemical
    [width=fit,
    height=fit]
  \startchemical
    \chemical[SIX,C,B,EB1,R6,RZ6][CH_3]
  \stopchemical
  \startchemical
    \chemical[SPACE,SPACE][]
  \stopchemical
  \startchemical
    \chemical[SIX,C,B,EB2,R6,RZ6][CH_3]
  \stopchemical
    \stopchemical
    \stopchemical
    \stopchemical
    \stopchemical
    \stopchemical
    \stopchemical
    \stopchemical
    \stopformula
```



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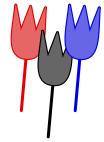


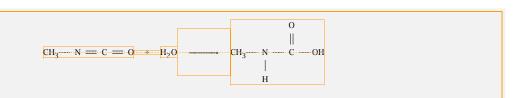
```
\startformula
  \setupchemical
    [width=fit,
     height=fit]
  \startchemical
    \chemical
      [ONE, ZO, SB1,
       MOV1,Z0,DB1,
       MOV1,Z0,DB1,
       MOV1,Z0]
      [CH_3,N,C,0]
  \stopchemical
  \startchemical
    \chemical[SPACE,PLUS,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, ZO, SB1,
       MOV1,Z0]
      [CH_3,OH]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, ZO, SB1,
       MOV1,Z0,SB13,Z3,
       MOV1,Z0,SB1,DB7,Z7,
       MOV1,Z0,SB1,
       MOV1, OFF1, Z0]
      [CH_3, N, H, C, 0, 0, CH_3]
  \stopchemical
\stopformula
```



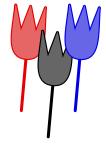
```
О
```

```
\startchemical[width=fit]
  \chemical
    [ONE,ZO,SB1,
    MOV1,Z03,SB13,
    MOV1,Z07,SB1,DB7,
    MOV1,Z03,SB13,
    MOV1,0FF1,Z0]
    [CH_3,N,H,C,O,N,H,CH_3]
\stopchemical
```

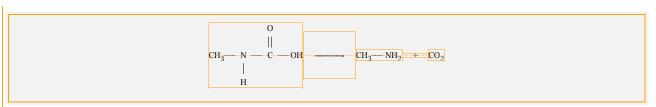




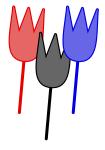
```
\startformula
 \setupchemical
    [width=fit,
     height=fit]
 \startchemical
    \chemical
      [ONE,ZO,SB1,
      MOV1,Z0,DB1,
       MOV1,Z0,DB1,
      MOV1,Z0]
      [CH_3,N,C,0]
 \stopchemical
 \startchemical
    \chemical[SPACE,PLUS,SPACE]
 \stopchemical
 \startchemical
    \chemical[ONE,Z0,][H_20]
 \stopchemical
 \startchemical
    \chemical[SPACE,GIVES,SPACE]
 \stopchemical
 \startchemical
    \chemical
      [ONE, ZO, SB1,
      MOV1,Z0,SB13,Z3,
       MOV1,Z0,SB1,DB7,Z7,
       MOV1,Z0
      [CH_3,N,H,C,0,OH]
 \stopchemical
\stopformula
```

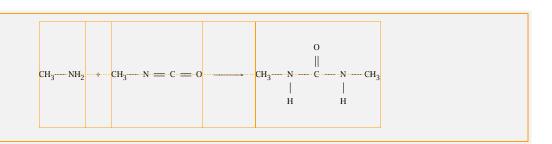


75

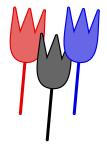


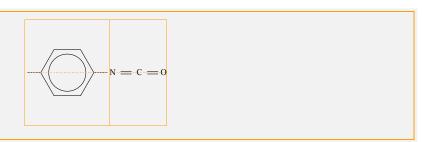
```
\startformula
 \setupchemical
    [width=fit,
     height=fit]
 \startchemical
    \chemical
      [ONE, ZO, SB1,
       MOV1,Z0,SB13,Z3,
       MOV1,Z0,SB1,DB7,Z7,
       MOV1,Z0]
      [CH_3, N, H, C, 0, OH]
 \stopchemical
 \startchemical
    \chemical[SPACE,GIVES,SPACE]
 \stopchemical
 \startchemical
    \chemical[ONE,Z0,SB1,Z1][CH_3,NH_2]
 \stopchemical
 \startchemical
    \chemical[SPACE,PLUS,SPACE]
 \stopchemical
 \startchemical
    \chemical[ONE,Z0][CO_2]
 \stopchemical
\stopformula
```



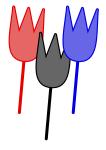


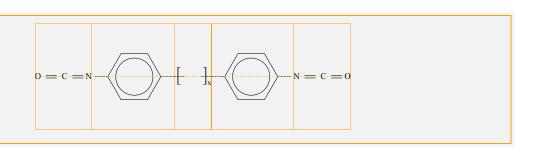
```
\startformula
  \setupchemical
    [width=fit]
  \startchemical
    \chemical[ONE, Z0, SB1, Z1][CH_3, NH_2]
  \stopchemical
  \startchemical
    \chemical[SPACE, PLUS, SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE,ZO,SB1,
       MOV1, Z0, DB1,
       MOV1,Z0,DB1,
       MOV1,Z0]
      [CH_3,N,C,0]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, ZO, SB1,
       MOV1, Z03, SB13,
       MOV1, Z07, SB1, DB7,
       MOV1, Z03, SB13,
       MOV1, OFF1, Z0]
      [CH_3,N,H,C,O,N,H,CH_3]
  \stopchemical
\stopformula
```



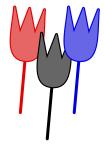


```
\startformula
  \setupchemical
    [width=fit]
  \startchemical
    \chemical[SIX,ROT2,B,C,R36,RZ36]
  \stopchemical
  \startchemical
    \chemical[ONE,Z0,DB1,MOV1,Z01,DB1][N,C,0]
  \stopchemical
  \stopformula
```

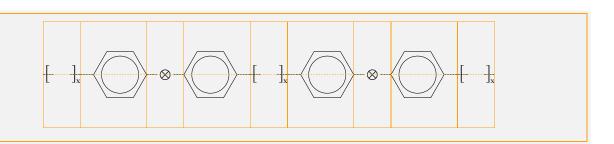




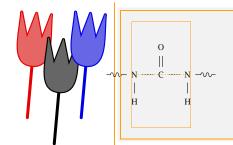
```
\startformula
 \setupchemical
    [width=fit]
 \startchemical
   \chemical[ONE,Z0,DB1,MOV1,Z01,DB1][0,C,N]
 \stopchemical
 \startchemical
   \chemical[SIX,ROT2,B,C,R36,RZ3]
 \stopchemical
 \startchemical
   \chemical[ONE,SB15,ZT5,ZT1][\[,\]{x}]
 \stopchemical
 \startchemical
   \chemical[SIX,ROT2,B,C,R36,RZ6]
 \stopchemical
 \startchemical
   \chemical[ONE,Z0,DB1,MOV1,Z01,DB1][N,C,0]
 \stopchemical
\stopformula
```



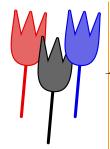
79



```
\startformula
 \setupchemical
    [width=fit]
 \startchemical
    \chemical[ONE,SB15,ZT5,ZT1][\[,\]\{x\}]
 \stopchemical
 \startchemical
    \chemical[SIX,ROT2,R3,B,C]
 \stopchemical
 \startchemical
    \chemical[ONE,SB15,Z0][{\bigotimes}]
 \stopchemical
 \startchemical
    \chemical[SIX,ROT2,R6,B,C]
 \stopchemical
 \startchemical
    \chemical[ONE,SB15,ZT5,ZT1][\[,\]{x}]
 \stopchemical
 \startchemical
    \chemical[SIX,ROT2,R3,B,C]
 \stopchemical
 \startchemical
    \chemical[ONE,SB15,Z0][{\bigotimes}]
 \stopchemical
 \startchemical
    \chemical[SIX,ROT2,R6,B,C]
 \stopchemical
 \startchemical
    \chemical[ONE,SB15,ZT5,ZT1][\[,\]{x}]
 \stopchemical
\stopformula
```

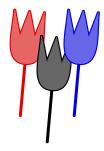


```
\startchemical[width=fit]
  \chemical
    [ONE,Z0,OE5,SB13,Z3,
         MOV1,Z0,SB1,DB7,Z7,
         MOV1,Z0,Z3,SB3,OE1]
    [N,H,C,O,N,H]
\stopchemical
```



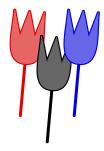
```
О
C N --- CH<sub>2</sub>--~-
```

```
\startchemical[width=fit]
  \chemical
    [ONE, Z0, OE5, OFF1, SB1,
     MOV1,Z0,SB13,Z3,
     MOV1,Z07,SB1,DB7,
     MOV1,Z03,SB13,
     MOV1,0FF1,Z0,0E1]
    [CH_2,N,H,C,O,N,H,CH_2]
\stopchemical
```



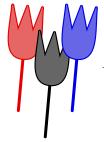
```
CH_2 = CH - C = CH_2
            CH_3
```

```
\startchemical[width=fit]
  \chemical
    [ONE,Z0,OFF1,DB1,
    MOV1,Z0,SB1,
    MOV1,Z0,SB3,Z3,DB1,
    MOV1,0FF1,Z0]
    [CH_2,CH,C,CH_3,CH_2]
\stopchemical
```

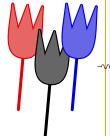


```
CH_3
```

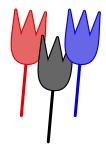
```
\startchemical[width=fit]
  \chemical
    [ONE,Z0,SB1,
    MOV1,0FF1,Z0,20FF1,SB1,
    MOV1,Z0,SB3,Z3,DB1,
    MOV1,0FF1,Z0]
    [CH_3,CH_2,C,CH_3,CH_2]
\stopchemical
```

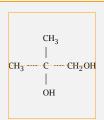


```
\startchemical[width=fit]
  \chemical
    [ONE,OFF1,Z0,OE5,OFF1,SB1,
        MOV1,Z0,SB137,Z37,
        MOV1,OFF1,Z0,20FF1,SB1,
        MOV1,OFF1,Z0,OFF1,DB1,
        MOV1,Z0,SB13,Z3,
        MOV1,OFF1,Z0,20FF1,SB1,
        MOV1,Z0,SB137,Z37,
        MOV1,OFF1,Z0,20FF1,SB1,
        MOV1,Z0,SB137,Z37,
        MOV1,OFF1,Z0,20FF1,SB1,
        MOV1,Z0,SB137,Z37,
        MOV1,OFF1,Z0,OE1]
    [CH_2,C,CH_3,CH_3,CH_2,CH,C,CH_3,CH_2,C,CH_3,CH_2,C,CH_3,CH_2,C,CH_3,CH_2,C,CH_3,CH_2]
    \stopchemical
```

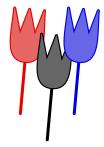


```
\startchemical[width=fit]
  \chemical
    [ONE, Z0, OE5, SB137, Z37,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1, Z0, SB13,
   SAVE,
     MOV1, OFF1, Z0, OFF1, SB1,
     MOV1,Z0,SB137,Z37,
     MOV1, OFF1, Z0, OFF1, SB1,
     MOV1,Z0,SB37,Z37,OE1,
   RESTORE,
     MOV3,Z0,SB1,DB3,Z13]
    [C,CH_3,CH_3,CH_2,CH_2,CH,CH_2,C,CH_3,CH_3,
     CH_2,C,CH_3,CH_3,C,CH_3,CH_2]
\stopchemical
```

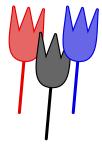




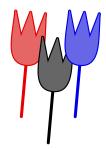
\startchemical[width=fit] \chemical [ONE,SB1357,Z0,Z1357] [C,CH_2OH,OH,CH_3,CH_3] \stopchemical



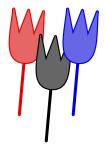
```
\definechemical
  [molecule]
  {\chemical
     [ONE, SB5, OFF1, Z0, 20FF1, SB1,
      MOV1, OFF1, Z0, 20FF1, SB1,
      MOV1, OFF1, Z0, 20FF1, SB1,
      MOV1, OFF1, Z0, SB3,
    SAVE,
      MOV3,Z0,SB3,
      MOV3,Z0,SB3,
      MOV3,Z0,SB3,
      MOV3,Z0,
    RESTORE,
      OFF1, SB1, MOV1]
     [CF_2,CF_2,CF_2,CF,0,CF_2,CF_2,S0_3H]}
\startchemical[width=fit]
  \chemical[molecule, molecule]
\stopchemical
```

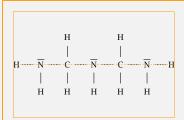


```
\startformula
 \setupchemical
    [height=2000,
    width=fit.
    scale=small,
    size=small,
    bottom=2000]
 \startchemical
   \chemical[ONE,SB1,SB3,SB5,Z0,Z1,Z3,Z5][N,\bf H,H,H]
 \stopchemical
 \startchemical
   \chemical[SPACE]
 \stopchemical
 \startchemical
    \chemical[ONE,DB3,Z0,Z3][\bf 0,CH_2]
 \stopchemical
 \startchemical
   \chemical[SPACE]
 \stopchemical
 \startchemical
   \chemical[ONE,SB1,SB3,SB5,Z0,Z1,Z3,Z5][N,\bf H,H,\bf H]
 \stopchemical
 \startchemical
   \chemical[SPACE]
 \stopchemical
 \startchemical
   \chemical[ONE,DB3,Z0,Z3][\bf 0,CH_2]
 \stopchemical
 \startchemical
   \chemical[SPACE]
 \stopchemical
 \startchemical
 \chemical[ONE,SB1,SB3,SB5,Z0,Z1,Z3,Z5][N,H,H,\bf H]
 \stopchemical
 \startchemical
   \chemical[SPACE,GIVES,SPACE]
 \stopchemical
 \startchemical
   \chemical
      [ONE, Z0, 20FF1, SB1,
      MOV1, OFF1, Z0, 20FF1, SB1,
      MOV1,SB1,SB3,Z0,Z3,
```

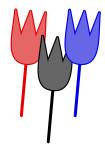


```
MOV1,OFF1,Z0,20FF1,SB1,
MOV1,OFF1,Z0]
[H_2N,CH_2,N,H,CH_2,NH_2]
\bottext{compound A}
\stopchemical
\startchemical
\chemical[SPACE,PLUS,SPACE]
\stopchemical
\startchemical
\chemical[ONE,Z0][2\,H_20]
\stopchemical
\stopformula
```

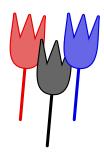




```
\startchemical[width=fit]
\chemical
  [ONE, Z035, SB135, EP7,
  MOV1,Z037,SB137,
  MOV1,Z03,SB13,EP7,
  MOV1,Z037,SB137,
  MOV1,Z013,SB13,EP7]
  [N,H,H,C,H,H,N,H,C,H,H,N,H,H]
\stopchemical
```

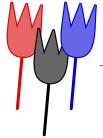


```
\startchemical[width=fit]
  \chemical
    [ONE, Z035, SB135, EP7,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1,Z03,SB13,EP7,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1,Z013,SB13,EP7]
    [N,H,H,CH_2,N,H,CH_2,N,H,H]
\stopchemical
```

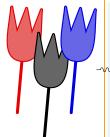


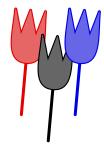


\startchemical[width=fit] \chemical[ONE,Z0,SB1,DB3,SB5,Z135][C,NH_2,0,H_2N] \stopchemical

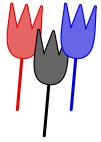


```
\startchemical[width=fit]
  \chemical
  [ONE, Z0, SB1, OE5,
      MOV1, Z0, SB1, Z3, DB3,
      MOV1, Z0, SB1,
      MOV1, OFF1, Z0, 20FF1, SB1,
      MOV1, Z0, SB1,
      MOV1, Z0, SB1,
      MOV1, Z0, SB1, Z3, DB3,
      MOV1, Z0, SB1,
      MOV1, OFF1, Z0, OE1]
  [NH, C, O, NH, CH_2, NH, C, O, NH, CH_2]
\stopchemical
```

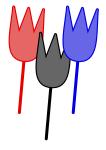




```
\startchemical[width=fit,scale=small,size=medium]
  \chemical
    [ONE,OFF1,Z0,OFF1,SB1,OE5,
     MOV1, OFF1, Z0, OFF1, SB1, Z3, DB3,
     MOV1, OFF1, Z0, OFF1, SB1,
     MOV1, OFF1, Z0, OFF1, SB1,
     MOV1, OFF1, Z0, OFF1, SB1, Z3, DB3,
     MOV1, OFF1, Z0, OFF1, SB1,
     MOV1, OFF1, Z0, OFF1, SB1,
     MOV1, OFF1, Z0, OFF1, SB1, Z3, DB3,
     MOV1, OFF1, Z0, OFF1, SB1,
     MOV1, OFF1, Z0, OFF1, SB1,
     MOV1, OFF1, Z0, OFF1, SB1, Z3, DB3,
     MOV1, OFF1, Z0, OFF1, OE1]
    [NH,C,O,NH,NH,C,O,NH,NH,C,O,NH,NH,C,O,NH]
\stopchemical
```

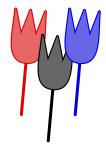


```
\startchemical[scale=small,width=fit,top=3000,bottom=3000]
\chemical[SIX,SB2356,DB14,Z2346,SR3,RZ3,-SR6,+SR6,-RZ6,+RZ6][C,N,C,C,H,H,H]
\chemical[PB:Z1,ONE,Z0,DIR8,Z0,SB24,DB7,Z27,PE][C,C,CH_3,O]
\chemical[PB:Z5,ONE,Z0,DIR6,Z0,SB24,DB7,Z47,PE][C,C,H_3C,O]
\chemical[SR24,RZ24][CH_3,H_3C]
\bottext{\crlf B} % ??
\stopchemical
```

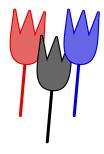


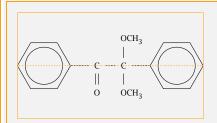
```
0 \qquad 0 \\ \parallel \qquad \parallel \\ H_2 C = O \longrightarrow NH_3 \longrightarrow 2 \\ H_3 C \longrightarrow C \longrightarrow CH_2 \longrightarrow C \longrightarrow CH_3 \longrightarrow verbinding B \longrightarrow 3 \\ H_2 O
```

```
\startformula
 \setupchemical
    [width=fit,
     size=small,
     scale=small]
 \startchemical
    \chemical[ONE, Z0, OFF1, DB1, Z1][H_2C, 0]
 \stopchemical
  \startchemical
    \chemical[SPACE,PLUS,SPACE]
 \stopchemical
  \startchemical
    \chemical[ONE,Z0][NH_3]
 \stopchemical
  \startchemical
    \chemical[SPACE, PLUS, SPACE]
 \stopchemical
  \startchemical
    \chemical[ONE,Z0][2\,\,]
 \stopchemical
 \startchemical
    \chemical
      [ONE, Z0, SB1, Z57, SB5, DB7,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1,Z0,SB1,Z17,DB7]
      [C,H_3C,0,CH_2,C,CH_3,0]
 \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
 \stopchemical
  \startchemical
    \chemical[ONE,Z0][verbinding B]
 \stopchemical
 \startchemical
    \chemical[SPACE, PLUS, SPACE][]
 \stopchemical
 \startchemical
    \chemical[ONE,Z0][3\,H_20]
 \stopchemical
\stopformula
```

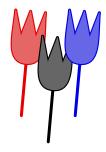


```
\startchemical[width=fit]
  \chemical
    [ONE, Z0, SB1, Z57, DB5, SB7,
    MOV1,Z0,SB1,Z7,DB7,
    MOV1,Z0,SB1,
     MOV1, OFF1, Z0, OFF1, SB1, ZT1, ZT5,
    MOV1,Z0,SB1,
    MOV1,Z0,SB1,Z7,DB7,
    MOV1,Z0,DB1,Z17,SB7]
    [C,H_2C,H,C,0,0,CH_2,\]{6},\[,0,C,0,C,CH_2,H]
\stopchemical
```

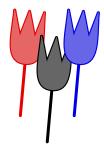




```
\startchemical[width=fit]
  \chemical
    [SIX,ROT2,B,C,R6,SUB1,
     ONE, Z0, Z3, SB15, DB3,
     MOV1,Z037,SB137,SUB1,
     SIX,ROT2,B,C,R3]
    [C,0,C,OCH_3,OCH_3]
\stopchemical
```

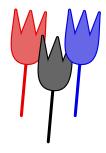




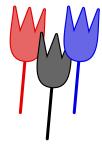


```
Н
```

```
\startchemical[width=fit]
\chemical
  [ONE, Z057, SB17, DB5,
  MOV1,Z0,SB1,
  MOV1, Z037, SB137,
  MOV1,0FF1,Z0,20FF1,SB1,
  MOV1, OFF1, Z0, 20FF1, SB1,
  MOV1,Z03,SB1,SB3,ES7,
  MOV1,Z0,SB1,
  MOV1,Z017,DB1,SB7]
  [C,H_2C,H,A,C,R,H,CH_2,CH_2,C,H,A,C,CH_2,H]
\stopchemical
```

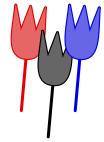


```
\startchemical[width=fit]
\chemical
[ONE,Z057,SB17,DB5,
MOV1,Z0,SB1,
MOV1,Z037,SB137,
MOV1,OFF1,Z0,20FF1,SB1,
MOV1,OFF1,Z0,20FF1,SB1,
MOV1,OFF1,Z0,20FF1,SB1,
MOV1,Z0,SB1,
MOV1,Z0,SB1,
MOV1,Z0,SB1,
MOV1,Z0,SB37,ES1]
[C,H_2C,H,A,C,R,H,CH_2,CH_2,CH_2,A,CH_2,C,H,H]
\stopchemical
```



```
HO - CH_2 - CH_2 - O - CH_2 - CH_2 - OH diethyleenglycol
```

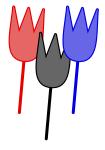
```
\startchemical[width=fit,bottom=1000]
\chemical
    [ONE,Z0,OFF1,SB1,
        MOV1,OFF1,Z0,20FF1,SB1,
        MOV1,OFF1,Z0,20FF1,SB1,
        MOV1,Z0,SB1,
        MOV1,OFF1,Z0,20FF1,SB1,
        MOV1,OFF1,Z0,20FF1,SB1,
        MOV1,Z0]
    [H0,CH_2,CH_2,O,CH_2,CH_2,OH]
    \bottext{diethyleenglycol}
\stopchemical
```



```
HO — CH<sub>2</sub> — CH<sub>2</sub> — OH

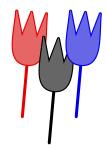
Sample A
```

```
\startchemical[width=fit,bottom=1000]
\chemical
  [ONE,Z0,OFF1,SB1,
    MOV1,OFF1,Z0,20FF1,SB1,
    MOV1,OFF1,Z0,20FF1,SB1,
    MOV1,OFF1,Z0]
  [H0,CH_2,CH_2,OH]
\bottext{Sample A}
\stopchemical
```

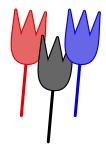


```
HO - CH_2 - CH_2 - O - CH_2 - CH_2 - OH - + H_2O - - 2 HO - CH_2 - CH_2 - OH
```

```
\startformula
  \setupchemical
    [width=fit]
  \startchemical
    \chemical
      [ONE, ZO, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, Z0, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1,Z0]
      [HO,CH_2,CH_2,O,CH_2,CH_2,OH]
  \stopchemical
  \startchemical
    \chemical[SPACE, PLUS, SPACE]
  \stopchemical
  \startchemical
    \chemical[ONE, Z0][H_20]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical[ONE,Z0,SPACE][2]
  \stopchemical
  \startchemical
    \chemical
      [ONE, ZO, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1,Z0
      [HO,CH_2,CH_2,OH]
  \stopchemical
\stopformula
```

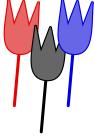


```
\startchemical[width=fit]
  \chemical
    [ONE, Z0, OFF1, SB1,
     MOV1,Z07,SB137,
   SAVE,
     MOV1, Z07, SB1, DB7,
     MOV1, Z07, SB17,
     MOV1,Z07,SB137,
   SAVE,
     MOV1, Z07, OE1, DB7,
   RESTORE,
     MOV3,SB3,Z0,
     MOV3,Z0,Z24,SB24,
   RESTORE,
     MOV3,,Z0,SB3,
     MOV3,Z0]
    [H_2N,C,H,C,0,N,H,C,H,C,0,
     CH_2,CH,CH_3,CH_3,CH_2,OH]
\stopchemical
```





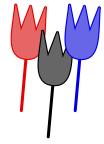
```
\startchemical[width=fit]
  \chemical
    [ONE,Z0,OFF1,SB1,
        MOV1,OFF1,Z0,OFF1,SB13,Z13]
    [H_2N,CH,COOH,R]
  \stopchemical
```



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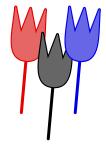
```
\begin{array}{c} \text{CH}_3 & \text{CH}_2 & \text{CH}_2 & \text{CH}_3 \\ \text{CH}_3 & \text{CH}_2 & \text{CH}_2 & \text{CH}_3 \\ \text{butaan} & \text{isobuteen} & \begin{array}{c} \text{CH}_3 \\ | \\ \text{CH}_3 & \text{CH}_3
```

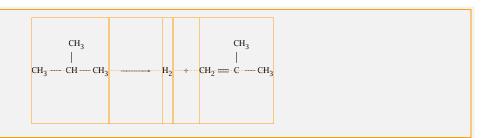
```
\startformula
 \startchemical[height=2000,width=fit]
    \chemical
      [ONE, Z0, OFF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0]
      [CH_3,CH_2,CH_2,CH_3]
    \bottext{butaan}
 \stopchemical
 \startchemical[height=2000,width=fit]
    \chemical
      [ONE,Z0,OFF1,DB1,
      MOV1, Z07, SB17,
       MOV1,0FF1,Z0]
      [CH_2,C,CH_3,CH_3]
   \bottext{isobuteen}
 \stopchemical
 \quad\quad\quad
 \startchemical[height=6000,width=fit]
    \chemical
      [ONE, ZO, SB1,
       MOV1, Z0, Z7, SB137,
       SAVE, MOV1, Z0, RESTORE,
       MOV3,Z0,SB3,
       MOV3,Z0]
      [CH_3,C,CH_3,CH_3,0,CH_3]
    \bottext{MTBE}
 \stopchemical
\stopformula
```



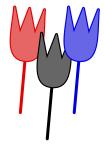
```
\begin{array}{c|c} & \text{CH}_3 \\ \text{CH}_3 - \text{CH}_2 - \text{CH}_2 - \text{CH}_3 \\ \end{array} \qquad \begin{array}{c|c} \text{CH}_3 \\ \text{CH}_3 - \text{CH} - \text{CH}_3 \\ \end{array}
```

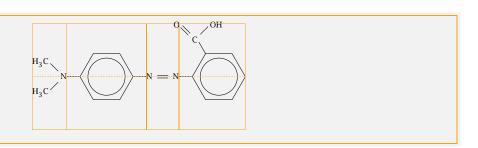
```
\startformula
  \setupchemical
    [height=2000,
     width=fit]
  \startchemical
    \chemical
      [ONE,Z0,OFF1,SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0]
      [CH_3,CH_2,CH_2,CH_3]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, Z0, OFF1, SB1,
       MOV1, Z07, SB17,
       MOV1, OFF1, Z0]
      [CH_3,CH,CH_3,CH_3]
  \stopchemical
\stopformula
```



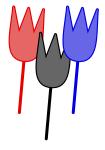


```
\startformula
 \setupchemical
    [width=fit]
 \startchemical
    \chemical
      [ONE,Z0,OFF1,SB1,
       MOV1, OFF1, Z0, OFF1, Z7, SB17,
       MOV1,0FF1,Z0]
      [CH_3,CH,CH_3,CH_3]
 \stopchemical
 \startchemical
    \chemical[SPACE,GIVES,SPACE]
 \stopchemical
 \startchemical
    \chemical[ONE,Z0,SPACE][H_2]
 \stopchemical
 \startchemical
    \chemical[SPACE,PLUS,SPACE]
 \stopchemical
 \startchemical
    \chemical
      [ONE, ZO, OFF1, DB1,
      MOV1,Z07,SB17,
       MOV1, OFF1, Z0]
      [CH_2,C,CH_3,CH_3]
 \stopchemical
\stopformula
```



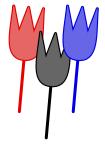


```
\startformula
 \setupchemical[width=fit]
 \startchemical
    \chemical[ONE, Z046, SB46,][N, H_3C, H_3C]
 \stopchemical
 \startchemical
    \chemical[SIX,ROT2,B,C,R36]
 \stopchemical
 \startchemical
    \chemical[ONE,Z0,DB1,MOV1,Z0][N,N]
 \stopchemical
 \startchemical
    \chemical
      [SIX,ROT2,B,C,R34,
       PB:RZ4,ONE,Z0,SB8,DB6,Z68,PE]
      [C,0,OH]
 \stopchemical
\stopformula
```

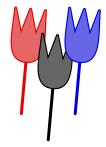


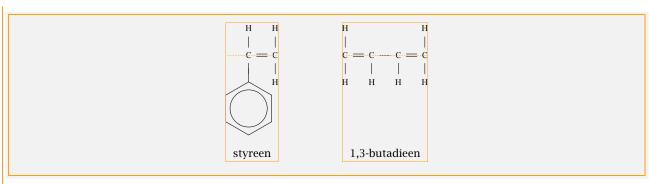
```
\begin{array}{c} \text{O} & \text{O} \\ \text{||} \\ \text{CH}_2 = \text{CH} - \text{CH}_2 - \text{CH}_2 - \text{C} - \text{OH} + \text{H}^{+} \\ \end{array}
```

```
\startformula
  \setupchemical
    [width=fit]
  \startchemical
    \chemical
      [ONE,Z0,OFF1,DB1,
       MOV1,0FF1,Z0,0FF1,SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1,Z0,SB1,DB7,Z1,Z7]
      [CH_2,CH,CH_2,CH_2,C,OH,O]
  \stopchemical
  \startchemical
    \chemical[SPACE,PLUS,SPACE]
  \stopchemical
  \startchemical
    \chemical[ONE,Z0][H^+]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, Z0, OFF1, SB1,
       MOV1, OFF1, Z0, OFF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1,Z0,SB1,DB7,Z1,Z7]
      [CH_3,\T{\oplus}CH,CH_2,CH_2,C,OH,O]
  \stopchemical
\stopformula
```

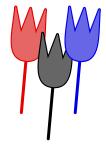


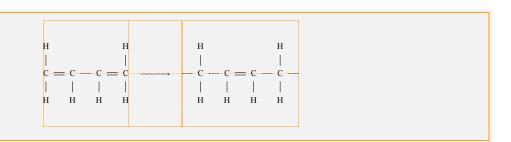
```
\startformula
  \setupchemical
    [width=fit]
  \startchemical
    \chemical
      [ONE, Z0, OFF1, SB1,
       MOV1, OFF1, Z0, OFF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1, OFF1, Z0, 20FF1, SB1,
       MOV1,Z0,SB1,DB7,Z1,Z7]
      [CH_3,\T{\oplus}CH,CH_2,CH_2,C,OH,O]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical[bottom=2500]
    \chemical
      [FIVE, ROT4, Z12345, SB, DR2, CRZ2, -SR4, +SR4, -RZ4, +RZ4]
      [H_2C~~,C,0,C,H_2C~~,0,CH_3,H]
    \bottext{compound B}
  \stopchemical
  \startchemical
    \chemical[SPACE,PLUS,SPACE]
  \stopchemical
  \startchemical
    \chemical[ONE,Z0][H^+]
  \stopchemical
\stopformula
```



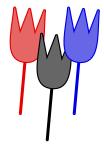


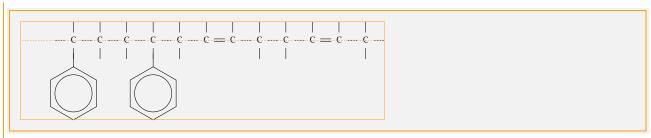
```
\startformula
  \setupchemical
    [width=fit,
     bottom=4000
  \startchemical
    \chemical
      [ONE, Z0, DB1, SB3, SB7, Z7,
       MOV1,Z0,SB3,SB7,Z3,Z7,
       MOV0, SUB2, SIX, B, R6, C]
      [C,H,C,H,H]
    \bottext{styreen}
  \stopchemical
  \quad\quad\quad\quad\quad
  \startchemical
    \chemical
      [ONE, Z0, DB1, SB3, SB7, Z3, Z7,
       MOV1,Z0,SB1,SB3,Z3,
       MOV1, Z0, DB1, SB3, Z3,
       MOV1,Z0,SB3,SB7,Z3,Z7]
      [C,H,H,C,H,C,H,C,H,H]
    \bottext{1,3-butadieen}
  \stopchemical
\stopformula
```



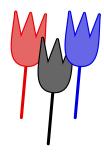


```
\startformula
  \setupchemical
    [width=fit]
  \startchemical
    \chemical
      [ONE, Z0, DB1, Z37, SB37,
       MOV1,Z03,SB13,
       MOV1,Z03,DB1,SB3,
       MOV1,Z037,SB37]
      [C,H,H,C,H,C,H,C,H,H]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, Z037, SB1357,
       MOV1,Z03,DB1,SB3,
       MOV1, Z03, SB13,
       MOV1, Z037, SB137]
      [C,H,H,C,H,C,H,C,H,H]
  \stopchemical
\stopformula
```



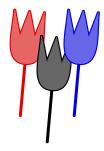


```
\setupchemical
  [height=fit]
\startchemical
  \chemical
    [ONE, Z0, SB1357,
   SAVE,
     SUB2, SIX, B, C, R6,
   RESTORE,
     MOV1, Z0, SB137,
     MOV1, Z0, SB137,
     MOV1,Z0,SB137,
   SAVE,
     SUB2, SIX, B, C, R6,
   RESTORE,
     MOV1, Z0, SB137,
     MOV1,Z0,DB1,SB7,
     MOV1, Z0, SB17,
     MOV1, Z0, SB137,
     MOV1, Z0, SB137,
     MOV1,Z0,DB1,SB7,
     MOV1,Z0,SB17,
     MOV1,Z0,SB137]
    [C,C,C,C,C,C,C,C,C,C,C]
\stopchemical
```





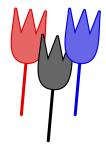
\setupchemical
 [height=fit]
\startchemical
 \chemical[NEWMAN,C,B123,Z123][C00^{-},C00^{-},^{-}00C]
\stopchemical



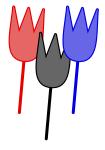
```
OH OH

OH OH
```

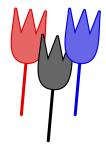
```
\setupchemical
    [width=fit]
\startchemical
    \chemical
    [ONE, Z0, SB14, DB6, Z46,
        MOV1, Z0, SB137, Z37,
        MOV1, Z0, SB137, Z37,
        MOV1, Z0, SB2, DB8, Z28]
    [C, ^-0,0,C,OH,H,C,OH,H,C,O^-,O]
\stopchemical
```



```
\setupchemical
    [width=fit]
\startchemical
    \chemical
    [ONE, Z0, SB14, DB6, Z46,
        MOV1, Z0, SB137, Z37,
        MOV1, Z0, SB137, Z37,
        MOV1, Z0, SB2, DB8, Z28]
    [C, H0, 0, C, 0^-, H, C, 0^-, H, C, OH, 0]
\stopchemical
```



```
\startformula
  \setupchemical
    [width=fit]
  \startchemical
    \chemical
      [ONE, Z0, OFF1, SB1,
       MOV1, OFF1, Z0, OFF1, SB13, Z3,
       MOV1, Z0, DB1, SB3, Z3,
       MOV1,Z0]
      [CH_3,CH,OH,C,OH,O]
  \stopchemical
  \startchemical
    \chemical[SPACE,EQUILIBRIUM,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, Z0, OFF1, SB1,
       MOV1, OFF1, Z0, OFF1, SB13, Z3,
       MOV1, Z0, DB1, SB3, Z3,
       MOV1,Z0]
      [CH_3, CH, OH, C, 0^-, 0]
   \stopchemical
   \startchemical
     \chemical[SPACE, PLUS, SPACE]
  \stopchemical
  \startchemical
    \chemical[ONE,OFF1,Z0][H^+]
  \stopchemical
\stopformula
```



```
\startchemical[width=fit]
\chemical

[ONE,Z0357,SB1357,

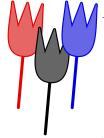
MOV1,Z037,SB137,

MOV1,Z037,SB137,

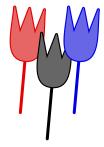
MOV1,Z037,SB137,

MOV1,Z0137,SB137]

[C,OH,H,H,C,OH,H,C,OH,H,C,OH,H,C,H,OH,H]
\stopchemical
```

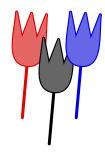


```
\setupchemical
  [width=fit,
   height=fit,
   bottom=2000
\startchemical
  \chemical
    [ONE,OFF1,Z0,OFF1,SB1,OE5,
     MOV1, OFF1, Z0, OFF1, SB1, Z3, SB3,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1, OFF1, Z0, OFF1, SB1, Z3, SB3,
     MOV1, OFF1, Z0, 20FF1, SB1,
     MOV1, OFF1, Z0, Z3, SB3, OE1]
    [CH_2,CH,CH_3,CH_2,CH,CH_3,CH_2,CH,CH_3]
  \bottext{poloymer A}
\stopchemical
\blank[3*big]
\startchemical
  \chemical
    [ONE, ZO, SB1, OE5,
     MOV1, OFF1, Z0, OFF1, SB1, Z3, SB3,
     MOV1, Z0, SB1, Z7, DB7,
     MOV1,Z0,SB1,
     MOV1, Z0, SB1, Z3, SB3,
     MOV1, OFF1, Z0, OFF1, SB1, Z7, DB7,
     MOV1, Z0, SB1,
     MOV1, OFF1, Z0, Z3, SB3, OE1]
    [0,CH,CH_3,C,0,0,CH,CH_3,C,0,0,CH,CH_3]
  \bottext{polymer B}
\stopchemical
```

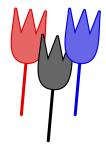


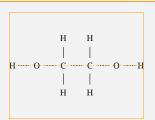
```
О
                                      О
                                                                                                                                                                       О
                                                                                                                                CH
                                                                                                                                                             CH-
                                                                                                                                                                              - OH
                                                                                                                                                                                               HO --- CH ---
CH-
                            CH-
                                                        CH-\\c
                                                                                 H<sub>2</sub>O
                                                                                                                                                                                                          |
CH<sub>3</sub>
CH<sub>3</sub>
                             CH_3
                                                         CH<sub>3</sub>
                                                                                                                                 CH_3
                                                                                                                                                              CH<sub>3</sub>
```

```
\startformula
 \setupchemical
    [width=fit,
     height=2750,
     scale=small,
     size=small]
  \startchemical
    \chemical
      [ONE, ZO, SB1, OE5,
       MOV1, OFF1, Z0, OFF1, SB1, Z3, SB3,
       MOV1,Z0,SB1,Z7,DB7,
       MOV1, Z0, SB1,
       MOV1, OFF1, Z0, OFF1, SB1, Z3, SB3,
       MOV1,Z0,SB1,Z7,DB7,
       MOV1, Z0, SB1,
       MOV1, OFF1, Z0, Z3, SB3, OE1]
      [0,CH,CH_3,C,0,0,CH,CH_3,C,0,0,CH,CH_3]
 \stopchemical
 \quad\quad
 \startchemical
    \chemical[SPACE, PLUS, SPACE]
 \stopchemical
 \startchemical
    \chemical[ONE,OFF1,Z0][H_20]
 \stopchemical
 \startchemical
    \chemical[SPACE,GIVES,SPACE]
 \stopchemical
 \quad\quad
 \startchemical
    \chemical
      [ONE, ZO, SB1, OE5,
       MOV1, OFF1, Z0, OFF1, SB1, Z3, SB3,
       MOV1,Z0,SB1,Z7,DB7,
       MOV1, Z0, SB1,
       MOV1, OFF1, Z0, OFF1, Z3, SB3, SB1,
       MOV1,Z0,SB1,DB7,Z17]
      [0,CH,CH_3,C,0,0,CH,CH_3,C,OH,0]
 \stopchemical
 \startchemical
    \chemical[SPACE,PLUS,SPACE]
  \stopchemical
```



\startchemical
 \chemical[ONE,OFF1,Z0,SB35,Z35,OE1][CH,CH_3,H0]
 \stopchemical
 \stopformula





```
\startchemical[width=fit]
\chemical

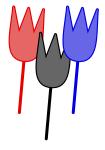
[ONE,Z05,SB15,

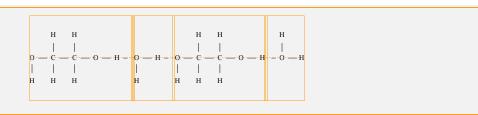
MOV1,Z037,SB137,

MOV1,Z037,SB137,

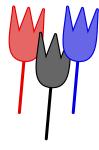
MOV1,Z01,SB1]

[O,H,C,H,H,C,H,H,O,H]
\stopchemical
```



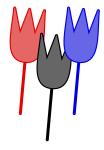


```
\startformula
  \setupchemical
    [width=fit,
     scale=small,
     size=small]
  \startchemical
    \chemical
      [ONE, Z03, SB13,
       MOV1, Z037, SB137,
       MOV1, Z037, SB137,
       MOV1,Z0,SB1,
       MOV1,Z0,SD1]
      [0,H,C,H,H,C,H,H,O,H]
  \stopchemical
  \startchemical[width=100]
    \chemical[SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, SB1, SB3, Z0, Z3, MOV1, Z0, SD1] [0, H, H]
  \stopchemical
  \startchemical[width=100]
    \chemical[SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, Z03, SB13,
       MOV1, Z037, SB137,
       MOV1, Z037, SB137,
       MOV1, Z0, SB1,
       MOV1,Z0
      [0,H,C,H,H,C,H,H,O,H]
  \stopchemical
  \startchemical[width=100]
    \chemical[SPACE]
  \stopchemical
  \startchemical
    \chemical[ONE,Z017,SB17,SD5][0,H,H]
  \stopchemical
\stopformula
```

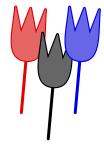


```
----CH<sub>3</sub> -----N-==-C-== O
```

```
\startchemical[height=2000]
  \chemical
    [ONE,DB1,SB5,Z0,Z5,
        MOV1,DB1,Z0,Z1]
    [N,CH_3,C,0]
  \stopchemical
```

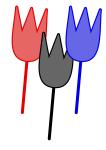


```
\startchemical
  [width=fit]
\chemical
   [ONE,OFF1,Z0,OFF1,Z5,SB135,
  SAVE,
    MOV3,Z0,SB3,
    MOV3,Z0,SB3,
    MOV3,Z013,DB1,SB3,
  RESTORE,
    MOV1, Z07, SB1, DB7,
    MOV1,Z07,SB17,
    MOV1, OFF1, Z0, OFF1, SB13,
  SAVE,
    MOV3,Z03,SB3,
  RESTORE,
    MOV1, Z07, SB1, DB7,
    MOV1, Z07, SB17,
    MOV1, OFF1, Z0, 20FF1, SB1,
    MOV1,Z017,SB1,DB7]
   [CH, H_2N, CH_2, CH_2, C, O, OH, C, O, N, H, CH, CH_2,
    SH,C,O,N,H,CH_2,C,OH,O]
\stopchemical
```



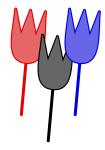
```
CH_3 - N = C = 0 + H - S additieprodukt
```

```
\def\Subs
   {\hbox
      {\setbox0=\hbox\framed{Glu\hairline Cys\hairline Gly}%
       \hskip-.5\wd0\lower.5\ht0\box0}
\startformula
 \setupchemical
    [width=fit,
     height=2000,
     size=big]
 \startchemical
    \chemical
      [ONE, Z05, DB1, SB5, MOV1, Z01, DB1] [N, CH_3, C, O]
 \stopchemical
 \startchemical
    \chemical[SPACE, PLUS, SPACE]
 \stopchemical
 \startchemical
    \chemical[ONE, Z035, SB35][S, \Subs, H]
 \stopchemical
 \startchemical
    \chemical[SPACE,GIVES,SPACE]
 \stopchemical
 \startchemical
    \chemical[ONE,Z0][additieprodukt]
 \stopchemical
\stopformula
```





```
\def\Subs
  {\hbox
     {\setbox0=\hbox\framed{Glu\hairline Cys\hairline Gly}%
      \hskip-.5\wd0\lower.5\ht0\box0}
\startformula
  \setupchemical
    [width=fit,
     height=1500]
  \startchemical
    \chemical
      [ONE, Z035, SB135,
      MOV1,Z01,DB1,SB3,
       MOV0, MOV1, MOV3, Z03, SB3]
      [N,H,CH_3,C,0,S,\Subs]
  \stopchemical
\stopformula
```



```
H H H

| | | |

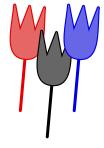
H—C—C—C—H

| | |

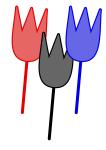
H C H

H | H
```

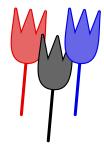
```
\startchemical[width=fit]
\chemical
[ONE,Z0357,SB1357,
MOV1,Z07,SB137,
SAVE,MOV1,Z0137,SB137,RESTORE,
MOV3,Z0234,SB234,]
[C,H,H,H,C,H,C,H,H,H,C,H,H,H]
\stopchemical
```

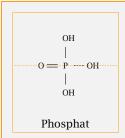


```
\startformula
  \setupchemical
    [width=fit,
     scale=small,
     size=small]
  \startchemical
    \chemical
      [ONE, ZO, SB1357, Z357,
       MOV1,Z0,SB137,Z37,
       MOV1,Z0,SB137,Z37,
       MOV1,Z0,SB137,Z37,
       MOV1, Z0, SB137, Z37,
       MOV1, Z0, SB137, Z37,
       MOV1,Z0,SB137,Z137,]
      [C,H,H,H,C,H,H,C,H,H,C,H,H,C,H,H,C,H,H,H]
  \stopchemical
  \startchemical
    \chemical[SPACE,GIVES,SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, ZO, SB1357, Z357,
       MOV1, Z0, SB137, Z37,
       MOV1, Z0, SB137, Z37,
       MOV1,Z0,SB137,Z37,
       MOV1,Z0,SB137,Z137]
      [C,H,H,H,C,H,H,C,H,H,C,H,H,C,H,H,H]
  \stopchemical
  \startchemical
    \chemical[SPACE, PLUS, SPACE]
  \stopchemical
  \startchemical
    \chemical
      [ONE, ZO, DB1, SB37, Z37,
       MOV1,Z0,SB37,Z37]
      [C,H,H,C,H,H]
  \stopchemical
\stopformula
```

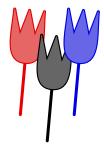


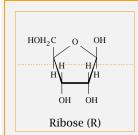
```
\startformula
\startchemical[width=fit]
\chemical
    [ONE,Z037,SB37,DB1,
        MOV1,Z07,SB17,
        MOV1,Z037,SB137,
        MOV1,Z0137,SB137]
        [C,H,H,C,H,C,H,H,C,H,H,H]
\stopchemical
\stopformula
```



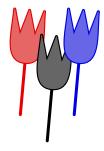


```
\definechemical[phosphat]
  {\chemical[ONE,SB137,DB5,Z01357][P,OH,OH,O,OH]}
\startchemical[height=6500,top=2000,bottom=2500,width=4000]
  \chemical[phosphat]
  \bottext{Phosphat}
\stopchemical
```



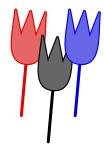


```
\definechemical[ribose]
  {\chemical[FIVE,FRONT,BB125,+SB3,-SB4,Z4][0]
   \chemical[FIVE,FRONT,+R1235,+RZ1235][H,H,\SR{HOH_2C},OH]
   \chemical[FIVE,FRONT,-R1235,-RZ1235][OH,OH,H,H]}
\startchemical[height=6500,top=2000,bottom=2500,width=4500]
   \chemical[ribose]
   \bottext{Ribose (R)}
\stopchemical
```



```
HOH<sub>2</sub>C O OH
H H
OH H
Desoxyribose (dR)
```

```
\definechemical[desoxyribose]
    {\chemical[FIVE,FRONT,BB125,+SB3,-SB4,Z4][0]
      \chemical[FIVE,FRONT,+R1235,+RZ1235][H,H,\SR{HOH_2C},OH]
      \chemical[FIVE,FRONT,-R1235,-RZ1235][H,OH,H,H]}
\startchemical[height=6500,top=2000,bottom=2500,width=4500]
      \chemical[desoxyribose]
      \bottext{Desoxyribose (dR)}
\stopchemical
```



```
NH<sub>2</sub>

N

N

N

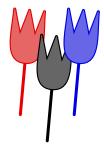
N

N

H

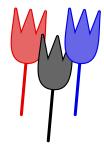
Cytosin (C)
```

```
\definechemical[cytosin]
{\chemical
       [SIX,B1,EB1,+SB2,-SB3,+SB4,-SB5,EB5,B6,Z3,Z5,SR3,ER4,R6,RZ3,RZ4,RZ6]
       [N,N,H,0,NH_2]}
\startchemical[height=6500,top=2000,bottom=2500,width=4500]
    \chemical[cytosin]
    \bottext{Cytosin (C)}
\stopchemical
```



```
O CH<sub>3</sub>
O N CH<sub>3</sub>
H Thymin (T)
```

```
\definechemical[thymin]
    {\chemical
        [SIX,B1,EB1,+SB2,-SB3,+SB4,-SB5,B6,Z35,R1,SR3,ER4,SR5,ER6,RZ13456]
        [N,N,CH_3,H,O,H,O,]}
\startchemical[height=6500,top=2000,bottom=2500,width=4500]
    \chemical[thymin]
    \bottext{Thymin (T)}
\stopchemical
```



```
O

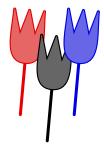
N

O

N

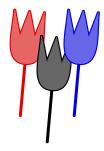
H

Uracil (U)
```



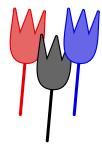
```
H<sub>2</sub>N N N H<sub>2</sub>H Guanin (G)
```

```
\definechemical[guanin]
    {\chemical[SIX,B6,B1,+SB2,-SB3,EB3,+SB4,-SB5,Z35][N,N]
        \chemical[SIX,R4,SR5,ER6,RZ456][\SR{H_2N},H,O]
        \chemical[SIX,MOV1,-SB1,EB1,+SB2,-SB3,EB4,-SS6,Z13,SR3,RZ3][N,N,H]}
\startchemical[height=6500,top=2000,bottom=2500,width=5000,left=2000]
        \chemical[guanin]
        \bottext{Guanin (G)}
\stopchemical
```



```
NH<sub>2</sub>
N
N
N
N
Adenin (A)
```

```
\definechemical[adenin]
{\chemical[SIX,B6,B1,+SB2,-SB3,EB3,+SB4,-SB5,EB5,Z35][N,N]
  \chemical[SIX,R6,RZ6][\SL{NH_2}]
  \chemical[SIX,MOV1,-SB1,EB1,+SB2,-SB3,EB4,-SS6,Z13,SR3,RZ3][N,N,H]}
\startchemical[height=6500,top=2000,bottom=2500,width=4500,left=1500]
  \bottext{Adenin (A)}
  \chemical[adenin]
\stopchemical
```



```
HOH<sub>2</sub>C 5 O OH

C4 1C

H 3 2 / H

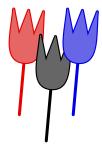
OH - C - C - H

H H

Desoxyribose (dR)
```

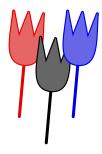
```
\definechemical[desoxyribose again]
    {\chemical[FIVE,ROT2,SB12345,Z12345][\TL{2}C,\TR{3}C,\R{4}C,0,\L{1}C]
    \chemical[FIVE,ROT2,+SR1235,+RZ1235][H,OH,\TR{5}{HOH_2C},H]
    \chemical[FIVE,ROT2,-SR1235,-RZ1235][H,H,H,OH]}
\startchemical[height=6500,top=2000,bottom=2500,width=5000]
    \chemical[desoxyribose again]
    \bottext{Desoxyribose (dR)}
\stopchemical
```

Na \longrightarrow Na⁺ + e⁻] \times 2

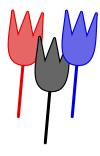


```
\startformula
  \chemical{Na}
  \chemical{GIVES}
  \chemical{Na^{+}}
  \chemical{PLUS}
  \chemical{e^{-}} \quad \rbrack \times 2
\stopformula
```



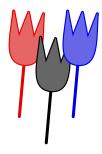


```
\startformula
  \chemical{2\,NH_{3}}
  \chemical{PLUS}
  \chemical{2\,e^{-}}
  \chemical{GIVES}
  \chemical{2\,NH_{2}^{-}}
  \chemical{PLUS}
  \chemical{H_{2}}
  \stopformula
```



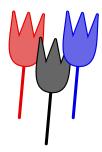
```
\startformula
  \chemical{2\,Na}
  \chemical{PLUS}
  \chemical{2\,NH_{3}}
  \chemical{GIVES}
  \chemical{2\,Na^{+}}
  \chemical{PLUS}
  \chemical{2\,NH_{2}^{-}}
  \chemical{PLUS}
  \chemical{PLUS}
  \chemical{H_{2}}
  \stopformula
```

$\mathrm{NH_2^-} \ + \ \mathrm{H_2O} \quad \longrightarrow \quad \mathrm{NH_3} \ + \ \mathrm{OH^-}$

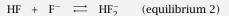


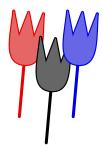
```
\startformula
  \chemical{NH_{2}^{-}}
  \chemical{PLUS}
  \chemical{H_{2}0}
  \chemical{GIVES}
  \chemical{NH_{3}}
  \chemical{PLUS}
  \chemical{OH^{-}}
\stopformula
```





```
\startformula
  \matrix
    {\chemical{HF}
     \quad
     \chemical{PLUS}
     \quad
     \chemical{H_{2}0}
     \quad
     \chemical{EQUILIBRIUM}
     \quad
     \chemical{H_{3}0^{+}}
     \quad
     \chemical{PLUS}
     \quad
     \chemical{F^{-}}}
                             &
     \quad
     \rm (equilibrium~1)
                             \cr}
\stopformula
```

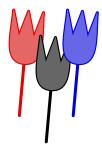




```
\startformula
  \matrix
    {\chemical{HF}}
    \quad
    \chemical{PLUS}
    \quad
    \chemical{F^{-}}
    \quad
    \chemical{EQUILIBRIUM}
    \quad
    \chemical{HF_{2}^{-}}

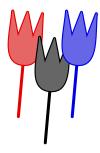
    \quad
    \chemical{HF_2^2}
    \quad
    \rm (equilibrium~2) \cr}
\stopformula
```





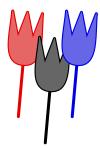
```
\startformula
  \chemical{C_{5}H_{10}\quad(g)}
  \quad
  \chemical{EQUILIBRIUM}
  \quad
  \chemical{C_{5}H_{8}\quad(g)}
  \quad
  \chemical{PLUS}
  \quad
  \chemical{H_{2}\quad(g)}
\stopformula
```



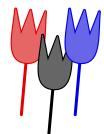


```
\startformula
  \chemical{2\,S_{2}0_{3}^{2-}}
  \quad
  \chemical{GIVES}
  \quad
  \chemical{S_{2}0_{6}^{2-}}
  \quad
  \chemical{PLUS}
  \quad
  \chemical{2\,e^{-}}
  \stopformula
```



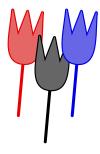


```
\startformula
  \chemical{S_{2}0_{3}^{2-}}
  \quad
  \chemical{PLUS}
  \quad
  \chemical{5},H_{2}0
  \quad
  \chemical{GIVES}
  \quad
 \chemical{2\,}SO_{4}^{2}}
  \quad
  \chemical{PLUS}
  \quad
  \chemical{10\,H^{+}}
  \quad
  \chemical{PLUS}{}
  \quad
  \chemical{8},e^{-}
\stopformula
```



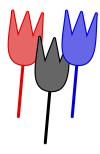
```
S_2O_3^{2-} \quad (aq) \quad + \quad 2\,H^+ \quad (aq) \quad \longrightarrow \quad SO_2 \quad (aq) \quad + \quad S \quad (s) \quad + \quad H_2O \quad (l) \qquad (reactie \ 1)
```

```
\startformula
  \matrix
    {\chemical{S_{2}0_{3}^{2}}\qquad ad(aq)}
     \quad
     \chemical{PLUS}
     \quad
     \chemical{2\,H^{+}}\quad (aq)
     \quad
     \chemical{GIVES}
     \quad
     \chemical{SO_{2}\quad(aq)}
     \quad
     \chemical{PLUS}
     \quad
     \chemical{S\quad(s)}
     \quad
     \chemical{PLUS}
     \quad
     \chemical{H_{2}0\qquad(1)}
     \quad
     \rm (reactie<sup>1</sup>)
                                  \cr}
\stopformula
```

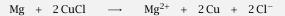


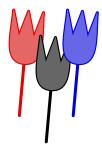
```
\startformula
  \chemical{FeTi0_3\quad(s)}
  \chemical{PLUS}
  \chemical{H_2\quad(g)}
  \chemical{GIVES}
  \chemical{Fe\quad(s)}
  \chemical{PLUS}
  \chemical{Ti0_2\quad(s)}
  \chemical{PLUS}
  \chemical{PLUS}
  \chemical{H_20\quad(g)}
  \stopformula
```

$2\,\text{NaN}_3 \quad (s) \qquad \longrightarrow \qquad 2\,\text{Na} \quad (s) \quad + \quad 3\,\text{N}_2 \quad (g)$

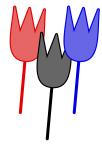


```
\startformula
  \chemical{2\,NaN_3\quad(s)}
  \chemical{GIVES}
  \chemical{2\,Na\quad(s)}
  \chemical{PLUS}
  \chemical{3\,N_2\quad(g)}
  \stopformula
```



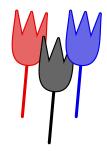


```
\startformula
  \chemical{Mg}
  \chemical{PLUS}
  \chemical{2\,CuCl}
  \chemical{GIVES}
  \chemical{Mg^{2+}}
  \chemical{PLUS}
  \chemical{2\,Cu}
  \chemical{PLUS}
  \chemical{PLUS}
  \chemical{2\,Cl^-}
  \stopformula
```



```
SO_4^{2-}+4H_2O+8e^- \Longrightarrow S^{2-}+8OH^-
                                 4\,\text{Fe} \rightleftharpoons 4\,\text{Fe}^{2+} + 8\,\text{e}^{-}
4 \text{ Fe+SO}_4^{2-} + 4 \text{ H}_2 \text{O} \Longrightarrow \text{S}^{2-} + 4 \text{ Fe}^{2+} + 8 \text{ OH}^-
                    Fe^{2+}+S^{2-}\longrightarrow FeS
```

```
\startformula
 \eqalign
    {\chemical{SO_4^{2-}}}
     \chemical{PLUS}
     \chemical{4}, H_20
     \chemical{PLUS}
     \chemical{8}, e^-
                                      &
     \chemical{EQUILIBRIUM}
     \cdot S^{S^{2-}}
     \chemical{PLUS}
     \chemical{8},OH^-
                                      \cr
     \chemical{4}, Fe
                                      &
     \chemical{EQUILIBRIUM}
     \chemical{4}, Fe^{2+}
     \chemical{PLUS}
     \chemical{8}, e^-
                                      \cr
     \chemical{4}, Fe}
     \chemical{PLUS}
     \chemical{SO_4^{2-}}
     \chemical{PLUS}
     \chemical{4}, H_20
                                      &
     \chemical{EQUILIBRIUM}
     \chemical{S^{2-}}
     \chemical{PLUS}
     \chemical{4}, Fe^{2+}
     \chemical{PLUS}
     \chemical{8},OH^-
                                      \cr
     \chemical{Fe^{2+}}
     \chemical{PLUS}
     \chemical{S^{2-}}
                                      &
     \chemical{GIVES}
     \chemical{FeS\downarrow}{dark} \cr}
\stopformula
```



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examples, manuals and code | www.ntg.nl/context

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