

# Formatted manual of singularcurve.lib

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# 1 Singular libraries

## 1.1 singularcurve\_lib

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**Library:** singularcurve.lib

**Purpose:** targets calculation of Zeta-function of a curve with ordinary double points over a finite field

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**Procedures:**

### 1.1.0.1 sumintvec

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve_lib`], page 1).

**Usage:** `sumintvec(e); e intvec`

**Assume:** none

**Return:** int: the sum of the entries of e

**Note:** none

**Example:**

```
LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
intvec e = 1,2,3;
int six = sumintvec(e);
six;
⇒ 6
```

### 1.1.0.2 list2ideal

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve_lib`], page 1).

**Usage:** `list2ideal(polylist); list polylist`

**Assume:** none

**Return:** ideal: the ideal  $I = \text{polylist}[1], \dots, \text{polylist}[n]$ ;

**Note:** this procedure can be used to sum ideals avoiding simplification

**Example:**

```

LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
ring r = 0, x, dp;
ideal I1 = x;
ideal I2 = 1;
list IL;
for(int i = 1; i<=size(I1); i++){
  IL = insert(IL, I1[i]);
}
for(int i = 1; i<=size(I2); i++){
  IL = insert(IL, I2[i]);
}
⇒ // ** redefining i **
ideal I3 = list2ideal(IL);
I1 + I2;
⇒ _[1]=1
I3;
⇒ I3[1]=1
⇒ I3[2]=x

```

### 1.1.0.3 multinomial

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve.lib`], page 1).

**Usage:**      `multinomial(N, iv):` int N, intvec iv

**Assume:**    none

**Return:**    bigint multinomial coefficient  $N!/(iv[1]! \cdots iv[n]!)$

**Note:**      needs `general.lib`; doesn't check  $N = \text{sumintvec}(iv)$

**Example:**

```

LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
LIB "general.lib";
intvec iv = 1,2,3;
bigint mulCoef = multinomial(6, iv);
mulCoef;
⇒ 60

```

### 1.1.0.4 ceilquot

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve.lib`], page 1).

**Usage:** `ceilquot(n1, n2); int n1, int n2`

**Assume:** none

**Return:** `ceil(decimal(n1/n2))`

**Note:** none

**Example:**

```
LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
ceilquot(1,2);
⇒ 1
ceilquot(2,2);
⇒ 1
ceilquot(-1,2);
⇒ 0
```

### 1.1.0.5 floorquot

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve.lib`], page 1).

**Usage:** `floorquot(n1, n2); int n1, int n2`

**Assume:** none

**Return:** `floor(decimal(n1/n2))`

**Note:** none

**Example:**

```
LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
floorquot(1,2);
⇒ 0
floorquot(2,2);
⇒ 1
floorquot(-1,2);
⇒ -1
```

### 1.1.0.6 padicOrder

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve.lib`], page 1).

**Usage:** `padicOrder(n, p)`: number  $n$ , bigint  $p$

**Assume:** basering is of characteristic 0

**Return:** returns the padic order of  $n$ : if  $n = p^k \cdot a/b$  then returns  $\text{order}(a) - \text{order}(b) + k$ ; returns 'Inf' if  $n == 0$

**Note:** needs `poly.lib`

**Example:**

```
LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
ring r = 0, x, dp;
LIB "poly.lib";
number n1 = number(1/2);
number n2 = number(98/3);
number n3 = number(0);
padicOrder(n1, 7);
⇒ 0
padicOrder(n2, 7);
⇒ 2
padicOrder(n3, 7);
⇒ Inf
```

### 1.1.0.7 padicApprox

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve.lib`], page 1).

**Usage:** `padicApprox(quot, p, N)`: number(quot), int  $p$ , int  $N$

**Assume:** basering is of characteristic 0

**Return:** returns the p-adic approximation of a rational number up to certain precision

**Note:**  $a$  is approximation of  $b$  to precision  $N \iff a - b = 0 \pmod{p^N}$ , i.e.  $a_i = 0$ ,  $i \geq N$ , include `poly.lib`

**Example:**

```
LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
```

```

⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
LIB "poly.lib";
ring r = 0, x, dp;
number n1 = number(0);
number n2 = number(1/2);
number n3 = 7*3 + 49*4;
padicApprox(n1, 7, 0);
⇒ 0
padicApprox(n2, 7, 7);
⇒ 411772
padicApprox(n2, 7, 4);
⇒ 1201
padicApprox(n3, 7, 3);
⇒ 217

```

### 1.1.0.8 reduceStep

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve.lib], page 1).

**Usage:**      `reduceStep(nom, jf, s):` poly nom, ideal jf, int s

**Assume:**    `nom = 0 mod jf`, basering of characteristic 0

**Return:**    the polynomial  $f$  s.t.  $f/F^{(s-1)}\Omega = \text{nom}/F^{(s)}\Omega$

**Note:**      none

**Example:**

```

LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
ring r = 0, (x,y,z), dp;
poly f = x^2 + y^2 + z^2;
ideal jf = jacob(ideal(f));
"2s-2:0->2::-2,0,2";
⇒ 2s-2:0->2::-2,0,2
"2s-2:3::4::s==3";
⇒ 2s-2:3::4::s==3
poly nom = x^2*y^2;
def reduced = reduceStep(nom, jf, 3);
reduced;
⇒ 1/4x2

```

### 1.1.0.9 reduceAll

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve.lib], page 1).

**Usage:**      `reduceAll(nom, jf, s):` poly nom, ideal jf, int s

**Assume:**    `nom == 0 mod jf`

**Return:** poly expression that expresses nom in terms of a base of  $R/jf$

**Note:**  $s$  is the degree of the denominator  $F_k^s$

**Example:**

```
LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
ring r = 0, (x,y,z), dp;
poly f = x2z3 + y2z3 + x5 + y5 + 7z5;
ideal jf = jacob(f);
poly nom = x^2 + y^2;
nom = nom*(f^10);
int s = (deg(nom)+3)/5;
⇒ // ** int division with '/': use 'div' instead in line >>int s = (deg(nom\
)+3)/5;<<
reduceAll(nom, jf, s);
⇒ x2+y2
```

### 1.1.0.10 extractIdeal

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve.lib`], page 1).

**Usage:** `extractIdeal(base, degrees)`: ideal base, intvec degrees

**Assume:** none

**Return:** the ideal containing only those generators of the input ideal of certain degrees

**Note:** none

**Example:**

```
LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
LIB "general.lib";
ring r = 0, (x,y,z), dp;
poly f = x2z3 + y2z3 + x5 + y5 + 7z5;
ideal jf = jacob(f);
ideal base = kbase(std(jf));
size(base);
⇒ 64
base = extractIdeal(base, intvec(2,7));
```

[illegible]



```

⇒ // ** redefining j **
⇒ // ** redefining j **
⇒ // ** redefining j **
⇒ // ** redefining j **
⇒ // ** redefining j **
base;
⇒ base[1]=z2
⇒ base[2]=yz
⇒ base[3]=xz
⇒ base[4]=y2
⇒ base[5]=xy
⇒ base[6]=x2
⇒ base[7]=z7
⇒ base[8]=yz6
⇒ base[9]=xz6
⇒ base[10]=xyz5
⇒ base[11]=xy3z3
⇒ base[12]=x3y3z

```

### 1.1.0.11 getMatrix

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve-lib], page 1).

**Usage:** `getMatrix(base, image)`: ideal base, list image

**Assume:** the generators of base are base vectors of a module; image contains the image of each generator represented by a linear combination of the generators of base

**Return:** the matrix representing the map  $\text{base}[i] \rightarrow \text{image}[i]$

**Note:** none

**Example:**

```

LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
LIB "general.lib";
ring r = 0, (x,y,z), dp;
ideal bf = x2,y2,z2, xy,xz,yz;
bf = sort(bf)[1];
matrix m1[6][6] = 1,2,3,4,5,6,7,8,9,10,11,12,1,2,3,4,5,6,7,8,9,10,11,12,1,2,3,4,5,6,7,8,9,10,11,12;
list image = list();
for(int col = 6; col>=1; col--){
  poly tmpoly = 0;
  for(int row = 1; row<=6; row++){
    tmpoly = tmpoly + m1[row, col]*bf[row];
  }
  image = insert(image, tmpoly);
}
⇒ // ** redefining tmpoly **
⇒ // ** redefining row **

```

[illegible]

```

⇒ m2[2,3]=9
⇒ m2[2,4]=10
⇒ m2[2,5]=11
⇒ m2[2,6]=12
⇒ m2[3,1]=1
⇒ m2[3,2]=2
⇒ m2[3,3]=3
⇒ m2[3,4]=4
⇒ m2[3,5]=5
⇒ m2[3,6]=6
⇒ m2[4,1]=7
⇒ m2[4,2]=8
⇒ m2[4,3]=9
⇒ m2[4,4]=10
⇒ m2[4,5]=11
⇒ m2[4,6]=12
⇒ m2[5,1]=1
⇒ m2[5,2]=2
⇒ m2[5,3]=3
⇒ m2[5,4]=4
⇒ m2[5,5]=5
⇒ m2[5,6]=6
⇒ m2[6,1]=7
⇒ m2[6,2]=8
⇒ m2[6,3]=9
⇒ m2[6,4]=10
⇒ m2[6,5]=11
⇒ m2[6,6]=12

```

### 1.1.0.12 inverseI

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve\_lib], page 1).

**Usage:**      `inverseI(int, int)`

**Assume:**    there is an  $m$  in  $\{2, \dots, q-1\}$  s.t.  $nm = n\_m q - 1 = -1 \bmod q$

**Return:**    `list(m, n_m)`

**Example:**

```

LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
ring r = 0, x, dp;
int a = 5;
int q = 7;
def result = inverseI(a,q);
if(typeof(b) == "string"){
//error handling
}
result;

```

```

⇒ [1]:
⇒ 4
⇒ [2]:
⇒ 3
a*result[1] == q*result[2] - 1;
⇒ 1

```

### 1.1.0.13 inverse

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve.lib], page 1).

**Usage:**     `inverse(int, int)`

**Assume:**   there is an  $m$  in  $\{2, \dots, q-1\}$  s.t.  $nm = n\_m q + 1 = 1 \bmod q$

**Return:**    `list(m, n_m)`

**Example:**

```

LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
ring r = 0, x, dp;
int a = 5;
int q = 7;
def result = inverseI(a,q);
if(typeof(b) == "string"){
//error handling
}
result;
⇒ [1]:
⇒ 4
⇒ [2]:
⇒ 3
a*result[1] == q*result[2] - 1;
⇒ 1

```

### 1.1.0.14 areEqualLists

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve.lib], page 1).

**Usage:**     `areEqualLists(list, list)`

**Assume:**    nothing

**Returns:**    0 if lists are not equal, 1 else

**Example:**

```

LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **

```

```

⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
ring r = 0, x, dp;
def l1 = list(1,2,3);
def l2 = list(1,2,4);
def l3 = list(1,2,3);
areEqualLists(l1,l2);
⇒ 0
areEqualLists(l1,l3);
⇒ 1

```

### 1.1.0.15 list2intvec

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve-lib], page 1).

**Usage:**     list2intvec(list)

**Assume:**   list contains int only

**Returns:**   an intvec holding the same values as the list

**Example:**

```

LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
ring r = 0, x, dp;
def l = list(1,2,3);
def v = list2intvec(l);
v;
⇒ 1,2,3
typeof(v);
⇒ intvec

```

### 1.1.0.16 sublist

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve-lib], page 1).

**Usage:**     sublist(list, start, end)

**Assume:**   nothing

**Returns:**   the sublist {list(start),...,list(end)}

**Example:**

```

LIB "singularcurve.lib";
⇒ // ** redefining Psi **
⇒ // ** redefining Psi **
⇒ // ** redefining ReduceAllForPoleOrder **

```

```

⇒ // ** redefining ReduceAllForPoleOrder **
⇒ // ** redefining D **
⇒ // ** redefining D **
⇒ // ** redefining C **
⇒ // ** redefining C **
⇒ // ** redefining Sigma **
⇒ // ** redefining Sigma **
ring r = 0,x,dp;
def l = list(1,2,3,4);
def sl = sublist(l, 2,3);
sl;
⇒ [1]:
⇒ 2
⇒ [2]:
⇒ 3

```

### 1.1.0.17 complexAbs

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve.lib], page 1).

**Usage:** complexAbs(number n)

**Assume:** basering with complex coefficient field resp. in subring of CC and atkins.lib has been imported

**Returns:** complex norm of n

### 1.1.0.18 qdivides

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve.lib], page 1).

**Usage:** qdivides(intvec, int)

**Assume:** nothing

**Returns:** 1, if  $q \mid \text{intvec}[i]$  for all  $i$ , 0 else

### 1.1.0.19 PolyMod

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve.lib], page 1).

**Usage:** PolyMod(poly f, bigint q)

**Assume:** f has coefficients that are integers or bigintegers

**Returns:**  $f \bmod q$

### 1.1.0.20 PolyPadicOrder

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve.lib], page 1).

**Usage:** PolyPadicOrder(poly f, bigint p)

**Assume:** f has rational coefficients

**Returns:**  $v_p(f) = \min \{ v_p(a), a \text{ coefficient of } f \}$

### 1.1.0.21 Psi

Procedure from library `singularcurve.lib` (see Section 1.1 [singularcurve.lib], page 1).

**1.1.0.22 ReduceAllForPoleOrder**

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve.lib`], page 1).

**1.1.0.23 D**

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve.lib`], page 1).

**1.1.0.24 C**

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve.lib`], page 1).

**1.1.0.25 Sigma**

Procedure from library `singularcurve.lib` (see Section 1.1 [`singularcurve.lib`], page 1).

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