

8 REFERENCES

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## APPENDIX 1: SOFTWARE, APPARATUS AND TESTS USED

### 1 Software

After investigating the ore dressing simulators available, as documented in section 3, it was decided to use Modsim, available from the University of the Witwatersrand. The first version of the package was obtained in February 1986. However, the package was not entirely debugged, and two further releases were obtained before September 1986.

Since extensive work had been done on Microsim by the University of the Witwatersrand in the meantime, it was decided to use Microsim rather than Modsim, because of Microsim's expandability to hydrometallurgical stream definitions and because of its superior structure. These features have been documented in section 3.

Microsim Version 1.0 was obtained in September 1986. However, this too was not completely debugged, and was superseded by Microsim Version 2.0 in November 1986. Microsim Version 3.0, with optimisation and design capabilities, was used from January 1989.

Models added to the package for this plant were

CRUSH5

CRUSH6\_7 (replaced CRUSH2\_3)

SCREEN5

ROSE1

ROSE2, as described in section 4.

In order to incorporate new models into Microsim, additional program modifications have to be made to files TYPEDEF.PAS, MODEL.DAT, and MODELS.PAS. These modifications will not be given. They serve the purpose of directing the simulator's executive program to the correct model and to allocate the correct model storage space. Furthermore, the simulator's database of form-filling input screens and default parameters must be updated. This is done through compiling and executing DBASE and modifying the file MICROSIM.MOD.

Also necessary for the running of Microsim and the incorporation of

new models are DOS Version 3 (or higher) and the TURBO Pascal Version 3.0 compiler.

## 2 Apparatus

All software was run on an IBM-AT microcomputer with 1MB memory, 1 floppy disk drive and a 40MB hard disk. An IBM EGA graphics screen was used for the flowsheet input and a Roland DXY-980A plotter for output of flowsheets, partition curves and size distributions. The printer used was a Fujitsu DX2200 printer.

## 3 Tests

At the iron ore mine the following tests were performed:

- screen analyses of material
- chemical analyses of ore for grade

At a pilot plant in Pretoria, impact work index tests were done.

## 4 Data gathered

Data gathered for this work were from records at the iron ore beneficiation plant.

Information on Nordberg crushers was obtained directly from Nordberg.

Information on Microsim was obtained from the Metallurgy department of the University of the Witwatersrand.

APPENDIX 2: CORRELATIONS FOR THE KARRA MODELS

The correlations for factors A through F for the Karra model (27), and A through H for the enhanced model are calculated as follows:

Once the following have been determined:

- Q=percentage oversize in feed
- R=percentage half size in feed
- T=theoretical undersize flow rate in tph
- U=bulk density in  $\text{kg/m}^3$

and the following parameters are known for one or two decks:

- screen area in  $\text{m}^2$ , S
- angle of inclination, theta
- square mesh size in mm, w
- length of aperture in mm, l
- wire (or poly) thickness in mm, d
- poly or wire deck

for one deck at a time, the effective aperture, h, in mm, and the factors A through F may be calculated from the empirical correlations of Karra and Nordberg (27,38):

$$h=(w+d)\cos(\text{theta}) - d$$

factor A:

$$\begin{aligned} \text{if } h < 50,8 \quad A &= 12,1286h^{0,3162} - 10,2991 \\ \text{else} \quad A &= 0,3388h + 14,4122 \end{aligned}$$

factor B:

$$\begin{aligned} \text{if } Q > 87, \quad B &= -0,012Q + 1,6 \\ \text{else} \quad B &= 0,0425Q + 4,275 \end{aligned}$$

factor C:

$$\begin{aligned} \text{if } R \leq 30, \quad C &= 0,012R + 0,7 \\ \text{if } 30 < R < 55, \quad C &= 0,1528R^{0,564} \\ \text{if } 55 \leq R < 80, \quad C &= 0,0061R^{1,37} \\ \text{if } R \geq 80, \quad C &= 0,05R - 1,5 \end{aligned}$$

factor D:

if top deck,  $D=1,0$   
if second deck,  $D=0,9$

factor E:

let  $t=1,26h$   
if  $t < 1$ ,  $E=1,0$   
if  $1 \leq t < 2$ ,  $E=t$   
if  $2 < t < 4$ ,  $E=1,5+0,25t$   
if  $4 \leq t < 6$ ,  $E=2,5$   
if  $6 < t \leq 10$ ,  $E=3,25-0,125t$   
if  $10 < t < 12$ ,  $E=4,5-0,25t$   
if  $12 \leq t \leq 16$ ,  $E=2,1-0,05t$   
if  $16 < t < 24$ ,  $E=1,5-0,125t$   
if  $24 \leq t \leq 32$ ,  $E=1,35-0,00625t$   
if  $t > 32$ ,  $E=1,15$

factor F:

$F=U/1602$

factor G (determined from poly decks used):

if wire deck,  $G=1,0$   
if poly deck,  $G=1/1,15$

factor H (given in Kelly and Spottiswood (29):

if  $1/w < 2$ ,  $H=1,0$   
if  $2 \leq 1/w < 3$ ,  $H=1,1$   
if  $3 \leq 1/w < 6$ ,  $H=1,4$   
if  $1/w \geq 6$ ,  $H=1,6$

## APPENDIX 3: PASCAL LISTINGS OF MODELS

The Pascal listings for new and modified Microsim models are given below. Models were written in accordance with the Microsim data structure as given in the Technical Reference Manual (50).

The crushing models CRUSH5 and CRUSH6\_7 are contained in the Microsim library file UNIT1.PAS CRUSH6\_7 replaces the general Whiten model CRUSH2\_3, of which it is a modification for haematite, whereas CRUSH5 is a new model for gyradisc crushing of haematite.

The screening models ROSE1, ROSE2 and SCREEN5 were added to the Microsim library file UNIT2.PAS. SCREEN5 is an enhancement of the original Karra model, SCREEN4. ROSE1 and ROSE2 are both new models.

## A3.1 Gyradisc model CRUSH5

```

-----
{ GYRADISC MODEL - Sishen test model
}
-----
OVERLAY PROCEDURE CRUSH5( var stm          : stms ;
                          point          : tounit;
                          type_num      : UNITYPES);

VAR
feedpt,prodpt,der2pt          : tor4;
feedh2o,prodh2o              : tor6;

g,nd,y1,y,ws,ws1,ak,an      : real ;

i,j,k,l,m                    : integer ;

BEGIN
  feedpt := solidselect(stm[point^.inout[2]]);
  prodpt := solidselect(stm[point^.inout[4]]);

  feedh2o := lixivselect(stm[point^.inout[2]]);
  prodh2o := lixivselect(stm[point^.inout[4]]);
  MAKESCRATCH(der2pt);

  prodh2o^.flow := feedh2o^.flow ;
  prodpt^.tonnes := 0.0 ;

  g := point^.param[1];

  WITH system DO
  begin
    FOR j:=1 TO ngc DO
      FOR k:=1 TO nsc DO
        FOR l:=1 TO n4c DO
          FOR m:=1 TO n5c DO
            SOLIDPUT(1,j,k,l,m,der2pt,0.0);

          FOR i:=1 TO ndc DO
            FOR j:=1 TO ngc DO
              FOR k:=1 TO nsc DO
                FOR l:=1 TO n4c DO
                  FOR m:=1 TO n5c DO
                    SOLIDPUT(1,j,k,l,m,der2pt,SOLIDVALUE(1,j,k,l,m,feedpt)+
                              SOLIDVALUE(1,j,k,l,m,der2pt));

```

```

nd := ndc -1 ;                y1 := 1.0 ;
FOR i := 1 to ndc DO
BEGIN (* for i *)
ws := 0.0 ;
IF i < ndc THEN ws := sqrt(size[i]*size[i+1])/g;

IF ws >= 0.15 THEN
begin
an := 2.77 ;
ak := 0.38 ;
end
ELSE
begin
an := 0.32 ;
ak := 250.0 ;
end;

ws1 := POW((ws/ak),an);
y := 1.0 ;
IF ws1 < 10 THEN y := 1.0 - EXP(-ws1);

FOR j := 1 TO ngc DO
FOR k := 1 to nsc DO
FOR l := 1 to n4c DO
FOR m := 1 to n5c DO
SOLIDPUT(i,j,k,l,m,prodpt,(y1-y)*SOLIDVALUE(i,j,k,l,m,der2pt));
y1:=y;
END; {For l}

FOR i:=1 TO ndc DO
FOR j:=1 TO ngc DO
FOR k:=1 TO nsc DO
FOR l:=1 TO n4c DO
FOR m:=1 TO n5c DO
prodpt^.tonnes := prodpt^.tonnes + SOLIDVALUE(i,j,k,l,m,prodpt) ;

KILLSCRATCH(der2pt);
END ;(* with system *)
END; (* crush5 *)

```

---

## A3.2 Whiten model CRUSH6\_7

```

-----}
{ CONE CRUSHER MODELS 6 and 7 : only for haemetite }
{ Standard Cone crusher model : Based on Whitens model }
{ Short Head crusher model : Based on Whitens model }
-----}
OVERLAY PROCEDURE CRUSH6_7( var stm : stms ;
                           point : tounit ;
                           type_num : UNITYPES);

VAR
  feedpt, prodpt, out2pt : tor4;
  feedh2o, prodh2o : tor6;
  c, ws, wsl, w : REAL;
  b, b1, bii : REAL;
  i, j, k, l, m, ll, ll1 : integer ;

PROCEDURE WHITEN(css, kk, k1, k2, k3, n, mm : REAL);
      {Common calculation part}
BEGIN
  WITH system DO
  BEGIN
    FOR i:=1 TO ndc DO
    BEGIN
      c:=1;
      IF (size[i]<k2) THEN
        c:=1-POW((size[i]-k2)/(k1-k2),k3);
      IF (size[i]<k1) THEN
        c:=0.0;
      wsl:=0.0; ws:=size[i];
      IF (i<ndc) THEN
        wsl:=SQRT(size[i]*size[i+1]);
      IF (i>1) THEN
        ws:=SQRT(size[i-1]*size[i]);

      FOR j:=1 TO ngc DO
        FOR k:=1 TO nsc DO
          FOR l:=1 TO n4c DO
            FOR m:=1 TO n5c DO
              BEGIN
                w:=SOLIDVALUE(i, j, k, l, m, feedpt); ll:=i-1;
                IF (ll<>0) THEN
                  BEGIN
                    FOR ll:=1 TO ll DO
                      BEGIN
                        b1:=(1-kk)*POW(wsl/size[ll],n)+kk*POW(wsl/size[ll],mm);
                        b:=(1-kk)*POW(ws/size[ll],n)+kk*POW(ws/size[ll],mm)-b1;
                        w:=w+b*SOLIDVALUE(ll, j, k, l, m, out2pt);
                      END; {For ll}
                    END; {If ll}
                    bii:=1-(1-kk)*POW(wsl/size[i],n)-kk*POW(wsl/size[i],mm);
                    w:=w/(1-bii*c); SOLIDPUT(i, j, k, l, m, out2pt, w*c);
                    SOLIDPUT(i, j, k, l, m, prodpt, w*(1-c));
                    prodpt^.tonnes := prodpt^.tonnes+SOLIDVALUE(i, j, k, l, m, prodpt) ;
                  END; {j, k, l, m}
                END; {For l}
              END; {with system}
            END; {Calculate procedure}
          -----}
        VAR
          css, kk, k1, k2, k3, n, mm : REAL;

        BEGIN
          feedpt := solidselect(stm[point^.inout[2]]);
          prodpt := solidselect(stm[point^.inout[4]]);
          feedh2o := lixivselect(stm[point^.inout[2]]);
          prodh2o := lixivselect(stm[point^.inout[4]]);

```

```

12.3 Single deck Ross efficiency model (ROSE)
MAKESCRATCH(out2pt);

prodh2o^.flow := feedh2o^.flow ;
prodpt^.tonnes := 0.0 ;

css:=point^.param[1];          kk:=point^.param[2];

CASE point^.MODEL OF
  2 : BEGIN
      k1:=0.653*css;      k2:=1.21*css;      k3:=2.0;
      n:=2.0;      mm:=0.535;
      END; {Standard Symons cone crusher for haemetite}
  3 : BEGIN
      k1:=0.944*css;      k2:=1.722*css+0.004826;      k3:=3.0;
      n:=2.000;      mm:=0.535;
      END; {Short head cone crusher for haemetite}
END;
WHITEN(css,kk,k1,k2,k3,n,mm);
KILLSCRATCH(out2pt);
END; {Procedure}
-----

```

## A3.3 Single deck Rose efficiency model ROSE1

```

-----}
{ SINGLE DECK ROSE MODEL }
-----}
OVERLAY PROCEDURE ROSE1 ( var stm          : stms   ;
                          point          : tounit  ;
                          type_num      : UNITYPES);

VAR
feedpt,outlpt,out2pt          : tor4;
feedh2o,outlh2o,out2h2o      : tor6;
eff,value,value2,d50,x,pof   : real  ;
i,j,k,l,m                    : integer ;

BEGIN
feedpt := solidselect(stm[point^.inout[2]]);
outlpt := solidselect(stm[point^.inout[4]]);
out2pt := solidselect(stm[point^.inout[5]]);
feedh2o := 1ixivselect(stm[point^.inout[2]]);
outlh2o := 1ixivselect(stm[point^.inout[4]]);
out2h2o := 1ixivselect(stm[point^.inout[5]]);
outlpt^.tonnes := 0.0 ;
out2pt^.tonnes := 0.0 ;
eff           := point^.param[2];
d50          := eff*point^.param[1];

WITH system DO
BEGIN
FOR i := 1 to ndc DO
  BEGIN
  x:=POW(size[i]/d50,5.846);
  IF (x<60) THEN pof:=1-EXP(-0.693*x)
  ELSE pof:=1;
  IF pof>1 THEN pof:=1
  ELSE IF pof<0 THEN pof:=0;
FOR j := 1 TO ngc DO
FOR k := 1 to nsc DO
FOR l := 1 to n4c DO
FOR m := 1 to n5c DO
  BEGIN
  value := 0.0 ;
  IF size[i] <= point^.param[1] THEN
  value := (1-pof)* solidvalue(i,j,k,l,m,feedpt);
  value2 := solidvalue(i,j,k,l,m,feedpt) - value;

  SOLIDPUT(i,j,k,l,m,outlpt,value);
  SOLIDPUT(i,j,k,l,m,out2pt,value2);

  outlpt^.tonnes := outlpt^.tonnes + value ;

  out2pt^.tonnes := out2pt^.tonnes + value2;
  END;
  END; (* for *)
END; (* with system *)
outlh2o^.flow := 0.95*feedh2o^.flow;
out2h2o^.flow := 0.05*feedh2o^.flow;
END; (* rose1 *)
-----}

```

## A3.4 Double deck Rose efficiency model ROSE2

```

-----}
{ DOUBLE DECK ROSE MODEL
}
-----}
OVERLAY PROCEDURE ROSE2 ( var stm          : stms   ;
                          point          : tounit  ;
                          type_num      : UNITYPES);

VAR
feedpt,out1pt,out2pt,out3pt      : tor4;
feedh2o,out1h2o,out2h2o,out3h2o : tor6;

eff1,eff2,d501,d502,x1,x2,pof1,pof2 : real  ;
value,value1,value2 ,value3         : real  ;

i,j,k,l,m                          : integer ;

BEGIN

feedpt := solidselect(stm[point^.inout[2]]) ;
out1pt := solidselect(stm[point^.inout[4]]) ;
out2pt := solidselect(stm[point^.inout[5]]) ;
out3pt := solidselect(stm[point^.inout[6]]) ;

feedh2o := lixivselect(stm[point^.inout[2]]) ;
out1h2o := lixivselect(stm[point^.inout[4]]) ;
out2h2o := lixivselect(stm[point^.inout[5]]) ;
out3h2o := lixivselect(stm[point^.inout[6]]) ;

out1pt^.tonnes := 0.0 ;
out2pt^.tonnes := 0.0 ;
out3pt^.tonnes := 0.0 ;

eff1          := point^.param[2];
eff2          := point^.param[4];
d501:=eff1*point^.param[1];
d502:=eff2*point^.param[3];

WITH system DO begin

FOR i := 1 to ndc DO
BEGIN
x1:=POW(size[i]/d501,5.846);
x2:=POW(size[i]/d502,5.846);
IF (x1<60) THEN pof1:=1-EXP(-0.693*x1)
ELSE pof1:=1;
IF pof1>1 THEN pof1:=1
ELSE IF pof1<0 THEN pof1:=0;
IF (x2<60) THEN pof2:=1-EXP(-0.693*x2)
ELSE pof2:=1;
IF pof2>1 THEN pof2:=1
ELSE IF pof2<0 THEN pof2:=0;
FOR j := 1 TO ngc DO
FOR k := 1 to nsc DO
FOR l := 1 to n4c DO
FOR m := 1 to n5c DO
BEGIN
value := solidvalue(i,j,k,l,m,feedpt) ;
IF size[i] < point^.param[1] THEN
value2 := (pof1) * value ;
IF size[i] >= point^.param[1] THEN
value2 := value;
IF size[i] < point^.param[3] THEN
value3 := (pof2)*(value-value2);
IF size[i] >= point^.param[3] THEN
value3 := (value-value2);

value1 := value-value2-value3;

```

```

solidput(i,j,k,l,m,out1pt,value1);
solidput(i,j,k,l,m,out2pt,value2);
solidput(i,j,k,l,m,out3pt,value3);

out1pt^.tonnes := out1pt^.tonnes + value1;
out2pt^.tonnes := out2pt^.tonnes + value2;
out3pt^.tonnes := out3pt^.tonnes + value3;
END;
END; (* for *)

```

```

out1h2o^.flow := feedh2o^.flow * 0.9 ;
out2h2o^.flow := feedh2o^.flow * 0.05;
out3h2o^.flow := feedh2o^.flow * 0.05;
END; (* with system *)
END; (* rose2 *)

```

## A3.5 Enhanced Karra Model SCREEN5

```

-----}
{ SCREEN MODEL 5
{ Enhanced Karra screen model for single or double screens
}
-----}
OVERLAY PROCEDURE SCREEN5(VAR stm : STMS;
                          point : TOUNIT;
                          type_num : UNITYPES);

PROCEDURE PARTICLE_SIZE(stream_ptr : TOR4;
                        VAR cumulative : SIZES);

VAR
  cumulats : REAL;
  i : INTEGER;

BEGIN
  cumulats:=0;
  FOR i:=1 TO system.NDC DO
  BEGIN
    cumulats:=cumulats+SIGNDC(i,stream_ptr);
    cumulative[i]:=100-cumulats;
  END;
END;
-----}
FUNCTION INTERPOLATE( x,y : SIZES;
                     n : INTEGER;
                     xx : REAL) : REAL;

VAR
  c : ARRAY[sizrange,1..3] OF REAL;
  i,n0,k : INTEGER;
  rm3,t1,rm2,rm1,rm4,t2,b : REAL;

BEGIN
  rm3:=(y[2]-y[1])/(x[2]-x[1]);          t1:=-rm3-(y[2]-y[3])/(x[2]-x[3]);
  rm2:=rm3+t1;                          rm1:=rm2+t1;
  n0:=n-2;

  FOR i:=1 TO N DO BEGIN
    IF (i>n0) THEN
      rm4:=rm3-rm2+rm3
    ELSE
      rm4:=(y[i+2]-y[i+1])/(x[i+2]-x[i+1]);
      t1:=ABS(rm4-rm3);                  t2:=ABS(rm2-rm1);
      b:=-t1+t2;
      IF (b<>0) THEN
        c[i,1]:=(t1*rm2+t2*rm3)/b
      ELSE
        c[i,1]:=0.5*(rm2+rm3);
      rm1:=rm2;    rm2:=rm3;    rm3:=rm4
    END;

    n0:=n-1;

    FOR i:=1 TO n0 DO BEGIN
      t1:=1/(x[i+1]-x[i]);                t2:=(y[i+1]-y[i])*t1;
      b:=(c[i,1]+c[i+1,1]-t2-t2)*t1;    c[i,3]:=b*t1;
      c[i,2]:=-b+(t2-c[i,1])*t1
    END;

    FOR i:=1 TO n-1 DO
      IF ((xx>=x[i]) AND (xx<x[i+1])) THEN k:=i;
      xx:=xx-x[k];
      interpolate:=c[k,3]*xx*xx*xx+c[k,2]*xx*xx+c[k,1]*xx+y[k]
    END;
  END;

```

```

-----}
FUNCTION PERCENTAGE_PASSING(cumulative : SIZES;           {Find % passing psize}
                             psize      : REAL) : REAL;   {psize in microns  }
VAR
  x,y      : SIZES;
  i        : INTEGER;
  dummy    : REAL;

BEGIN
  psize:=psize/1E6;           {Convert to metres}
  WITH system DO
  BEGIN
    FOR i:=1 TO ndc DO
    BEGIN
      x[i]:=size_class[ndc-i+1];
      y[i]:=cumulative[ndc-i+1];
    END;
    IF psize>size_class[1] THEN
    BEGIN
      dummy:=psize/size_class[1]*cumulative[1];
      IF dummy>100 THEN dummy:=100;
    END
    ELSE
      dummy:=INTERPOLATE(x,y,ndc,psize);
    END;
  percentage_passing:=dummy;
END;
-----}
PROCEDURE MODCALC_ONE_DECK(feedpt,ofpt,ufpt : TOR4;
                            lw,ht,area,poly : REAL;           {Ht in mm}
                            deck : INTEGER;
                            water : REAL);
VAR
  i,j,k,l,m      : INTEGER;
  a,b,c,d,e,f,kk,xn,p,q,r,t,g,h : REAL;
  undersize_theory,svm,sga,u,pof,d50,x : REAL;
  cumulative      : SIZES;

BEGIN
  PARTICLE_SIZE(feedpt,cumulative);           {Get cumulative in feed}
  q:=100-PERCENTAGE_PASSING(cumulative,ht*1000);  {% oversize in feed  }
  r:=PERCENTAGE_PASSING(cumulative,ht*500);      {% half size in feed  }
  xn:=PERCENTAGE_PASSING(cumulative,1250*ht)-
      PERCENTAGE_PASSING(cumulative,750*ht);     {% near size in feed  }
  undersize_theory:=(1-q/100)*feedpt^.TONNES*3.6; {Theoretical underflow TPH}
  p:=ht;
  t:=1.26*ht;

      {Get constants A,B,C,D,E,F,G,H}
  IF (p>=50.8) THEN a:=-14.4122+0.3388*p
  ELSE a:=-12.1286*POW(p,0.3162)-10.2991;

  IF (q>87) THEN b:=4.275+0.0425*q
  ELSE b:=1.6-0.012*q;

  IF (r>=80) THEN c:=0.05*r-1.5
  ELSE
    IF ((r<80) AND (r>=55)) THEN c:=0.0061*POW(r,1.37)
    ELSE
      IF ((r<55) AND (r>=30)) THEN c:=0.1528*POW(r,0.564)
      ELSE c:=0.012*r+0.7;

  d:=1.1-0.1*deck;
  {g = polydeck factor for reduced area}
  {h = factor for nonsquare apertures}
  IF poly=1.0 THEN g:=1./1.15
  ELSE g:=1.0;
  IF lw<2.0 THEN h:=1.0
  ELSE
    IF lw<3.0 THEN h:=1.1

```

```

ELSE
  IF 1w<6 THEN h:=1.4
  ELSE h:=1.6;
IF (water>0) THEN      {Wet screening}
BEGIN
  IF (t>32) THEN e:=-1.15
  ELSE
    IF ((t<=-32) AND (t>=24)) THEN e:=1.35-0.00625*t
    ELSE
      IF ((t<24) AND (t>16)) THEN e:=1.5-0.0125*t
      ELSE
        IF ((t<=16) AND (t>=12)) THEN e:=2.1-0.05*t
        ELSE
          IF ((t<12) AND (t>10)) THEN e:=4.5-0.25*t
          ELSE
            IF ((t<=10) AND (t>6)) THEN e:=3.25-0.125*t
            ELSE
              IF ((t<=6) AND (t>=4)) THEN e:=2.5
              ELSE
                IF ((t<4) AND (t>2)) THEN e:=1.5+0.25*t
                ELSE
                  IF ((t<=2) AND (t<=-1)) THEN e:=t
                  ELSE
                    e:=1.0;

    sga:=-Ore_Density(feedpt);
    if sga > 0.0 then svm:=1/(1000*sga)
      else svm:=0;
    u:=-sga*0.6*1000;
    f:=u/1602;
  END
ELSE
  BEGIN                                  {Dry screening}
    e:=1;                                f:=1;
  END;

kk:=undersize_theory/area/(a*b*c*d*e*f*g*h);

d50:=0.975*ht*POW(kk,-0.148)*POW((1-xn/100),0.511);

d50:=d50*1e-03;                          {050 in meters}

ufpt^.TONNES:=0.0;      ofpt^.TONNES:=0.0;
WITH system DO
BEGIN
  FOR i:=1 TO ndc DO
  BEGIN
    x:=POW(size[i]/d50,5.846);
    IF (x<60) THEN pof:=1-EXP(-0.693*x)
    ELSE pof:=1;
    IF pof>1 THEN pof:=1
    ELSE IF pof<0 THEN pof:=0;
    FOR j:=1 TO ngc DO
    FOR k:=1 TO nsc DO
    FOR l:=1 TO n4c DO
    FOR m:=1 TO n5c DO
    BEGIN
      SOLIDPUT(i,j,k,l,m,ofpt,pof*SOLIDVALUE(i,j,k,l,m,feedpt));
      SOLIDPUT(i,j,k,l,m,ufpt,SOLIDVALUE(i,j,k,l,m,feedpt)
        -SOLIDVALUE(i,j,k,l,m,ofpt));
      ufpt^.TONNES:=ufpt^.TONNES+SOLIDVALUE(i,j,k,l,m,ufpt);
      ofpt^.TONNES:=ofpt^.TONNES+SOLIDVALUE(i,j,k,l,m,ofpt);
    END;
  END;
END;
END;
{-----}
VAR
feedpt,out1pt,out2pt,out3pt,dummyfeedpt : TOR4;
feedh2o,out1h2o,out2h2o,out3h2o       : TOR6;

```

```

i, j, k, l, m                                : INTEGER;
area, ht, theta, water, lw, poly             : REAL;
deck                                          : INTEGER;

BEGIN
feedpt := solidselect(stm[point^.inout[2]]) ;
out1pt := solidselect(stm[point^.inout[4]]) ;
out2pt := solidselect(stm[point^.inout[5]]) ;

IF (type_num=19) THEN
BEGIN
out3pt := solidselect(stm[point^.inout[6]]) ;
out3h2o := lixivselect(stm[point^.inout[6]]) ;
END;

feedh2o := lixivselect(stm[point^.inout[2]]) ;
out1h2o := lixivselect(stm[point^.inout[4]]) ;
out2h2o := lixivselect(stm[point^.inout[5]]) ;

MAKESCRATCH(dummyfeedpt);

area:=point^.param[1];          theta:=point^.param[2]/180*3.1416;
WITH point^ DO
BEGIN
ht:=(param[3]+param[5])*COS(theta)-param[5];
lw:=param[4]/param[3];
poly:=param[6];
END;
water:=feedh2o^.FLOW*3.6;

MODCALC_ONE_DECK(feedpt, out2pt, dummyfeedpt, lw, ht, area, poly, 1, water);

WITH system DO
CASE type_num OF
15 : BEGIN
FOR i:=-1 TO ndc DO
FOR j:=-1 TO ngc DO
FOR k:=-1 TO nsc DO
FOR l:=-1 TO n4c DO
FOR m:=-1 TO n5c DO
SOLIDPUT(i, j, k, l, m, out1pt, SOLIDVALUE(i, j, k, l, m, dummyfeedpt));
out1pt^.TONNES:=dummyfeedpt^.TONNES;
out1h2o^.FLOW:=0.95*feedh2o^.FLOW;
out2h2o^.FLOW:=0.05*feedh2o^.FLOW;
END;
19 : BEGIN

WITH point^ DO
BEGIN
ht:=(param[7]+param[9])*COS(theta)-param[9];
lw:=param[8]/param[7];
poly:=param[10];
END;
MODCALC_ONE_DECK(dummyfeedpt, out3pt, out1pt, lw, ht, area, poly, 2, water);
out1h2o^.flow := feedh2o^.flow * 0.9 ;
out2h2o^.flow := feedh2o^.flow * 0.05;
out3h2o^.flow := feedh2o^.flow * 0.05;
END;

END;

KILLSCRATCH(dummyfeedpt);
END;
-----

```