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GRAVPOLY: A modification of a three-dimensional gravity
modelling program

) by

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Abstract

The three-dimensional gravity modelling program of Donald Plouff (Plouff, 1975) has been made operational on a Honeywell Multics 68/80 computer. Modifications were made to input and output procedures to make the program compatible with U.S. Geological Survey (USGS) standard geophysical input and output formats.

Introduction

This program calculates the gravity effect of 3D models that consist of elements that form polygons in plan view (Plouff, 1976). The elements of models are bounded by two horizontal-planes and a series of intersecting vertical planes. The calculations use an exact formula rather than using a numerical integration approach of previous Talwani (1960) programs. Thus, values can be obtained on or inside the model. A least-squares comparison between the calculated anomalies and an observed field can be used to determine the best density contrast.

Description

Model input to the program consists of x, y, and z coordinates of polygon corners. A maximum of 50 polygons can be specified and the total number of corners for all polygons cannot exceed 900. Parameter input such as plotting variables, file names, and so forth are entered from a command file in namelist format. Fieldpoint locations and optionally observed values at these locations are read in from a disk file. This file can be user-formatted random data (see Appendix B) or it can be in USGS geophysical grid format (see Appendix A). If the height of the fieldpoint locations are not on a level plane, a second USGS grid-format file can be read from a disk that will contain the elevations of the fieldpoint locations.

Output from the program consists of printed information about the bodies

and fieldpoint coordinates and values. Contoured output may be directed to a line printer or terminal, or grid-format files can be created that can be used as direct input to a contouring program.

Program Usage

Program variables and body coordinate data are obtained from an ASCII segment on disk created by the user prior to running the program. This segment is read by Fortran unit 9, and attachment and detachment of the segment are accomplished internally in the program. The program variables are read from the first part of the segment in namelist format. The segment therefore must start with the characters "&parms" followed by any number of namelist variables explained below and followed with the character "&". After the above character, the body coordinate information explained below is read from the segment.

Fieldpoint information in either a user-specified format or in a USGS geophysics grid format is read from a separate segment created on disk by the user before running the program. The maximum number of fieldpoints allowed is 1,000. This segment is read by Fortran unit 10 and is attached and detached internally in the program. With user-formatted input, the height of each fieldpoint location is specified; this is not the case with grid files. In order to allow this flexibility with grid files, a second file that contains the heights of the fieldpoint locations may be read in from disk. This file must have the same header information as the location file, that is, it must be registered. This optional segment is read by Fortran unit 11 and is attached and detached by the program.

Following are explanations of the namelist variables, grouped under usage subheadings.

File information

ifile - the name of the file that contains the fieldpoint information.

It must be enclosed in quotes with a maximum length of 50 characters.

ifile2 - the name of a grid file that contains the heights of the fieldpoint locations. It must be enclosed in quotes with a maximum length of 50 characters. This file is not necessary if the parameter "height" below is used.

name - a character variable that states the type of data that will be read from "ifile". Default is the characters "gridded", which means that the fieldpoint data will be read as a standard USGS geophysics grid file. Any other characters (maximum of 8) enclosed in quotes will assumed a user-formatted file.

height - the height of the fieldpoint grid above the same datum that is used to reference the body heights, for example, sea level. Positive values are upward. This parameter is used only when name = "gridded" and ifile2 is not specified. Default is zero.

ifmt - the format to be used when reading a user-formatted file for fieldpoint data. Four variables are read for each fieldpoint location in the following order: x coordinate, y coordinate, observed value, and height. The parameter must be enclosed in quotes, for example, "(4F10.2)". A maximum of 80 characters can be used. Default is "(v)" for list directed input.

Plotting information and scaling

- I. iplotr - a number that determines which plotter device is to be used. The number 9, which is the default, directs all plotting to an output file on disk, which is named with the characters "print." as a prefix to

the command file name. One of the numbers listed below will direct the plot of the body corners to that device.

0 = Calcomp 7900

1 = Tektronix 4010

2 = Hewlett-Packard 7202

3 = Hewlett-Packard 7203

4 = Tektronix 4014 low resolution

5 = Tektronix 4014 high resolution

6 = General vector output, for example, electrostatic printer/plotter

ibody - a nonzero number will create a plot of body corners and fieldpoint locations. Default is zero.

iobs - a nonzero number and `iplotr = 9` will create printer contours for gridded input fieldpoint data. Default is zero.

icalc - a nonzero number and `iplotr = 9` will create printer contours of the calculated data. A nonzero number will create a USGS geophysical grid file with the characters "calc." as a prefix to the command file name, which can be contoured later by a contouring program. Default is zero. (This parameter is applicable only with gridded input fieldpoint data).

ires - a nonzero number and `iplotr = 9` will create printer contours of the residual data values (observed-calculated). A nonzero number will create a standard USGS geophysics grid file with the characters "res." as a prefix to the command file name, which can be contoured later by a contouring program. Default is zero. (This parameter is applicable only with gridded input fieldpoint data).

idplot - 40 characters of identification that is printed at the top of printed output. This parameter must be enclosed in quotes and the

default is blanks.

dc - the contour interval to be used if printer plots are requested.

Default is zero.

xscale - the scaling factor to be used only when plotting body

coordinates on a plotter device other than a printer. It is specified in coordinate units per inch, for example, 62500, if inches are the units of measurement (see "unix" below). Default is zero.

unix - the units of measurement in the x and y directions. It is a four character variable which must be enclosed in quotes. Options available are "feet", "mile", "kilf", "kilm" and "metr". Default is "kilm".

uniz - the units of measurement in the z direction. It has the same options as "unix" above. Default is "feet".

maxcol - a number that controls the horizontal width of printer plots. If later printing is desired on a data terminal then 79 should be used. Default is 130 for plotting on a line printer.

Calculation information

lsqs - a number that determines if fieldpoint (observed) values will be read or not or whether a least-squares comparison will be made between the calculated and observed values.

0 = no fieldpoint (observed) values will be used. It must be kept in mind that it is always necessary to read fieldpoint locations in order to define the grid points at which the model values will be calculated

1 = fieldpoint values will be read and used to calculate residual values (observed-calculated).

7 = a least-squares comparison between the observed and calculated

values will be made to determine the best density contrast datum - a number to be added to the calculated anomaly to convert to the observed anomaly values. This is nearly always zero, which is the default.

Body information

Following the namelist variables in the command segment is parameter information for each body of the model. A total of 50 bodies is permitted and each body has the following format (see Examples of Program Execution):

Five parameter values precede the body coordinates. The values are read from one line in v format. The order of the values are as follows. 1) The number of corners of the body. 2) A print switch for body information. A zero means no additional printout. A positive number will print the contribution of this body at each field point. A negative number, used for debugging, will print the values of key parameters for each polygonal side of the body (integrals V_1-V_6). 3) The density contrast in g/cm^3 . A positive contrast will produce positive anomalies. A hole in a body can be produced by using a density contrast of opposite sign from that of the parent body. 4) The height of the top of body. The positive direction is upward. 5) The height of the bottom of the body and again the positive direction is upward.

The body coordinate corners follow the above parameter line in x and y pairs. The progression of corners is clockwise as viewed from above. The first pair is not repeated at the end of the list and the total number of pairs must equal the first number of the preceding parameter line. A hole may be removed from a large body by treating the hole as a separate body with the same density contrast as the parent body but with input progressing counterclockwise from corner to corner.

Program Execution

To run the program type the characters: >online>RegPgms>gravpoly. The program will then type the statement: "enter command file name:". The user then types in the name of the command segment that he has previously created on disk and program execution continues. If no error messages are generated and the program terminates normally, a message is printed that states that the output is in a file that is named with the characters "print." as a prefix to the command file name. The user then can observe the results of the program by printing this file, which resides in the user's working directory on disk.

If a plot of body corners and fieldpoint locations is requested on any of the offline plotting devices, then the program will be interrupted by the plotting system to ask the user questions.

References

- Evenden, G. J., and Wahl, R. P., 1977, Basic plotting system: U.S. Geological Survey unpublished documentation.
- Plouff, Donald, 1975, Derivation of formulas and Fortran programs to compute gravity anomalies of prisms: U.S. Geological Survey Report, 90 p. available from National Technical Information Service No. PB-243-526, U.S. Department of Commerce, Springfield, VA 22161.
- Plouff, Donald, 1976, Gravity and magnetic fields of polygonal prisms and application to magnetic terrain corrections: *Geophysics*, v. 41, no. 4, p. 727-741.
- Talwani, Manik, and Ewing, Maurice, 1960, Rapid computation of gravitational attraction of three-dimensional bodies of arbitrary shape: *Geophysics*, v. 25, no. 1, p. 203-225.

Command Segment Examples

1. The following example will calculate a best density contrast between the observed and calculated values.

```

&parms uniz="kilm",iobs=1,ifile="gravpoly.test1",dc=5,
name="formatted",idplot="gravpoly test 1",icalc=1,ires=1,
naxcol=40,ibody=1,lsas=7,3
4 0 1 0 -1
5.657 0 2.828 2.828 5.657 5.657 8.485 2.828
4 0 1 -1 -2
5.657 0 2.828 2.828 5.657 5.657 8.485 2.828

```

2. This example will perform a forward calculation from the two model bodies specified below. The fieldpoint values are read from one file and the elevations of the fieldpoint locations are read from a second file. Output is placed in a disk file.

```

&parms unix="kilm",uniz="kilm",xscale=5,idplot="hood",
naxcol=90,lsas=1,ibody=1,icalc=1,ires=1,dc=2,
ifile="hdresre",ifile2="hdfc",3
5 0 0.6 2.6 -0.5
-14 44 -14 42 -17 42 -17 44 -14 45
5 0 0.4 -0.5 -3.0
-18 46 -13 47 -12 41 -19 38 -21 40

```

3. This example will perform the same calculations as example 2. Contoured output is not placed in a disk file but two grid files of the calculated and residual values are created.

```

&parms unix="kilm",uniz="kilm",xscale=5,idplot="hood",
iplotr=1,lsas=1,ibody=1,icalc=1,ires=1,dc=2,
ifile="hdresre",ifile2="hdfc",3
5 0 0.6 2.6 -0.5
-14 44 -14 42 -17 42 -17 44 -14 45
5 0 0.4 -0.5 -3.0
-18 48 -13 47 -12 41 -19 38 -21 40

```

Appendix B

Output Example 1

The following output was produced from command example 1.

gravity effect of polyhedral prisms
 for gravpoly test 1
 units of distance measurement are in kiln and heights are in kiln
 body 1 between denths 0.000 and -1.000 and has 4 vertices. density = 1.000
 5.657 0.000 2.828 2.828 5.657 5.657 8.485 2.828
 body 2 between denths -1.000 and -2.000 and has 4 vertices. density = 1.000
 5.657 0.000 2.828 2.828 5.657 5.657 8.485 2.828

observed values		z	q
0.0000	0.0000	0.0000	1.0000
1.4140	0.0000	0.0000	3.0000
2.8280	0.0000	0.0000	5.0000
4.2430	0.0000	0.0000	10.0000
5.6570	0.0000	0.0000	20.0000
0.0000	1.4140	0.0000	2.0000
1.4140	1.4140	0.0000	5.0000
2.8280	1.4140	0.0000	10.0000
4.2430	1.4140	0.0000	15.0000
5.6570	1.4140	0.0000	50.0000
0.0000	2.8280	0.0000	3.0000
1.4140	2.8280	0.0000	5.0000
2.8280	2.8280	0.0000	15.0000
4.2430	2.8280	0.0000	45.0000
5.6570	2.8280	0.0000	50.0000

least-squares density is 0.959 (std 0.062) with correlation coefficient= 0.97 datum= 1.21 mgal.

standard deviation= 4.26 and average difference= 1.21 milligals

	x	y	z	q	residual
1	0.000	0.000	0.000	0.85	0.15
2	1.414	0.000	0.000	1.64	1.36
3	2.828	0.000	0.000	3.42	1.58
4	4.243	0.000	0.000	7.33	2.67
5	5.657	0.000	0.000	15.46	4.54
6	0.000	1.414	0.000	1.10	0.90
7	1.414	1.414	0.000	2.56	2.44
8	2.828	1.414	0.000	7.33	2.67
9	4.243	1.414	0.000	27.00	-12.00
10	5.657	1.414	0.000	42.88	7.12
11	0.000	2.828	0.000	1.22	1.78
12	1.414	2.828	0.000	3.15	1.85
13	2.828	2.828	0.000	13.45	-0.45
14	4.243	2.828	0.000	42.88	2.12
15	5.657	2.828	0.000	43.56	1.44

Output Example 2

The following output was produced from command example 2.

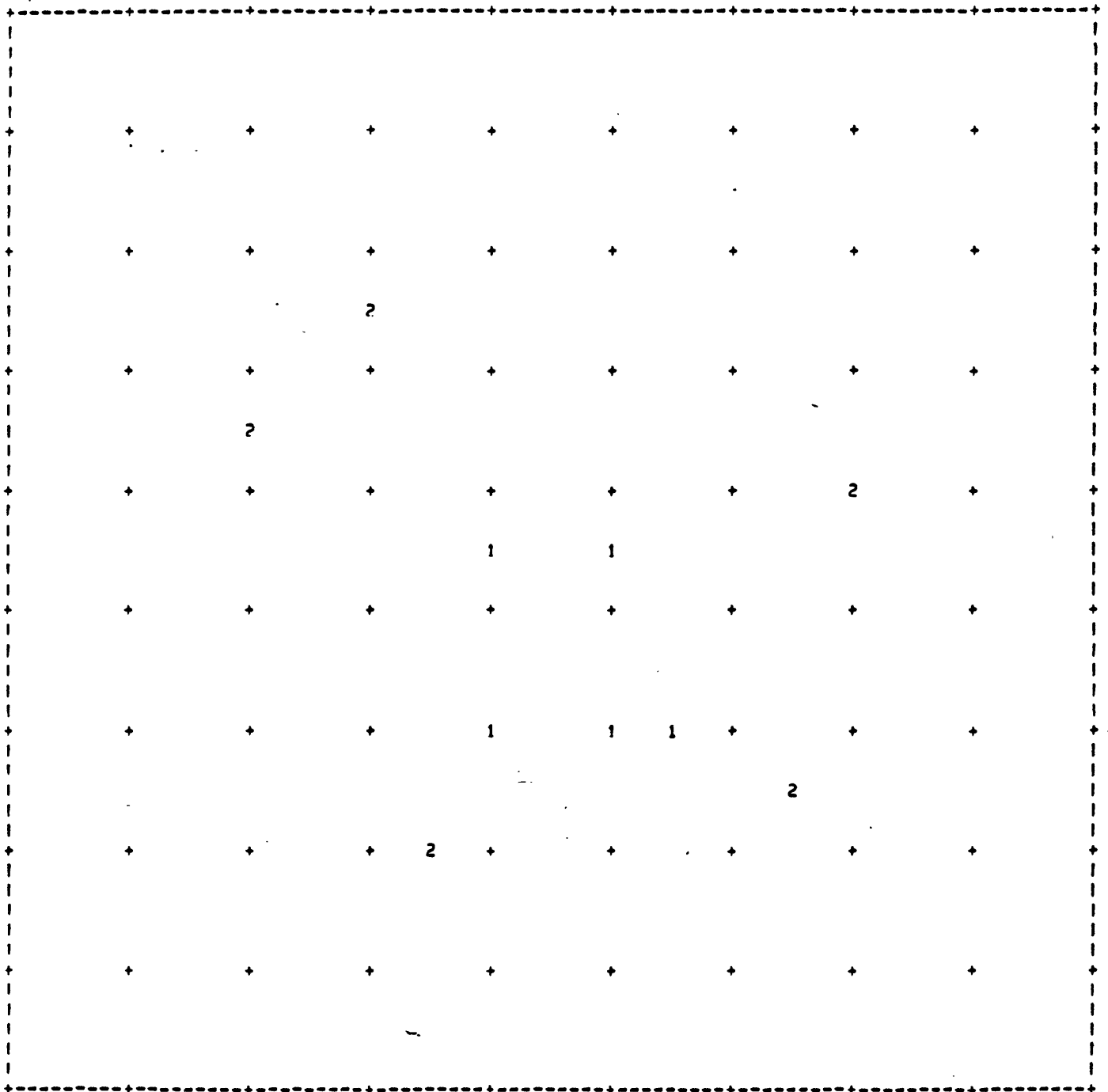
gravity effect of polyhedral prisms for		units of distance measurement are in kilm and heights are in kilm		-0.500 and has 5 vertices. density =		0.600	
body 1	between depths	2.600 and	-14.000	42.000	-17.000	42.000	-17.000
body 2	between depths	-0.500 and	-13.000	47.000	-3.000 and has 5 vertices. density =	41.000	-19.000
x	observed values	z	y	u			
-26.0000	34.0000	0.7032	-4.4752				
-24.0000	34.0000	0.8063	-5.0963				
-22.0000	34.0000	1.0032	-4.7125				
-20.0000	34.0000	1.1685	-5.0729				
-18.0000	34.0000	1.2915	-5.7330				
-16.0000	34.0000	1.4336	-3.7549				
-14.0000	34.0000	1.5534	41.8552				
-12.0000	34.0000	1.2578	0.9125				
-10.0000	34.0000	1.3793	-0.2056				
-8.0000	34.0000	1.2469	-4.1507				
-26.0000	36.0000	1.1045	-6.0303				
-24.0000	36.0000	1.0740	-3.3738				
-22.0000	36.0000	1.0710	-5.1146				
-20.0000	36.0000	1.3257	-6.2990				
-18.0000	36.0000	1.5777	-4.5202				
-16.0000	36.0000	1.6541	-2.3822				
-14.0000	36.0000	1.6054	-4.3734				
-12.0000	36.0000	1.5400	-3.5456				
-10.0000	36.0000	1.3500	-2.6900				
-8.0000	36.0000	1.2354	-1.9731				
-26.0000	38.0000	1.2555	3.8396				
-24.0000	38.0000	1.3862	5.0040				
-22.0000	38.0000	1.3163	0.2649				
-20.0000	38.0000	1.4803	0.7490				
-18.0000	38.0000	1.6247	0.6795				
-16.0000	38.0000	2.0185	-2.5144				
-14.0000	38.0000	2.0043	-2.3035				
-12.0000	38.0000	1.7091	-3.7793				
-10.0000	38.0000	1.5532	-1.9242				
-8.0000	38.0000	1.5604	0.9555				
-26.0000	40.0000	1.1749	6.1622				
-24.0000	40.0000	1.0047	4.4865				
-22.0000	40.0000	1.1950	1.1651				
-20.0000	40.0000	1.5276	3.1051				
-18.0000	40.0000	2.0512	4.0331				
-16.0000	40.0000	2.6200	3.5525				
-14.0000	40.0000	2.3423	2.6886				
-12.0000	40.0000	1.8463	-1.4664				
-10.0000	40.0000	1.7129	-1.0795				
-8.0000	40.0000	1.5006	2.2547				
-26.0000	42.0000	0.7608	1.0344				
-24.0000	42.0000	1.0243	0.8072				
-22.0000	42.0000	1.0529	1.2495				
-20.0000	42.0000	1.4979	2.3965				
-18.0000	42.0000	2.0366	6.0353				
-16.0000	42.0000	2.4138	9.8200				
-14.0000	42.0000	2.5052	7.7781				
-12.0000	42.0000	2.0769	2.6175				

Output Example 2--Continued

-10.0000	42.0000	1.6770	4.4229
-8.0000	42.0000	1.3685	3.0072
-26.0000	44.0000	1.0325	3.1538
-24.0000	44.0000	0.8599	0.3559
-22.0000	44.0000	0.9701	-0.7647
-20.0000	44.0000	1.2288	0.6161
-18.0000	44.0000	1.7800	3.3673
-16.0000	44.0000	2.1822	10.9393
-14.0000	44.0000	1.9485	6.4238
-12.0000	44.0000	1.8721	0.5234
-10.0000	44.0000	1.5460	3.7946
-8.0000	44.0000	1.3610	5.9164
-26.0000	46.0000	0.8376	1.0846
-24.0000	46.0000	1.0251	-1.5601
-22.0000	46.0000	1.1882	-0.4104
-20.0000	46.0000	1.1681	0.3335
-18.0000	46.0000	1.4215	3.0304
-16.0000	46.0000	1.7419	3.1077
-14.0000	46.0000	1.4516	0.2309
-12.0000	46.0000	1.5282	0.0212
-10.0000	46.0000	1.4067	1.6231
-8.0000	46.0000	1.2071	1.2785
-26.0000	48.0000	1.1950	0.2554
-24.0000	48.0000	1.2186	-7.5915
-22.0000	48.0000	1.0251	-7.7261
-20.0000	48.0000	1.1310	-2.5403
-18.0000	48.0000	1.4392	0.7354
-16.0000	48.0000	1.3835	0.4184
-14.0000	48.0000	1.3401	-0.4183
-12.0000	48.0000	1.2443	0.2504
-10.0000	48.0000	1.2734	2.1052
-8.0000	48.0000	1.1459	-0.8801
-26.0000	50.0000	1.1200	3.8535
-24.0000	50.0000	1.2394	1.0059
-22.0000	50.0000	0.7588	-2.2630
-20.0000	50.0000	0.9185	-3.4962
-18.0000	50.0000	1.3691	-1.2682
-16.0000	50.0000	1.1329	-0.6621
-14.0000	50.0000	1.0653	-2.4461
-12.0000	50.0000	0.9946	-2.1838
-10.0000	50.0000	1.0258	-0.5849
-8.0000	50.0000	0.9843	-2.7911
-26.0000	52.0000	1.1602	4.4701
-24.0000	52.0000	0.9325	3.3534
-22.0000	52.0000	0.7125	-1.9564
-20.0000	52.0000	0.9177	-4.2629
-18.0000	52.0000	1.3534	-0.8839
-16.0000	52.0000	1.2964	0.0481
-14.0000	52.0000	1.1620	-1.7044
-12.0000	52.0000	1.0913	-2.8486
-10.0000	52.0000	0.8489	-3.3899
-9.0000	52.0000	0.7865	-4.6056

expect plot of body corners x= -26.0 to 34.0 to 52.0

Output Example 2--Continued



standard deviation= 6.61 and Average difference= -4.85 milligals

Output Example 2—Continued

	x	y	z	q	residual
1	-26.000	34.000	0.793	0.49	-4.97
2	-24.000	34.000	0.806	0.74	-5.84
3	-22.000	34.000	1.093	1.12	-5.83
4	-20.000	34.000	1.169	1.49	-6.56
5	-18.000	34.000	1.291	1.74	-7.47
6	-16.000	34.000	1.434	1.76	-5.51
7	-14.000	34.000	1.553	1.49	-3.35
8	-12.000	34.000	1.258	1.15	-0.24
9	-10.000	34.000	1.379	0.90	-1.11
10	-8.000	34.000	1.247	0.63	-4.78
11	-6.000	36.000	1.104	0.76	-6.79
12	-4.000	36.000	1.074	1.21	-4.54
13	-2.000	36.000	1.071	1.97	-7.06
14	-20.000	36.000	1.326	3.05	-9.35
15	-18.000	36.000	1.578	3.71	-8.23
16	-16.000	36.000	1.654	3.56	-5.94
17	-14.000	36.000	1.605	2.90	-7.26
18	-12.000	36.000	1.540	2.12	-5.66
19	-10.000	36.000	1.550	1.38	-4.07
20	-8.000	36.000	1.235	0.69	-2.86
21	-6.000	38.000	1.255	1.05	2.79
22	-4.000	38.000	1.386	1.91	3.09
23	-2.000	38.000	1.516	3.57	-3.31
24	-20.000	38.000	1.480	6.26	-5.51
25	-18.000	38.000	1.625	7.63	-6.95
26	-16.000	38.000	2.019	7.18	-6.70
27	-14.000	38.000	2.004	5.74	-8.04
28	-12.000	38.000	1.709	3.82	-7.60
29	-10.000	36.000	1.553	2.28	-4.20
30	-8.000	36.000	1.560	1.37	-0.41
31	-6.000	40.000	1.175	1.22	4.94
32	-4.000	40.000	1.005	2.28	2.20
33	-2.000	40.000	1.195	5.23	-4.06
34	-20.000	40.000	1.526	10.03	-6.93
35	-18.000	40.000	2.057	12.23	-8.19
36	-16.000	40.000	2.620	12.23	-8.67
37	-14.000	40.000	2.542	10.54	-7.65
38	-12.000	40.000	1.846	6.57	-8.04
39	-10.000	40.000	1.713	3.48	-4.56
40	-8.000	40.000	1.501	1.78	0.47
41	-6.000	42.000	0.761	1.13	-0.10
42	-4.000	42.000	1.024	2.45	-1.64
43	-2.000	42.000	1.053	5.54	-4.29
44	-20.000	42.000	1.408	11.29	-8.89
45	-18.000	42.000	2.037	16.31	-10.26
46	-16.000	42.000	2.814	20.74	-16.92
47	-14.000	42.000	2.505	21.00	-13.31
48	-12.000	42.000	2.071	9.33	-6.71
49	-10.000	42.000	1.677	4.41	0.01
50	-8.000	42.000	1.569	2.03	0.97
51	-6.000	44.000	1.035	1.18	1.97
52	-4.000	44.000	0.860	2.12	-1.77
53	-2.000	44.000	0.970	4.63	-5.40
54	-20.000	44.000	1.229	10.12	-9.50
55	-18.000	44.000	1.780	16.28	-12.91
56	-16.000	44.000	2.182	28.64	-17.90
57	-14.000	44.000	1.949	21.27	-14.84
58	-12.000	44.000	1.872	9.76	-9.23

Output Example 2--Continued

59	-10.000	44.000	1.547	4.44	-0.64
60	-8.000	44.000	1.361	2.07	3.85
61	-26.000	46.000	0.838	0.95	0.13
62	-24.000	46.000	1.025	1.81	-3.37
63	-22.000	46.000	1.188	3.62	-4.03
64	-20.000	46.000	1.168	7.36	-7.02
65	-18.000	46.000	1.422	12.65	-9.62
66	-16.000	46.000	1.742	14.87	-11.76
67	-14.000	46.000	1.452	13.10	-12.87
68	-12.000	46.000	1.528	7.56	-7.54
69	-10.000	46.000	1.407	3.62	-2.00
70	-8.000	46.000	1.207	1.74	-0.47
71	-26.000	48.000	1.195	0.87	-0.62
72	-24.000	48.000	1.219	1.44	-9.03
73	-22.000	48.000	1.025	2.38	-10.11
74	-20.000	48.000	1.131	4.48	-7.03
75	-18.000	48.000	1.439	7.47	-6.74
76	-16.000	48.000	1.383	8.75	-8.33
77	-14.000	48.000	1.340	7.34	-7.76
78	-12.000	48.000	1.244	4.54	-4.29
79	-10.000	48.000	4.273	2.48	-0.37
80	-8.000	48.000	1.146	1.34	-2.22
81	-26.000	50.000	1.120	0.65	3.20
82	-24.000	50.000	1.239	1.03	-0.03
83	-22.000	50.000	0.759	1.38	-3.64
84	-20.000	50.000	0.919	2.30	-5.80
85	-18.000	50.000	1.369	3.54	-4.81
86	-16.000	50.000	1.133	3.82	-4.49
87	-14.000	50.000	1.065	3.27	-5.72
88	-12.000	50.000	0.995	2.28	-4.46
89	-10.000	50.000	1.026	1.47	-2.05
90	-8.000	50.000	0.984	0.91	-3.70
91	-26.000	52.000	1.160	0.50	3.97
92	-24.000	52.000	0.933	0.64	2.71
93	-22.000	52.000	0.713	0.83	-2.79
94	-20.000	52.000	0.918	1.24	-5.50
95	-18.000	52.000	1.353	1.75	-2.63
96	-16.000	52.000	1.296	1.85	-1.81
97	-14.000	52.000	1.162	1.64	-3.34
98	-12.000	52.000	1.091	1.28	-4.13
99	-10.000	52.000	0.849	0.85	-4.24
100	-8.000	52.000	0.786	0.58	-5.19

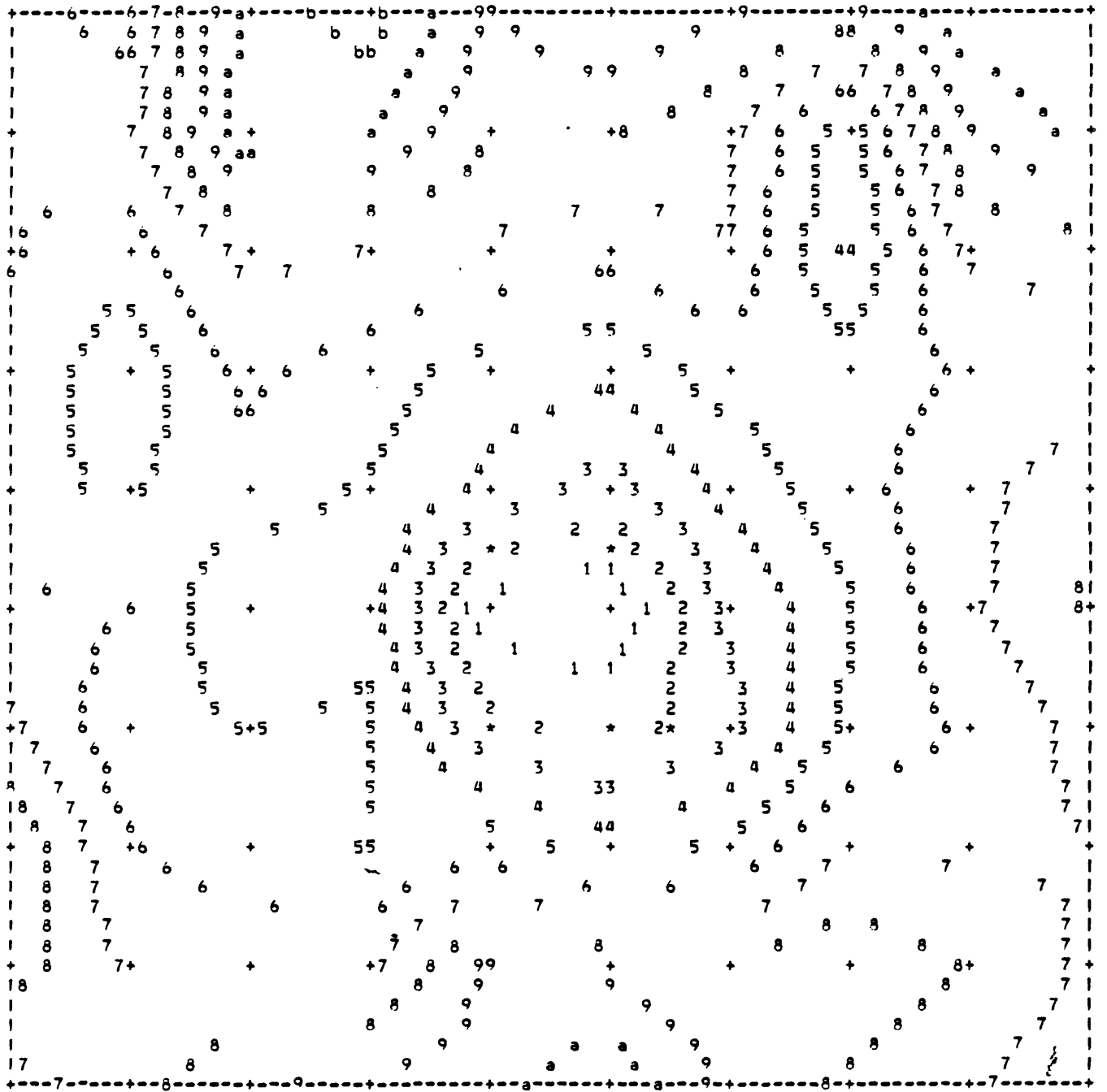
exact plot of calculated values of 10 by 10(100) array. x= -26.0 to -8.0 y= 34.0 to 52.0 contoured at 2.0 milligals

2 4.0000
3 6.0000
4 8.0000
5 10.0000
6 12.0000
7 14.0000
8 16.0000
9 18.0000
a 20.0000
b 22.0000
c 24.0000
d 26.0000
e 28.0000
+ fieldpoint

* corner of body

expect plot of residual values of 10 by 10(100) array. x= -26.0 to -8.0 y= 34.0 to 52.0 contoured at 2.0 milligals

Output Example 2--Continued



symbol value (milliads). y increases to right and x increases downward
 1 -16.0000

Output Example 2--Continued

2	-14.0000
3	-12.0000
4	-10.0000
5	-8.0000
6	-6.0000
7	-4.0000
8	-2.0000
9	0.0000
a	2.0000
b	4.0000
+	fieldpoint

* corner of body

Output Example 3

The following output was produced using command example 3 and a contouring program to produce the plots of the observed, calculated, and residual values.

gravity effect of unlygonal prisms for bond		units of distance measurement are in kilm and heights are in kilm		-0.500 and has 5 vertices. density =		47.000 and has 5 vertices. density =	
body 1	between depths	2.600 and	-14.000	47.000	-17.000	42.000	-17.000
body 2	between depths	-0.500 and	-13.000	47.000	-3.000 and has 5 vertices. density =	-12.000	-19.000
-1A.000	4B.000	-13.000	47.000	47.000	41.000	40.000	3A.000
-26.0000	34.0000	0.7932	-4.4752	0.600			
-24.0000	34.0000	0.8963	-5.0463	44.000			
-22.0000	34.0000	1.0032	-4.7125	0.400			
-20.0000	34.0000	1.1685	-5.0729	3A.000			
-18.0000	34.0000	1.2915	-5.7330				
-16.0000	34.0000	1.4336	-3.7549				
-14.0000	34.0000	1.3534	-1.8552				
-12.0000	34.0000	1.257A	0.9125				
-10.0000	34.0000	1.3793	-0.2056				
-8.0000	34.0000	1.2469	-0.1507				
-6.0000	34.0000	1.1045	-6.0303				
-4.0000	36.0000	1.0740	-3.3738				
-2.0000	36.0000	1.0710	-5.1146				
-20.0000	36.0000	1.3257	-6.2990				
-18.0000	36.0000	1.5777	-0.5202				
-16.0000	36.0000	1.6541	-2.3822				
-14.0000	36.0000	1.6054	-0.3734				
-12.0000	36.0000	1.5400	-3.5456				
-10.0000	36.0000	1.3500	-2.6900				
-8.0000	36.0000	1.2354	-1.9731				
-6.0000	3A.0000	1.2555	3.8396				
-4.0000	3A.0000	1.3A62	5.0040				
-2.0000	3A.0000	1.3163	0.2649				
-20.0000	3A.0000	1.4A03	0.7490				
-18.0000	3A.0000	1.8247	0.6795				
-16.0000	3A.0000	2.0185	-2.5144				
-14.0000	3A.0000	2.0043	-2.3035				
-12.0000	3A.0000	1.7091	-3.7793				
-10.0000	3A.0000	1.5532	-1.9242				
-8.0000	3A.0000	1.5604	0.9555				
-6.0000	40.0000	1.1749	6.1622				
-4.0000	40.0000	1.0047	4.4865				
-2.0000	40.0000	1.1950	1.1651				
-20.0000	40.0000	1.5276	3.1051				
-18.0000	40.0000	2.0572	4.0331				
-16.0000	40.0000	2.6200	3.5525				
-14.0000	40.0000	2.3423	2.6886				
-12.0000	40.0000	1.8463	-1.4664				
-10.0000	40.0000	1.7129	-1.0795				
-8.0000	40.0000	1.5006	2.2547				
-6.0000	42.0000	0.760A	1.0344				
-4.0000	42.0000	1.0243	0.8072				
-2.0000	42.0000	1.0529	1.2485				
-20.0000	42.0000	1.4979	2.3965				
-18.0000	42.0000	2.0366	4.0553				
-16.0000	42.0000	2.813A	9.8200				
-14.0000	42.0000	2.5052	7.7781				
-12.0000	42.0000	2.0709	2.6175				

Output Example 3--Continued

-10.0000	42.0000	1.6770	-4.4229
-8.0000	42.0000	1.3685	3.0072
-26.0000	44.0000	1.0325	3.1538
-24.0000	44.0000	0.8590	0.3559
-22.0000	44.0000	0.9701	-0.7647
-20.0000	44.0000	1.2288	0.6161
-18.0000	44.0000	1.7800	3.3673
-16.0000	44.0000	2.1822	10.9393
-14.0000	44.0000	1.9485	6.4238
-12.0000	44.0000	1.8721	0.5234
-10.0000	44.0000	1.5469	3.7946
-8.0000	44.0000	1.3610	5.9164
-26.0000	46.0000	0.8376	1.0846
-24.0000	46.0000	1.0251	-1.5601
-22.0000	46.0000	1.1882	-0.4104
-20.0000	46.0000	1.1681	0.3335
-18.0000	46.0000	1.4215	3.0304
-16.0000	46.0000	1.7419	3.1077
-14.0000	46.0000	1.4516	0.2309
-12.0000	46.0000	1.5282	0.0212
-10.0000	46.0000	1.4067	1.6231
-8.0000	46.0000	1.2071	1.2785
-26.0000	48.0000	1.1950	0.2554
-24.0000	48.0000	1.2186	-7.5915
-22.0000	48.0000	1.0251	-7.7261
-20.0000	48.0000	1.1310	-2.5403
-18.0000	48.0000	1.4392	0.7354
-16.0000	48.0000	1.3835	0.4184
-14.0000	48.0000	1.3401	-0.4183
-12.0000	48.0000	1.2443	0.2504
-10.0000	48.0000	1.2734	2.1052
-8.0000	48.0000	1.1459	-0.8801
-26.0000	50.0000	1.1200	3.8535
-24.0000	50.0000	1.2394	1.0059
-22.0000	50.0000	0.7588	-2.2630
-20.0000	50.0000	0.9185	-3.4962
-18.0000	50.0000	1.3691	-1.2682
-16.0000	50.0000	1.1329	-0.6621
-14.0000	50.0000	1.0653	-2.4461
-12.0000	50.0000	0.9946	-2.1838
-10.0000	50.0000	1.0258	-0.5849
-8.0000	50.0000	0.9843	-2.7911
-26.0000	52.0000	1.1602	4.4701
-24.0000	52.0000	0.9325	3.3534
-22.0000	52.0000	0.7125	-1.9564
-20.0000	52.0000	0.9177	-4.2629
-18.0000	52.0000	1.3534	-0.8839
-16.0000	52.0000	1.2964	0.0481
-14.0000	52.0000	1.1620	-1.7044
-12.0000	52.0000	1.0913	-2.4486
-10.0000	52.0000	0.8489	-3.3899
-8.0000	52.0000	0.7865	-4.6056

standard deviation= 6.61 and average difference= -4.85 milligrams

	x	y	z	q	residual
1	-26.000	34.000	0.793	0.49	-4.97
2	-24.000	34.000	0.896	0.74	-5.84

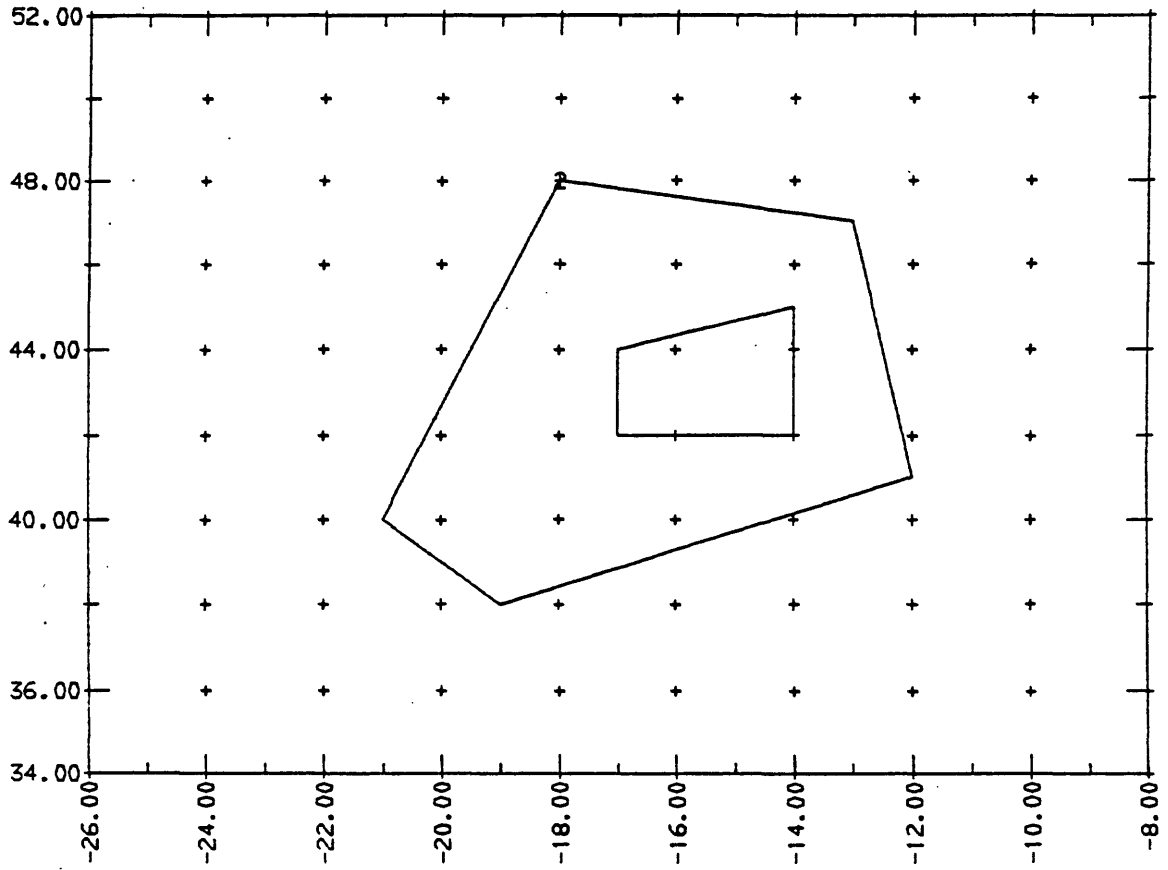
Output Example 3--Continued

3	-22.000	34.000	1.093	1.12	-5.83
4	-20.000	34.000	1.169	1.49	-6.56
5	-18.000	34.000	1.291	1.74	-7.47
6	-16.000	34.000	1.434	1.76	-5.51
7	-14.000	34.000	1.553	1.49	-3.35
8	-12.000	34.000	1.258	1.15	-0.24
9	-10.000	34.000	1.379	0.90	-1.11
10	-8.000	34.000	1.247	0.63	-4.78
11	-26.000	36.000	1.104	0.76	-6.79
12	-24.000	36.000	1.074	1.21	-4.59
13	-22.000	36.000	1.071	1.97	-7.08
14	-20.000	36.000	1.326	3.05	-9.35
15	-18.000	36.000	1.578	3.71	-8.23
16	-16.000	36.000	1.654	3.56	-5.94
17	-14.000	36.000	1.605	2.90	-7.28
18	-12.000	36.000	1.540	2.12	-5.66
19	-10.000	36.000	1.350	1.38	-4.07
20	-8.000	36.000	1.235	0.89	-2.86
21	-26.000	38.000	1.255	1.05	2.79
22	-24.000	38.000	1.386	1.91	3.09
23	-22.000	38.000	1.316	3.57	-3.31
24	-20.000	38.000	1.480	6.26	-5.51
25	-18.000	38.000	1.825	7.63	-6.95
26	-16.000	38.000	2.019	7.18	-9.70
27	-14.000	38.000	2.004	5.74	-8.04
28	-12.000	38.000	1.709	3.82	-7.60
29	-10.000	38.000	1.553	2.28	-4.20
30	-8.000	38.000	1.560	1.37	-0.41
31	-26.000	40.000	1.175	1.22	4.94
32	-24.000	40.000	1.005	2.28	2.20
33	-22.000	40.000	1.195	5.23	-4.06
34	-20.000	40.000	1.528	10.03	-6.93
35	-18.000	40.000	2.057	12.23	-8.19
36	-16.000	40.000	2.620	12.23	-8.67
37	-14.000	40.000	2.342	10.34	-7.65
38	-12.000	40.000	1.846	6.57	-8.04
39	-10.000	40.000	1.713	3.48	-4.56
40	-8.000	40.000	1.501	1.78	0.47
41	-26.000	42.000	0.761	1.13	-0.10
42	-24.000	42.000	1.024	2.45	-1.64
43	-22.000	42.000	1.053	5.54	-4.29
44	-20.000	42.000	1.498	11.29	-8.89
45	-18.000	42.000	2.037	16.31	-10.28
46	-16.000	42.000	2.814	26.74	-16.92
47	-14.000	42.000	2.505	21.09	-13.31
48	-12.000	42.000	2.071	9.33	-6.71
49	-10.000	42.000	1.677	4.41	0.01
50	-8.000	42.000	1.369	2.03	0.97
51	-26.000	44.000	1.033	1.18	1.97
52	-24.000	44.000	0.860	2.12	-1.77
53	-22.000	44.000	0.970	4.63	-5.40
54	-20.000	44.000	1.229	10.12	-9.50
55	-18.000	44.000	1.780	16.28	-12.91
56	-16.000	44.000	2.182	28.84	-17.90
57	-14.000	44.000	1.949	21.27	-14.84
58	-12.000	44.000	1.872	9.76	-9.23
59	-10.000	44.000	1.547	4.44	-0.64
60	-8.000	44.000	1.361	2.07	3.85
61	-26.000	46.000	0.838	0.95	0.13
62	-24.000	46.000	1.025	1.61	-3.37

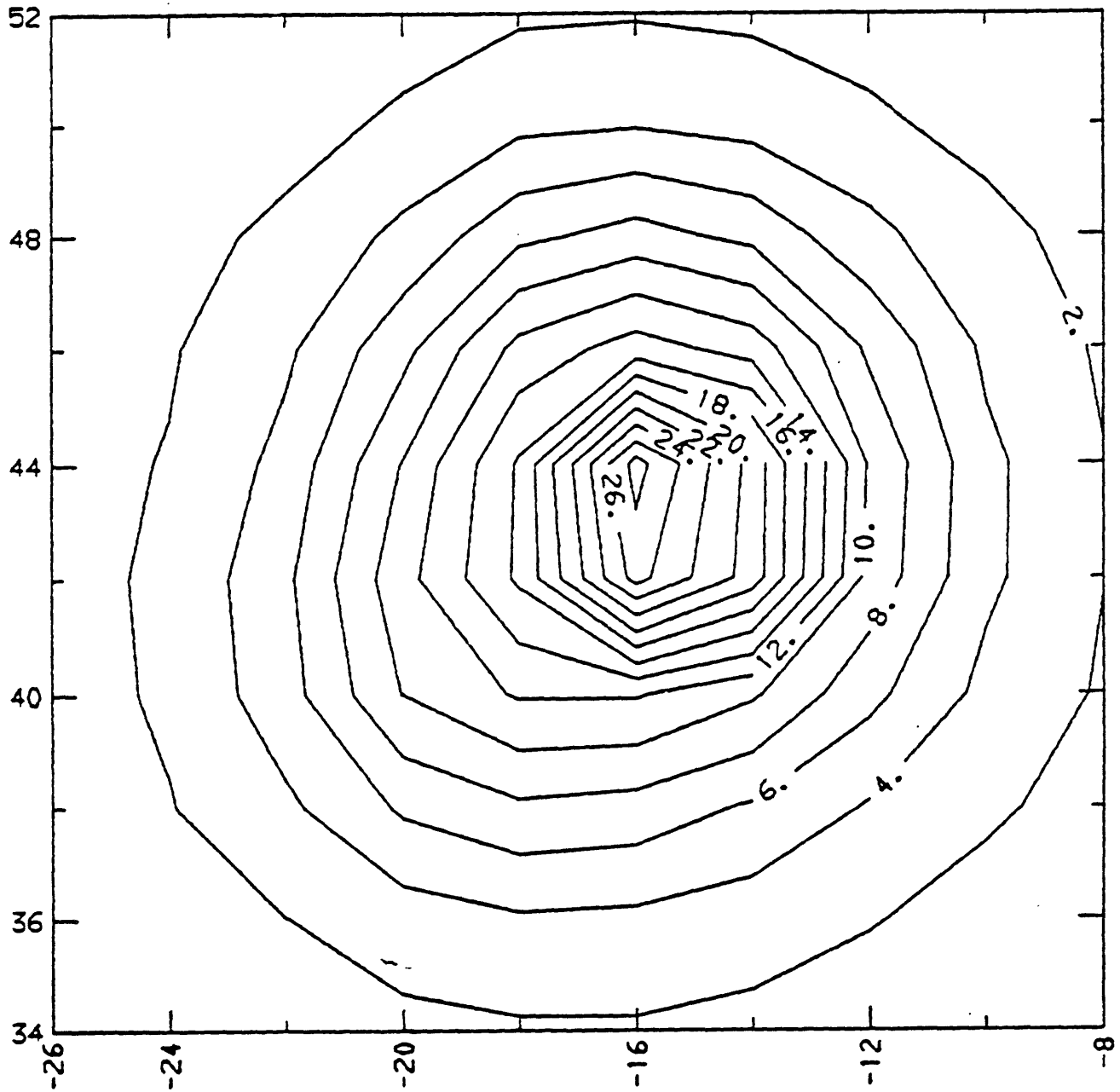
Output Example 3--Continued

63	-22.000	46.000	1.188	3.62	-4.03
64	-20.000	46.000	1.168	7.36	-7.02
65	-18.000	46.000	1.422	12.65	-9.62
66	-16.000	46.000	1.742	14.87	-11.76
67	-14.000	46.000	1.452	13.10	-12.87
68	-12.000	46.000	1.528	7.56	-7.54
69	-10.000	46.000	1.407	3.62	-2.00
70	-8.000	46.000	1.207	1.74	-0.47
71	-26.000	48.000	1.195	0.87	-0.62
72	-24.000	48.000	1.219	1.44	-9.03
73	-22.000	48.000	1.025	2.38	-10.11
74	-20.000	48.000	1.131	4.48	-7.03
75	-18.000	48.000	1.439	7.47	-6.74
76	-16.000	48.000	1.383	8.75	-8.33
77	-14.000	48.000	1.340	7.34	-7.76
78	-12.000	48.000	1.244	4.54	-4.29
79	-10.000	48.000	1.273	2.48	-0.37
80	-8.000	48.000	1.146	1.34	-2.22
81	-26.000	50.000	1.120	0.65	3.20
82	-24.000	50.000	1.239	1.03	-0.03
83	-22.000	50.000	0.759	1.38	-3.64
84	-20.000	50.000	0.919	2.30	-5.80
85	-18.000	50.000	1.369	3.54	-4.81
86	-16.000	50.000	1.133	3.82	-4.49
87	-14.000	50.000	1.065	3.27	-5.72
88	-12.000	50.000	0.995	2.28	-4.46
89	-10.000	50.000	1.026	1.47	-2.05
90	-8.000	50.000	0.984	0.91	-3.70
91	-26.000	52.000	1.160	0.50	3.97
92	-24.000	52.000	0.933	0.64	2.71
93	-22.000	52.000	0.713	0.83	-2.79
94	-20.000	52.000	0.918	1.24	-5.50
95	-18.000	52.000	1.353	1.75	-2.63
96	-16.000	52.000	1.296	1.85	-1.81
97	-14.000	52.000	1.162	1.64	-3.34
98	-12.000	52.000	1.091	1.28	-4.13
99	-10.000	52.000	0.849	0.85	-4.24
100	-8.000	52.000	0.786	0.58	-5.19

Output Example 3--Continued

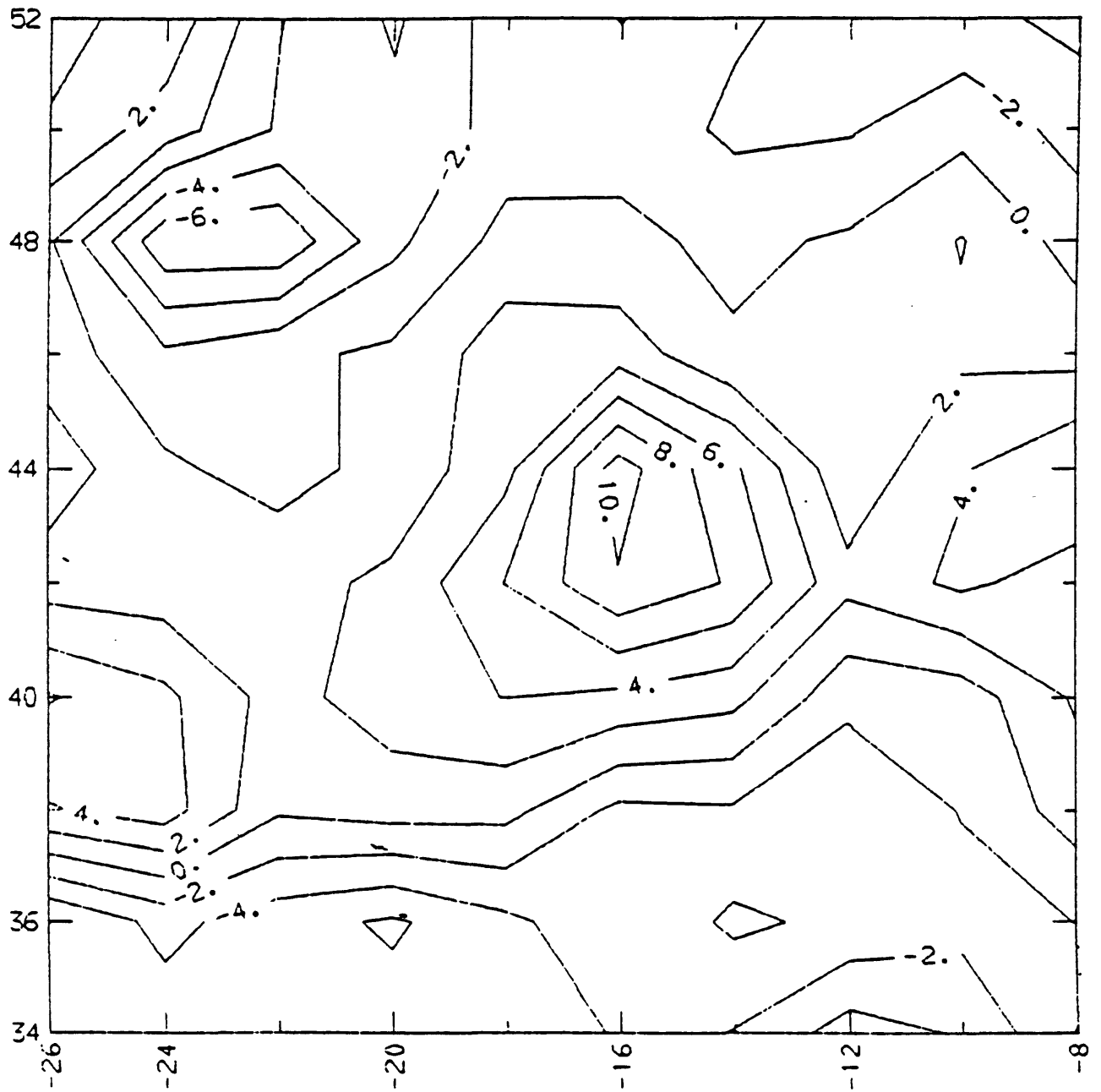


Body corners and fieldpoint locations



Calculated values

Output Example 3.--Continued



Observed values

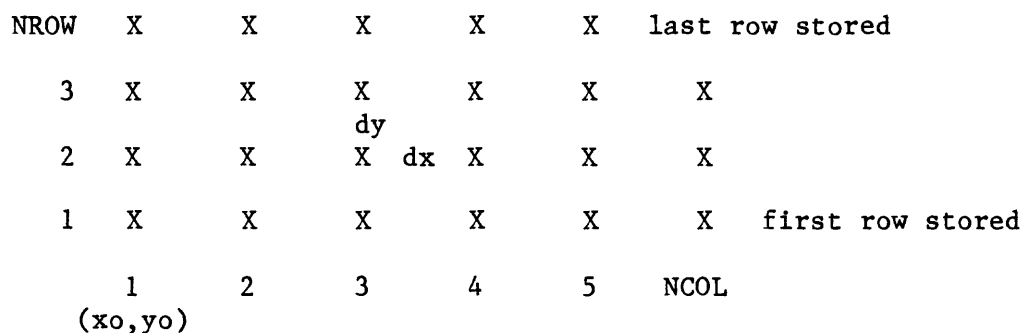
Appendix E

Standard Grid

The grid can be any one of many types: (a) a rectangular grid with equal spacing in the x- and y-directions, (b) a rectangular grid with constant but unequal x- and y-intervals, (c) a rectangular grid with varying distances between grid positions in either the x- or y-direction or both, and d) a quadrilateral grid that consists of connecting quadrilaterals whose interior angles do not exceed 180 degrees.

The file of the gridded data consists of two basic parts: (1) a header record and optionally, a following record that contains the x-coordinates for each column, and (2) a series of data records, each containing the column values for one row.

The following diagram shows the relationship of the grid elements in the usual case where dx and dy are positive.



A. Header record (23 words long)

id: 56 ASCII characters of identification (14 words).

pgm: 8 ASCII characters of creation program identification (2 words).

ncol: number of columns of data (integer, 1 word).

nrow: number of rows of data (integer, 1 word).

nz: number of words per data element (integer, 1 word). For single

precision use 1, double precision or complex use 2, double precision

complex use 4. For quadrilateral grids this value is 3.

xo: position of first column of data (real, 1 word).

dx: equal spacing interval of columns (real, 1 word).

If equal to zero, then coordinates for each column are in the following data record; otherwise the following record consists of data for row one.

yo: position of first row (real, 1 word).

dy: equal spacing interval of rows (real, 1 word). If equal to zero, the coordinate for each row is the first word of each data record row.

B. Column coordinate record, present only if dx of header record is equal to zero. Record consists of ncol real words specifying the coordinates of each data column in monotonic order. If nz=3 then this record is present but the values are meaningless.

C. Data record. Each data record contains one row of real data items. The total record length is ncol times (nz plus 1) words. For quadrilateral data the sequence of data values is x, y, and z. The first word contains the row coordinate if dy of the header record is zero, else the value is a dummy. Again, the row coordinates should be in monotonic sequence, if specified.

In general, i/o for this standard file can be stated in fortran as:

```
dimension g(iz,ix,iy),id(14),pgm(2),x(ix),y(iy)
read or write (..) id,pgm,ncol,nrow,nz,xo,dx,yo,dy
if (dx.eq.0.) read or write (..) (x(i),i=1,ncol)
if (dy.ne.0.) go to 15
do 10 j=1,nrow
10 read or write (..) y(j),((g(k,i,j),k=1,nz),i=1,ncol)
go to 25
15 do 20 j=1,nrow
```

```
20   read or write (..) dum,((g(k,i,j),k=1,nz),i=1,ncol)
```

```
25   continue
```

In the usual case where dx and dy are constant and nz=1. the code simplifies to:

```
dimension g(ix,iy),id(14),pgm(2)
```

```
read or write(..) id,pgm,ncol,nrow,nz,xo,dx,yo,dy
```

```
do 10 j=1,nrow
```

```
10   read or write (..) dum,(g(i,j),i=1,ncol)
```


Appendix F

Example of User-Formatted File

0.	0.	1.	0
1.414	0.	3.	0
2.828	0.	5.	0
4.243	0.	10.	0
5.657	0.	20.	0
0.	1.414	2.	0
1.414	1.414	5.	0
2.828	1.414	10.	0
4.243	1.414	15.	0
5.657	1.414	50.	0
0.	2.828	3.	0
1.414	2.828	5.	0
2.828	2.828	15.	0
4.243	2.828	45.	0
5.657	2.828	50.	0

Appendix G

External Subroutines

The following subroutines are called externally by the program:

Basic Plot System (Evdenden and Wahl, 1977)

char (x, y, a, nc, icode, size, theta, xoff, yoff) - writes
characters on the plot

endpt (ie) - terminates plot

line (x, y, n, icon, ipn) - draws one or more lines

neatl () - draws a box around data area

pltset (iplotr, xbd, ybd, isl) - initializes plot system

scale (dxp, dyp, xp, yp, nopts, ier) - defines the plot area
and units

xaxis (dxp, dyp, xp, del, ip, size, fmt, nfmt) - draws and
labels x axis

yaxis (dyp, dyp, xp, del, ip, size, fmt, nfmt) - draws and
labels y axis

where:

a = character string to be plotted

del = data unit interval of tick marks

dxp = array defining plot size (see fig. C1)

dyp = array defining plot size (see fig. C1)

fmt = Fortran format for labels

icode = 1 means x, y are arrays, 3 means x, y are single elements

icon = 0 start new line, 1 continue old line

ie = 10 element array containing diagnostics

ier = set to 0 to indicate no problems

ip = interval between tick mark labels

ipn = line type, 0=solid, 1-7 dashed lines
 iplotr = plotter number
 isl = 5 element array, isl(1)=1 inches, 2 cm
 n = dimension of x, y arrays in line call
 nc = if icode is 0 or 1, dimension of x and y arrays, otherwise
 length of a
 nfmt = length of format
 nopts = length of xp and yp arrays
 size = character size
 theta = rotation of char in radians
 x = x location (array of dimension n)
 y = y location (array of dimension n)
 xbd = returned maximum x size for selected plotter
 ybd = returned maximum y size for selected plotter
 xoff = x offset for center of label
 yoff = y offset for center of label
 xp = array defining plot size (see fig. C1)
 yp = array defining plot size (see fig. C1)
 scale3(xmin, xmax, nxl, xminp, xmaxp, xdel)

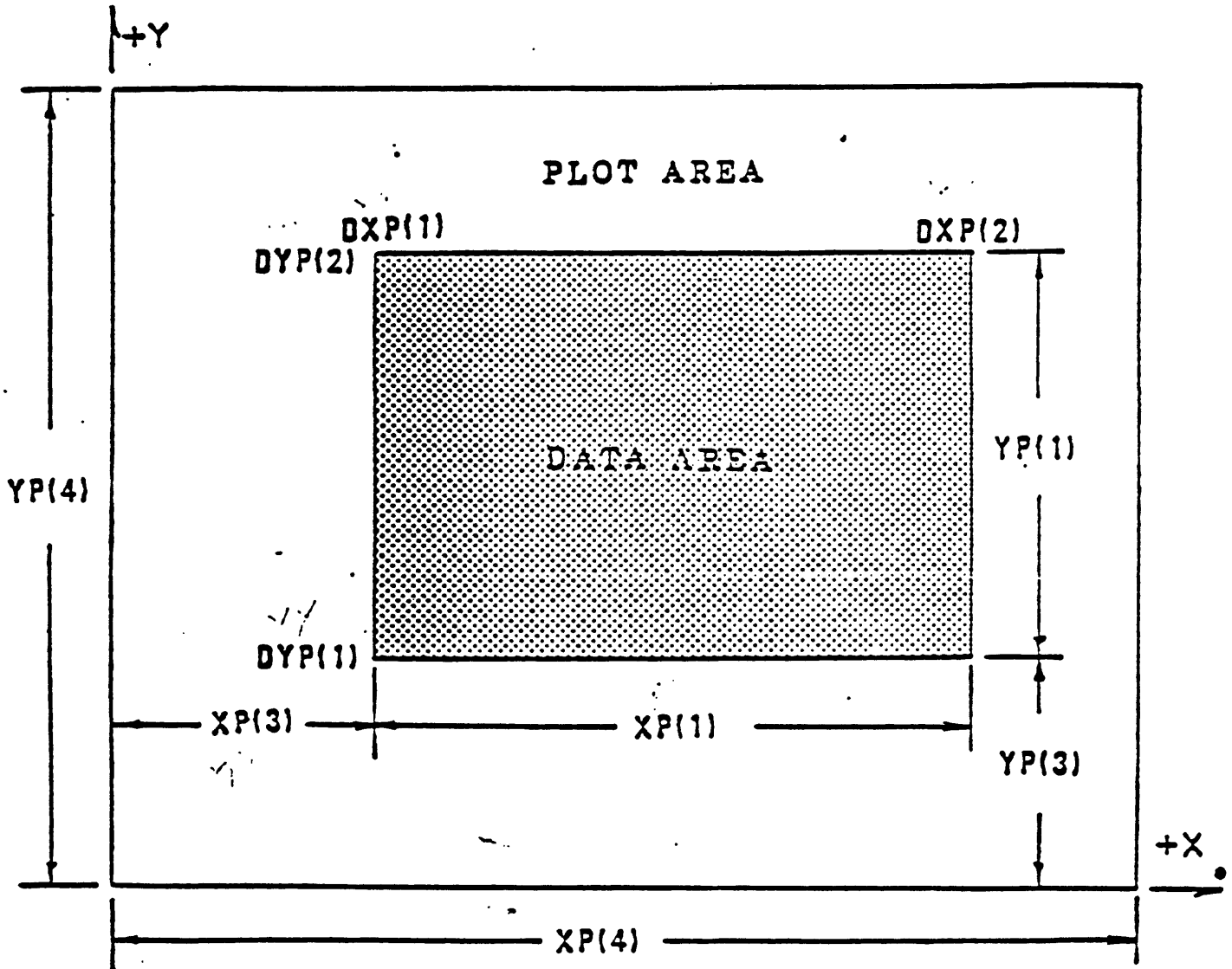
where:

xmin = minimum x value of data (input)
 xmax = maximum x value of data (input)
 nxl = number of divisions along x axis (input)
 xminp = minimum whole number of x value (output)
 xmaxp = maximum whole number of x value (output)
 xdel = increment for labelling x axis (output)

Figure C1. DXP,DYP,XP,YP arrays and their relation to the plot area.

DXP and DYP are in data units.

XP and YP are in plot units (inches).



```

c*****
c program to input exact talwani polygonal prisms for 3d gravity-plouff
c modified by r.godson for denver multics 5/81
c
  character*4 unix,uniz
  character*8 name
  character*18 title
  character*40 idplot
  character*50 ifile,ifile2,cfile
  character*54 rfile
  character*55 cfile
  character*56 pfile
  character*80 ifmt
  common/st/cfile,rfile,pfile
  common/graving/iplotr,ibody,ifile,ifile2,lsqs,
  1 datum,xscale,dc,unix,uniz,name,height,
  2 ifmt,iplot,iobs,icalc,ires
  common /field/r(1000)
  common /fpts/ xf(1000),yf(1000),zf(1000)
  common /bodk/ mm,nlim,jbody,short,nend
  common /trans/ inpt,iout,jin,iprint
  dimension a(1000),b(1000),dstnc(900)
c dimension limits for 3 parameters. corners, bodies, fieldpoints.
c one more than dimension of total number of fieldpoints
  data ncorn,nbod,npt/900,50,1001/
c
c   get input parameters
c
  call min($220)
  10 write (iout,20) idplot
  20 format ('1 gravity effect of polygonal prisms',/,
  1 1x, ' for ',5x,a40)
  facx=0.0
  facz=0.0
  call unit(unix,uniz,facx,facz)
  if (facx .eq. 0.0) go to 200
  zfac=1.0/facz
  ofac=6.670*facx
  30 short=1.0e20
c
c   read body coordinates
c
  call bodg(facz,nbod,dstnc,ncorn,ofac)
c mm (common) is total number of layers
  if (mm .eq. 0) go to 200
c
c   read fieldpoint locations and maybe values
c
  call ptsm(nk,$220)
  if (nk .lt. 1) go to 200
c
c   check if printer plot of bodies & fieldpoints
c   desired if fieldpoints are in arid form
c
  title="body corners"
  if(ibody.ne.0.and.iplotr.eq.9.and.name.eq."gridded")
  1 call cplot(dc,a,1,0,title,$365)
c
c   check if location plot of bodies & fieldpoints on
c   plotting device wanted

```

```

c      if(ibody.ne.0.and.iplotr.ne.9) call grav3dplot(nk,$220)
c
c      check if plot of observed values desired
c
      title="observed values"
      if(iobs.ne.0.and.iplotr.eq.9.and.name.eq."gridded")
1 call cplot(dc,r,nk,1,title,$365)
      do 40 k=1,nk
40  zf(k)=zf(k)*zfac
50  call qtalw(nk,mm,short,dstnc,a)
      if(datum .ne. 0.) write (iout,60) datum
60  format (/,1x,f10.2," maal is added to gravity values")
      if (lsas .eq. 0) go to 130
      do 70 k=1,nk
70  r(k)=r(k)-datum
      if (lsas .ne. 7) go to 90
      ier=0
      call qlsas(nk,a,ier,datum)
      if (ier .ne. 0) go to 200
90  std=0.0
      avd=0.0
      do 100 k=1,nk
      errk=r(k)-a(k)
      b(k)=errk
      avd=avd+errk
100  std=std+errk*errk
110  if (nk .lt. 2) go to 130
      std=sqrt(std/(nk -1))
      avd=avd/nk
      write (iout,120) std,avd
120  format (///,5x,"standard deviation=",
1 f8.2," and average difference=",f8.2," milligals")
130  write (iout,140)
140  format (/,15x,"x",9x,"y",9x,"z",9x,"a",4x,"residual")
      do 180 k=1,nk
      x=xf(k)
      y=yf(k)
      zfk=zf(k)*zfac
      grav=a(k)+datum
      a(k)=grav
      if (lsas .eq. 0) go to 160
      write (iout,150) k,y,x,zfk, grav,b(k)
150  format (i6,3x,3f10.3,2f9.2)
      do to 180
160  write (iout,170) k,y,x,zfk, grav
170  format (i6,3x,3f10.3,f9.2)
180  continue
c
c      contour plots
c
340  title="calculated values"
      if (icalc.ne.0.and.iplotr.eq.9.and.name.eq."aridded")
1 call cplot (dc,a,nk,1,title,$365)
      title="residual values"
      if (ires.ne.0.and.iplotr.eq.9.and.lsas.ne.0.and.name.eq."gridded")
1 call cplot (dc,b,nk,1,title,$365)
      if(icalc.ne.0.and.iplotr.lt.9.and.name.eq."gridded")
1 call stdout(1,a)
      if (ires.ne.0.and.iplotr.lt.9.and.lsas.ne.0.and.name.eq."aridded")

```

```

1 call stdout(1,a)
  if(ires.ne.0.and.iplotr.lt.9.and.lsas.ne.0.and.name.eq."aridded")
1 call stdout(2,b)
  do to 220
365 write(iprint,375)
375 format(" contour interval(dc) is 0.")
200 write(iprint,210)
210 format(" error - program abort")
220 close(jin)
  close(iout)
  write(iprint,230) pfile
230 format(/," output is in file ",a56)
  stop
  end
c*****
  subroutine min(*)
c
c  namelist input parameter subroutine
c
c  r.aodson = 3/81
c
  character*4 unix,uniz
  character*8 name
  character*40 idplot
  character*50 ifile,ifile2,cfile
  character*54 rfile
  character*55 cfile
  character*56 pfile
  character*80 ifmt
  common/st/cafile,rfile,pfile
  common/oravinp/iplotr,ibody,ifile,ifile2,lsas,
1 datum,xscale,dc,unix,uniz,name,height,
2 ifmt,idplot,iobs,icalc,ires
  common/contq/nrow,ncol,dy,dx,xmin,ymin,xmax,ymax
  common /vec/cde,cie,sde,sie,ex,ey,ez,earth,cdci,sdci
  common /trans/ inpt,iout,jin,iprint
  common/widect/naxcol
  namelist/parms/iplotr,ibody,ifile,ifile2,lsas,
1 datum,xscale,dc,unix,uniz,name,height,
2 ifmt,idplot,iobs,icalc,ires,naxcol
  inpt=5
  iout=8
  iprint=6
  jin=9
  iplotr=9
  naxcol=130
  ibody=0
  iobs=0
  icalc=0
  ires=0
  idplot=" "
  ifile=" "
  ifile2=" "
  lsas=0
  datum=0.
  xscale=0.
  dc=0.
  unix="kiln"
  uniz="feet"
  name="aridded"

```

```

height=0.
ifmt="(v)"
write(iprint,10)
10 format(" enter command file name: ",$)
read(inpt,20) cfile
20 format(v)
open(jin,file=cfile,form="formatted")
read(jin,parms,end=30)
if(lsas.eq.0) ires=0
if(ifile.eq." ") go to 50
if(iplotr.lt.0.or.iplotr.gt.9) iplotr=9
c
c   encode output file names
c
   encode(pfile,25) cfile
25 format("p rint.",a50)
   encode(cfile,26) cfile
26 format("c alc.",a50)
   encode(rfile,27) cfile
27 format("r es.",a50)
   open(iout,file=pfile,form="formatted",carriage=.true.)
   return
30 write(iprint,40)
40 format(" end of file reading namelist command file")
   go to 70
50 write(iprint,60)
60 format(" no observed data file name specified")
70 return 1
   end
c*****
   subroutine bodq(facz,nbod,dstnc,ncorn,afac)
c
c   reads body corner coordinates
c
c   r.qodson = 4/81
c
   character*4 unix,uniz
   character*8 name
   character*40 idplot
   character*50 ifile,ifile2,cfile
   character*80 ifmt
   dimension dstnc(1)
   common/mplot/ymin,xmin,ymax,xmax
   common/conta/nrow,ncol,dy,dx,yo,xo,xma,yma
   common/aravino/iplotr,ibody,ifile,ifile2,lsas,
1 datum,xscale,dc,unix,uniz,name,height,
2 ifmt,idplot,iobs,icalc,ires
   common /bod/ rho(50),ifprnt(50),zee(50),zm(50)
c see ncorn dimension limit in main program
   common /side/xb(900),yb(900), ll(50)
   common /bodk/ mm,nlim,jbody,short,nend
   common /trans/ inpt,iout,jin,iprint
   nend=0
   mm=0
   xmin=1.0e38
   xmax=-1.0e38
   ymin=1.0e38
   ymax=-1.0e38
10 read(jin,20,end=180)num,ifprnt(mm+1),dencon,top,bottom
20 format(v)

```



```

    if(1sas.ea.7) dencon=1.
    nst=nend+1
    nend=nend+num
    mm=mm+1
    if(mm.at.nbod) go to 130
    if(nend.at.ncorn) go to 70
    ll(mm)=num
    zee(mm)=top*facz
    zm(mm)=bottom*facz
    rho(mm)=dencon*gfac
    write(iout,30) mm,top,bottom,num,dencon
30  format (' body',i3,' between depths',f11.3,' and',f11.3,
1    ' and has',i4,' vertices. density = ',f8.3)
40  if (top .at. bottom) go to 90
50  write (ioint,60) top,bottom
60  format ('0top of body is below bottom',2f7.2)
    go to 170
70  write(ioprint,80) mm,nend,ncorn
80  format (' body',i3,' includes sufficient corners',i4,' to exceed'
1    ' limitation of',i5,' for all bodies.')
    go to 170
c  coordinates of corners of body east,north (x,y) pairs.
c  body has at least 3 corners
90  read(jin,20,end=150) (yb(j),xb(j),j=nst,nend)
    write(iout,100) (yb(j),xb(j),j=nst,nend)
100 format (5(5x,2f9.3))
    do 105 j=nst,nend
        if(xb(j).lt.xmin) xmin=xb(j)
        if(xb(j).gt.xmax) xmax=xb(j)
        if(yb(j).lt.vmin) vmin=yb(j)
        if(yb(j).gt.vmax) vmax=yb(j)
105 continue
110 np=nend+1
    xb(np)=xb(nst)
    yb(np)=yb(nst)
    do 120 j=nst,nend
        dist=(xb(j+1)-xb(j))**2+(yb(j+1)-yb(j))**2
        if (dist .eq. 0.0) go to 120
        dist=sqrt(dist)
        if (dist .lt. short) short=dist
120 dstnc(j)=dist
    go to 10
130 write (ioint,140) mm
140 format (5x,'number of bodies=',i3,' exceeds limitation.')
    go to 170
150 write(ioprint,160) mm
160 format(" end of file reading x,y pairs for body no. ",i3)
170 mm=0
180 return
    end
c*****
    subroutine unit(unix,uniz,face,facez)
c  conversion of horizontal (x) and vertical (z) distances-3d magnetic
character*4 unix,uniz,feet,kilm,mile,kilf,metr,jblank,
1  lunx,lunz
    common /trans/ inpt,iout,jin,ioprint
    data kilm, feet, mile, kilf, metr,jblank
1  /'kilm','feet','mile','kilf','metr',' /
    write (iout,10) unix,uniz
10  format (5x,'units of distance measurement are in ',a4,

```

```

1  and heights are in ',a4)
20 if (unix .eq. mile) facx=1.609347
   if (uniz .eq. mile) facz=1.609347
   if (unix .eq. feet) facx=3.048006e-4
   if (uniz .eq. feet) facz=3.048006e-4
   if (unix .eq. kilf) facx=3.048006e-1
   if (uniz .eq. kilf) facz=3.048006e-1
   if (unix .eq. kilm) facx=1.0
   if (uniz .eq. kilm) facz=1.0
   if (unix .eq. metr) facx=1.0e-3
   if (uniz .eq. metr) facz=1.0e-3
   if (facx .ne. 0.0) go to 40
   write(iprint,30)
30 format(" unix specification is incorrect")
   return
40 if (facz .ne. 0.0) go to 50
   facx=0.0
   return
c multiplying factor for heights to convert to distance units.
c negative sign to compensate for depths + downward later 'calc'.
50 facz=-facz/facx
   return
end
c*****
   subroutine atalw(nk,mm,short,dstnc,a)
c exact solution for talwani 3d gravity in single precision (7 figures)
   common /side/xb(900),yb(900), ll(50)
   common /bod/ rho(50),ifprnt(50),zee(50),zm(50)
   common /fpts/ xp(1000),yp(1000),zp(1000)
   common /trans/ inpt,iout,jdisk,page
   dimension a(1),dstnc(1)
c 1/1000 length of shortest non-zero side
   test=short*1.0e-3
   do 240 k=1,nk
     x=xp(k)
     y=yp(k)
     z=zp(k)
     grav=0.0
     ne =0
c loop of bodies
     do 230 m=1,mm
c number of vertices
       ns =ne +1
       lk= ll(m)
       z1=zee(m)-z
       z2=zm (m)-z
       ne =ns +lk-1
       af=rho(m)
       jfprnt=ifprnt(m)
       a=0.0
c to give clockwise positive
       h1=abs(z1)
       h2=abs(z2)
       zz=z1+z2
       if (zz)10,230,20
c case where most of mass is above fieldpoint (minus a)
       10 af=-af
       20 if (h1 .lt. h2) go to 30
         t=h2
         h2=h1

```

```

    h1=t
30 x1=xb(ns)-x
    v1=yb(ns)-y
c last point is repeat of first
    xf=x1
    yf=y1
    r1s=x1*x1+y1*y1
    r1sf=r1s
    hh1=h1*h1
    hh2=h2*h2
    dh=h2-h1
    r11= sqrt(r1s+hh1)
    r12= sqrt(r1s+hh2)
    r11f=r11
    r12f=r12
    ana=0.0
    t1=0.0
    t2=0.0
    w1=0.0
    w2=0.0
    w3=0.0
    n=ns
c loop of all sides
do 210 l=1,lk
    nm=n
    if (l .ne. lk) go to 40
    x2=xf
    y2=yf
    r2s=r1sf
    r21=r11f
    r22=r12f
    go to 50
40 n=n+1
    x2=xb(n)-x
    y2=yb(n)-y
    r2s=x2*x2+y2*y2
    r22= sqrt(r2s+hh2)
    r21= sqrt(r2s+hh1)
50 ds=dstnc(nm)
    if (test .at. ds) go to 190
c same sign as definition used for magnetics. opposite last gravity
    p=(x1*y2-x2*y1)/ds
    dx=x2-x1
    dy=y2-y1
c test if perpendicular distance to line segment is less than 1/1000
c of shortest segment length
    ap=abs(p)
    if (ap .lt. test) go to 190
    d1= (x1*dx+y1*dy)/ds
    d2= (x2*dx+y2*dy)/ds
    dd1= abs(d1)
    dd2= abs(d2)
    if (dd1 .lt. test) d1=0.0
    if (dd2 .lt. test) d2=0.0
    fnum=ap*ds
    den=x1*x2+y1*y2
    if (den) 60,100,110
60 a1=3.141593
    if (n .at. 0.0) go to 80
    a1=-a1

```

```

70 den=-den
80 call tandet(fnum/den,w3,ana)
90 ana=ana+a1
  go to 120
100 a1=1.570796
  if (p .lt. 0.0) a1=-1.570796
  go to 90
110 a1=0.0
  if (p .gt. 0.0) go to 80
  go to 70
120 if (d1 .eq. 0.0) go to 130
  call tandet(-d1*h1/(p*r11),w1,t1)
  call tandet( d1*h2/(p*r12),w2,t2)
130 if (d2 .eq. 0.0) go to 140
  call tandet( d2*h1/(p*r21),w1,t1)
  call tandet(-a2*h2/(p*r22),w2,t2)
140 ph=p*p+hh1
  rz1=r11+d1
  if (d1 .ae. 0.0) go to 160
  if (d2 .at. 0.0) go to 150
  a1=(r21-d2)/(r11-d1)
  go to 180
150 dtest=ph/(d1*d1)
  if (dtest .lt. 0.01) rz1=dd1*0.5*dtest*(1.0-0.25*dtest)
160 rz2=r21+d2
  if (d2 .ae. 0.0) go to 170
  dtest=ph/(d2*d2)
  if (dtest .lt. 0.01) rz2=dd2*0.5*dtest*(1.0-0.25*dtest)
170 a1=rz1/rz2
180 a1=p*a1+a1*(r22+d2)/(r12+d1)
  a=a-a1
190 if (jfornt .lt. 0) write (iout,200) 1,p,d1,d2,ana,a1,w1,w2
200 format (i4,7e16.8)
  x1=x2
  y1=y2
  r1s=r2s
  r12=r22
  r11=r21
210 continue
  t1=t1+ atan(w1)
  t2=t2+ atan(w2)
  ana= ana+atan(w3)
  q=a1*(ana*dh+a+n1*t1+h2*t2)
  if (jfornt .ne. 0) write (iout,220) k,a,ana ,t1,t2
220 format (i4,f8.2,' mqa1',15x,3f9.3)
230 grav=grav+a
240 a(k)=grav
  return
  end
c*****
  subroutine tandet(f,zze,sume)
c program for arctangent addition. negative f for subtraction.
  a1=zze+f
  a2=1.0d0-zze*f
  if (a2 .ne. 0.0 ) go to 40
  if (a1) 10,20,20
  10 sume=sume-1.570796
  go to 30
  20 sume=sume+1.570796
  30 zze=0.0

```

```

    return
40 zze=a1/a2
    if (a2 .at. 0.0 ) return
    if (a1) 50,60,70
50 sume=sume-3.141593
60 return
70 sume=sume+3.141593
    return
end
c*****
subroutine alsas(kk,a,ier,datum)
c least-squares calculation of density regression
real*8 fn,sx,sy,sxx,sxy,syv
integer nm(4)/'n','x','s','y'/
common /field/r(1000)
common /trans/ inpt,iout,jdisk,page
dimension a(1)
l=1
if (kk .lt. 3) go to 40
fn=kk
sx =0.0d0
sy =0.0d0
syv=0.0d0
sxy=0.0d0
sxx=0.0d0
do 10 k=1,kk
rk=r(k)
ak=a(k)
sx =sx +ak
sy =sy +rk
syv=syv+rk*rk
sxy=sxy+ak*rk
10 sxx=sxx+ak*ak
xx=sxx-sx*sx/fn
l=2
if (xx .le. 0.0) go to 40
xy=sxy-sx*sy/fn
yy=svv-sy*sy/fn
tm=xy/xx
tb=(sy-tm*sx)/fn
l=4
if (vy .le. 0.0) go to 40
coer=xy/sqrt(xx*vy)
stdm=yv*(1.0-coer*coer)/(xx*(kk-2))
l=3
if (stdm .lt. 0.0) go to 40
stdm=sqrt(stdm)
tb=tb+datum
write (iout,20) tm,stdm,coer,tb
20 format (//,' least-squares density is',f7.3,' (std',f6.3,') with',
1 ' correlation coefficient=', f6.2,3x,' datum=',f7.2,
2 ' maal.')
```

```

do 30 k=1,kk
30 a(k)=tm*a(k)
return
40 ier=1
write (iout,50) nm(1)
50 format ('0stop. density can be calculated. cutoff ',a1)
return
end

```

```

c*****
      subroutine cplot(pc,a,nt,jk,title,*)
c   jk must be zero for first call of cplot. later calls use revised xb/yb
c       ci---the initial contour value (calculated)
c       pc,dc---the increment in the contour values (read in)
c       a(y,x)---the input array containing the data to be contoured.
c   index for equivalent 1d array is n*(x-1)+y
c       m---the final value of x-index in a-array
c       n---the final value of y-index in a-array
c       limits of plot are a(yi,xi) to a(n,m)
c   nx is number of intervals in x-direction (right) less than 41.
c   ny is number of intervals in y-direction (down) less than 41.
      character*18 title
      integer ont(131),sym(35),c,iy(41),jx(41)
      common /side/xb(900),yb(900),ll(50)
      common /fpts/xf(1000),yf(1000),zf(1000)
      common /conta/ m,n, dx,dy, xmn,ymn,xmx,ymx
      common /trans/ inpt,iout,jin,ioprint
      common /widect/ naxcol
      common /bodk/ mm,nlim,jbody,short,nend
      dimension a(1),xa(900),ya(900),lll(50)
      data sym/'1','2','3','4','5','6','7','8','9','a','b','c','d','e',
1  'f','g','h','i','j','k','l','m','n','o','p','q','r','s','t',
2  'u','v','w','x','y','z'//,ihor/'-'/,ivert/'|'/,      jplus/'+'/'
c   needs carriage with 132 columns available or change maxcol
      data maxcol,iblack,jex/131,' ','*'/
      if(pc.ea.0.) return 1
      dc=pc
      if (m .lt. 2) go to 650
      if (m .at. 41) go to 650
      if (n .lt. 2) go to 650
      if (n .at. 41) go to 650
      maxcol=maxcol+1
      if (naxcol .ea. 0 .or. naxcol .at. 130) maxcol=131
      nx=m-1
      ny=n-1
      if(jk.ea.0) go to 60
      write (iout,10) title,n,m,nt,ymn,ymx,xmn,xmx,dc
10  format (///,' expect plot of "a18," of "i3," ov "i3,
1  "( "i3,") arra',
2  "y. x=",f8.1," to",f8.1," y=",f8.1," to",f8.1," contoured at",
3  f7.1," milligals",//)
      fmin= 1.0e20
      fmax=-1.0e20
      do 20 i=1,nt
      aij=a(i)
      if (aij .gt. fmax) fmax=aij
      if (aij .lt. fmin) fmin=aij
20  continue
30  fc=fmin/dc
      ic=fc
      if(amod(fc,1.).ne.0.) ic=ic+1
      if (ic .lt. 0) ic=ic-1
      jc=fmax/dc
      if (jc .lt. 0) jc=jc-1
      ncont=jc-ic+1
      if (ncont .lt. 36) go to 50
      write (iout,40) dc
40  format (5x,"requested contour interval of",f9.4, " is doubled",
1  " because more than 35 contours are produced.")

```

```

dc=2.0*dc
do to 30
50 cd=1.0/dc
   ci=dc*ic
   cf=dc*ic
c
60 width=maxcol-1
   ux=nx/width
c number of columns in one grid interval
   fjinc=1.0/ux
   fjincp=0.5+fjinc
   colefm=1.0-fjinc
c maintains true proportions with x expanded to 130 columns
   uy= dx*ux/( dy*0.6)
   if (uy .lt. 1.0) do to 80
   uy=1.0/uy
   write (iout,70) uy
70 format(5x,"y-distances are exaggerated by",f9.4," for contouring")
   uy=1.0
c number of columns per unit distance
80 dlx=fjinc/dx
   fiinc=1.0/uy
   dlv=fiinc/dy
   row=1.0
   do 90 k=1,ny
   row=row+fiinc
90 iy(k)=row+0.5
   lastrw=iy(ny)
   col=1.0
   do 100 k=1,m
   jx(k)=col+fjinc
100 col=col+fjinc
   if (lk .ne. 0) do to 300
c
c   plot body(s) corners
c
   write (iout,110) title,ymn,ymx,xmn,xx
110 format (///," expect plot of 'a18,
1 'x=",f8.1," to",f8.1," y=",f8.1," to",f8.1,/)
   do 120 i=1,nend
   xa(i)=xb(i)
   ya(i)=yb(i)
120 continue
   do 130 i=1,mm
   lll(i)=ll(i)
130 continue
   nb=nend
   jr=1
   do 280 jrow=1,lastrw
   do 140 l=1,maxcol
   pnt(l)=iblack
140 continue
   ie=0
   do 180 i=1,mm
   is=ie+1
   ie=ie+lll(i)
   if(lll(i).eq.0) do to 180
150 ky=1.5001+dlv*(ya(is)-ymn)
   if(ky.ne.jrow) do to 170
   kx=1.5001+dlx*(xa(is)-xmn)
   if(kx.lt.1.or.kx.gt.maxcol) do to 170

```

```

      pnt(kx)=sym(i)
      nb=nb-1
      do 160 k=is,nb
        kk=k+1
        xa(k)=xa(kk)
        va(k)=va(kk)
160    continue
        is=is-1
        ie=ie-1
        lll(i)=lll(i)-1
170    is=is+1
        if(is.le.ie) go to 150
180    continue
        itest=0
        if(jrow .eq. lastrw) go to 220
        if(jrow .eq.      1) go to 220
        if(jrow .eq. iy(jr)) go to 210
        if(pnt(maxcol) .eq. iblank) pnt(maxcol)=ivert
        if(pnt(      1) .eq. iblank) pnt(      1)=ivert
190    write (iout,200) (pnt(k),k=1,maxcol)
200    format ( 1x,131a1)
        go to 280
210    itest=1
        jr=1+jr
220    do 230 k=1,m
        kx=jx(k)
        if(pnt(kx) .eq. iblank) pnt(kx)=jplus
230    continue
        if(itest .eq. 1) go to 190
        do 240 k=1,maxcol
        if(pnt(      k) .eq. iblank) pnt(      k)=ihor
240    continue
        if(jrow .eq. lastrw) go to 290
        if(maxcol.eq.131) go to 260
        write(iout,250)
250    format("1")
        write (iout,200) (pnt(k),k=1,maxcol)
        go to 280
260    write(iout,270) (pnt(k),k=1,maxcol)
270    format("1",131a1)
280    continue
290    write (iout,200) (pnt(k),k=1,maxcol)
        return
c next statement used if only first body corners to be plotted
c otherwise all could be plotted
c nb is total number of corners of first body of model here
300    nb= ll(1)
        nf=1
        l=0
310    do 320 k=nf,nb
        kk=k
        dum=xb(k)
        if(dum .lt. xmn) go to 330
        if(dum .gt. xmx) go to 330
        kx=1.5001+d1x*(dum-xmn)
        dum=yb(k)
        if(dum .lt. ymn) go to 330
        if(dum .gt. ymx) go to 330
        ky=1.5001+d1y*(dum-ymn)
        l=l+1

```



```

    xa(1)=kx
320 ya(1)=ky
    do to 350
330 nf=kk+1
    write (iout,340) vb(kk),xh(kk)
340 format (5x,"body point x=",f10.3," y=",f10.3," is outside",
    1 " plot boundary.")
    if (kk.lt. nb) do to 310
350 nb=1
360 cm=ci-dc
    do 370 i=1,nt
c changes values inputted as 'a' (restored later).
370 a(i)=cd*(a(i)-cm)
    ir=1
c loop of rows from top to bottom
    do 580 jrow=1,lastrw
    row=1.0+uy*(jrow-1)
c major grid index (i) increases once each fiinc rows
    i=row
    ful=row-i
    ind=0
    kt=i-n
    do 380 l = 1,maxcol
380 pnt(l) = iblank
    coleft=coleftm
c loop of grid intervals from left to right
    do 460 j=1,nx
    coleft=coleft+fjinc
    if (jrow .eq. lastrw) go to 390
    kt=kt+n
    kb=kt+n
    dum = a(kt)
    zl = dum + ful*(a(kt+1) -dum)
    dum = a(kb)
    zr = dum + ful*(a(kb+1) -dum)
    go to 400
390 ind=ind+n
    zl=a(ind)
    kb=ind+n
    zr=a(kb)
400 den = zr-zl
    if (abs(den) .at. 0.000001) go to 420
    izc=zl
    fl=izc
    if (fl .ne. zl) go to 460
    left=coleft+0.5
    irt=coleft+fjinco
c loop of contours in this grid interval
    do 410 l=left,irt
410 pnt(l)=sym(izc)
    go to 460
420 if (zl .at. zr) go to 430
    fmin=zl
    fmax=zr
    go to 440
430 fmin=zr
    fmax=zl
440 lmin=fmin+0.9999
    lmax=fmax
    if (lmax .eq. 0) go to 460

```

```

    if (lmax .lt. lmin) go to 460
    ratio=fjinc/den
    do 450 izc=lmin,lmax
    c=coleft+ratio*(izc-zl)
450 pnt(c)=sym(izc)
460 continue
    if (nb .eq. 0) go to 510
    row=ir0w
    nf=1
470 do 480 k=nf,nb
    kk=k
    if (ya(k) .eq. row) go to 490
480 continue
    go to 510
490 nb=nb-1
    kx=xa(kk)+0.5
    pnt(kx)=jex
    if (kk .at. nb) go to 510
    nf=kk
    do 500 k=nf,nb
    kp=k+1
    xa(k)=xa(kp)
500 ya(k)=ya(kp)
    go to 470
510 itest=0
    if (jrow .eq. lastrw) go to 540
    if (jrow .eq. 1) go to 540
    if (jrow .eq. iy(jr)) go to 530
    if (pnt(maxcol) .eq. iblank) pnt(maxcol)=ivert
    if (pnt( 1) .eq. iblank) pnt( 1)=ivert
520 write (iout,200) (pnt(k),k=1,maxcol)
    go to 580
530 itest=1
    jr=1+jr
540 do 550 k=1,m
    kx=jx(k)
    if (pnt(kx) .eq. iblank) pnt(kx)=jplus
550 continue
    if (itest .eq. 1) go to 520
    do 560 k=1,maxcol
    if (pnt( k) .eq. iblank) pnt( k)=ihor
560 continue
    if (jrow .eq. lastrw) go to 590
    if(maxcol.eq.131) go to -570
    write(iout,250)
    write (iout,200) (pnt(k),k=1,maxcol)
    go to 580
570 write(iout,270) (pnt(k),k=1,maxcol)
580 continue
590 write (iout,200) (pnt(k),k=1,maxcol)
    write (iout,600)
600 format('/',, ' symbol',8x,
    1 'value (milliads). y increases',
    2 ' to right and x increases downward')
    zc = ci-dc
    do 610 i=1,ncont
    zc=zc+dc
610 write (iout,620) sym(i),zc
620 format(' ',3x,a1,6x,f14.4)
    write (iout,630)

```

```

630 format(4x,"+",6x,"fieldpoint",14x,"*",6x,"corner of body")
do 640 i=1,nt
640 a(i)=cm+dc*a(i)
return
650 write (ioprnt,660) m,n
660 format("0","grid of",i4," by",i4," (x,y) is too large (41) or ",
1 "too small for printer contour.")
return
end
c*****
subroutine ptsm(nots,*)
c
c input data subroutine for user formatted or gridded data
c r.aodson = 3/81
c
character*4 unix,uniz
character*8 name,date1
character*40 idplot
character*50 ifile,ifile2,cfile
character*64 id,id1
character*80 ifmt
common/mplot/ymin,xmin,ymax,xmax
common/fpts/xf(1000),yf(1000),zf(1000)
common/aravinp/iplotr,ibody,ifile,ifile2,lsqs,
1 datum,xscale,dc,unix,uniz,name,height,
2 ifmt,idplot,iobs,icalc,ires
common/conta/nrow,ncol,dv,dx,yo,xo,xma,yma
common/field/r(1000)
common/trans/ inpt,iout,jin,ioprnt
dimension z(1001),z2(1001)
data in/10/,in2/11/,nrcmax/1000/
write(iout,10)
10 format(//,14x,"observed values",/,
1 7x,"x",10x,"y",10x,"z",10x,"a",/)
if(name.eq."gridded") go to 50
c
c user formatted data
c
open(in,file=ifile,form="formatted")
i=1
nz=1
20 read(in,ifmt,end=40) yf(i),xf(i),r(i),zf(i)
write(iout,30) yf(i),xf(i),zf(i),r(i)
30 format(1x,4f11.4)
if(xf(i).lt.xmin) xmin=xf(i)
if(xf(i).gt.xmax) xmax=xf(i)
if(yf(i).lt.ymin) ymin=yf(i)
if(yf(i).gt.ymax) ymax=yf(i)
i=i+1
go to 20
40 k=i-1
go to 230
c
c gridded data
c
50 open(in,file=ifile)
read(in,end=190) id,ncol,nrow,nz,xo,dx,yo,dy
if((ncol*nrow).gt.nrcmax) go to 110
if(dx.eq.0..or.dy.eq.0.) go to 130
if(nz.ne.1) go to 170

```

```

if(ifile2.ea." ") go to 60
open(in2,file=ifile2)
read(in2,end=190) id1,ncol1,nrow1,nz1,xo1,dx1,yo1,dy1
if(ncol.ne.ncol1.or.nrow.ne.nrow1) go to 90
if(dx.ne.dx1.or.dy.ne.dy1.or.xo.ne.xo1.or.yo.ne.yo1) go to 90
if(nz.ne.nz1) go to 90
60 mcol=ncol+1
xol=xo-(2*dx)
yol=yo-dy
k=0
do 80 i=1,nrow
call rowin(z,mcol,in)
if(ifile2.ne." ") call rowin(z2,mcol,in2)
do 70 j=2,mcol
k=k+1
r(k)=z(j)
if(ifile2.ne." ") height=z2(j)
zf(k)=height
yf(k)=xol+(j*dx)
xf(k)=yol+(i*dy)
write(iout,30) yf(k),xf(k),zf(k),r(k)
70 continue
80 continue
if(xf(1).lt.xmin) xmin=xf(1)
if(xf(k).gt.xmax) xmax=xf(k)
if(yf(1).lt.ymin) ymin=yf(1)
if(yf(k).gt.vmax) ymax=yf(k)
xma=xf(k)
vma=yf(k)
go to 230
90 write(iprint,100)
100 format(" parameters on two input standard files do not match")
go to 260
110 write(iprint,120) nrcmax
120 format(" number of grid points .at. allowed maximum of ",i4)
go to 260
130 write(iprint,140)
140 format(" present implementation does not allow unequal",
1 " grid intervals")
go to 260
150 write(iprint,160) nrcmax
160 format(" the number of columns * number of rows requested is",
1 " greater than the maximum allowed of ",i5)
go to 260
170 write(iprint,180)
180 format(" only single precision values allowed")
go to 260
190 write(iprint,200)
200 format(" end of file while reading input data")
go to 260
210 write(iprint,220) k,nrcmax
220 format(" number of formatted data values",i5,
1 " is greater than the allowed number of ",i5)
go to 260
230 npts=k
if(npts.lt.1) go to 240
ierr=0
go to 270
240 write(iprint,250)
250 format(" number of field points less than 1")

```

```

260 close(in)
   if(ifile2.ne." ") close(in2)
   return 1
270 close(in)
   if(ifile2.ne." ") close(in2)
   return
   end
c*****
   subroutine rowin(z,mcol,in)
   dimension z(mcol)
   read(in) z
   return
   end
c*****
   subroutine arav3dplot(npts,*)
c
c   plotting routine for body corners & fieldpoints
c   r.aodson - 4/81
c
   external char(descriptors),pplot(descriptors),araph2(descriptors),
1  xaxis(descriptors),yaxis(descriptors)
   character*4 unix,uniz
   character*8 name
   character*40 idplot
   character*50 ifile,ifile2,cfile
   character*80 ifmt
   common/mplot/xmin,ymin,xmax,ymax
   common/fpts/xf(1000),yf(1000),zf(1000)
   common/aravinp/iplotr,ibody,ifile,ifile2,lsqs,
1  datum,xscale,dc,unix,uniz,name,height,
2  ifmt,idplot,iobs,icalc,ires
   common /bodk/ mm,nlim,jbody,short,nend
   common /side/xh(900),yb(900), ll(50)
   common /trans/ inpt,iout,jin,iprint
   common/contq/nrow,ncol,dv,dx,yo,xo,vma,xma
   character*1 revx,svm(50)
   dimension xd(2),yd(2),xs(4),ys(4),jdim(1),
1  ifmt1(2),nscale(5),xxb(900),yyb(900)
   data jdim(1)/0/,nscale/1,0,2,0,2/,ifmt1/("f10",".1) "/,
1  nsh1/6/,nsv1/10/,nvl/12/,
2  nx/2/,nv/2/,vexaa/1./,
3  sym/"1","2","3","4","5","6","7","8","9","a","b","c","d","e",
4  "f","g","h","i","j","k","l","m","n","o","p","q","r","s",
5  "t","u","v","w","x","y","z",15*"*/
c
c
c   denver plotting system
c
10  if(xscale.eq.0.) go to 170
   if(iplotr.eq.1.or.iplotr.eq.4) go to 20
   nx1=(xmax-xmin)/xscale*2.+5
   if(nx1.lt.2) nx1=2
   ny1=nx1
   go to 30
20  nx1=24
   ny1=17
30  call scale3(xmin,xmax,nx1,xminp,xmaxp,xdel)
   call scale3(ymin,ymax,ny1,yminp,ymaxp,ydel)
   xd(1)=xminp
   xd(2)=xmaxp

```

```

    vd(1)=vminp
    vd(2)=vmaxp
    xs(1)=(xmaxp-xminp)/xscale
    ys(1)=(ymaxp-yminp)/yscale
    if(ys(1).gt.18.) vs(1)=18.
    do to 60
40  nyl=(ymax-ymin)/yscale*vexaq*4.+5
    if(nyl.lt.2) nyl=2
50  call scale3(vmin,ymax,nyl,vminp,vmaxp,vdel)
    vd(1)=vmaxp
    vd(2)=vminp
    xs(1)=(xmaxp-xminp)/yscale
    ys(1)=(ymaxp-yminp)/yscale*vexaq
60  call pltset(iplotr,xs(4),vs(4),jdim)
    vyss=vs(4)
    xs(2)=0.
    ys(2)=0.
    xs(3)=2.
    vs(3)=2.
    if((xs(1)+xs(3)+2).at.xs(4)) do to 70
    if((ys(1)+vs(3)+2).at.vs(4)) do to 70
    do to 100
70  xs(4)=amin1(xs(4),13.)
80  vs(4)=amin1(vs(4),10.)
90  xs(3)=1.0
    vs(3)=1.0
    xs(1)=xs(4)-xs(3)-.2
    ys(1)=vs(4)-vs(3)-.2
    do to 110
100 xs(4)=xs(1)+xs(3)+.2
    ys(4)=vs(1)+vs(3)+.2
110 call scale(xd,yd,xs,ys,4,icode)
    if(icode.lt.0) do to 150
    call neat1
    call xaxis(xd,yd,xs,xdel,2,.08,"(f10.2)",10)
    call yaxis(vd,yd,ys,ydel,2,.08,"(f10.2)",10)
    if(iplotr.ne.1.and.iplotr.ne.4)
1  call char(2.,.1,idplot,40,3,.15,0.,0.,0.)
120 ie=0
c
c  plot body corners
c
    do 140 i=1,mm
    is=ie+1
    ie=ie+11(i)
    call char(vb(is),xb(is),sym(i),1,0,.08,0.,0.,0.)
    m=0
    do 130 k=is,ie
    m=m+1
    xxb(m)=xb(k)
    vyb(m)=vb(k)
130 continue
    xxh(m+1)=xb(is)
    vyb(m+1)=vb(is)
    call line(vyb,xxh,m+1,0,0)
140 continue
c
c  plot fieldpoints
c
    do 145 i=1,npts

```

```

    call char(vf(i),xf(i),"+",1,0,.08,0.,0.,0.)
145 continue
    call endot(idim)
    return
150 write(iprint,160)
160 format(" unable to scale plotter")
    go to 190
170 write(iprint,180)
180 format(" xscale is zero")
190 return 1
    end
c*****
    subroutine stdout(isw,z)
c
c    subroutine to create standard file
c    r.godson - 4/81
c
    character*32 id1(2),id2(2)
    character*54 file2
    character*55 file1
    character*56 file3
    dimension z(1)
    common/st/file1,file2,file3
    common/gravinp/inlotr,ibody,ifile,ifile2,lsqs,
1 datum,xscale,dc,unix,uniz,name,height,
2 ifmt,idplot,iobs,icalc,ires
    common/conta/nrow,ncol,dy,dx,xmin,vmin,xmax,vmax
    data iout/11/,nz/1/,dum/0./,
1 id1/"gravpoly calculated values from ",
2 "model          "/,
3 id2/"gravpoly residual values(observe",
4 "d-calculated   "/
    if(isw.eq.1) open(iout,file=file1)
    if(isw.eq.2) open(iout,file=file2)
    if(isw.eq.1) write(iout) id1,ncol,nrow,nz,vmin,dx,xmin,dy
    if(isw.eq.2) write(iout) id2,ncol,nrow,nz,vmin,dx,xmin,dy
    is=1
    ie=ncol
    do 10 i=1,nrow
    write(iout) dum,(z(j),j=is,ie)
    is=ie+1
    ie=ie+ncol
10 continue
    close(iout)
    return
    end

```