

Appendix A

Thermo-physical Properties of the S355 Grade Steel

T (°C)	Fraction liquid (wt%)	Density (g/cm ³)	Thermal conductivity [W/(m K)]	Electrical resistivity (10e-6 Ohm m)	Young's modulus (GPa)
1600	100	6.91218	34.91665	1.31434	0
1580	100	6.92858	34.54952	1.31412	0
1560	100	6.94487	34.18239	1.3139	0
1540	100	6.96107	33.81526	1.31367	0
1520	100	6.97717	33.44813	1.31344	0
1513.7	99.99996	6.98222	33.33249	1.31337	1.13E-24
1500	46.85386	7.13721	36.00007	1.20672	6.38733
1486.24	25.80748	7.20396	37.09981	1.16187	25.52687
1483.93	10.7205	7.25349	35.0414	1.2285	46.73048
1480	6.46744	7.2678	35.1107	1.22334	57.04138
1465.4	0.19412	7.29098	35.10308	1.21341	76.17767
1465.22	4.91E-05	7.29142	35.10561	1.2132	76.81795
1460		7.29426	35.03964	1.21183	77.25785
1420		7.31594	34.53488	1.20117	80.62455
1400		7.32673	34.28274	1.19571	82.30757
1360		7.34827	33.77882	1.18453	85.67304
1340		7.35903	33.52701	1.17882	87.35554
1300		7.38052	33.02361	1.16711	90.72018
1280		7.39127	32.772	1.16112	92.40234
1240		7.41278	32.26891	1.14885	95.76642
1200		7.43433	31.76595	1.13619	99.13026
1180		7.44512	31.5145	1.12971	100.81211
1140		7.46675	31.01165	1.11642	104.17572
1120		7.47759	30.76024	1.10962	105.85749
1080		7.49932	30.25745	1.09567	109.22097

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T (°C)	Fraction liquid (wt%)	Density (g/cm ³)	Thermal conductivity [W/(m K)]	Electrical resistivity (10e-6 Ohm m)	Young's modulus (GPa)
1040		7.52113	29.75469	1.08125	112.5844
1020		7.53206	29.50331	1.07385	114.2661
980		7.55399	29.00056	1.05868	117.6295
940		7.57599	28.49781	1.04296	120.99288
900		7.59809	27.99507	1.02669	124.35625
880		7.60917	27.7437	1.01833	126.03793
840		7.6314	27.24096	1.00115	129.4013
822.87		7.64095	27.02563	0.99359	130.84191
820		7.63921	27.19745	0.98473	131.75148
780		7.62239	29.24646	0.88223	141.69614
760		7.61956	30.00841	0.8435	145.34667
720		7.62059	31.21428	0.77952	151.26621
706.93		7.62221	31.54512	0.7612	152.93539
700		7.61719	32.55451	0.73238	155.69517
693.09		7.61436	33.32865	0.71029	157.88905
660		7.62712	33.61721	0.68007	160.57609
640		7.63474	33.8191	0.66152	162.19246
600		7.64982	34.2856	0.62394	165.40907
580		7.65727	34.55049	0.60497	167.00971
540		7.67204	35.14422	0.56687	170.19637
520		7.67936	35.47328	0.5478	171.78252
480.87		7.69356	36.1795	0.5106	174.87229
480		7.69389	36.20218	0.5097	174.94655
440		7.70932	37.31181	0.46827	178.40557
431.86		7.71246	37.54104	0.4601	179.09951
420		7.71658	37.80282	0.44923	180.02249
380		7.73035	38.74111	0.41305	183.11825
360		7.73716	39.24242	0.39529	184.65566
320		7.7506	40.3099	0.36051	187.70846
300		7.75724	40.87627	0.34353	189.2233
260		7.77033	42.07472	0.31045	192.22824
240		7.77678	42.70698	0.29438	193.71764
200		7.7895	44.03823	0.26323	196.66875
180		7.79577	44.73759	0.24816	198.13017
140		7.80812	46.2054	0.21907	201.02563
120		7.81421	46.97471	0.20505	202.46062
80		7.82621	48.58644	0.17808	205.30793
68.52		7.82962	49.06718	0.1706	206.1196

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T (°C)	Fraction liquid (wt%)	Density (g/cm ³)	Thermal conductivity [W/(m K)]	Electrical resistivity (10e-6 Ohm m)	Young's modulus (GPa)
40		7.83811	50.27003	0.15262	208.11896
25		7.84248	50.93719	0.14341	209.1696

T (°C)	Poisson's ratio	Liquid viscosity (mPa s)	Enthalpy (J/g)	Specific heat [J/(g K)]	Latent heat (J/g)
1600	0.49992	5.66944	1341.7645	0.82688	
1580	0.49992	5.84298	1325.2278	0.82684	
1560	0.49992	6.0258	1308.6915	0.82682	
1540	0.49992	6.21856	1292.1556	0.82679	
1520	0.49992	6.42199	1275.7131	0.81806	
1513.7	0.49992	6.48841	1270.5687		50.97787
1500	0.42097	6.44337	1131.8002	6.05028	96.93005
1486.24	0.3882	6.40566	1072.0753		128.23126
1483.93	0.39375	6.40141	1015.0818		150.54388
1480	0.38815	6.38227	1001.5291	2.97105	167.01298
1465.4	0.37924	6.32939	975.70594		211.36935
1465.22	0.37898	6.32857	975.06483		241.82898
1460	0.37868		971.38423	0.70351	
1420	0.37633		943.38307	0.69556	
1400	0.37515		929.49892	0.69173	
1360	0.3728		901.9554	0.68431	
1340	0.37162		888.29294	0.6807	
1300	0.36925		861.18047	0.67365	
1280	0.36807		847.72861	0.67016	
1240	0.36571		821.03032	0.66333	
1200	0.36334		794.59762	0.65679	
1180	0.36216		781.47837	0.65355	
1140	0.35979		755.43174	0.64711	
1120	0.35861		742.50364	0.64393	
1080	0.35624		716.83646	0.63761	
1040	0.35387		691.41987	0.63129	
1020	0.35269		678.80514	0.62816	
980	0.35032		653.76217	0.62191	
940	0.34796		628.96727	0.61568	
900	0.34559		604.41988	0.60943	

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T (°C)	Poisson's ratio	Liquid viscosity (mPa s)	Enthalpy (J/g)	Specific heat [J/(g K)]	Latent heat (J/g)
880	0.3444		592.23888	0.60632	
840	0.34204		568.06213	0.60011	
822.87	0.34102		557.78226		
820	0.3401		554.9352	0.99814	
780	0.32991		514.98116	0.99243	
760	0.32655		495.06587	1.00289	
720	0.32173		452.16563	1.0585	
706.93	0.32051		438.48967		
700	0.31745		421.67167	2.16441	
693.09	0.3153		408.16114		
660	0.31401		377.71663	0.8885	
640	0.31324		360.24442	0.85654	
600	0.3117		327.04394	0.80251	
580	0.31093		311.19806	0.77943	
540	0.30939		280.78854	0.73914	
520	0.30862		266.15472	0.72132	
480.87	0.30712		238.49024		
480	0.30708		237.85883	0.72682	
440	0.3052		209.02962	0.70229	
431.86	0.30482		203.33605		
420	0.30436		195.61239	0.64602	
380	0.30281		170.2417	0.62038	
360	0.30203		157.93091	0.6083	
320	0.30047		134.01322	0.5859	
300	0.29969		122.39202	0.57492	
260	0.29813		99.79758	0.55366	
240	0.29735		88.81615	0.54331	
200	0.29579		67.46756	0.52339	
180	0.29501		57.09232	0.51364	
140	0.29345		36.9104	0.49522	
120	0.29267		27.08885	0.48655	
80	0.29112		7.95299	0.46994	
68.52	0.29067		2.58538		
40	0.28956			0.45329	
25	0.28897			0.44669	

Appendix B

Thermo-physical Properties of the C45 Grade Steel

T (°C)	Fraction liquid (wt)	Density (g/cm ³)	Thermal conductivity [W/(m K)]	Electrical resistivity (10e-6 Ohm m)	Young's modulus (GPa)
1600	1	6.91435	35.34082	1.29856	0
1550	1	6.95495	34.4199	1.29771	0
1505	1	6.99096	33.59108	1.29691	0
1500	1	6.99493	33.49898	1.29682	0
1495	1	6.99889	33.40689	1.29673	0
1494.79	1	6.99906	33.40304	1.29672	0
1490	0.87841	7.03448	33.97108	1.27159	0.0165101
1489.88	0.87555	7.03532	33.9845	1.271	0.0181056
1489.88	0.87553	7.03534	33.98456	1.271	0.0181014
1485	0.7273	7.07865	34.14199	1.26163	0.0941919
1480	0.60769	7.11483	34.26067	1.25369	0.84088
1475	0.51111	7.14508	34.34632	1.24699	2.86007
1470	0.43136	7.17097	34.40652	1.24125	6.27462
1465	0.36426	7.19354	34.44713	1.23623	10.88971
1460	0.30683	7.2136	34.47117	1.23182	16.42427
1455	0.25701	7.23165	34.48333	1.22783	22.61462
1450	0.21335	7.24802	34.48592	1.22419	29.25093
1445	0.17473	7.26307	34.48068	1.22082	36.18725
1440	0.14025	7.277	34.46906	1.21768	43.32377
1435	0.10925	7.28997	34.45225	1.21472	50.59053
1430	0.0811754	7.30214	34.43121	1.2119	57.94115
1425	0.0556055	7.31363	34.40654	1.20921	65.33844
1420	0.0321876	7.32452	34.37859	1.20663	72.758
1415	0.0106378	7.33489	34.34723	1.20416	80.18024
1412.42	0.0061221	7.34013	34.32971	1.20293	84.06567

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T (°C)	Fraction liquid (wt)	Density (g/cm ³)	Thermal conductivity [W/(m K)]	Electrical resistivity (10e-6 Ohm m)	Young's modulus (GPa)
1405		7.34391	34.23823	1.20084	84.69597
1400		7.34646	34.17663	1.19942	85.12081
1370		7.36177	33.80816	1.19075	87.66882
1340		7.3771	33.44172	1.18182	90.21507
1310		7.39246	33.07744	1.17262	92.75958
1280		7.40785	32.71543	1.16313	95.30232
1250		7.42327	32.35579	1.15334	97.8433
1220		7.43871	31.99862	1.14324	100.38251
1190		7.45417	31.64402	1.13283	102.91995
1160		7.46967	31.29205	1.12208	105.45561
1130		7.48519	30.9428	1.11099	107.98949
1100		7.50075	30.59631	1.09955	110.52159
1070		7.51633	30.25262	1.08775	113.05189
1040		7.53194	29.91177	1.07557	115.5804
1000		7.5528	29.46171	1.05874	118.94894
960		7.57372	29.01666	1.0412	122.31426
920		7.59469	28.57653	1.02294	125.67635
900		7.6052	28.35825	1.01354	127.35618
860		7.62625	27.92515	0.99416	130.71339
830		7.64208	27.60321	0.97913	133.22916
800		7.65795	27.28359	0.96366	135.74307
770		7.67385	26.96612	0.94775	138.25514
765		7.6765	26.91341	0.94506	138.67363
760		7.67915	26.86075	0.94235	139.09208
755		7.68181	26.80814	0.93963	139.51047
750		7.68446	26.75559	0.93689	139.92881
745		7.68712	26.70308	0.93415	140.3471
740		7.68978	26.65063	0.93139	140.76534
735		7.69244	26.59822	0.92862	141.18352
730		7.6951	26.54587	0.92584	141.60166
725		7.69776	26.49356	0.92304	142.01974
720		7.70042	26.4413	0.92023	142.43777
715		7.70308	26.38908	0.91741	142.85575
710		7.70575	26.33691	0.91458	143.27367
705		7.70841	26.28478	0.91173	143.69154
700		7.71108	26.23269	0.90887	144.10937
695		7.71374	26.18064	0.906	144.52713
690		7.71641	26.12863	0.90312	144.94485

(continued)

(continued)

T (°C)	Fraction liquid (wt)	Density (g/cm ³)	Thermal conductivity [W/(m K)]	Electrical resistivity (10e-6 Ohm m)	Young's modulus (GPa)
685		7.71908	26.07666	0.90022	145.36252
680		7.72175	26.02473	0.89731	145.78013
675		7.72442	25.97284	0.89438	146.19769
670		7.72709	25.92098	0.89145	146.61519
665		7.72977	25.86915	0.8885	147.03265
660		7.73244	25.81736	0.88553	147.45005
655		7.73511	25.7656	0.88256	147.8674
650		7.73779	25.71387	0.87957	148.2847
645		7.74047	25.66217	0.87657	148.70194
640		7.74314	25.6105	0.87355	149.11913
635		7.74582	25.55885	0.87053	149.53627
630		7.7485	25.50723	0.86749	149.95336
615		7.75655	25.35252	0.85828	151.2043
610		7.75923	25.30099	0.85519	151.62118
600		7.7646	25.19799	0.84896	152.45477

T (°C)	Poisson's ratio	Liquid viscosity (mPa s)	Enthalpy (J/g)	Specific heat [J/(g K)]	Latent heat (J/g)
1600	0.49968	5.44574	1351.003		
1550	0.49968	5.87455	1309.562	0.8288	
1505	0.49968	6.31224	1272.547	0.814	
1500	0.49968	6.36427	1268.489	0.8116	
1495	0.49968	6.41703	1264.442	0.8094	
1494.79	0.49968	6.41926	1264.274	0.80371	
1490	0.48196	6.40787	1229.413	7.2764	28.871
1489.88	0.48155	6.40771	1228.587		29.554
1489.88	0.48154	6.40771	1228.638		29.502
1485	0.46441	6.37848	1184.669	8.9488	67.637
1480	0.45033	6.35021	1148.982	7.1374	98.081
1475	0.43876	6.32358	1119.988	5.7988	122.357
1470	0.42905	6.29814	1095.888	4.82	142.121
1465	0.42075	6.27418	1075.471	4.0834	158.489
1460	0.41354	6.25124	1057.88	3.5182	172.255
1455	0.40719	6.22958	1042.516	3.0728	183.969
1450	0.40155	6.20922	1028.956	2.712	194.021

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T (°C)	Poisson's ratio	Liquid viscosity (mPa s)	Enthalpy (J/g)	Specific heat [J/(g K)]	Latent heat (J/g)
1445	0.39648	6.18996	1016.866	2.418	202.715
1440	0.39189	6.17154	1005.992	2.1748	210.288
1435	0.38771	6.15421	996.137	1.971	216.921
1430	0.38386	6.13801	987.143	1.7988	222.76
1425	0.38031	6.1227	978.883	1.652	227.919
1420	0.37701	6.10826	971.257	1.5252	232.494
1415	0.37393	6.09474	964.18	1.4154	236.562
1412.42	0.3724	6.06324	960.689	1.35151	238.52
1405	0.37197		955.515	0.69759	
1400	0.37169		952.036	0.6958	
1370	0.36997		931.247	0.691	
1340	0.36825		910.595	0.6864	
1310	0.36654		890.077	0.682	
1280	0.36482		869.693	0.6776	
1250	0.3631		849.442	0.6732	
1220	0.36139		829.321	0.6688	
1190	0.35967		809.328	0.6648	
1160	0.35795		789.463	0.6604	
1130	0.35624		769.721	0.6564	
1100	0.35452		750.102	0.6522	
1070	0.3528		730.6	0.6484	
1040	0.35109		711.212	0.6448	
1000	0.3488		685.531	0.64	
960	0.34651		660.03	0.6356	
920	0.34422		634.687	0.632	
900	0.34308		622.067	0.6304	
860	0.34079		596.902	0.6284	
830	0.33908		578.066	0.6278	
800	0.33736		559.217	0.6288	
770	0.33564		540.288	0.633	
765	0.33536		537.119	0.6338	
760	0.33507		533.878	0.6482	
755	0.33479		530.629	0.6498	
750	0.3345		527.391	0.6476	
745	0.33421		524.166	0.645	
740	0.33393		520.952	0.6428	
735	0.33364		517.748	0.6408	
730	0.33336		514.556	0.6384	

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T (°C)	Poisson's ratio	Liquid viscosity (mPa s)	Enthalpy (J/g)	Specific heat [J/(g K)]	Latent heat (J/g)
725	0.33307		511.374	0.6364	
720	0.33278		508.203	0.6342	
715	0.3325		505.042	0.6322	
710	0.33221		501.891	0.6302	
705	0.33193		498.749	0.6284	
700	0.33164		495.617	0.6264	
695	0.33135		492.495	0.6244	
690	0.33107		489.382	0.6226	
685	0.33078		486.277	0.621	
680	0.3305		483.182	0.619	
675	0.33021		480.095	0.6174	
670	0.32993		477.017	0.6156	
665	0.32964		473.947	0.614	
660	0.32935		470.886	0.6122	
655	0.32907		467.832	0.6108	
650	0.32878		464.787	0.609	
645	0.3285		461.75	0.6074	
640	0.32821		458.72	0.606	
635	0.32792		455.698	0.6044	
630	0.32764		452.684	0.6028	
615	0.32678		443.686	0.5984	
610	0.32649		440.701	0.597	
600	0.32592		434.753	0.5942	

Appendix C

Complete Source Code: Steady Heat Flow

As an example, let's consider the area presented in Fig. C.1, which is a cross-section of e.g. a rolled strand with dimensions of 0.5×0.5 m. The left, right and top surface of the cast strand are heated by the medium with a temperature 500°C . On the other hand the bottom surface of the cast strand is cooled by the medium with a temperature 20°C . The heat transfer coefficient for all boundary conditions assumes the same value of $\alpha = 100(\text{W}/\text{m}^2\text{K})$, and heat transfer coefficients of $\lambda_x = \lambda_y = 78(\text{W}/\text{mK})$. The power of heat sources referred to the volume unit in the analysed case $q_{vol} = 0(\text{W}/\text{m}^3)$.

Below find a complete print-out of the source code written in the C++ language, simulating the steady state condition of heat transfer in the cross-section. The comments placed within the code will enable the reader to analyse the numerical code individually, and will be helpful during potential modifications implemented by the user, or when creating original solutions. The program code may be

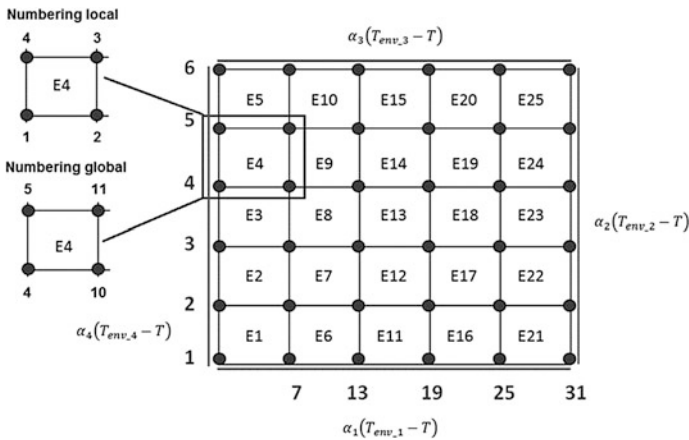


Fig. C. 1 A cross-section of a cast strand with defined third type boundary conditions

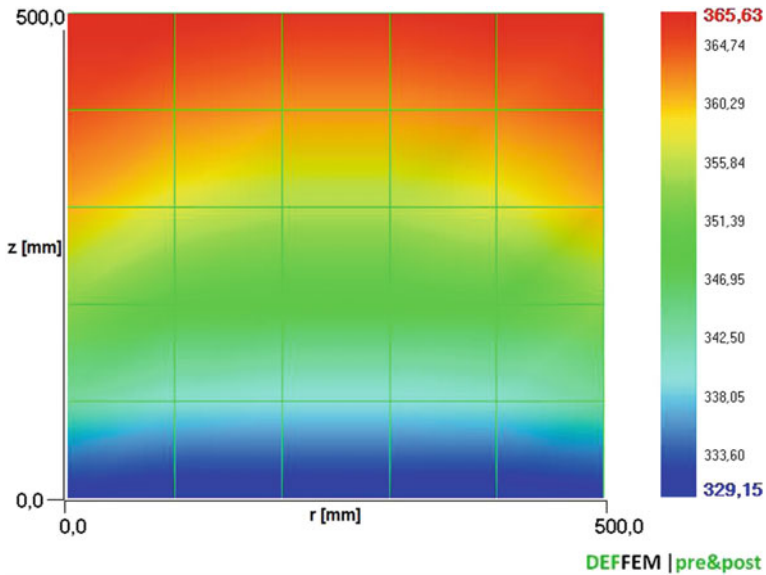


Fig. C.2 The temperature field distribution for the case analysed (steady state condition)

compiled with any C++ compiler (e.g. Dev-C++ 5.11). The final result will be the temperature field distribution for the steady state condition (test data) (Fig. C.2).

```
#include <iostream>
#include <fstream>

main()
{
float a[220][220], x_w[251], zamiana, stala, b, max;
int nlh, nlv, nw, ne, ie[1000][5];
int l, licznik, k, i, j, u, ii, jj;
float r, h, lambdaX, lambdaY, moc_zrodel, strumien;
float t_czynnikal, t_czynnika2, t_czynnika3, t_czynnika4;
float alfa1, alfa2, alfa3, alfa4;
float dx, dy, dv, dh, wx[1000], wy[1000], x[1000], y[1000], xe[500], ye[500];
float Kc_e[5][5], Kalfa_e[5][5];
float Kc[251][251], Kalfa[251][251], K[251][251], f[251][2];
float fq_duze[251][2], fq_duze_e[5][2], falfa_e[5][2], falfa[251][2];
float x1, x2, x3, x4, y1, y2, y3, y4;
float czlon_lambda_x, czlon_lambda_y, czlon_alfa_bok23, czlon_alfa_bok41;
float czlon_alfa_bok12, czlon_alfa_bok34;
float czlon_fq_duze, czlon_falfa, czlon_q;
int n, imax, spr=1;
```

```

//----- INPUT PARAMETERS-----
//nlh-the number of horizontal lines in the mesh
//nlv-the number of vertical lines in the mesh
//nw,ne-number of mesh nodes, number of elements
//h,r-height, width section
//dv,dh-increases the length of elements for h and r
//dx,dy-operating variables
//wx[],wy[]-tables containing the coordinates of the mesh
//x[],y[]-coordinates x and y nodes of the mesh
//----- INPUT DATA-----
nlh=6;
nlv=6;
h=500;
r=500;
h=h/1000;
r=r/1000;           //conversion of centimeters to meters
lambdaX=78;        //thermal conductivity, x-direction, W/(m*K)
lambdaY=78;        //thermal conductivity, y-direction, W/(m*K)
alfa1=100;         //heat transfer coefficient, bottom section, W/(m2*K)
alfa2=100;         //heat transfer coefficient, right section, W/(m2*K)
alfa3=100;         //heat transfer coefficient, top section, W/(m2*K)
alfa4=100;         //heat transfer coefficient, left section, W/(m2*K)
moc_zrodel=0;     //power sources of heat per unit volume, (W/m3)
t_czynnika1=20;   //
the temperature of the cooling medium, bottom section, [st. C]
t_czynnika2=500;  //
the temperature of the cooling medium, right section, [st. C]
t_czynnika3=500;  //
the temperature of the cooling medium, top section, [st. C]
t_czynnika4=500;  //
the temperature of the cooling medium, left section, [st. C]
//-----
dx=0;
dy=0;
dv=h/(nlh-1);
dh=r/(nlv-1);
nw=nlh*nlv;
ne=(nlh-1)*(nlv-1);
std::cout<<"Parameters the generated mesh:"<<std::endl;
std::cout<<"Number of elements:"<<" "<<ne<<std::endl;
std::cout<<"Number of nodes:"<<" "<<nw<<std::endl;
std::cout<<std::endl;
std::cout<<"ENTER-START CALCULATION";
std::cin.get();
std::cin.get();

```

```

//----- DETERMINATION OF THE MESH COORDINATES -----
for(int i=1;i<=nlv;i++)
{
wx[i]=dx;
dx=dx+dh;
}
for(int j=1;j<=nlh;j++)
{
wy[j]=dy;
dy=dy+dv;
}
//----- DETERMINING COORDINATES OF A GRID NODES -----
licznik=1;
l=0;
for(int i=1;i<=nw;i++)
{
x[i]=wx[licznik];
l++;
if(l==nlh)
{
licznik++;
l=0;
}}
licznik=1;
for(int i=1;i<=nw;i++)
{
y[i]=wy[licznik];
licznik++;
if(licznik==nlh+1)
{
licznik=1;
}}
//----- DETERMINATION OF MATRIX CONNECTION -----
k=1;
i=1;
for(int w=1;w<=ne;w++)
{
ie[w][1]=i;
ie[w][2]=i+nlh;
ie[w][3]=1+i+nlh;
ie[w][4]=1+i;
k++;
i++;
if(k==nlh)
{

```

```

k=1;
i=i+1;
}}
//-----
//Calculation of the thermal conductivity of the matrix
//Kc_e for one single element and global matrix conductivity Kc
//-----
for (k=1;k<=250;k++)
{
for (l=1;l<=250;l++)
{
Kc[k][l]=0;
}}
//----- the beginning of the loop elements -----
for (u=1;u<=ne;u++)
{
for (j=1;j<=4;j++)
{
xe[j]=x[ie[u][j]];
ye[j]=y[ie[u][j]];
}
x1=xe[1];x2=xe[2];x3=xe[3];x4=xe[4]; //determining coordinates
of a single element
y1=ye[1];y2=ye[2];y3=ye[3];y4=ye[4];
for (k=1;k<=4;k++) //resetting single array Kc_e
{
for (l=1;l<=4;l++)
{
Kc_e[k][l]=0;
}
czlon_lambda_x= (lambdaX/6)*((y4-y1)/((x2-x1)));
czlon_lambda_y= (lambdaY/6)*((x2-x1)/((y4-y1)));
Kc_e[1][1]=(czlon_lambda_x)*2+(czlon_lambda_y)*2;
Kc_e[1][2]=(czlon_lambda_x)*-2+(czlon_lambda_y)*1;
Kc_e[1][3]=(czlon_lambda_x)*-1+(czlon_lambda_y)*-1;
Kc_e[1][4]=(czlon_lambda_x)*1+(czlon_lambda_y)*-2;
Kc_e[2][1]=(czlon_lambda_x)*-2+(czlon_lambda_y)*1;
Kc_e[2][2]=(czlon_lambda_x)*2+(czlon_lambda_y)*2;
Kc_e[2][3]=(czlon_lambda_x)*1+(czlon_lambda_y)*-2;
Kc_e[2][4]=(czlon_lambda_x)*-1+(czlon_lambda_y)*-1;
Kc_e[3][1]=(czlon_lambda_x)*-1+(czlon_lambda_y)*-1;
Kc_e[3][2]=(czlon_lambda_x)*1+(czlon_lambda_y)*-2;
Kc_e[3][3]=(czlon_lambda_x)*2+(czlon_lambda_y)*2;
Kc_e[3][4]=(czlon_lambda_x)*-2+(czlon_lambda_y)*1;
Kc_e[4][1]=(czlon_lambda_x)*1+(czlon_lambda_y)*-2;

```

```

Kc_e[4][2]=(czlon_lambda_x)*-1+(czlon_lambda_y)*-1;
Kc_e[4][3]=(czlon_lambda_x)*-2+(czlon_lambda_y)*1;
Kc_e[4][4]=(czlon_lambda_x)*2+(czlon_lambda_y)*2;
std::cout<<std::endl;
for(i=1;i<=4;i++) //the creation of a global matrix Kc
{
ii=ie[u][i];
for(j=1;j<=4;j++)
{
jj=ie[u][j];
if(jj>0)
Kc[ii][jj]=Kc[ii][jj]+Kc_e[i][j];
}}
//----- the end of the loop elements -----
//Calculation of matrix thermal conductivity taking into account
the convective boundary
conditions Kalfa_e for single element and global matrix Kalfa
//-----
for(k=1;k<=250;k++)
{
for(l=1;l<=250;l++)
{
Kalfa[k][l]=0;
}}
//----- the beginning of the loop elements -----
for(u=1;u<=ne;u++)
{
for(j=1;j<=4;j++)
{
xe[j]=x[ie[u][j]];
ye[j]=y[ie[u][j]];
}
x1=xe[1];x2=xe[2];x3=xe[3];x4=xe[4]; //determining coordinates
of a single element
y1=ye[1];y2=ye[2];y3=ye[3];y4=ye[4];
for(k=1;k<=4;k++) //resetting single array Kalfa_e
{
for(l=1;l<=4;l++)
{
Kalfa_e[k][l]=0;
}}
if(u==21||u==22||u==23||u==24||u==25) //boundary condition,
right section
{
//exchange heat on the side 2-3 according to the local numbering nodes

```

```

czlon_alfa_bok23= (alfa2*(y4-y1))/6;
Kalfa_e[1][1]= 0;
Kalfa_e[1][2]= 0;
Kalfa_e[1][3]= 0;
Kalfa_e[1][4]= 0;
Kalfa_e[2][1]=0;
Kalfa_e[2][2]=czlon_alfa_bok23*2;
Kalfa_e[2][3]=czlon_alfa_bok23*1;
Kalfa_e[2][4]=0;
Kalfa_e[3][1]=0;
Kalfa_e[3][2]=czlon_alfa_bok23*1;
Kalfa_e[3][3]=czlon_alfa_bok23*2;
Kalfa_e[3][4]=0;
Kalfa_e[4][1]=0;
Kalfa_e[4][2]=0;
Kalfa_e[4][3]=0;
Kalfa_e[4][4]=0;
}
if (u==1 || u==2 || u==3 || u==4 || u==5)           //boundary condition, left
section
{
//exchange heat on the side 4-1 according to the local numbering nodes
czlon_alfa_bok41= (alfa4*(y4-y1))/6;
Kalfa_e[1][1]= czlon_alfa_bok41*2;
Kalfa_e[1][2]= 0;
Kalfa_e[1][3]= 0;
Kalfa_e[1][4]= czlon_alfa_bok41*1;
Kalfa_e[2][1]=0;
Kalfa_e[2][2]=0;
Kalfa_e[2][3]=0;
Kalfa_e[2][4]=0;
Kalfa_e[3][1]=0;
Kalfa_e[3][2]=0;
Kalfa_e[3][3]=0;
Kalfa_e[3][4]=0;
Kalfa_e[4][1]=czlon_alfa_bok41*1;
Kalfa_e[4][2]=0;
Kalfa_e[4][3]=0;
Kalfa_e[4][4]=czlon_alfa_bok41*2;
}
if (u==1 || u==6 || u==11 || u==16 || u==21)       //boundary condition,
bottom section
{
//exchange heat on the side 1-2 according to the local numbering nodes
czlon_alfa_bok12= (alfa1*(x2-x1))/6;

```



```

Kalfa_e[1][1]=czlon_alfa_bok12*2;
Kalfa_e[1][2]=czlon_alfa_bok12*1;
Kalfa_e[1][3]=0;
Kalfa_e[1][4]=0;
Kalfa_e[2][1]=czlon_alfa_bok12*1;
Kalfa_e[2][2]=czlon_alfa_bok12*2;
Kalfa_e[2][3]=0;
Kalfa_e[2][4]=0;
Kalfa_e[3][1]=0;
Kalfa_e[3][2]=0;
Kalfa_e[3][3]=0;
Kalfa_e[3][4]=0;
Kalfa_e[4][1]=0;
Kalfa_e[4][2]=0;
Kalfa_e[4][3]=0;
Kalfa_e[4][4]=0;
}
if (u==5 || u==10 || u==15 || u==20 || u==25) //boundary condition,
top section
{
//exchange heat on the side 3-4 according to the local numbering nodes
czlon_alfa_bok34= (alfa3*(x2-x1))/6;
Kalfa_e[1][1]=0;
Kalfa_e[1][2]=0;
Kalfa_e[1][3]=0;
Kalfa_e[1][4]=0;
Kalfa_e[2][1]=0;
Kalfa_e[2][2]=0;
Kalfa_e[2][3]=0;
Kalfa_e[2][4]=0;
Kalfa_e[3][1]=0;
Kalfa_e[3][2]=0;
Kalfa_e[3][3]=czlon_alfa_bok34*2;
Kalfa_e[3][4]=czlon_alfa_bok34*1;
Kalfa_e[4][1]=0;
Kalfa_e[4][2]=0;
Kalfa_e[4][3]=czlon_alfa_bok34*1;
Kalfa_e[4][4]=czlon_alfa_bok34*2;
}
if (u==7 || u==8 || u==9 || u==12 || u==13 || u==14 || u==17 || u==18 || u==19)
{
Kalfa_e[1][1]=0;Kalfa_e[1][2]=0;Kalfa_e[1][3]=0;Kalfa_e[1][4]=0;
Kalfa_e[2][1]=0;Kalfa_e[2][2]=0;Kalfa_e[2][3]=0;Kalfa_e[2][4]=0;
Kalfa_e[3][1]=0;Kalfa_e[3][2]=0;Kalfa_e[3][3]=0;Kalfa_e[3][4]=0;
Kalfa_e[4][1]=0;Kalfa_e[4][2]=0;Kalfa_e[4][3]=0;Kalfa_e[4][4]=0;
}

```

```

}
for (i=1; i<=4; i++)          //the creation of a global matrix Kalfa
{
  ii=ie[u][i];
  for (j=1; j<=4; j++)
  {
    jj=ie[u][j];
    if (jj>0)
    Kalfa[ii][jj]=Kalfa[ii][jj]+Kalfa_e[i][j];
  }
}
//----- the end of the loop elements -----
//-----
//Determination of the final matrix K = Kc+Kalfa
//-----
for (k=1; k<=nw; k++)          //matrix addition Kc i Kalfa
{
  for (l=1; l<=nw; l++)
  {
    K[k][l]=Kc[k][l]+Kalfa[k][l];
  }
}
//-----
//Calculation of the free vector fq_duze_e for single element
//and global vector taking into account
//the power sources of heat per unit volume
//-----
for (k=1; k<=250; k++)          //resetting vector fq_duze
{
  for (l=1; l<=2; l++)
  {
    fq_duze[k][l]=0;
  }
}
//----- the beginning of the loop elements -----
for (u=1; u<=ne; u++)
{
  for (j=1; j<=4; j++)
  {
    xe[j]=x[ie[u][j]];
    ye[j]=y[ie[u][j]];
  }
  x1=xe[1]; x2=xe[2]; x3=xe[3]; x4=xe[4]; //determining coordi-
nates of a single element
  y1=ye[1]; y2=ye[2]; y3=ye[3]; y4=ye[4];
  for (k=1; k<=4; k++)          //resetting vector fq_male_e for sin-
gle element
  {

```

```

for (l=1;l<=1;l++)
{
fq_duze_e[k][l]=0;
}
czlon_fq_duze= ((moc_zrod1)*((x2-x1)*(y4-y1)))/4;
fq_duze_e[1][1]= czlon_fq_duze*1;fq_duze_e[2][1]= czlon_fq_duze*1;
fq_duze_e[3][1]= czlon_fq_duze*1;fq_duze_e[4][1]= czlon_fq_duze*1;
for(i=1;i<=4;i++)      //the creation of a global fq_duze
{
ii=ie[u][i];
if(jj>0)
fq_duze[ii][1]=fq_duze[ii][1]+fq_duze_e[i][1];
}
//----- the end of the loop elements -----
//-----
//Calculation of the free vector falfa_e for single element
//and global vector falfa taking into account convection heat transfer
//-----
for (k=1;k<=250;k++)      //resetting vector falfa
{
for (l=1;l<=2;l++)
{
falfa[k][l]=0;
}
//----- the beginning of the loop elements -----
for(u=1;u<=ne;u++)
{
for(j=1;j<=4;j++)
{
xe[j]=x[ie[u][j]];
ye[j]=y[ie[u][j]];
}
x1=xe[1];x2=xe[2];x3=xe[3];x4=xe[4]; //determining coordinates of a single element
y1=ye[1];y2=ye[2];y3=ye[3];y4=ye[4];
for (k=1;k<=4;k++)      //resetting vector falfa_e for single element
{
for (l=1;l<=1;l++)
{
falfa_e[k][l]=0;
}
}
if (u==21 || u==22 || u==23 || u==24 || u==25) //boundary condition, right section
{
czlon_falfa= ((alfa2*t_czynnika2)*(y4-y1))/2;

```

```

falfa_e[1][1]=0;
falfa_e[2][1]=czlon_falfa*1;
falfa_e[3][1]=czlon_falfa*1;
falfa_e[4][1]=0;
}
if (u==1 || u==2 || u==3 || u==4 || u==5) //boundary condition, left section
{
czlon_falfa= ((alfa4*t_czynnika4)*(y4-y1))/2;
falfa_e[1][1]=czlon_falfa*1;
falfa_e[2][1]=0;
falfa_e[3][1]=0;
falfa_e[4][1]=czlon_falfa*1;
}
if (u==1 || u==6 || u==11 || u==16 || u==21) //boundary condition, bot-
tom section
{
czlon_falfa= ((alfa1*t_czynnika1)*(x2-x1))/2;
falfa_e[1][1]=czlon_falfa*1;
falfa_e[2][1]=czlon_falfa*1;
falfa_e[3][1]=0;
falfa_e[4][1]=0;
}
if (u==5 || u==10 || u==15 || u==20 || u==25) //boundary condition, top sec-
tion
{
czlon_falfa= ((alfa3*t_czynnika3)*(x2-x1))/2;
falfa_e[1][1]=0;
falfa_e[2][1]=0;
falfa_e[3][1]=czlon_falfa*1;
falfa_e[4][1]=czlon_falfa*1;
}
if (u==7 || u==8 || u==9 || u==12 || u==13 || u==14 || u==17 || u==18 || u==19)
{
falfa_e[1][1]=0; falfa_e[2][1]=0;
falfa_e[3][1]=0; falfa_e[4][1]=0;
}
for(i=1; i<=4; i++) //the creation of a global falfa
{
ii=ie[u][i];
if(jj>0)
falfa[ii][1]=falfa[ii][1]+falfa_e[i][1];
}}
//----- the end of the loop elements -----
//-----
// Determination of the End vector f= falfa+fq_duze

```

```

//-----
for (k=1;k<=nw;k++)          //adding a vector
{
for (l=1;l<=1;l++)
{
f[k][l]=fq_duze[k][l]+falfa[k][l];
}}
//-----
// The solution of equations using Gaussian elimination
//-----

    for (i=0;i<nw;i++)
        {
            for (j=0;j<nw;j++)
                {
                    a[i][j]=a[i][j]+K[i+1][j+1];
                }
            a[i][nw]=f[i+1][1];
        }
i=-1;
while (i<nw-1 && spr)
    {
        i++;
        max=a[i][i]*a[i][i];
        imax=i;
        for (j=i+1;j<nw;j++)
            {
                if (max<a[j][i]*a[j][i])
                    {
                        max=a[j][i]*a[j][i];
                        imax=j;
                    }
                &#160;
            }
        if (imax!=i)
            {
                for (j=0;j<nw+1;j++)
                    {
                        zamiana=a[imax][j];
                        a[imax][j]=a[i][j];
                        a[i][j]=zamiana;
                    }
            }
        for (j=i+1;j<nw;j++)
            {
                if (!a[i][i]) spr=0;
                stala=a[j][i]/a[i][i];
            }
    }

```

```

                                for(k=i;k<nw+1;k++) a[j][k]=a[j][k]-sta-
la*a[i][k];
                                }
                                }
                                for (i=nw-1;i>-1;i -----)
                                {
                                b=0;
                                for (j=nw-1;j>i;j--- ) b=b+a[i][j]*x_w[j];
                                x_w[i]=(a[i][nw]-b)/a[i][i];
                                if(!a[i][i]) spr=0;
                                }
std::ofstream stream("Output.txt" );
    if (spr)
    {
        std::cout<<std::endl;
        std::cout<<"Results"<<std::endl;

                                for (i=0;i<nw;i++)
                                {
                                std::cout<<"X["<<i+1<<"] = "<<x_w[i]<<std::endl;
                                stream<<1000*x[i+1]<<"\t"<<1000*y[i+1]<<"\t"<<x_w[i]<<std:::
endl;
                                }
                                }
    else
    {
        std::cout<<" Method Gaussian used in the algorithm solver ";
        std::cout<<"\n it does not give an unambiguous solu-
tion in this case ";
        std::cout<<"\n or the equations does not give an unambigu-
ous solution \n";
    }
getchar();
}

```

Appendix D

Subroutine: Gauss Method

```
PROGRAM gauss_program
IMPLICIT NONE
INTEGER, PARAMETER :: rozmiar = 3
DOUBLE PRECISION, DIMENSION(rozmiar,rozmiar) :: A
DOUBLE PRECISION, DIMENSION(rozmiar) :: b
! Predefined matrix A
A(1,1)=-1.0
A(1,2)=2.0
A(1,3)=1.0
A(2,1)=1.0
A(2,2)=-3.0
A(2,3)=-2.0
A(3,1)=3.0
A(3,2)=-1.0
A(3,3)=-1.0
! Predefined vector b
b(1)=-1
b(2)=-1
b(3)=4
! Call Gauss method
CALL Gauss_method(A,b,rozmiar)
END PROGRAM gauss_program

SUBROUTINE Gauss_method(A,b,n)
INTEGER :: n, i_max, Maks_bezw_calk
DOUBLE PRECISION, DIMENSION(n,n) :: A, AMatrix
DOUBLE PRECISION, DIMENSION(n) :: b, bVector, xVector
AMatrix = A
bVector = b
```

```

xVector(:) = -1
do i = 1, n
  if (AMatrix(i,i) == 0) then
    i_max = Maks_bezw_calk(AMatrix(i:n,i), (n-i+1)) + i - 1
    if(AMatrix(i_max,i) /= 0) then
      CALL Podmien_wiersz(i, i_max, AMatrix, n)
      CALL Podmien_elem_vec(i, i_max, bVector, n)
    else
      if (AMatrix(Maks_bezw_calk(AMatrix(:,i),n),i) /= 0) then
        STOP 'Impossible to calculate the roots'
      else
        xVector(i) = 0
      end if
    end if
  end if
end if
  if (AMatrix(i,i) /= 0) then
    do j = (i+1), n
      do k = (i+1), n
        AMatrix(j,k) = AMatrix(j,k) - AMatrix(i,k) * ( AMatrix(j,
i) / AMatrix(i,i) )
      end do
      bVector(j) = bVector(j) - bVector(i) * ( AMatrix(j,i) / AMatrix(i,
i) )
      AMatrix(j,i) = 0
    end do
  end if
end do
do i = n, 1, -1
  if(xVector(i) /= 0) then
    xVector(i) = bVector(i) - SUM( xVector((i+1):n) * AMatrix(i, (i+1):
n) )
    if(AMatrix(i,i) /= 0) xVector(i) = xVector(i) / AMatrix(i,i)
  end if
end do
print*, 'xVector='
print*, xVector
END SUBROUTINE Gauss_method

SUBROUTINE Podmien_wiersz(r1, r2, A, n)

INTEGER :: r1, r2, n
DOUBLE PRECISION, DIMENSION(n,n) :: A
do i = 1, n
  CALL Zamiana( A(r1,i) , A(r2,i) )
end do

```



```
END SUBROUTINE Podmien_wiersz
SUBROUTINE Podmien_elem_vec(i1, i2, A, n)
INTEGER :: n, i1, i2
DOUBLE PRECISION :: swap
DOUBLE PRECISION, DIMENSION(n) :: A
swap = A(i1)
A(i1) = A(i2)
A(i2) = swap
END SUBROUTINE Podmien_elem_vec
FUNCTION Maks_bezw_calk(A, n)
INTEGER :: Maks_bezw_calk, n
DOUBLE PRECISION, DIMENSION(n) :: A
Maks_bezw_calk = 1
do i = 2, n
if (ABS(A(i)) > ABS(A(Maks_bezw_calk))) Maks_bezw_calk = i
end do
END FUNCTION Maks_bezw_calk
SUBROUTINE Zamiana(v1, v2)
DOUBLE PRECISION :: v1, v2, swap
swap = v1
v1 = v2
v2 = swap
END SUBROUTINE Zamiana
```

Appendix E

Subroutine: Transformation and Integration (3D)

```
subroutine TRANSFORM_INTEGRATION
use globalvariables !in this module define global variables
implicit REAL*8 (A-H,O-Z)
integer(kind=8), dimension( 8) :: node_local

!WFACTOR - Weighting factor (each Gaussian point)
!COORD - Coordinates
!PKSI - Derivative of shape function by KSI(each Gaussian point)
!PETA - Derivative of shape function by ETA(each Gaussian point)
!PZETA - Derivative of shape function by ZETA(each Gaussian point)
!NO_ELEMENTS - Number of elements in solution domain
!LAMBDA - Thermal conductivity factor W/mK
!DPX - Derivative of shape function by X(each Gaussian point)
!DPY - Derivative of shape function by Y(each Gaussian point)
!DPZ - Derivative of shape function by Z(each Gaussian point)

COORD(1)= -0.5773502692D+00
COORD(2)= +0.5773502692D+00
WFACTOR(1)= +1.0000000000D+00
WFACTOR(2)= +1.0000000000D+00
do kpnt= 1, 2
do jpnt= 1, 2
do ipnt= 1, 2
```

!Shape function

```

SHAPE_FUN(ipnt,jpnt,kpnt,1)= 0.125 * (1.d0 - COORD(ipnt)) *
&(1.d0 - COORD(jpnt)) * (1.d0 - COORD(kpnt))
SHAPE_FUN(ipnt,jpnt,kpnt,2)= 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 - COORD(jpnt)) * (1.d0 - COORD(kpnt))
SHAPE_FUN(ipnt,jpnt,kpnt,3)= 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 + COORD(jpnt)) * (1.d0 - COORD(kpnt))
SHAPE_FUN(ipnt,jpnt,kpnt,4)= 0.125 * (1.d0 - COORD(ipnt)) *
&(1.d0 + COORD(jpnt)) * (1.d0 - COORD(kpnt))
SHAPE_FUN(ipnt,jpnt,kpnt,5)= 0.125 * (1.d0 - COORD(ipnt)) *
&(1.d0 - COORD(jpnt)) * (1.d0 + COORD(kpnt))
SHAPE_FUN(ipnt,jpnt,kpnt,6)= 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 - COORD(jpnt)) * (1.d0 + COORD(kpnt))
SHAPE_FUN(ipnt,jpnt,kpnt,7)= 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 + COORD(jpnt)) * (1.d0 + COORD(kpnt))
SHAPE_FUN(ipnt,jpnt,kpnt,8)= 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 + COORD(jpnt)) * (1.d0 + COORD(kpnt))

```

!Derivative of shape function

```

PKSI(jpnt, kpnt, 1) = - 0.125 * (1.d0 - COORD(jpnt)) *
&(1.d0 - COORD(kpnt))
PKSI(jpnt, kpnt, 2) = + 0.125 * (1.d0 - COORD(jpnt)) *
&(1.d0 - COORD(kpnt))
PKSI(jpnt, kpnt, 3) = + 0.125 * (1.d0 + COORD(jpnt)) *
&(1.d0 - COORD(kpnt))
PKSI(jpnt, kpnt, 4) = - 0.125 * (1.d0 + COORD(jpnt)) *
&(1.d0 - COORD(kpnt))
PKSI(jpnt, kpnt, 5) = - 0.125 * (1.d0 - COORD(jpnt)) *
&(1.d0 + COORD(kpnt))
PKSI(jpnt, kpnt, 6) = + 0.125 * (1.d0 - COORD(jpnt)) *
&(1.d0 + COORD(kpnt))
PKSI(jpnt, kpnt, 7) = + 0.125 * (1.d0 + COORD(jpnt)) *
&(1.d0 + COORD(kpnt))
PKSI(jpnt, kpnt, 8) = - 0.125 * (1.d0 + COORD(jpnt)) *
&(1.d0 + COORD(kpnt))
PETA(ipnt, kpnt, 1) = - 0.125 * (1.d0 - COORD(ipnt)) *
&(1.d0 - COORD(kpnt))
PETA(ipnt, kpnt, 2) = - 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 - COORD(kpnt))
PETA(ipnt, kpnt, 3) = + 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 - COORD(kpnt))
PETA(ipnt, kpnt, 4) = + 0.125 * (1.d0 - COORD(ipnt)) *
&(1.d0 - COORD(kpnt))
PETA(ipnt, kpnt, 5) = - 0.125 * (1.d0 - COORD(ipnt)) *

```

```

&(1.d0 + COORD(kpnt))
PETA(ipnt,kpnt,6)= - 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 + COORD(kpnt))
PETA(ipnt,kpnt,7)= + 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 + COORD(kpnt))
PETA(ipnt,kpnt,8)= + 0.125 * (1.d0 - COORD(ipnt)) *
&(1.d0 + COORD(kpnt))
PZETA(ipnt,jpnt,1)= - 0.125 * (1.d0 - COORD(ipnt)) *
&(1.d0 - COORD(jpnt))
PZETA(ipnt,jpnt,2)= - 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 - COORD(jpnt))
PZETA(ipnt,jpnt,3)= - 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 + COORD(jpnt))
PZETA(ipnt,jpnt,4)= - 0.125 * (1.d0 - COORD(ipnt)) *
&(1.d0 + COORD(jpnt))
PZETA(ipnt,jpnt,5)= + 0.125 * (1.d0 - COORD(ipnt)) *
&(1.d0 - COORD(jpnt))
PZETA(ipnt,jpnt,6)= + 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 - COORD(jpnt))
PZETA(ipnt,jpnt,7)= + 0.125 * (1.d0 + COORD(ipnt)) *
&(1.d0 + COORD(jpnt))
PZETA(ipnt,jpnt,8)= + 0.125 * (1.d0 - COORD(ipnt)) *
&(1.d0 + COORD(jpnt))
enddo
enddo
enddo

```

```
do i= 1, NO_ELEMENTS
```

```
!Global number for each element
```

```

node1= CONNECT_MATRIX(i,1)
node2= CONNECT_MATRIX(i,2)
node3= CONNECT_MATRIX(i,3)
node4= CONNECT_MATRIX(i,4)
node5= CONNECT_MATRIX(i,5)
node6= CONNECT_MATRIX(i,6)
node7= CONNECT_MATRIX(i,7)
node8= CONNECT_MATRIX(i,8)

```

```
write (*,*) 'Global numbers for element:', i
```

```
write (*,*) '-----'
```

```

write (*,*) node1
write (*,*) node2
write (*,*) node3
write (*,*) node4

```

```

write (*,*) node5
write (*,*) node6
write (*,*) node7
write (*,*) node8
!XYZ coordinates for each element
X1= XYZ (node1,1)
X2= XYZ (node2,1)
X3= XYZ (node3,1)
X4= XYZ (node4,1)
X5= XYZ (node5,1)
X6= XYZ (node6,1)
X7= XYZ (node7,1)
X8= XYZ (node8,1)
Y1= XYZ (node1,2)
Y2= XYZ (node2,2)
Y3= XYZ (node3,2)
Y4= XYZ (node4,2)
Y5= XYZ (node5,2)
Y6= XYZ (node6,2)
Y7= XYZ (node7,2)
Y8= XYZ (node8,2)
Z1= XYZ (node1,3)
Z2= XYZ (node2,3)
Z3= XYZ (node3,3)
Z4= XYZ (node4,3)
Z5= XYZ (node5,3)
Z6= XYZ (node6,3)
Z7= XYZ (node7,3)
Z8= XYZ (node8,3)

write (*,*) 'Coordinates of nodes of element:', i
write (*,*) '-----'
write (*,*) X1,Y1,Z1
write (*,*) X2,Y2,Z2
write (*,*) X3,Y3,Z3
write (*,*) X4,Y4,Z4
write (*,*) X5,Y5,Z5
write (*,*) X6,Y6,Z6
write (*,*) X7,Y7,Z7
write (*,*) X8,Y8,Z8

call JACOBIAN (iflaga1, iflaga2,detJ, DPX, DPY, DPZ, PKSI, PETA,
& PZETA, X1, X2, X3, X4, X5, X6, X7, X8,

```

```
& Y1, Y2, Y3, Y4, Y5, Y6, Y7, Y8,
& Z1, Z2, Z3, Z4, Z5, Z6, Z7, Z8 )
```

!Integration

```
wspij= 0.d0
```

```
do iii= 1, 8
```

```
do jjj= 1, 8
```

```
do kpn= 1, 2
```

```
do jpn= 1, 2
```

```
do ipn= 1, 2
```

```
segment= abs(detJ(ipn,jpn,kpn)) *WFACTOR(ipn) *WFACTOR(jpn) *WFACTOR(kpn)
```

```
wspij= wspij + LAMBDA * (DPX(ipn,jpn,kpn,iii) *
```

```
&DPX(ipn,jpn,kpn,jjj)
```

```
&+DPY(ipn,jpn,kpn,iii) *DPY(ipn,jpn,kpn,jjj)+DPZ(ipn,jpn,kpn,iii) *
```

```
&DPZ(ipn,jpn,kpn,jjj)) *segment
```

```
enddo
```

```
enddo
```

```
enddo
```

```
enddo
```

```
enddo
```

```
enddo
```

```
return
```

```
end
```

subroutine JACOBIAN (iflaga1, iflaga2,detJ, PX, DPY, DPZ, PKSI,
&PETA, PZETA,X1, X2, X3, X4, X5, X6, X7, X8, Y1, Y2, Y3, Y4, Y5,
&Y6, Y7, Y8,Z1, Z2, Z3, Z4, Z5, Z6, Z7, Z8)

```
implicit REAL*8 (A-H,O-Z)
```

```
dimension detJ(2,2,2)
```

```
dimension PKSI(2,2,8), PETA(2,2,8), PZETA(2,2,8)
```

```
dimension DPX(2,2,2,8), DPY(2,2,2,8), DPZ(2,2,2,8)
```

```
do kpnt= 1, 2
```

```
do jpnt= 1, 2
```

```
do ipnt= 1, 2
```

```
DPX(ipnt,jpnt,kpnt,1)=0.d0
```

```
DPX(ipnt,jpnt,kpnt,2)=0.d0
```

```

DPX(ipnt, jpnt, kpnt, 3) = 0.d0
DPX(ipnt, jpnt, kpnt, 4) = 0.d0
DPX(ipnt, jpnt, kpnt, 5) = 0.d0
DPX(ipnt, jpnt, kpnt, 6) = 0.d0
DPX(ipnt, jpnt, kpnt, 7) = 0.d0
DPX(ipnt, jpnt, kpnt, 8) = 0.d0
DPY(ipnt, jpnt, kpnt, 1) = 0.d0
DPY(ipnt, jpnt, kpnt, 2) = 0.d0
DPY(ipnt, jpnt, kpnt, 3) = 0.d0
DPY(ipnt, jpnt, kpnt, 4) = 0.d0
DPY(ipnt, jpnt, kpnt, 5) = 0.d0
DPY(ipnt, jpnt, kpnt, 6) = 0.d0
DPY(ipnt, jpnt, kpnt, 7) = 0.d0
DPY(ipnt, jpnt, kpnt, 8) = 0.d0
DPZ(ipnt, jpnt, kpnt, 1) = 0.d0
DPZ(ipnt, jpnt, kpnt, 2) = 0.d0
DPZ(ipnt, jpnt, kpnt, 3) = 0.d0
DPZ(ipnt, jpnt, kpnt, 4) = 0.d0
DPZ(ipnt, jpnt, kpnt, 5) = 0.d0
DPZ(ipnt, jpnt, kpnt, 6) = 0.d0
DPZ(ipnt, jpnt, kpnt, 7) = 0.d0
DPZ(ipnt, jpnt, kpnt, 8) = 0.d0

```

DPXDPKSI =

```

& + PKSI(jpnt, kpnt, 1) * X1 + PKSI(jpnt, kpnt, 2) * X2
& + PKSI(jpnt, kpnt, 3) * X3 + PKSI(jpnt, kpnt, 4) * X4
& + PKSI(jpnt, kpnt, 5) * X5 + PKSI(jpnt, kpnt, 6) * X6
& + PKSI(jpnt, kpnt, 7) * X7 + PKSI(jpnt, kpnt, 8) * X8

```

DPYDPKSI =

```

& + PKSI(jpnt, kpnt, 1) * Y1 + PKSI(jpnt, kpnt, 2) * Y2
& + PKSI(jpnt, kpnt, 3) * Y3 + PKSI(jpnt, kpnt, 4) * Y4
& + PKSI(jpnt, kpnt, 5) * Y5 + PKSI(jpnt, kpnt, 6) * Y6
& + PKSI(jpnt, kpnt, 7) * Y7 + PKSI(jpnt, kpnt, 8) * Y8

```

DPZDPKSI =

```

& + PKSI(jpnt, kpnt, 1) * Z1 + PKSI(jpnt, kpnt, 2) * Z2
& + PKSI(jpnt, kpnt, 3) * Z3 + PKSI(jpnt, kpnt, 4) * Z4
& + PKSI(jpnt, kpnt, 5) * Z5 + PKSI(jpnt, kpnt, 6) * Z6
& + PKSI(jpnt, kpnt, 7) * Z7 + PKSI(jpnt, kpnt, 8) * Z8

```

DPXDPETA =

```

& + PETA(ipnt, kpnt, 1) * X1 + PETA(ipnt, kpnt, 2) * X2
& + PETA(ipnt, kpnt, 3) * X3 + PETA(ipnt, kpnt, 4) * X4
& + PETA(ipnt, kpnt, 5) * X5 + PETA(ipnt, kpnt, 6) * X6
& + PETA(ipnt, kpnt, 7) * X7 + PETA(ipnt, kpnt, 8) * X8

```

DPYDPETA =

```

& + PETA(ipnt, kpnt, 1) * Y1 + PETA(ipnt, kpnt, 2) * Y2
& + PETA(ipnt, kpnt, 3) * Y3 + PETA(ipnt, kpnt, 4) * Y4
& + PETA(ipnt, kpnt, 5) * Y5 + PETA(ipnt, kpnt, 6) * Y6
& + PETA(ipnt, kpnt, 7) * Y7 + PETA(ipnt, kpnt, 8) * Y8
DPZDPETA =
& + PETA(ipnt, kpnt, 1) * Z1 + PETA(ipnt, kpnt, 2) * Z2
& + PETA(ipnt, kpnt, 3) * Z3 + PETA(ipnt, kpnt, 4) * Z4
& + PETA(ipnt, kpnt, 5) * Z5 + PETA(ipnt, kpnt, 6) * Z6
& + PETA(ipnt, kpnt, 7) * Z7 + PETA(ipnt, kpnt, 8) * Z8
DPXDPZETA =
& + PZETA(ipnt, jpnt, 1) * X1 + PZETA(ipnt, jpnt, 2) * X2
& + PZETA(ipnt, jpnt, 3) * X3 + PZETA(ipnt, jpnt, 4) * X4
& + PZETA(ipnt, jpnt, 5) * X5 + PZETA(ipnt, jpnt, 6) * X6
& + PZETA(ipnt, jpnt, 7) * X7 + PZETA(ipnt, jpnt, 8) * X8
DPYDPZETA =
& + PZETA(ipnt, jpnt, 1) * Y1 + PZETA(ipnt, jpnt, 2) * Y2
& + PZETA(ipnt, jpnt, 3) * Y3 + PZETA(ipnt, jpnt, 4) * Y4
& + PZETA(ipnt, jpnt, 5) * Y5 + PZETA(ipnt, jpnt, 6) * Y6
& + PZETA(ipnt, jpnt, 7) * Y7 + PZETA(ipnt, jpnt, 8) * Y8
DPZDPZETA =
& + PZETA(ipnt, jpnt, 1) * Z1 + PZETA(ipnt, jpnt, 2) * Z2
& + PZETA(ipnt, jpnt, 3) * Z3 + PZETA(ipnt, jpnt, 4) * Z4
& + PZETA(ipnt, jpnt, 5) * Z5 + PZETA(ipnt, jpnt, 6) * Z6
& + PZETA(ipnt, jpnt, 7) * Z7 + PZETA(ipnt, jpnt, 8) * Z8
detJ(ipnt, jpnt, kpnt) = DPXDPKSI * (DPYDPETA*DPZDPZETA-DPZDPETA*
&DPYDPZETA) + DPYDPKSI * (DPZDPETA*DPXDPZETA-DPXDPETA*DPZDPZETA) +
&DPZDPKSI * (DPXDPETA*DPYDPZETA-DPYDPETA*DPXDPZETA)

segment= 1.d0 / detJ(ipnt, jpnt, kpnt)
w11= segment * ( DPYDPETA*DPZDPZETA - DPZDPETA*DPYDPZETA )
w12= segment * ( DPZDPKSI*DPYDPZETA - DPYDPKSI*DPZDPZETA )
w13= segment * ( DPYDPKSI*DPZDPETA - DPZDPKSI*DPYDPETA )
w21= segment * ( DPZDPETA*DPXDPZETA - DPXDPETA*DPZDPZETA )
w22= segment * ( DPXDPKSI*DPZDPZETA - DPZDPKSI*DPXDPZETA )
w23= segment * ( DPZDPKSI*DPXDPETA - DPXDPKSI*DPZDPETA )
w31= segment * ( DPXDPETA*DPYDPZETA - DPYDPETA*DPXDPZETA )
w32= segment * ( DPYDPKSI*DPXDPZETA - DPXDPKSI*DPYDPZETA )
w33= segment * ( DPXDPKSI*DPYDPETA - DPYDPKSI*DPXDPETA )

detJ(ipnt, jpnt, kpnt) = dabs(detJ(ipnt, jpnt, kpnt))

DPX(ipnt, jpnt, kpnt, 1) = w11*PKSI(jpnt, kpnt, 1) +
&w12*PETA(ipnt, kpnt, 1) + w13*PZETA(ipnt, jpnt, 1)

```



```

DPX(ipnt, jpnt, kpnt, 2) = w11*PKSI(jpnt, kpnt, 2) +
&w12*PETA(ipnt, kpnt, 2) + w13*PZETA(ipnt, jpnt, 2)
DPX(ipnt, jpnt, kpnt, 3) = w11*PKSI(jpnt, kpnt, 3) +
&w12*PETA(ipnt, kpnt, 3) + w13*PZETA(ipnt, jpnt, 3)
DPX(ipnt, jpnt, kpnt, 4) = w11*PKSI(jpnt, kpnt, 4) +
&w12*PETA(ipnt, kpnt, 4) + w13*PZETA(ipnt, jpnt, 4)
DPX(ipnt, jpnt, kpnt, 5) = w11*PKSI(jpnt, kpnt, 5) +
&w12*PETA(ipnt, kpnt, 5) +w160;w13*PZETA(ipnt, jpnt, 5)
DPX(ipnt, jpnt, kpnt, 6) = w11*PKSI(jpnt, kpnt, 6) +
&w12*PETA(ipnt, kpnt, 6) + w13*PZETA(ipnt, jpnt, 6)
DPX(ipnt, jpnt, kpnt, 7) = w11*PKSI(jpnt, kpnt, 7) +
&w12*PETA(ipnt, kpnt, 7) + w13*PZETA(ipnt, jpnt, 7)
DPX(ipnt, jpnt, kpnt, 8) = w11*PKSI(jpnt, kpnt, 8) +
&w12*PETA(ipnt, kpnt, 8) + w13*PZETA(ipnt, jpnt, 8)
DPY(ipnt, jpnt, kpnt, 1) = w21*PKSI(jpnt, kpnt, 1) +
&w22*PETA(ipnt, kpnt, 1) + w23*PZETA(ipnt, jpnt, 1)
DPY(ipnt, jpnt, kpnt, 2) = w21*PKSI(jpnt, kpnt, 2) +
&w22*PETA(ipnt, kpnt, 2) + w23*PZETA(ipnt, jpnt, 2)
DPY(ipnt, jpnt, kpnt, 3) = w21*PKSI(jpnt, kpnt, 3) +
&w22*PETA(ipnt, kpnt, 3) + w23*PZETA(ipnt, jpnt, 3)
DPY(ipnt, jpnt, kpnt, 4) = w21*PKSI(jpnt, kpnt, 4) +
&w22*PETA(ipnt, kpnt, 4) + w23*PZETA(ipnt, jpnt, 4)
DPY(ipnt, jpnt, kpnt, 5) = w21*PKSI(jpnt, kpnt, 5) +
&w22*PETA(ipnt, kpnt, 5) + w23*PZETA(ipnt, jpnt, 5)
DPY(ipnt, jpnt, kpnt, 6) = w21*PKSI(jpnt, kpnt, 6) +
&w22*PETA(ipnt, kpnt, 6) + w23*PZETA(ipnt, jpnt, 6)
DPY(ipnt, jpnt, kpnt, 7) = w21*PKSI(jpnt, kpnt, 7) +
&w22*PETA(ipnt, kpnt, 7) + w23*PZETA(ipnt, jpnt, 7)
DPY(ipnt, jpnt, kpnt, 8) = w21*PKSI(jpnt, kpnt, 8) +
&w22*PETA(ipnt, kpnt, 8) + w23*PZETA(ipnt, jpnt, 8)
DPZ(ipnt, jpnt, kpnt, 1) = w31*PKSI(jpnt, kpnt, 1) +
&w32*PETA(ipnt, kpnt, 1) + w33*PZETA(ipnt, jpnt, 1)
DPZ(ipnt, jpnt, kpnt, 2) = w31*PKSI(jpnt, kpnt, 2) +
&w32*PETA(ipnt, kpnt, 2) + w33*PZETA(ipnt, jpnt, 2)
DPZ(ipnt, jpnt, kpnt, 3) = w31*PKSI(jpnt, kpnt, 3) +
&w32*PETA(ipnt, kpnt, 3) + w33*PZETA(ipnt, jpnt, 3)
DPZ(ipnt, jpnt, kpnt, 4) = w31*PKSI(jpnt, kpnt, 4) +
&w32*PETA(ipnt, kpnt, 4) + w33*PZETA(ipnt, jpnt, 4)
DPZ(ipnt, jpnt, kpnt, 5) = w31*PKSI(jpnt, kpnt, 5) +
&w32*PETA(ipnt, kpnt, 5) + w33*PZETA(ipnt, jpnt, 5)
DPZ(ipnt, jpnt, kpnt, 6) = w31*PKSI(jpnt, kpnt, 6) +
&w32*PETA(ipnt, kpnt, 6) + w33*PZETA(ipnt, jpnt, 6)
DPZ(ipnt, jpnt, kpnt, 7) = w31*PKSI(jpnt, kpnt, 7) +
&w32*PETA(ipnt, kpnt, 7) + w33*PZETA(ipnt, jpnt, 7)

```

```
DPZ(ipnt, jpnt, kpnt, 8) = w31*PKSI(jpnt, kpnt, 8) +  
&w32*PETA(ipnt, kpnt, 8) + w33*PZETA(ipnt, jpnt, 8)  
enddo  
enddo  
enddo  
return  
end
```

Appendix F

Function Calculating Geometry for Stereo Presentation

```
void Stereo
(float x_min, float x_max, float y_min, float y_max, float z_near, float z_far,
float z_zp, float dist, float eye)
{
//Determination of the position of the clipping planes
x_min=-6.0;
x_max=6.0;
y_min=-4.8;
y_max=4.8;
z_near=6.0;
z_far=-20.0;
//The position of the plane of zero parallax
z_zp=0.0;
//The distance between the projection center and the parallax plane
dist=10.5;
//Half the distance between the eyes, positive to the
right eye, negative to the left eye
eye=-0.3;
float dx = x_max - x_min;
float dy = y_max - y_min;
float xmid = (x_max + x_min) / 2.0;
float ymid = (y_max + y_min) / 2.0;
float clip_near = dist + z_zp - z_near;
float clip_far = dist + z_zp - z_far;
float n_over_d = clip_near / dist;
float topw = n_over_d * dy / 2.0;
float bottomw = -topw;
float rightw = n_over_d * (dx / 2.0 - eye);
```

```
float leftw = n_over_d * (-dx / 2.0 - eye);
glMatrixMode(GL_PROJECTION);
glLoadIdentity();
glFrustum(leftw, rightw, bottomw, topw, clip_near, clip_far);
glTranslatef(-xmid - eye, -ymid, -z_zp - dist);
}
```