

Residual Stress Capability

AFGROW European Workshop

ZHAW

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Discussion

- This method is used to approximate the effect of local residual stresses by using a 2-D Gaussian integration method to calculate “residual” stress intensity values to be added to applied K_s (superposition)
- Although it does not make physical sense to calculate K -values for compression, the resulting net K is used to determine crack growth
- This capability is only provided as an Engineering tool for the purpose of estimating the net effect of local residual stresses

Typical Cold Worked Hole Case

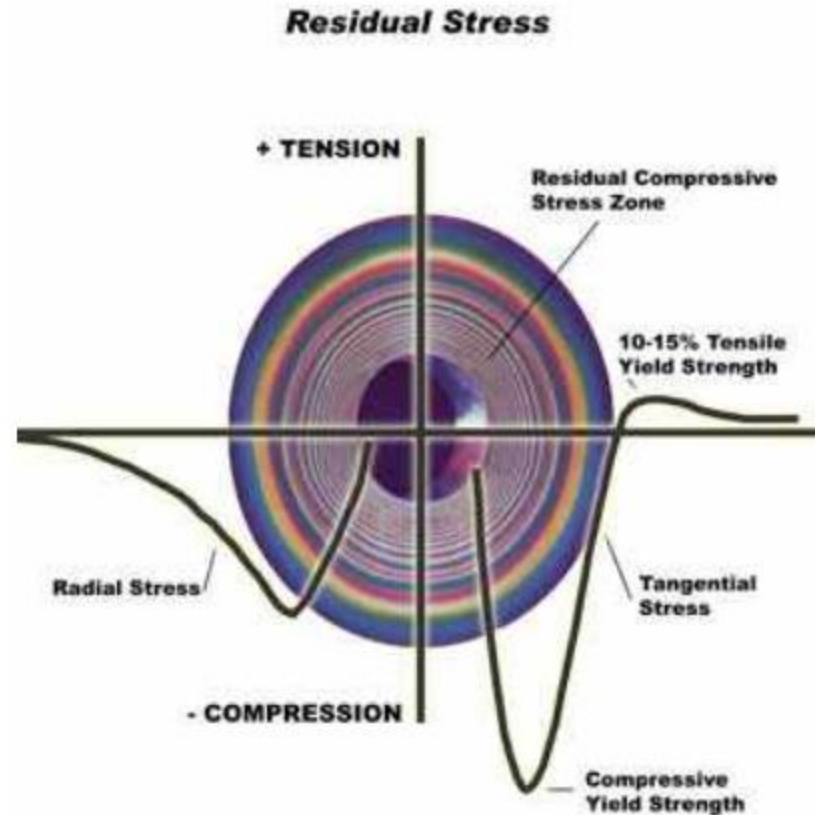
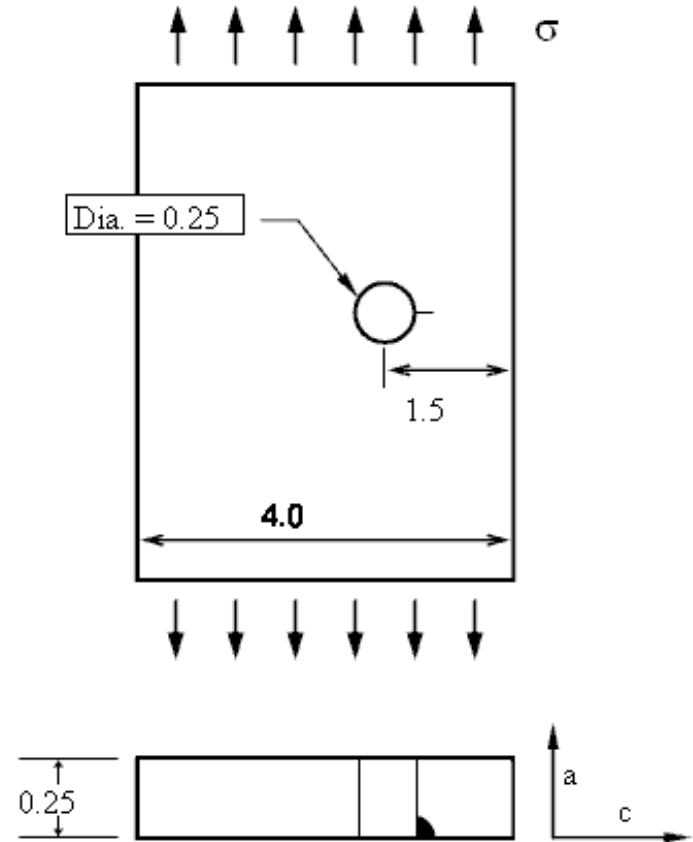


Image from: Fatigue Technologies , Inc.

Example Case in the User's Guide

Specimen Geometry:	Corner Crack at an Offset Hole in a Plate	
Dimensions:	W = 4.0 in., T = 0.25 in., Dia. = 0.25 in.	
Hole Offset:	B = 1.5 in.	
Initial Crack Size:	c = 0.05 in., a = 0.05 in.	
Material:	7050-T74 Plate (Harter T-Method)	
Stress Spectrum:	16 ksi to 0 ksi	1 Cycle
	12 ksi to 8 ksi	1000 Cycles
Retardation Model:	Generalized Willenborg Model, SOLR = 2.8	




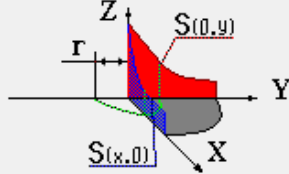
Residual Stress Input

AFGROW uses integration points for each crack growth direction at user-defined radial distances from the crack origin. The residual K value for each crack is determined using the linear interpolation between points.

This can be an additional source for error if the crack shape is not relatively close to circular, but it is not practical to anticipate crack shape changes in advance.

Residual Stresses

 AFGROW offers the option to model the effect of residual stresses on crack growth by reading in a table of residual stresses as a function of crack length, than AFGROW uses these values to generates a table of 'Residual Stress Intensity Factors' (SIF).



$S(x,y)$ - value of a stress in Z axis direction;
 r - distance from the center point of the crack along X or Y axis;

Select type of Data:

Stress Residual K

Enter stress and 'r'

Number of Sets: 7

Set	r	$S(r,0)$	$S(0,r)$
1	0	-2.4	-2.4
2	0.02	-1.2	-2.4
3	0.04	0	-2.4
4	0.1	0.4	-2.4
5	0.25	0.35	-2.4
6	0.5	0.3	0

Generate SIF table using:

Gauss Integration Weight Function

File

Open

Save

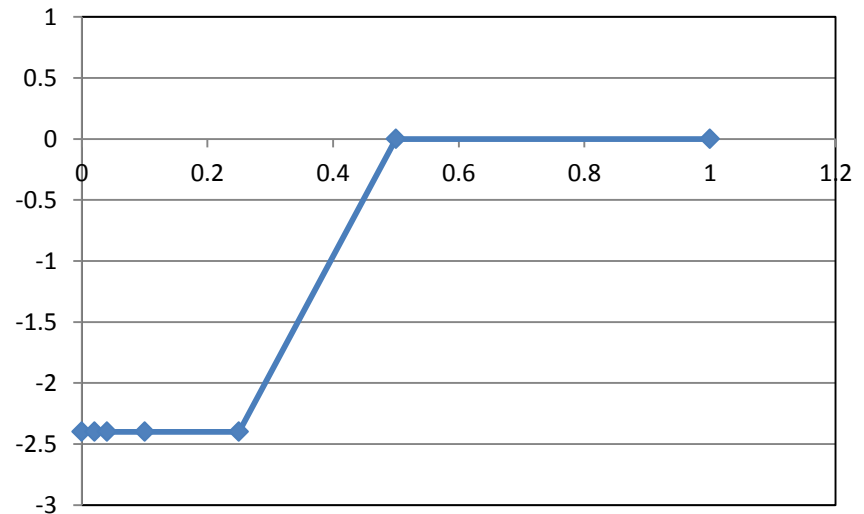
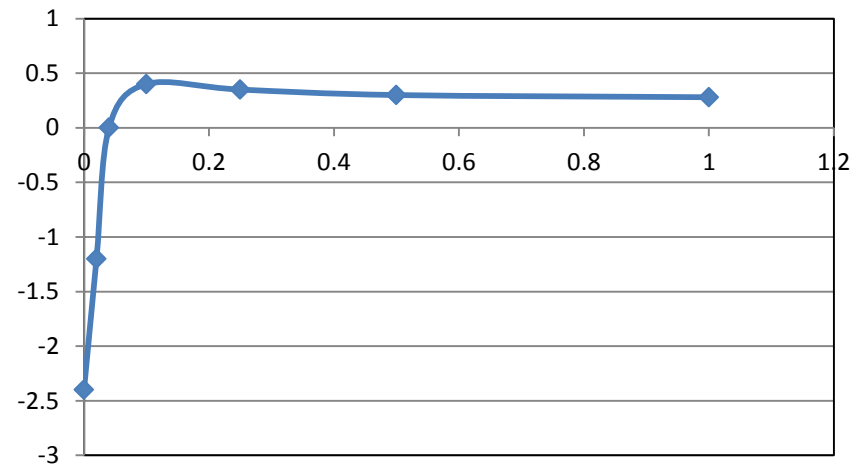
OK Cancel No Stresses

Assumed Residual Stress Distributions

X-Direction Y-Direction

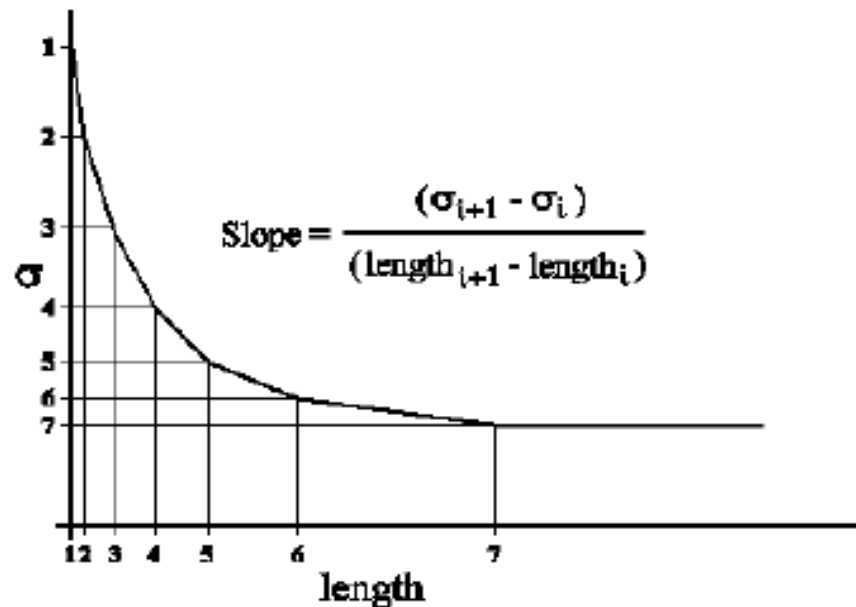
r	Residual Stress (r,0)	Residual Stress (0,r)
0.000	-2.40	-2.40
0.020	-1.20	-2.40
0.040	0.00	-2.40
0.100	0.40	-2.40
0.250	0.35	-2.40
0.500	0.30	0.00
1.000	0.28	0.00

Note: The distribution in the thickness direction must go to zero for radial distances > thickness. Otherwise, the Gaussian integration will include the effect of the distribution on the resulting residual K estimate in the width direction.



Issue

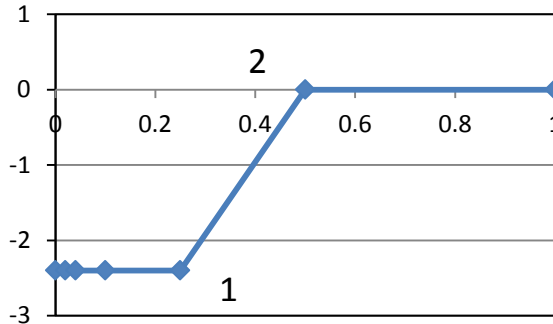
The Gaussian integration method is somewhat sensitive to the change in slope between integration points



Reasonable results have been obtained with the Gaussian method for:

$$\frac{dSlope}{dr} \leq |600|$$

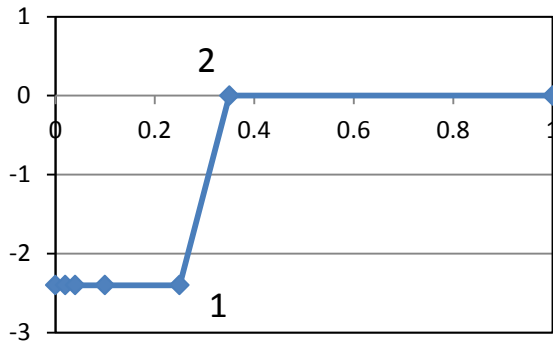
Examples



$$dS_1/dR = 38.4$$

$$dS_2/dR = -19.2$$

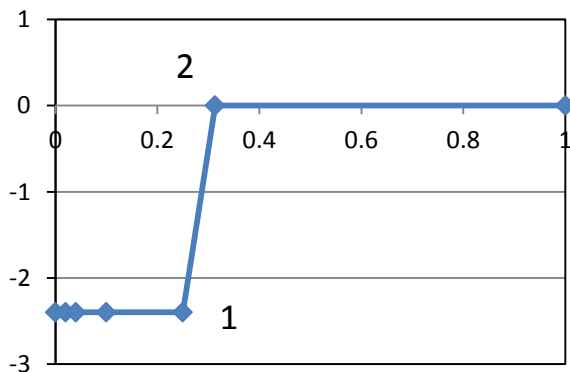
r	σ_a	$K_a \text{ resid}$	σ_c	$K_c \text{ resid}$
0.000	-2.400	-0.076	-2.400	-0.076
0.020	-2.400	-0.733	-1.200	-0.518
0.040	-2.400	-0.940	0.000	-0.254
0.100	-2.400	-1.365	0.400	-0.059
0.250	-2.400	-2.482	0.350	-0.226
0.500	0.000	-0.264	0.300	0.132
1.000	0.000	0.002	0.280	0.429



$$dS_1/dR = 240$$

$$dS_2/dR = -36.9$$

r	σ_a	$K_a \text{ resid}$	σ_c	$K_c \text{ resid}$
0.000	-2.400	-0.076	-2.400	-0.076
0.020	-2.400	-0.733	-1.200	-0.518
0.040	-2.400	-0.940	0.000	-0.254
0.100	-2.400	-1.262	0.400	-0.021
0.250	-2.400	-2.920	0.350	-0.368
0.350	0.000	-0.831	0.300	-0.188
1.000	0.000	-0.025	0.280	0.402

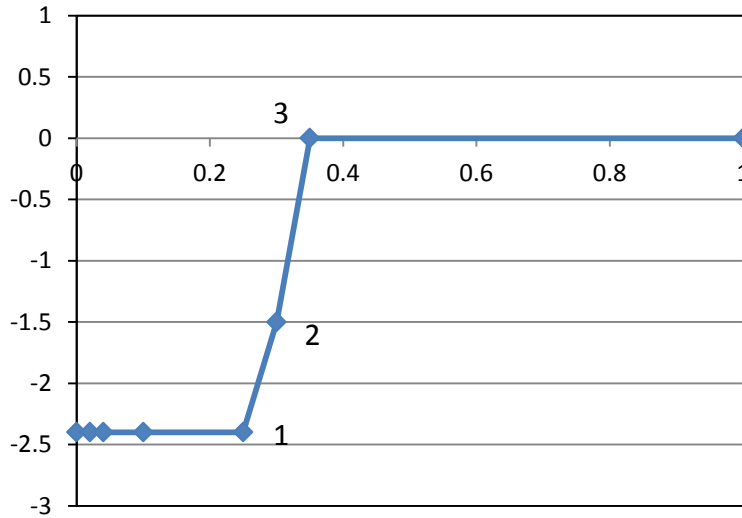


$$dS_1/dR = 599$$

$$dS_2/dR = -55.2$$

r	σ_a	$K_a \text{ resid}$	σ_c	$K_c \text{ resid}$
0.0000	-2.400	-0.076	-2.400	-0.076
0.0200	-2.400	-0.733	-1.200	-0.518
0.0400	-2.400	-0.940	0.000	-0.254
0.1000	-2.400	-1.144	0.400	0.021
0.2500	-2.400	-3.345	0.350	-0.503
0.3133	0.000	-1.336	0.300	-0.385
1.0000	0.000	-0.051	0.280	0.377

Adding Transition Points

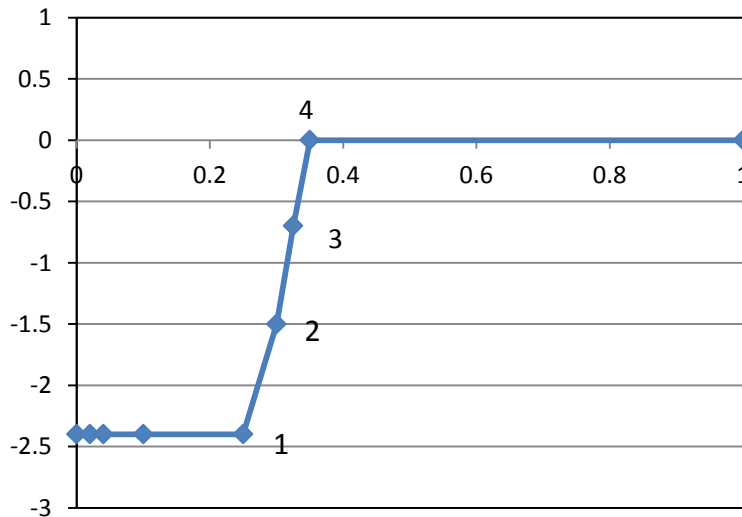


$$dS_1/dR = 360$$

$$dS_2/dR = 240$$

$$dS_3/dR = -46$$

r	σ_a	K_a resid	σ_c	K_c resid
0.000	-2.400	-0.076	-2.400	-0.076
0.020	-2.400	-0.733	-1.200	-0.518
0.040	-2.400	-0.940	0.000	-0.254
0.100	-2.400	-1.271	0.400	-0.025
0.250	-2.400	-2.256	0.350	-0.097
0.300	-1.500	-2.131	0.325	-0.133
0.350	0.000	-0.981	0.300	-0.094
1.000	0.000	-0.018	0.280	0.413



$$dS_1/dR = 360$$

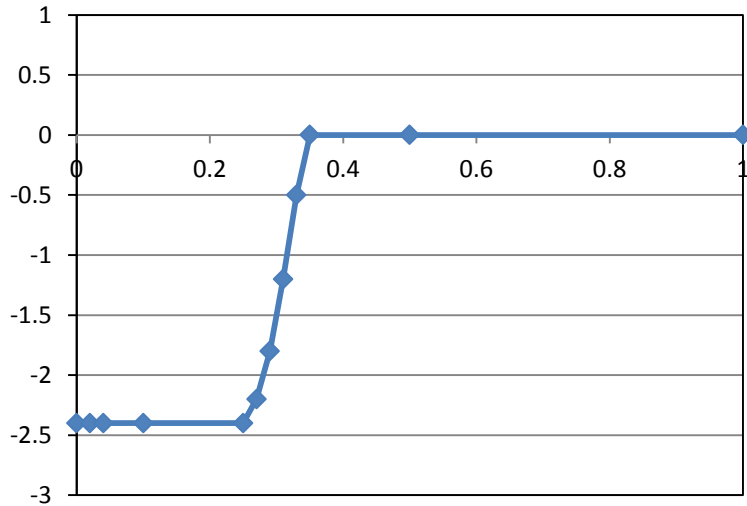
$$dS_2/dR = 560$$

$$dS_3/dR = -160$$

$$dS_4/dR = -43$$

r	σ_a	K_a resid	σ_c	K_c resid
0.000	-2.400	-0.076	-2.400	-0.076
0.020	-2.400	-0.733	-1.200	-0.518
0.040	-2.400	-0.940	0.000	-0.254
0.100	-2.400	-1.271	0.400	-0.025
0.250	-2.400	-1.993	0.350	0.011
0.300	-1.500	-2.035	0.325	-0.045
0.325	-0.700	-1.632	0.310	-0.054
0.350	0.000	-1.072	0.300	-0.042
1.000	0.000	-0.013	0.280	0.418

Using More Integration Points

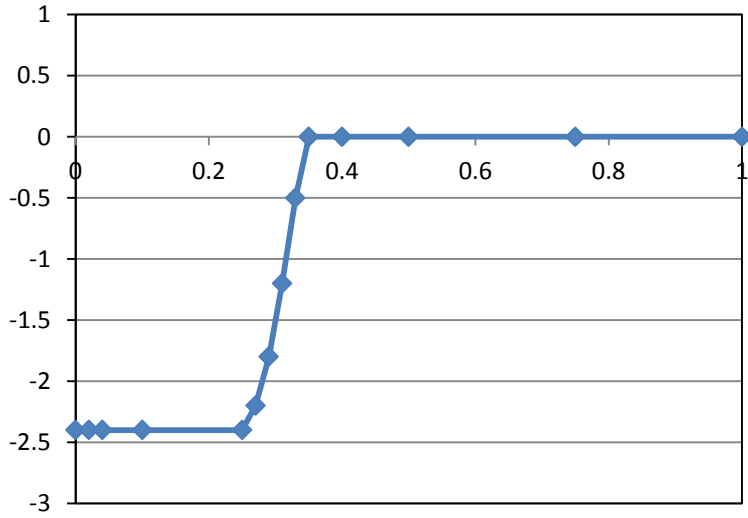


Added 5 new points to transition back to zero, keeping $dS/dr < |600|$

Point	dS/dr
1	500
2	500
3	500
4	250
5	-500
6	-166.667

r	σ_a	K_a resid	σ_c	K_c resid
0.000	-2.400	-0.076	-2.400	-0.076
0.020	-2.400	-0.733	-1.200	-0.518
0.040	-2.400	-0.940	0.000	-0.254
0.100	-2.400	-1.316	0.400	-0.042
0.250	-2.400	-1.718	0.350	0.122
0.270	-2.200	-1.834	0.340	0.094
0.290	-1.800	-1.827	0.330	0.069
0.310	-1.200	-1.689	0.320	0.050
0.330	-0.500	-1.352	0.310	0.043
0.350	0.000	-0.942	0.300	0.046
0.500	0.000	-0.287	0.300	0.180
1.000	0.000	0.010	0.280	0.441

How About a Few More....

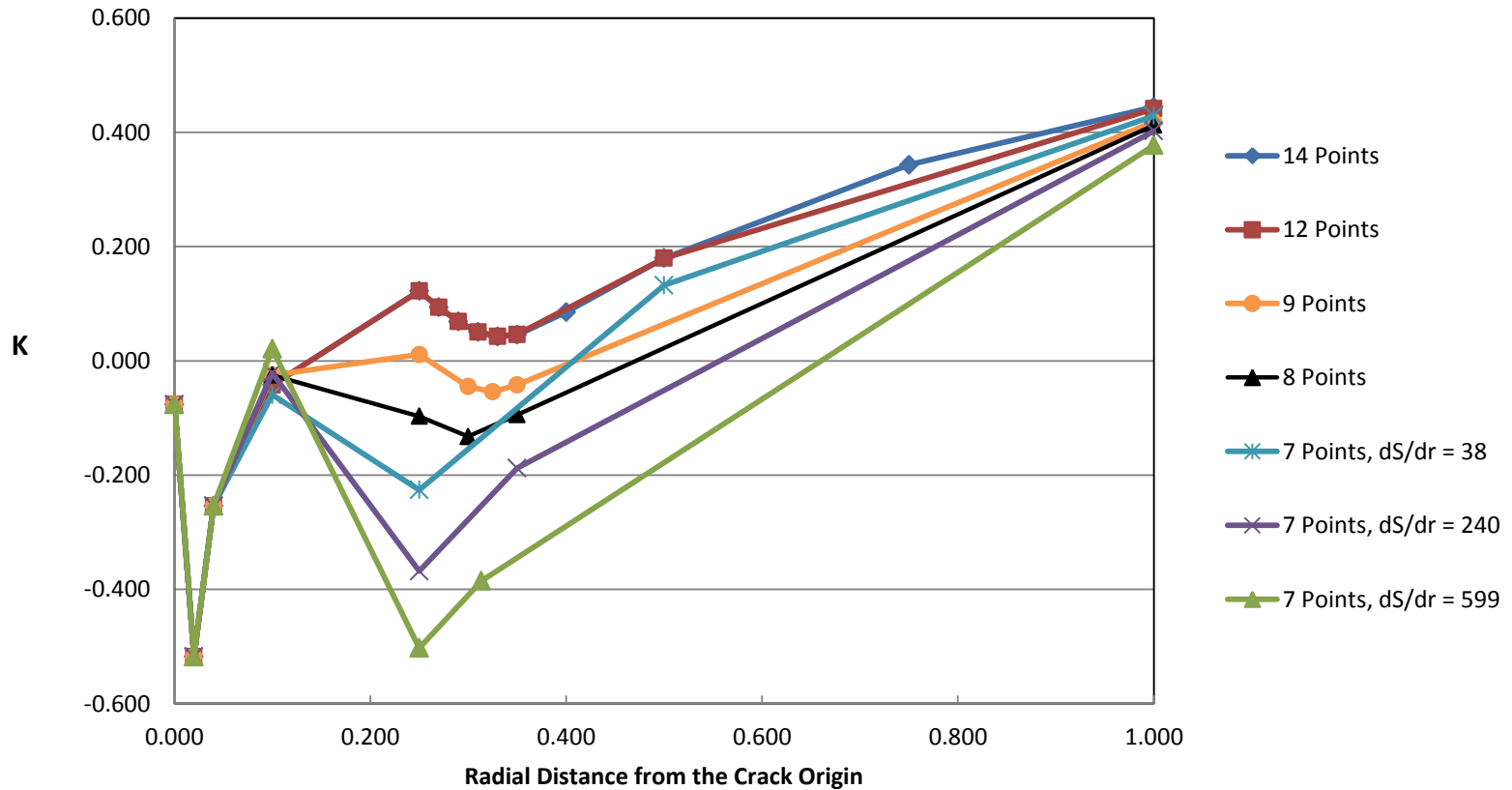


Just added two more points after the transition to zero in the y-direction (no difference in the slope changes)

r	σ_a	K_a resid	σ_c	K_c resid
0.000	-2.400	-0.076	-2.400	-0.076
0.020	-2.400	-0.733	-1.200	-0.518
0.040	-2.400	-0.940	0.000	-0.254
0.100	-2.400	-1.316	0.400	-0.042
0.250	-2.400	-1.718	0.350	0.122
0.270	-2.200	-1.834	0.340	0.094
0.290	-1.800	-1.827	0.330	0.069
0.310	-1.200	-1.689	0.320	0.050
0.330	-0.500	-1.371	0.310	0.042
0.350	0.000	-0.953	0.300	0.045
0.400	0.000	-0.580	0.300	0.085
0.500	0.000	-0.290	0.300	0.180
0.750	0.000	-0.067	0.290	0.343
1.000	0.000	0.011	0.280	0.444

Summary, continued

$K_{c,resid}$



Conclusions

- Keep slope changes between integration points below $|600|$
- Use multiple points (>3) to make the transition to $s = 0$ in the y -direction for $r > \text{thickness}$
- Minimize the transition distance as much as possible to minimize the error caused by the transition points.
- It is generally helpful to include multiple points for uniform stress distributions to decrease potential interpolation errors in the resulting stress intensities.

Remember

The beta correction capability uses the same Gaussian integration method, so the same rules apply with one major exception:

The stress ratio values used for the y -direction for $r > t$ must be set to 1.0 (not 0.0). This is because the beta correction calculations take a ratio of the calculated K for the input stress distributions divided by the K calculated for a unit stress distribution. Using a stress ratio of 1.0 beyond the thickness, means that there will be no net effect on the calculations in the x -direction for $r > t$.

There is a document on the AFGROW Web page (**Using AFGROW's Beta Correction Capability**) that demonstrates the use of the beta correction capability for a plate with out of plane bending.