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Q200312100011

Scientific Notebooks No. 069: Stochastic
Project (02/26/1993 through 12/16/1996)

S149

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Stochastic

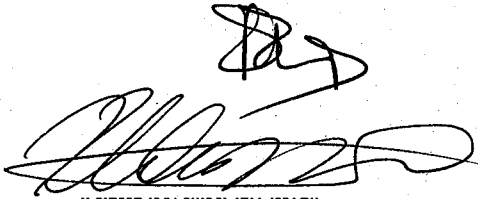
Project notebook #2.

Names

Sitakanta Mohanty

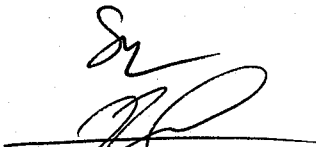
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Signatures





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All work is described
in the Project Plan.



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2/26/93

1

Bugs in the module relpmvg1.f were identified and taken care of. Test case was run w/ $C_1 = 1.0$, $C_2 = 0.1$, $\alpha_1 = \alpha_2 = 0.025$, $\beta_1 = 0.1$, $\beta_2 = 0.01$, $n = 3.7$, $a_{inc} = 0.001$

When ~~soe1~~ step increments in saturation was changed for ~~soe1~~ to ~~soe2~~ ~~after the~~ ~~soe1~~ approach 0, some results in rk_{w1} & rk_{w2} were obtained at two steps. That was fixed.

Test cases were also run with $C_1, \alpha_1, \beta_1 > C_2, \alpha_2, \beta_2$. We decided that it is better to keep $C_1, \alpha_1, \beta_1 < C_2, \alpha_2, \beta_2$

3/1/93

Module relpmvg4.f (a refined version of relpmvg1.f) was incorporated into the code fracvar3.f that determines uncat conductivity after generating the fracture system. relpmvg4.f is capable of calculating uncat. conductivity of binary systems only. Therefore fracvar3.f gives the warning that it can calculate only for binary ~~for~~ conductivity field

3/2/93 `fracvar3.f` was upgraded to `fracvar4.f` to see if all variables are compatible for with implicit double precision. The Fortran compiler is an old one and can not produce double precision random numbers using `rand48`.

~~frac~~ `relprmdist1.f` was created to calculate uncat conductivity for a dist. of matrix conductivity (~~sat~~). This implement method A (see prog)

3/3/93 `relprmdist1.f` was modified to `relprmdist2.f` to implement method B (see program).

3/4/93 Bugs in `relprmdist1.f` & `relprmdist2` were fixed. Fine intervals were taken in avg saturation (`soemini`) ~~is a~~ `cin-dist2.f`) & `soemin` (minimum perm) ~~data~~ is (`cin-dist1.f`) to get the whole range of relationship between S_{oe} & R_{wk} . See α & R_{wk} .

3/5/93

- `relprmdist1.f` module was combined with the main program for `fracvar3.f` ~~now~~ \rightarrow `fracvar5.f`
- Changes were made to `relprmdist1.f` to ~~incorporate~~ generate files with (for `satd.k`, α & β arrays) which later would be used as input files to BIGFLOW.

Pages 3 and 4 Are Intentionally
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3/25

Running program mdmuf under the title
mdmv.p. input files INPUT & IN DAT.

- Gives segmentation violation problem when a field of size $16 \times 16 \times 16$ is used on sebastian.
- Therefore, currently I ~~forced~~ am trying it on ~~the~~ vax 8700.

3/25

work
done
for
Gordon

POFLOW version 2.4.1 was tested on sparc machine (sebastian). The particular version that was used did not reproduce the results for prob1 in the manual (NUREG/CR-5991)

Need to get a new version and test it on VAX before transferring to Sparc. The above problem may answer questions regarding the strange answers we are obtaining on a one-D unsat flow problem.

3/26

A code "variogram3d.f" was developed to determine the variograms in three different directions of a 3-D field. This is a modification of variogram2d.f which was formerly named as variogram.f. The field was generated by using GSLIB program tb3dm.f (a turning bad method).

The data were specified for a $32 \times 32 \times 32$ size field, with an isotropic correlation length = 16. Nugget = 0. The correlation lengths ~~determined by the program~~ ^{specified by the program} "variogram3d.f" were very close to 16 though they were slightly different in different directions as one would expect.

4/19/93 & 4/20/93

Fracture generator code fracvar6.f was modified: the program now can create the alfa, beta & ksar files ^{for} with any matrix-fracture composite. Previously it was doing it only for a distribution of matrix conductivities in the subroutine relpmdist2. A subroutine was also added to generate the headerfile for BIGFLOW files mentioned above. Ashok is supposed to add the exact format to the header file. The α, β values for the homogeneous background now can be assigned in the main program. The bug in the ~~fr~~ postscript file generator of fracvar6.f has also been fixed.

4/26/93 — 4/30/93

A 2-D nnm2d.f and nnm2da.f codes were written to be able to use nnm method ~~as a~~ using regular fortran. The approach ~~was~~ of Baker was used. The programs ^{are} not complete yet. The programs can handle conditioning points.

Pages 1 through 7 of this Scientific Notebook were reviewed for compliance with QAP-001 in response to Corrective Action Request 94-02. Corrections and clarifications were made as appropriate. In some cases, the date of a change will reflect the date of this review rather than the date of the original Scientific Notebook entry.

Randy Falk
SWRC-QA
12/09/94

Project was closed December 1994 because it reached completion.

12/14/96

JS

**ADDITIONAL INFORMATION FOR SCIENTIFIC NOTEBOOK No.: 065, 067,
068, and 069**

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File Types: (.exe, .bat, .zip, etc.)	Various
Remarks: (computer runs, etc.)	Media contains: data and output files relative to the stochastic project.