

INTERNATIONAL ATOMIC ENERGY AGENCY
UNITED NATIONS EDUCATIONAL, SCIENTIFIC AND CULTURAL ORGANIZATION



INTERNATIONAL CENTRE FOR THEORETICAL PHYSICS
34100 TRIESTE (ITALY) • P.O.B. 586 • MIRAMARE - STRADA COSTIERA 11 • TELEPHONE: 2240-1
CABLE: CENTRATOM - TELEX 460892-I

SECOND SCHOOL ON ADVANCED TECHNIQUES
IN COMPUTATIONAL PHYSICS
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SMR.282/ 10

M.C.D. GENERATORS

F.JAMES
F.BERMEJO

EXERCISES
M.C.D. GENERATORS

F. James

F. Bermejo

1.) STUDY OF R.N. GENERATORS

a.) The mid-squares method, proposed by J. von Neuman was apparently the first algorithm proposed for the generation of pseudorandom numbers, and proceeds as follows:

Suppose we wish to generate four-digit integers and the last number generated was ABCD. Let's assume that we are using 8-digit decimals and that in successive steps, we just take the four middle digits as the number we want. To obtain the next number in the sequence, we square the last one and use the middle four digits of the product, i.e.

$$\begin{array}{l} \text{XYABCDZW} \\ \text{squared} = \underline{\text{RWEFGD YY}} \\ \text{squared} = \underline{\text{SVHLMOTP}} \\ \text{squared (etc.)} \end{array}$$

Study the statistical quality of this RNG by means of:

- a-1) Its dependence upon the "seed" (the first number).
- a-2) The repetition cycle (How long before the sequence ABCD is repeated).
- b.) The Fibonacci method consists in adding two preceding numbers and taking the remainder when the sum is taken with respect to some modulus i.e.:

$$X_i = (X_{i-1} + X_{i-2}) \bmod m$$

b-1) Write a code to study the quality of this RNG.

b-2) Consider the recursive formula,

$$X_i = (aX_{i-1} + c) \bmod m$$

where a, c and m are parameters which depend upon the particular word-length and determine the statistical quality of the generator. For c=0, the method is called a "pure multiplicative congruential" generator. Write a code to generate RN's using this method. (Hint: use a modulus $m=2^k - 1$). Test the quality of the distribution by means of the χ^2 statistic. (Hint: Partition the outcomes in m-subsets: $X < 0.10, 0.10 < X < 0.2$ etc. and compute the number of occurrences f_i . If we denote the expected frequency as f_1 , the computed quantity,

$$\chi^2 = \sum_{i=1}^m \frac{(f_i - f_1)^2}{f_1}$$

should approach the χ^2 -distribution for $m-1$ degrees of freedom. The approximation tends to be reasonable for large numbers of cells (300).

c.) The combination of two pseudorandom sequences X and Y to produce a third Z tends to reduce nonrandomness. Code the algorithms:

$$\begin{aligned} \text{a.) set } Z &= (X + Y) \bmod m \\ \text{b.) use } X &= 171X \bmod 30269 \\ &Y = 172Y \bmod 30307 \end{aligned}$$

and compare its quality with that of the preceding one by χ^2 testing.

2.) Make a Gaussian RNG by means of the central-limit theorems in the following way; consider a set of V_1, \dots, V_n random variables and set $N = \sum V_i$. As $n \rightarrow N$, tends to be normally-distributed. Since we have to settle for a "finite", a convenient way to obtain N is:

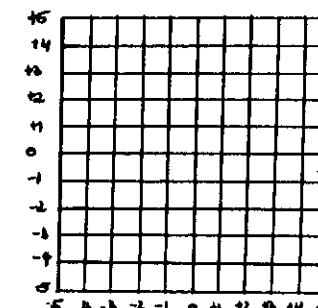
$$N = \sum_{i=1}^{12} V_i - 6$$

where the V_i are uniformly-generated pseudorandom numbers. Code the above equation and compute the expectation value and variance of the obtained distribution.

3.) Study the properties of the distribution obtained by dividing two independent Gaussian-distributed variables X, Y, (i.e. $Z = X/Y$). Compute the expectation and variance of Z numerically. Take X and $Y = N(1,1)$.

4.) 2D RANDOM WALK

Consider a planar square lattice like:



Suppose you are initially located at the lattice point (0,0). A "random walk" is simulated if at each lattice point a random choice is made about the direction of your next step.

a.) Code an algorithm to perform such a walk starting from an arbitrary point.

b.) Suppose that every time you reach the boundaries you get reflected (unable to leave the lattice) unless you are (-2,5) or (4,5). Compute the number of steps required to exit (try at least 10 simulations).

c.) Self-avoiding-walk.

Suppose that at each lattice point the random choice is subject to the restrictions that:

- you will never visit a lattice point more than once.
- all the boundaries are closed.

Write a program to simulate such a movement, bearing in mind that the simulation will stop by either achieving a maximum number of steps (200) or by getting stuck into a point (unable to move without violating the constraints a & b).

5.) BRUTE FORCE INTEGRATION.

Compute the value of the definite integral,

$$I = \int_0^1 2x^2 dx$$

by means of a program which determines the area under the function $f(x) = 2x^2$. The way to proceed would be to generate random numbers and check if the associated point lays below $f(x)$ (i.e. P is below $f(x)$ $0 < X < 1$, $0 < Y < 2$).

Generate at least 5000 events

6.) Code the Adaptive Integration algorithm given in the lectures (use the scheme, the abscissas and weights given separately).

MURKIN QUADRATURE. A GOOD ALGORITHM:

R₆₅ ON ONE INTERVAL, USE R₆₅:
 $G_n = \text{GAUSS RULE with } n \text{ points}$
 $I = G_6$
 $\Delta = |G_6 - G_5|$

THEN:

1. APPLY R₆₅ TO TOTAL INTERVAL
2. IF Δ TOO BIG GO TO 3
ELSE FINISHED
3. DIVIDE IN HALVES THE INTERVAL WITH THE LARGEST Δ
4. APPLY R₆₅ TO BOTH HALVES
5. SET $I = \sum_{\text{ALL INTERVALS}} I_i$
 $\Delta = \sum \Delta_i$
6. GO TO 2

Table 25.4 ABSCISSAS AND WEIGHT FACTORS FOR GAUSSIAN INTEGRATION

$$\int_{-1}^{+1} f(x) dx = \sum_{i=1}^n w_i f(x_i)$$

Abscissas - x_i ; (Zeros of Legendre Polynomials)

$\pm x_i$	w_i	$\pm x_i$	w_i
$n=2$		$n=8$	
0.57735 0.2691 0.89626	1.00000 0.00000 0.00000	0.18343 0.4424 0.9450	0.34268 0.7833 0.8362
0.77459 0.6692 0.41483	0.00000 0.00000 0.00000	0.52553 2.0999 1.6289	0.3170 0.6458 0.77087
0.33998 1.0435 0.48556	0.00000 0.00000 0.00000	0.79666 0.4774 1.3427	0.2234 1.0244 5.3174
0.86113 0.3115 0.94053	0.34785 0.48451 0.7454	0.96028 0.9564 0.9736	0.10122 0.5362 0.0376
$n=4$		$n=9$	
0.33846 0.9101 0.56683	0.45214 0.51548 0.52546	0.00000 0.00000 0.00000	0.33023 0.93550 0.1260
0.90617 0.98159 0.38664	0.33692 0.56189	0.32245 0.4224 0.3809	0.31234 0.70770 0.00003
$n=5$		$n=10$	
0.00000 0.00000 0.00000	0.56888 0.68888 0.88889	0.14887 0.33889 0.81631	0.29552 0.42247 1.4753
0.23861 0.91860 0.83197	0.7862 0.6704 0.93366	0.43339 0.53941 0.2947	0.26926 0.67193 0.09996
0.66120 0.93064 0.62625	0.36076 0.15730 0.48139	0.67940 0.95682 0.99024	0.21908 0.63625 1.9882
0.91246 0.5142 0.3152	0.17132 0.4923 0.79170	0.8606 0.33666 0.88085	0.14945 0.3491 0.50581
$n=6$		$n=12$	
0.00000 0.00000 0.00000	0.46791 0.39245 0.72691	0.97390 0.65285 0.17172	0.06667 0.13443 0.08888
0.45801 0.6776 0.57227	0.62643 0.748447	0.58731 0.79542 0.8617	0.20316 0.74267 0.20664
0.61787 0.6244 0.26447	0.53895	0.12462 0.671 0.5553	0.16007 0.83285 0.43346
0.75540 0.4083 0.5003	0.17132 0.4923 0.79170	0.96990 0.26741 0.9405	0.10607 0.874810 0.892863
0.86563 1.0223 0.87831	0.73880	0.90411 0.72563 0.70475	0.04717 0.53363 0.6512
0.74153 1.0855 0.99394	0.27970 0.53914 0.69277	0.98156 0.06342 0.46719	0.02715 0.2594 0.1754 0.04852
0.94910 0.79123 0.42759	0.91294 0.49661 0.60793		
$n=7$		$n=16$	
0.00000 0.00000 0.00000	0.41795 0.91836 0.74469	0.36783 1.4989 0.98180	0.23349 0.25365 0.31355
0.40584 0.51513 0.77397	0.88183 0.0505 0.51119	0.16915 0.62193 0.95002	0.16915 0.62193 0.95002
0.74153 1.0855 0.99394	0.27970 0.53914 0.69277	0.76990 0.26741 0.9405	0.16007 0.83285 0.43346
0.94910 0.79123 0.42759	0.91294 0.49661 0.60793	0.98156 0.06342 0.46719	0.04717 0.53363 0.6512
$n=8$		$n=20$	
0.09501 0.25098 0.37637	0.40185	0.15275 0.31871 0.30725	0.350698
0.28160 0.35507 0.79250	0.13230	0.49617 0.28864 0.72603	0.76788
0.45801 0.6776 0.57227	0.36342	0.18260 0.3150 0.44923	0.58867
0.61787 0.6244 0.26447	0.58004	0.16915 0.62193 0.95002	0.58867
0.75540 0.4083 0.5003	0.17132	0.14945 0.58088 0.16576	0.72081
0.86563 1.0223 0.87831	0.73880	0.12462 0.671 0.5553	0.672052
0.91223 0.4282 0.51525	0.80568	0.95513 0.8116 0.82492	0.784810
0.91223 0.4282 0.51525	0.80568	0.06225 0.34239 0.38647	0.62225
0.96397 1.9272 0.77913	0.79123	0.06267 0.20483 0.3109	0.035570
0.99312 0.85991 0.85094	0.924786	0.04060 0.14298 0.0086	0.941331
$n=24$			
0.06405 0.68928 0.62605	0.626083	0.12793 0.01953 0.46172	0.156974
0.19111 0.86674 0.73616	0.309159	0.12563 0.70563 0.46828	0.266121
0.311504 0.26796 0.96163	0.314367	0.12563 0.70563 0.46828	0.266121
0.431319 0.35076 0.26045	0.138487	0.12167 0.04729 0.27803	0.501329
0.54542 0.4713 0.88819	0.535659	0.11550 0.56680 0.53725	0.601353
0.64689 0.36519 0.36975	0.56252	0.10744 0.42701 0.1565	0.34783
0.74012 0.41915 0.78554	0.364244	0.09781 0.6521 0.4113	0.888270
0.82000 0.19859 0.73902	0.921954	0.08619 0.01615 0.1913	0.25917
0.88641 0.5270 0.0401	0.04213	0.07314 0.6814 0.1080	0.05734
0.97472 0.97472 0.45520	0.2772	0.05929 0.85849 1.5436	0.780746
0.93827 0.45520 0.2772	0.758524	0.04427 0.7388 1.7419	0.806169
0.99518 0.72199 0.97021	0.360160	0.02853 0.13886 0.2833	0.663181
0.99518 0.72199 0.97021	0.360160	0.01234 0.1227 0.99987	0.19547

Compiled from P. Davis and P. Rabinowitz, *Abscissas and weights for Gaussian quadratures of high order*, J. Research NBS 56, 35-37, 1956; RP2645; P. Davis and P. Rabinowitz, *Additional abscissas and weights for Gaussian quadratures of high order*. Values for $n=64$, 80, and 96, J. Research NBS 60, 613-614, 1958, RP2675; and A. N. Lowan, N. Davids, and A. Levenson, *Table of the zeros of the Legendre polynomials of order 1-16 and the weight coefficients for Gauss' mechanical quadrature formula*, Bull. Amer. Math. Soc. 48, 739-743, 1942 (with permission).