

# Laplace

Va=	0															
Vb=	100															
	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
0	0.4	0.8	1.2	1.6	2.1	2.5	2.9	3.2	3.6	3.9	4.2	4.5	4.7	4.8	4.9	
0	0.8	1.6	2.5	3.3	4.1	4.9	5.7	6.5	7.2	7.9	8.5	9.1	9.5	9.7	9.9	
0	1.2	2.5	3.7	5	6.2	7.4	8.6	9.8	11	12	13	14	14	15	15	
0	1.6	3.3	5	6.6	8.3	10	12	13	15	16	17	19	19	20	20	
0	2.1	4.1	6.2	8.3	10	12	15	17	19	21	22	24	25	26	26	
0	2.5	4.9	7.4	10	12	15	18	20	23	25	27	29	31	32	32	
0	2.9	5.7	8.6	12	15	18	21	24	27	30	32	35	37	38	39	
0	3.2	6.5	9.8	13	17	20	24	27	31	35	38	41	44	45	46	
0	3.6	7.2	11	15	19	23	27	31	35	40	44	48	51	53	54	
0	3.9	7.9	12	16	21	25	30	35	40	45	50	55	60	62	64	
0	4.2	8.5	13	17	22	27	32	38	44	50	57	64	70	73	74	
0	4.5	9.1	14	19	24	29	35	41	48	55	64	73	82	85	87	
0	4.7	9.5	14	19	25	31	37	44	51	60	70	82	100	100	100	
0	4.8	9.7	15	20	26	32	38	45	53	62	73	85	100	100	100	
0	4.9	9.9	15	20	26	32	39	46	54	64	74	87	100	100	100	

