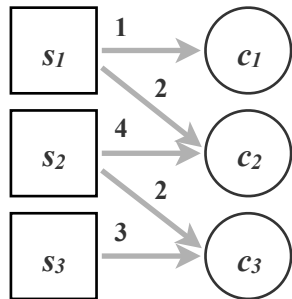


M

	c_1	c_2	c_3	...
s_1	1	2	0	...
s_2	0	4	2	...
s_2	0	0	3	...
...



P

	c_1	c_2	c_3	...
c_1	0	0.16	0.03	...
c_2	0.05	0	0.5	...
c_3	0.01	0.03	0	...
...

