



REPORT

- Preliminary measures
- + introduction
- + PSD(r0, L0) plot
- + ccl on influence of r0 and L0

+ (more to come...)

-> For next time: read Aime (Sec. 1 & Sec. 2) and Maire (Chap.1)...

Chapitre 1

Introduction

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Teaching astronomical speckle techniques

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Abstract

This paper gives an introduction to speckle techniques developed for high angular-resolution imagery in astronomy. The presentation is focussed on fundamental aspects of the techniques of Labeyrie and Weigelt. The formalism used is that of Fourier optics and statistical optics, and corresponds to graduate level. Several new approaches of known results are presented. An operator formalism is used to identify similar regions of the bispectrum. The relationship between the bispectrum and the phase closure technique is presented in an original geometrical way. Effects of photodetection are treated using simple Poisson statistics. Realistic simulations of astronomical speckle patterns illustrate the presentation. Jérôme Maire, PhD thesis (in French), chap.1

Long Telescopes may cause Objects to appear brighter and larger than short ones can do, but they cannot be so formed as to take away the confusion of the Rays which arises from the Tremors of the Atmosphere.

I. Newton, 1717 Optics, Sec. Ed., Book I, Part I, Prop. VIII



-> Perturbed wavefront generation

The well-known FFT method allows us to generate phase screens $\varphi(\vec{r})$, where \vec{r} is the two-dimensional position within the phase screen, assuming usually either a Kolmogorov or a von Karman spectrum $\Phi_{\varphi}(\vec{\nu})$, where $\vec{\nu}$ is the two-dimensional spatial frequency, from which is computed the modulus of $\tilde{\varphi}(\vec{\nu})$, the Fourier transform of $\varphi(\vec{r})$. Assuming the near-field approximation and small phase perturbation [3], the von Karman/Kolmogorov spectrum is given by

$$\Phi_{arphi}(ec{
u}) = 0.0229 r_0^{-rac{5}{3}} igg(
u^2 + rac{1}{\mathcal{L}_0^2} igg)^{-rac{11}{6}}, \qquad (1)$$

where r_0 is the Fried parameter and \mathcal{L}_0 is the wavefront outer scale (infinite for the Kolmogorov model). Within the framework of the classical FFT-based technique, a turbulent phase screen $\varphi_L(\vec{r})$ of physical length L is obtained by taking the inverse FFT of $\tilde{\varphi}_L(\vec{\nu})$, the modulus of which is obtained from Eq. (1) by applying the definition of the power spectrum, which is

$$\begin{split} \Phi_{\varphi}(\vec{\nu}) &= \lim_{L \to \infty} \left(\frac{\langle |\tilde{\varphi}_L(\nu)|^2 \rangle}{L^2} \right) \\ \Rightarrow &|\tilde{\varphi}_L(\nu)| \simeq L r_0^{-\frac{5}{6}} \sqrt{0.0228} \left(\nu^2 + \frac{1}{\mathcal{L}_0^2} \right)^{-\frac{11}{12}}, \quad (2) \end{split}$$

and which phase is random and uniformly distributed.

(From Carbillet & Riccardi, sec. 2: read it as well...)

(the same manipulation as before is applied here in order to obtain the numerical formulation here below.)

The obtained phase screen is thus numerically written

$$\begin{split} \varphi_{L}(i,j) &= \sqrt{2}\sqrt{0.0228} \left(\frac{L}{r_{0}}\right)^{\frac{5}{6}} \bigg\{ \text{FFT}^{-1} \bigg[\left(k^{2} + l^{2} + \left(\frac{L}{\mathcal{L}_{0}}\right)^{2}\right)^{-\frac{11}{12}} \exp\{\imath\theta(k,l)\} \bigg] \bigg\}, \end{split}$$
(3)

where *i* and *j* are the indices in the direct space, *k* and *l* are the indices in the FFT space, {} stands for either real part of or imaginary part of, 1 is the imaginary unit, and θ is the random uniformly distributed phase (between $-\pi$ and π). The factor $\sqrt{2}$ comes from the fact that here we use both the real and imaginary parts of the original complex generated FFT phase screens, which are independent one from the other [4]. This kind of phase screen suffers, however, from the lack of spatial frequencies lower than the inverse of the necessarily finite length *L* of the simulated array.

```
function wfgeneration, dim, length, L0, r0, lambda, SEED=seed
 wave-front (wf) generation following von Karman model
  (infinite L0 -Kolmogorov model- not allowed here).
                                                       Θ
                                                                                                            000
                                                                                                           nm
                                                                                                     4.4E+03
        = wf linear dimension [px],
  dim
                                                                                                    3.5E+03
                                                            10
  length = wf physical length [m],
                                                                                                    2.7E+03
        = wf outer-scale [m],
  L0
                                                                                                     1.8E+03
        = random generation seed (OPTIONAL),
  seed
                                                                                                    9.3E+02
        = Fried parameter at wavelength 'lambda' [m],
  r0
                                                         Ξ
                                                                                                    6.6E+01
  lambda = wavelength at which r0 is defined.
                                                                                                    -7.9E+02
                                                                                                    -1.7E+03
                                                            -5
 Marcel Carbillet [marcel.carbillet@unice.fr],
                                                                                                    -2.5E+03
  lab. Lagrange (UCA, OCA, CNRS), Feb. 2013.
                                                                                                    -3.4E+03
                                                           -10
                                                                                                    -4.2E+03
                                                                -20
                                                                       -10
                                                                               0
                                                                                      10
                                                                                             20
 Last modification: Feb. 2018.
                                                                              [m]
phase = (randomu(seed,dim,dim)-.5) * 2*!PI
                                             ; rnd uniformly distributed phase
                                                                                  wf generation:
                                             ; (between -PI and +PI)
rr = dist(dim)
                                                                                  generate a cube
modul = (rr^2+(length/L0)^2)^{(-11/12.)}
                                             ; von Karman model
screen = fft(modul*exp(complex(0,1)*phase), /INVERSE)
                                                                                  of statistically
                                             : compute wf
screen *= sqrt(2)*sqrt(.0228)*(length/r0)^(5/6.)*lambda/(2*!PI)
                                                                                  independent wf
                                             ; proper normalization of wf
screen -= mean(screen)
                                             ; force mean to zero
                                                                                   (typically 100)...
                                             ; deliver 2 independent wf:
return, screen
                                             ; float(screen) & imaginary(screen)
                                                                                   => compute mean
end
```

rms for different

 r_0, L_0

[IDL> .r wfgeneration % Compiled module: WFGENERATION. [IDL> wf=wfgeneration(128,2.,27.,.1,500E-9,SEED=seed) % Compiled module: DIST. [IDL> wf1=float(wf) [IDL> wf2=imaginary(wf) [IDL> tvscl, [wf1,wf2] [IDL> wf=wfgeneration(128,2.,27.,.1,500E-9,SEED=seed) [IDL> wf1=float(wf) [IDL> wf1=float(wf) [IDL> wf2=imaginary(wf) [IDL> tvscl, [wf1,wf2] IDL>

```
function compute_rms, cube
; cube: cube of wavefronts (square wf, no pupil!)
n_wf = (size(cube))[3]
rms = fltarr(n_wf)
for i=0,n_wf-1 do begin
   toto = moment(cube[*,*,i], SDEV=dummy)
   rms[i] = dummy
endfor
rms_moy = mean(rms)
return, rms_moy
end
```



```
function wfcube, dim, length, L0, r0, lambda, n_wf
 use:
           = 128L
                       ; [px] wf dimension
 dim
                       ; [m] wf physical dimension
 length
           = 27.
= .1
: L0
                     ; [m] outerscale
                       ; [m] Fried parameter
 r0
 lambda
           = 500E - 9
                       : [m] r0 wavelength
 n wf
           = 100L
                       ; nb of generated wf
 print, wfcube(dim,length,L0,r0,lambda,n wf,filename,SEED=seed)
 -> prints the rms value
 sub-routines needed:
 wfgeneration.pro, calcul_rms.pro
 Marcel Carbillet [marcel.carbillet@unice.fr],
 lab. Lagrange (UCA, OCA, CNRS), Feb. 2018.
: Last modification: Feb. 2018
cube = fltarr(dim, dim, n_wf)
for i=0, n_wf/2-1 do begin
   wf = wfgeneration(dim, length, L0, r0, lambda, SEED=seed)
   cube[*,*,2*i] = float(wf)
   cube[*,*,2*i+1] = imaginary(wf)
endfor
rms = compute_rms(cube)
return, rms
end
```

(IDL - 3)

; call with: IDL> @Exo2 Diam =1.0 r0 =0.3 N = 10	<pre>; call with: IDL> .rn Exo2_main Diam =1.0 r0 =0.3 N = 10</pre>
<pre>J = (N+1)*(N+2)/2-1 Noll = .2944*J^(-sqrt(3)/2)*(Diam/r0)^(5./3) S = exp(-Noll) ; see result with: IDL> print, S</pre>	<pre>J = (N+1)*(N+2)/2-1 Noll = .2944*J^(-sqrt(3)/2)*(Diam/r0)^(5./3) S = exp(-Noll)</pre>
<i>batch</i> : all variables are accessible.	<pre>end ; see result with: IDL> print, S</pre>
	<i>main</i> : idem (« .r » : run ; « .rn » : run new).
<pre>; call with: IDL> .rn Exo2_proc ; IDL> Exo2_proc, Diam, r0, N, S ; with, e.g: Diam=1.0, r0=0.3, N=10, S undefined</pre>	
pro Exo2_proc, Diam, r0, N, S	; call with: IDL> .rn Exo2_func
<pre>J = (N+1)*(N+2)/2-1 Noll = .2944*J^(-sqrt(3)/2)*(Diam/r0)^(5./3) S = exp(-Noll)</pre>	<pre>; IDL> print, Exo2_func(Diam, r0, N) ; with, e.g: Diam=1.0, r0=0.3, N=10 function Exo2_func, Diam, r0, N</pre>
end ; see result with: IDL> print, S	<pre>J = (N+1)*(N+2)/2-1 Noll = .2944*J^(-sqrt(3)/2)*(Diam/r0)^(5./3) S = exp(-Noll)</pre>
<i>procedure</i> : (input/output) parameters are accessible, but variables defined within	return, S end
the procedure are not.	<i>function</i> : no output parameters, inside variables not

accessible, result of the function returned.