

ZPX World Coordinate System

The ZPX World Coordinate System is a non-standard system for evaluating celestial coordinates from the image pixel coordinates. It follows the the FITS conventions for a zenithal polynomial projection (ZPN) but adds an additional two-dimensional polynomial distortion term to the evaluation. This discussion concentrates on the non-ZPN distortion extension and assumes the reader understands the FITS WCS conventions including applying a zenithal polynomial projection. The reference for the FITS WCS standard for undistorted celestial coordinates systems is [Representations of celestial coordinates in FITS](#) Calabretta, M. R., and Greisen, E. W., *Astronomy & Astrophysics*, **395**, 1077-1122, 2002. ([PDF](#), [HTML](#)).

One thing to note is that generally the radial "pin-cushion" distortion included in the ZPN projection is generally fixed and possibly defined by known optical design terms. For NOAO Mosaic data, each WCS calibration exposure has different distortion terms from fitting the refraction and other small effects but the calibration does not change the radial polynomial coefficients.

The ZPN World Coordinate System projection has a FITS keyword representation as illustrated in figure 1.

Figure 1: Sample header with ZPX WCS projection

```
WCSAXES =                2 / Number of WCS axes
CTYPE1   = 'RA---ZPX'      / Coordinate type
CTYPE2   = 'DEC--ZPX'      / Coordinate type
CRVAL1   =    320.687374999995 / Coordinate reference value
CRVAL2   =    36.9085555555556 / Coordinate reference value
CRPIX1   =    4167.56175625891 / Coordinate reference pixel
CRPIX2   =    4120.25894749731 / Coordinate reference pixel
CD1_1    = -5.2588308681025E-8 / Coordinate matrix
CD2_1    = -7.2772379161132E-5 / Coordinate matrix
CD1_2    = -7.2753930850119E-5 / Coordinate matrix
CD2_2    = -1.8637632244742E-8 / Coordinate matrix
WAT0_001 = 'system=image'   / Coordinate system
WAT1_001 = 'wtype=zpx axtype=ra projp0=0. projp1=1. projp2=0. projp3=337.74 proj'
WAT1_002 = 'p4=0. projp5=632052. lngcor = "3. 3. 3. 2. 0.001876397956622823 0.29'
WAT1_003 = '99113930557312 0.1542460039112511 0.3032873851581314 1.9247409545894'
WAT1_004 = '95E-5 -1.348328290485618E-5 1.414186703253352E-4 -1.792784764381400E'
WAT1_005 = '-4 -1.276226238774833E-4 4.339217671825231E-4 "'
WAT2_001 = 'wtype=zpx axtype=dec projp0=0. projp1=1. projp2=0. projp3=337.74 pro'
WAT2_002 = 'jp4=0. projp5=632052. latcor = "3. 3. 3. 2. 0.001876397956622823 0.2'
WAT2_003 = '999113930557312 0.1542460039112511 0.3032873851581314 9.963957331149'
WAT2_004 = '402E-5 -1.378185066830135E-4 1.559892401479664E-4 -8.280442729203771'
WAT2_005 = 'E-4 3.966701903249366E-4 0.001678960379199465 "'
```

The WCSAXES keyword (possible seen as WCSDIM) will always be 2. That this is a ZPX projection is indicated by the CTYPE keywords. These keywords also indicate that the first image axis corresponds to RA and the second to DEC.

The ZPX projection is evaluated as follows.

1. Compute the first order standard coordinates ξ and η from the linear part of the solution stored in CRPIX and the CD matrix.

$$\begin{aligned}\xi &= \text{CD1_1} * (x - \text{CRPIX1}) + \text{CD1_2} * (y - \text{CRPIX2}) \\ \eta &= \text{CD2_1} * (x - \text{CRPIX1}) + \text{CD2_2} * (y - \text{CRPIX2})\end{aligned}$$

2. Add the non-linear part of the projection using the coefficients in the WAT keywords as described below.

$$\begin{aligned}\xi' &= \xi + \text{lngcor}(\xi, \eta) \\ \eta' &= \eta + \text{latcor}(\xi, \eta)\end{aligned}$$

3. Apply the zenithal polynomial projection to ξ' and η' as described in *Calabretta and Greisen* where the P_m coefficients in that paper are given by the `projpm` coefficients in the WAT keywords. The reference tangent point given by the CRVAL values lead to the final RA and DEC in degrees. Note that the units of ξ , η , `lngcor`, and `latcor` are also degrees.

The coefficients for the zenithal polynomial (the P_m) and the non-linear polynomial distortion functions `lngcor( $\xi$ , $\eta$ )` and `latcor( $\xi$ , $\eta$ )` are stored as FITS keywords under the indexed `WATj_nnn` keywords. The `j` refers to the image axis and the `nnn` give a sequence number. The cards for a particular image axis are sorted by the sequence number and then concatenated together into one long string. Take care not to add spaces between the concatenated strings since the coefficients may be split across strings.

The long string for each image axis is composed of a set of keyword/value pairs where the value is quoted if it contains whitespace. Figure 2 shows how the WAT keywords in figure 1 would be decomposed into parameters and coefficients.

Figure 2: Decomposing the WAT keywords from figure 1

AXIS 1	AXIS 2
-----	-----
wtype=zpx	wtype=zpx
axtype=ra	axtype=dec
projp0=0.	projp0=0.
projp1=1.	projp1=1.
projp2=0.	projp2=0.
projp3=337.74	projp3=337.74
projp4=0.	projp4=0.
projp5=632052.	projp5=632052.
lngcor=	latcor=
3.	3.
3.	3.
3.	3.
2.	2.
0.001876397956622823	0.001876397956622823
0.2999113930557312	0.2999113930557312
0.1542460039112511	0.1542460039112511
0.3032873851581314	0.3032873851581314
1.924740954589495E-5	9.963957331149402E-5
-1.348328290485618E-5	-1.378185066830135E-4
1.414186703253352E-4	1.559892401479664E-4
-1.792784764381400E-4	-8.280442729203771E-4
-1.276226238774833E-4	3.966701903249366E-4
4.339217671825231E-4	0.001678960379199465

The list of coefficients are interpreted as follows.

1. The first number is the function type encoded as 1=chebyshev, 2=legendre, 3=polynomial. The example has a function of type 3 which is the simple polynomial.
2. The next two numbers represent the "order" of the function in ξ and η . The order is the one less than the highest polynomial power. The powers are represented below by m and n such as $m = 0$ to $\xi_{\text{order}} - 1$ and $n = 0$ to $\eta_{\text{order}} - 1$. In the example the orders are 4 which means cubic polynomials ($m=0$ to 3 and $n=0$ to 3).
3. The next (fourth) number specifies the type of cross-terms encoded as 0=no cross-terms, 1=full cross-terms, 2=half-cross terms. The cross-terms are terms of $\xi^m \eta^n$ where m and n are non-zero. Full cross-terms mean that both m and n will go to their maximum values independently while half-cross terms mean that $m + n$ will only go to the maximum of $\xi_{\text{order}} - 1$ and $\eta_{\text{order}} - 1$.
4. The next 4 numbers describe the region of validity of the fits in ξ and η space, e.g. ξ_{min} , ξ_{max} , η_{min} , η_{max} . They are used to compute normalized values for ξ and η used in the chebyshev and legendre polynomial functions:

```

xin = (2 * xi - (ximax + xmin)) / (ximax - xmin)
etan = (2 * eta - (etamax + etamin)) / (etamax - etamin)

```

5. The remaining terms are the coefficients of the polynomial terms. The functions are evaluated by summing polynomial terms $P_{mn}(xi,eta)$ multiplied by the coefficients C_{mn} as

```

lngcor(xi,eta) = sum (Cmn * Pmn(xi,eta))
latcor(xi,eta) = sum (Cmn * Pmn(xi,eta))

```

Representing the coefficients as C_{mn} for the polynomials P_{mn} , where m and n are the powers of xi and eta , they are ordered as

```

C00
C10
C20
C30
...
C01
C11
C21
C31
...
C02
C12
C22
C32
...
C03
C13
C23
C33
...

```

In the example with the half cross-terms and orders of 4 the ten coefficients would be C_{00} , C_{10} , C_{20} , C_{30} , C_{01} , C_{11} , C_{21} , C_{02} , C_{12} , and C_{03} .

The polynomials P_{mn} are defined below. The chebyshev and legendre polynomials are defined iteratively as functions of the normalized coordinates defined earlier.

```

Pmn = xi ** m * eta ** n      (polynomial)

Pmn = Pm(xin) * Pn(etan)      (chebyshev)

P0(xin) = 1.0
P1(xin) = xin
Pm+1(xin) = 2.0 * xin * Pm(xin) - Pm-1(xin)

P0(etan) = 1.0
P1(etan) = etan
Pn+1(etan) = 2.0 * etan * Pn(etan) - Pn-1(etan)

Pmn = Pm(xin) * Pn(etan)      (legendgre)

P0(xin) = 1.0
P1(xin) = xin
Pm+1(xin) = ((2m+1) * xin * Pm(xin) - m * Pm-1(xin)) / (m+1)

P0(etan) = 1.0
P1(etan) = etan
Pn+1(etan) = ((2n+1) * etan * Pn(etan) - n * Pn-1(etan)) / (n+1)

```

In the example with with a simple polynomial the functions would be given as follows.

```
lngcor/latcor = C00
+ C10 * xi      + C20*xi^2      + C30 * xi^3
+ C01 * eta     + C02*eta^2     + C03 * eta^3
+ C11 * xi*eta  + C21 * xi^2*eta + C12 * xi*eta^2
```