# A perturbation theory for the shape of central force orbits 

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#### Abstract

The two body central force orbit can be solved exactly only for the gravitational and simple harmonic oscillator potentials. When one discusses nonlinear oscillators, the trajectory in space-time can be found by various kinds of perturbative techniques- one of the most prominent ones being the Lindstedt-Poincare perturbation theory. In this work we show that a Lindstedt-Poincare like perturbation theory can be set up for the shape of a general central force orbit by working round a circular orbit. One also gets an answer for spatial frequency by this process. The effectiveness of our technique is checked against numerical simulations.


## Introduction

We consider the dynamics of a particle of mass ' $m$ ' moving in a central force field where the force is taken to be of the form $F=-m \lambda r^{-n}$, where $n$ is any number such that bound orbits exist. The distance of the particle from the center of force is ' $r$ ' and ' $\lambda$ ' is the interaction strength. The conservation of the angular momentum (magnitude ' $l$ ' per unit mass) restricts the particle to a plane. In terms of the polar co-ordinates ' $r$ ' and ' $\theta$ ', we have $(u=1 / r)$

$$
\begin{equation*}
\frac{d^{2} u}{d \theta^{2}}+u=\frac{\lambda}{l^{2}} u^{n-2} \tag{1}
\end{equation*}
$$

Our perturbation theory is set up around the circular orbit characterized by $u_{0}=\left(\frac{\lambda}{l^{2}}\right)^{\frac{1}{3-n}}$. The energy of the orbit is $E_{c}=\frac{1}{2} l^{2} u_{0}^{2} \frac{n-3}{n-1}$. Appropriate modifications are necessary for $n=1$. The deviation $u_{1}$ from the circular orbit defines the dimensionless quantity $X=\frac{u_{1}}{u_{0}}$. The variable $X$ satisfies the dynamics

$$
\begin{equation*}
\frac{d^{2} X}{d \theta^{2}}+(3-n) X=\sum_{k=2}^{\infty}{ }^{n-2} C_{k} X^{k} \tag{2}
\end{equation*}
$$

The energy is expressed in terms of $X$ as

$$
\begin{equation*}
\Delta E=E-E_{c}=\frac{1}{2} l^{2} u_{0}^{2}\left[\left(\frac{d X}{d \theta}\right)^{2}+2 X+X^{2}-\frac{2}{n-1}\left\{(1+X)^{n-1}-1\right\}\right] \tag{3}
\end{equation*}
$$

We have thus reduced the orbit equation formally to an anharmonic oscillator equation with coordinate $X$ and timelike variable $\theta$. The order of perturbation theory is determined by how many powers of $X$ is retained. In some ways this is another example of a traditional perturbation theory being used in an unexpected situation [1].

## Results and Discussion

The orbit upto second order in $\epsilon$ is (initial conditions suitably chosen)

$$
\begin{equation*}
u=\left(\frac{\lambda}{l^{2}}\right)^{\frac{1}{3-n}}\left[1-\epsilon^{2} \frac{n-2}{4}+\epsilon \cos (\sqrt{3-n} \theta)+\epsilon^{2} \frac{n-2}{12} \cos (2 \sqrt{3-n} \theta)\right] \tag{4}
\end{equation*}
$$

One gets a spatial frequency $\Omega=\sqrt{3-n}$ within this order. We get corrections to this as we go to higher order. Note, $\epsilon$ is the order of amplitude of $X$ and hence is the perturbation parameter. Our results agree with the exact solutions for $n=2$ and $n=-1$. The comparison between our perturbation theory result and the numerically obtained trajectory, spatial frequency is shown in Figure 1 .


Figure 1: (a) Plot of $\frac{1}{r}$ as a function of $\theta$ for $n=1, \epsilon=0.5, u_{0}=1$ (b) Plot of $\Omega$ as a function of $\epsilon$ for $n=2.5$

## References

[1] T Shah, R.Chattopadhyay, K.Vaidya, S.Chakraborty (2015) Conservative Perturbation Theory for Non Conservative Systems. Phys Rev E 92062927

