Application of the *t*-model of optimal prediction to the estimation of the rate of decay of solutions of the Euler equations in two and three dimensions

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Abstract

The "t-model" for dimensional reduction is applied to the estimation of the rate of decay of solutions of the Burgers equation and of the Euler equations in two and three space dimensions. The model was first derived in a statistical mechanics context, but here we analyze it purely as a numerical tool and prove its convergence. In the Burgers case the model captures the rate of decay exactly, as was already previously shown. For the Euler equations in two space dimensions, the model preserves energy as it should. In three dimensions, we find a power law decay in time and observe a temporal intermittency.

1 Introduction

Despite the rapid increase in available computational power there are still many systems which cannot be studied numerically without prior simplification. In earlier work [1, 2], we and others have derived methods for reducing the number of variables one has to solve for in complex problems, based on statistical projections. A special case, a long memory model called the "t-model", was thought to be particularly applicable to problems in fluid dynamics [3, 4], where temporal correlations decay slowly. An earlier application [5] of the t-model to the estimation of the rate of decay of solutions of the Burgers equation yielded remarkably accurate results.

In the present paper we use the t-model equations to reduce the number of variables in spectral methods and prove its convergence as the number of Fourier components increases . We then apply it to the estimation of the rate of decay of solutions of Euler's equations in two and three space dimensions. We do not address the claim implicit in earlier work, that the t-model equations may yield acceptable results even when the number of variables remains finite. The results we obtain are, however, surprisingly accurate, and a full analysis may well have to go through some version of the arguments on the basis of which the t-model was originally derived. Note that unlike previous damping methods for allowing spectral calculations to proceed to significant time spans (e.g. [6, 7, 8, 9, 10, 11, 12], the t-model equations contain no adjustable parameters and is guaranteed to remain stable.

The paper is organized as follows. In Section 2.1 we present the derivation of the *t*-model. In Section 2.2 we prove some results about its behavior for systems that conserve the L_2 norm of the solution and construct numerical methods that respect these properties. In Section 3, the *t*-model for the 3D Euler equations is constructed. In Section 4, we apply the *t*-model to the 1D inviscid Burgers equation and the 2D and 3D Euler equations and discuss how the numerical results compare to the known theoretical results.

2 The *t*-model

We begin with a system of ordinary differential equations

$$\frac{d}{dt}v(t) = f(v(t), w(t)), v(0) = x$$
(1)

$$\frac{d}{dt}w(t) = g(v(t), w(t)), w(0) = y$$
(2)

Here $x, v \in \mathbb{R}^n$, $y, w \in \mathbb{R}^m$ and $f : \mathbb{R}^n \times \mathbb{R}^m \to \mathbb{R}^n$, $g : \mathbb{R}^n \times \mathbb{R}^m \to \mathbb{R}^m$, and t is time. We think of v as the slow (resolved) variables and of w as the fast (unresolved) variables.

We assume that the system (1)-(2) conserves energy and that the energy is given by

$$E = \frac{1}{2}(\|v\|^2 + \|w\|^2).$$
(3)

Here ||v|| and ||w|| are the norms corresponding to the inner products $(v, v') = \sum_{i=1}^{n} v_i v'_i$ and $(w, w') = \sum_{i=1}^{m} w_i w'_i$. It follows from the conservation

of energy that

$$(v, f(v, 0)) = 0 (4)$$

$$||f(v,0)||^2 + (v, Df_v(v,0) \cdot f(v,0)) = 0$$
(5)

$$||g(v,0)||^2 + (v, Df_w(v,0) \cdot g(v,0)) = 0$$
(6)

for all $v \in \mathbb{R}^n$. Here the $n \times m$ matrix $D_w f(v, 0)$ consists of the derivatives of f(v, w), evaluated at w = 0.

To establish (4) we differentiate both sides of (3) with respect to t and use (1),(2). This gives

$$\frac{d}{dt}E = (v, f(v, w)) + (w, g(v, w)) = 0.$$
(7)

Since v, w can be given any initial values we see that (7) is an identity in v, w. In particular it holds when w = 0 so (v, f(v, 0)) = 0 for all v.

To prove (5) we use a variational argument. Let $a \in \mathbb{R}^n$. Since (4) remains true when we replace v by $v + \epsilon a$ it follows from Taylor's formula that

$$0 = (v + \epsilon a, f(v + \epsilon a, 0)) = (v, f(v, 0)) + \epsilon [(a, f(v, 0)) + (v, D_v f(v, 0) \cdot a)] + O(\epsilon^2).$$

But (v, f(v, 0)) = 0 so dividing by ϵ , letting $\epsilon \to 0$ and setting a = f(v, 0) yields (5).

The proof of (6) is similar. Let $w = \epsilon b = \epsilon g(v, 0)$. Using Taylor's formula in (7) we obtain

$$0 = (v, f(v, \epsilon b)) + (\epsilon b, g(v, \epsilon b)) = (v, f(v, 0)) + \epsilon [(v, D_w f(v, 0) \cdot b) + (b, g(v, 0))] + O(\epsilon^2).$$

To get (6) we divide by ϵ and let $\epsilon \to 0$.

2.1 Derivation of the *t*-model

In this section we will derive and analyze approximations for systems which conserve energy and which can be written as (1)-(2).

Let y = 0. Since the energy is conserved, w(t) = O(t). Expanding f(v, w) around w = 0 we see that

$$\frac{d}{dt}v(t) = f(v(t), 0) + D_w f(v(t), 0) \cdot w(t) + O(t^2)$$
(8)

with y = 0, equation (2) gives

$$w(t) = \int_0^t g(v(\tau), w(\tau)) d\tau$$

= $\int_0^t g(v(\tau), 0) d\tau + \int_0^t O(\tau) d\tau$
= $\int_0^t g(v(t), 0) d\tau + \int_0^t O(t - \tau) d\tau + O(t^2)$
= $tg(v(t), 0) + O(t^2)$

Inserting the expression for w in (8) and disregarding the $O(t^2)$ terms we arrive at the *t*-model

$$\frac{d}{dt}v(t) = f(v(t), 0) + tD_w f(v(t), 0) \cdot g(v(t), 0).$$
(9)

It is called the *t*-model because it contains the factor t, and because it can be derived—for the cases we are interested in—as the zero variance limit of the *t*-damping equations studied by Chorin, Hald and Kupferman [1].

2.2 Properties of the *t*-model and associated numerical methods

The energy for a solution of the *t*-model is not constant, but decreases. Indeed, it follows from (4),(6),(9) that

$$\frac{d}{dt}\frac{1}{2}\|v\|^2 = (v, f(v, 0)) + t(v, D_w f(v, 0) \cdot g(v, 0))$$
$$= -t\|g(v, 0)\|^2.$$
(10)

Thus the last term in (9) acts as a (non-linear) viscosity term. Similar results have been obtained for the *t*-damping method applied to Hamiltonian systems, see [2].

To solve eq.(9) we look for numerical methods where the energy decreases in each time step. Let F(v(t), t) denote the right-hand side of eq.(9) and consider Runge-Kutta methods of the form

$$k_i = F\left(v^n + h\sum_{j=1}^s a_{ij}k_j, t_n + hc_i\right)$$
$$v^{n+1} = v^n + h\sum_{i=1}^s b_i k_i$$

with $t_n = nh$, n = 0, 1, ... Set $V_i = v^n + h \sum_{j=1}^s a_{ij} k_j$ for i = 1, ..., s.

Theorem Let $b_i a_{ij} + b_j a_{ji} - b_i b_j = 0$ for i, j = 1, ..., s and assume that

$$\sum_{i=1}^{s} b_i = 1 \quad , \quad \sum_{i=1}^{s} b_i c_i = \frac{1}{2} \tag{11}$$

with $b_i, c_i \ge 0$. There is a \tilde{V} in the convex hull of V_1, \ldots, v_s such that

$$\frac{1}{2} \|v^{n+1}\|^2 - \frac{1}{2} \|v^n\|^2 = -h\left(t_n + \frac{h}{2}\right) \|g(\tilde{V}, 0)\|^2$$

Remark A numerical method that satisfies the assumptions in the theorem will be symplectic and at least second order. The simplest example is the implicit midpoint rule. It has $s = b_1 = 1$ and $a_{11} = c_1 = \frac{1}{2}$. Methods of higher order (4,5,6,8) can be found in [13] [p.207, p.209, p317]. **Proof** We begin by expanding $||v^{n+1}||^2$. After adding and subtracting

Proof We begin by expanding $||v^{n+1}||^2$. After adding and subtracting $h \sum_j a_{ij} k_j$ we get

$$(v^{n+1}, v^{n+1}) = (v^n, v^n) + h \sum_{i} b_i(k_i, v^n + h \sum_{j} a_{ij}k_j) + h \sum_{j} b_j(v^n + h \sum_{i} a_{ji}k_i, k_j) - h^2 \sum_{ij} [b_i a_{ij} + b_j a_{ji} - b_i b_j](k_i, k_j)$$

The last sum vanishes. Using the definitions of v_i and k_i yields

$$\|v^{n+1}\|^2 - \|v^n\|^2 = h \sum_i b_i \left(F(V_i, t_n + hc_i), V_i\right) + h \sum_j b_j \left(V_j, F(V_j, t_n + hc_j)\right).$$

Now V_i are real, so the two sums are equal. Consequently

$$\|v^{n+1}\|^2 - \|v^n\|^2 = 2h \sum_j b_j \left(V_j, f(V_j, 0) + (t_n + jc_j) D_w f(V_j, 0) \cdot g(V_j, 0)\right)$$

= $-2h \sum_j b_j (t_n + hc_j) \|g(V_j, 0)\|^2$

where we have used (4),(6). Let

$$G_0 = \min \|g(V_j, 0)\|^2 = \|g(V_{j_0}, 0)\|^2$$

$$G_1 = \max \|g(V_j, 0)\|^2 = \|g(V_{j_1}, 0)\|^2.$$

Using (11) we conclude that

$$-2h\left(t_n + \frac{h}{2}\right)G_1 \le \|v^{n+1}\|^2 - \|v^n\|^2 \le -2h\left(t_n + \frac{h}{2}\right)G_0.$$

Set $\tilde{V} = \theta V_j + (1 - \theta) V_j$. Since g(V, 0) is continuous there is a $\Theta \in [0, 1]$ such that

$$||v^{n+1}||^2 - ||v^n||^2 = -2h\left(t_n + \frac{h}{2}\right)||g(\tilde{V}, 0)||^2.$$

This completes the proof.

3 The *t*-model for the Euler equations

The Euler equations describe the flow of an incompressible, inviscid fluid in two or three dimensions. Here we look at flows in a cube with periodic boundary conditions and consider two kinds of approximations. First we use the Fourier method to obtain approximate solutions of Euler's equations. This leads to a large system of ordinary differential equations. Secondly we use the *t*-model to reduce the number of variables. The questions are: Does this method converge and what are the numerical results?

The Euler equations for 3 dimensional flows are

$$\partial_t u + (u \cdot \nabla) u = -\nabla p \tag{12}$$

$$\nabla \cdot u = 0 \tag{13}$$

where u = u(x,t) is the velocity, p is the pressure and t is time. We look for 2π periodic solutions and use the notation $u \cdot \nabla = \sum_j u^j \partial_j$ where $u = (u^1, u^2, u^3)^T$, T means transpose and $\partial_j u = \partial u / \partial x_j$ for j = 1, 2, 3. By taking the divergence of (12) and using (13) we see that

$$-\Delta p = \nabla \cdot (u \cdot \nabla)u$$

where $\Delta = \sum_j \partial_j^2$. Set $\int p = 0$. We can then solve for p and conclude from (12) that

$$\partial_t u = -\left[(u \cdot \nabla)u + \nabla \cdot (-\Delta)^{-1} \nabla \cdot (u \cdot \nabla)u\right].$$
(14)

Next we expand u in a Fourier series

$$u(x,t) = \sum_{k} u_k(t)e^{ikx}.$$
(15)

It follows from (13) that $k \cdot u_k \equiv 0$ for $k \neq 0$. By combining (12) and (15) and proceeding formally we get for $k \neq 0$

$$\frac{d}{dt}u_k = -i\sum_{p+q=k}k \cdot u_p A_k u_q \tag{16}$$

where $A_k = I - kk^T/|k|^2$. (For more details, see e.g. [14]). Since *u* is real, $u_{-k} = \bar{u}_k$. Finally we assume that $u_0 \equiv 0$. As $\int u$ is constant in time this amounts to a restriction of the initial data.

We get the Fourier method by setting $u_k \equiv 0$ for $|k|_{\infty} > m$ and for all time and setting $k \cdot u_k = 0$ and $u_{-k} = \bar{u}_k$ for t = 0. In this way (16) becomes a closed system of ordinary differential equations that conserves energy and yields solutions that are incompresible. To express (16) in the form (1)–(2) we set $v_k = u_k$ if $|k|_{\infty} \leq n$, and $w_k = u_k$ if $n < |k|_{\infty} \leq m$. Let $F = \{|k|_{\infty} \leq n\}$ and $G = \{n < |k|_{\infty} \leq m\}$. If $k \in F$ then

$$\frac{dv_k}{dt} = -i \bigg[\sum_{\substack{p+q=k\\p\in F, q\in F}} k \cdot v_p A_k v_q + \sum_{\substack{p+q=k\\p\in F, q\in G}} k \cdot v_p A_k v_q + \sum_{\substack{p+q=k\\p\in G, q\in F}} k \cdot w_p A_k v_q + \sum_{\substack{p+q=k\\p\in G, q\in G}} k \cdot w_p A_k w_q \bigg].$$

The equation for dw_k/dt is the same, except that $k \in G$. Following the derivation of (9) we obtain the *t*-model for Euler's equations

$$\frac{d}{dt}v_k = -i\left[\sum_{\substack{p+q=k\\p\in F, q\in F}} k \cdot v_p A_k v_q + \sum_{\substack{p+q=k\\p\in F, q\in G}} k \cdot v_p A_k(-i)t \sum_{\substack{r+s=q\\r\in F, s\in F}} q \cdot v_r A_q v_s \right] + \sum_{\substack{p+q=k\\p\in G, q\in F}} k \cdot (-i)t \sum_{\substack{r+s=p\\r\in F, s\in F}} p \cdot v_r A_p v_s A_k v_q\right]$$

If $k \cdot v_k = 0$ and $v_{-k} = \bar{v}_k$ at t = 0, it holds for all time. Moreover, a direct calculation shows that the energy decays, i.e.

$$\frac{d}{dt}\frac{1}{2}\sum_{k\in F}|v_k|^2 = -t\sum_{k\in G}|\sum_{\substack{p+q=k\\p\in F, q\in F}}k\cdot v_pA_kv_q|^2.$$
(18)

This is the complex analogue of (10). Finally we compare the solution u(x,t) of Euler's equations with the solutions generated by the *t*-model. Let \mathbb{T}^d be

the torus of length 2π in \mathbb{R}^d with d = 2, 3 and set

$$v(x,t) = \sum_{k \in F} v_k(t) e^{ik \cdot x}$$

Theorem Assume that $u(t) \in H^s(\mathbb{T}^d)$ for $s \geq 3$, d = 2, 3 and $0 \leq t \leq T$. Then

$$||u(t) - v(t)||_{L^2(\mathbb{T}^d)} \le \frac{\operatorname{const}(T)}{n^{s-1}} \cdot \max_{0 \le t \le T} ||u||_{H^s(\mathbb{T}^d)}.$$

Proof Will be presented elsewhere.

Thus if the solution of Euler's equations is smooth for $t \leq T$ then the energy is constant, and the energy of the *t*-model should converge to the energy of Euler's equations. In our numerical experiments this holds for d = 2 and fails for d = 3. This suggests that the solutions of the Euler equation lose smoothness when d = 3, and may develop singularities.

4 Numerical results

In this section we present numerical results of the application of the t-model to the 1D inviscid Burgers equation and the 2D and 3D Euler equations. The equations of motion for the Fourier modes were solved by a Runge-Kutta-Fehlberg method ([13]) with the tolerance set to 10^{-10} . Note that due to the quadratic nonlinearity and the form of the t-model term (see Eq. 17), the right hand side of the equation for each Fourier mode in the t-model contains interactions with Fourier modes of at most double the wavevector. So, the ratio of the number of unresolved modes in G to the resolved modes in F is 1.

The different terms appearing in the right hand side of the equations for the reduced model can be computed in real space using the Fast Fourier Transform (FFT). Since for a reduced model of size N in each spatial direction we include N unresolved modes in each direction, the arrays involved in the FFTs should be of size 2N. The fact that the *t*-model term can be computed using the FFT makes the numerical implementation of the *t*-model computationally efficient. Moreover, the FFT calculations involved are dealiased by construction and thus no extra (e.g. 3/2 rule) dealiasing is needed. For a calculation involving N modes in each direction, i.e. N/2positive and N/2 negative, we perform FFTs of size 2N, i.e. N positive and N negative modes. But we are interested only on the right hand side of the equations for the first N/2 modes. This means (see [14]) that to avoid aliasing (in the *t*-model term) we need for the total number of modes M used

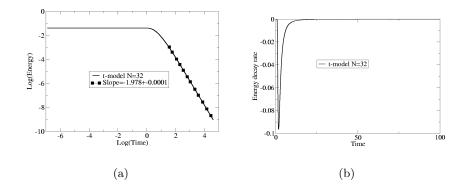


Figure 1: (a) Energy evolution of the *t*-model with N = 32 modes for the inviscid Burgers equation. (b) Evolution of the energy decay rate.

to satisfy the following inequality: $-N - N/2 \ge N/2 - 1 - M$ which yields $M \ge 2N$. But 2N is exactly how many modes we use in the FFTs, and thus the *t*-model term calculation through FFTs is dealiased by construction.

In order to study the asymptotic decay rate of the energy in the resolved modes, one has to evolve the system for long times. We evolved each case up to time t=100, so that we have enough points to perform an accurate estimate of the decay rate exponent. The need to perform calculations for long times, prevented us from conducting numerical experiments of larger size (within reasonable time) for the 3D Euler equations on a single processor workstation. However, the fact that the *t*-model term can be computed using the FFT means that the implementation of the model in any existing parallel spectral Navier-Stokes code is straightforward and we expect to report results of such simulations in the near future.

Figures 1, 2 and 3 present results of the application of the *t*-model to the 1D inviscid Burgers equation, the 2D Euler equations and the 3D Euler equations respectively. We present results for the evolution with time of the energy in the resolved modes and for the rate of energy decay (see Eq. 18). The numerical experiments are for resolved sets F (and the corresponding sets G) of size N = 32, $N = 32^2$, and $N = 32^3$ for the 1D, 2D and 3D cases respectively.

For the 1D inviscid Burgers equation $u_t + uu_x = 0$, the initial condition is $u_0(x) = \sin x$, which gives rise to a single shock wave at time T = 1. Until the moment of the formation of the shock wave the energy E is practically constant and soon after the well known t^{-2} energy decay law [15] is estab-

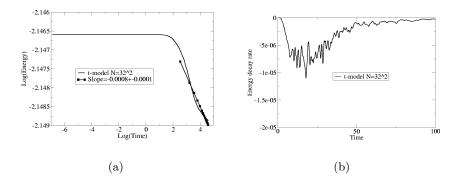


Figure 2: (a) Energy evolution of the *t*-model with $N = 32^2$ modes for the 2D Euler equations (note the extremely slow decay reflected in the very small slope of the linear fit). (b) Evolution of the energy decay rate. Again, note the very small values of the energy decay rate.

lished. The energy decay rate shown in Figure 1(b) reaches its peak around time $\tau = 1.219$ after which it starts decreasing. The t^{-2} energy decay regime is established soon after. The estimated exponent of the energy decay is estimated as -1.9781 ± 0.0001 , using about 15000 points. It is interesting that the right energy decay law is captured with only N = 32 Fourier modes.

For the 2D Euler equations the situation is drastically different. We present results for an incompressible, isotropic random initial condition with energy spectrum $E(k) = \exp(-2k)$ for the resolved modes and zero for the unresolved modes. After 100 units of time, the energy has decayed by 0.2%. In other words, the smooth initial condition does not lose its smoothness. In fact, as one can see in Figure 2(a), after the insignificant energy decay, the energy establishes a plateau which signifies the absence of drain of energy out of the resolved range of modes. A linear fit (in log-log coordinates) of the energy evolution gives a slope of -0.0008 ± 0.0001 , where we used about 15000 points.

The qualitative difference between the 2D and 3D case is striking. For the 3D Euler equations we use as initial condition the Taylor-Green vortex given by

$$u^{1}(x,0) = \sin(x_{1})\cos(x_{2})\cos(x_{3}),$$

$$u^{2}(x,0) = -\cos(x_{1})\sin(x_{2})\cos(x_{3}),$$

$$u^{3}(x,0) = 0$$

Note that the Taylor-Green initial condition is smooth having nonzero values

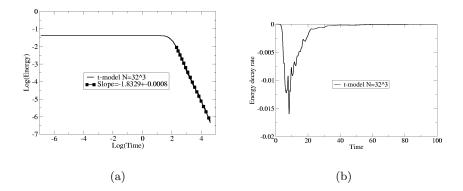


Figure 3: (a) Energy evolution of the *t*-model with $N = 32^3$ modes for the 3D Euler equations. (b) Evolution of the energy decay rate.

only for the Fourier modes with $k_i = \pm 1$, i = 1, 2, 3. A lot of numerical work (see e.g. [16] and references therein) has been devoted to the investigation of whether the solutions of the 3D Euler equations with the Taylor-Green initial condition blow up in finite time. All calculations show a rapid increase in the value of the maximum vorticity, but are hampered by the fact that they run out of resolution around time T = 5. Note that the first peak of the energy decay rate that we find is around time $\tau = 5.17$.

More interestingly, the decay of the energy appears to be organized in a collection of spikes of diminishing strength. This organization of the energy decay is reminiscent of the phenomenon of intermittency, i.e. bursts of activity followed by intervals of relative inaction on the part of the flow. Of course, the phenomenon of intermittency is not only of temporal nature, but has a spatial manifestation too. This is exhibited as concentration of the highest vorticity in small regions of the flow. The trend we observe in the decay of the energy seems to assign a specific purpose to the vorticity. Starting from a smooth initial condition, we have a steepening of the gradients in the field. This means that smaller scales are excited until the vorticity producing mechanism runs out of steam. Then we enter a period of relative inaction, until there is a restart of the mechanism of steepening. Energy is transferred again to the smaller scales and so forth. This scenario continues until there is no energy left in the large scales. After that, the flow just disintegrates and eventually comes to a halt. The purpose of vorticity mentioned above is to regulate the transfer of energy to the small scales [17]. This is reminiscent of the picture suggested by Moffatt, Kida and Okhitani [18] of the vortex structures acting as the "sinews of turbulence".

The slope of the linear fit (in log-log coordinates) for the energy decay is -1.8329 ± 0.0008 , where we have used about 15000 points. Estimates on the energy decay law for the 3D Euler equations are rare (see [19] and references therein). Moreover, all the estimates concern the infinite space case while we use periodic boundary conditions. For the infinite space case, under the assumption of complete self-preservation, i.e. self-similarity for all scales from 0 to ∞ , one finds that the energy should decay as t^{-1} [19]. If the assumption of complete self-preservation is not satisfied, the energy is expected to decay as $t^{-\alpha}$ where $\alpha > 1$. Note that the assumption of complete self-preservation is violated for the case of periodic boundary conditions and so the exponent of the energy decay should be larger than 1. However, it is not clear how the assumptions can be modified for this case. In 1D, the change of boundary conditions from infinite to periodic changed the exponent of energy decay from 1 to 2 [15]. It is not clear that this is also the case for 3D. If it is, then the numerical estimate -1.8329 for α becomes more plausible.

5 Conclusions

The problem of constructing reduced models for the Euler equations has been, and still is, a great challenge for scientific computing. The *t*-model proposed here should be considered a first step in deriving models directly from the equations without *ad hoc* approximations. It is based on numerical and physical observations about the behavior of the solution (more sophisticated reduced models for the Euler equations were constructed and simulated in [20]). Following [5], we tested the model on the 1D inviscid Burgers equation for an initial condition that gives rise to a shockwave. The model captures the right time of formation of the shock and the right rate of decay of the energy of the solution. For the 2D Euler equations, the model preserves the energy as it should since the solution remains smooth for all times. The numerical results for the Taylor-Green vortex for the 3D Euler equations produce rates of decay compatible with current thinking, and suggest that the solution loses its smoothness in finite time.

The terms appearing in the reduced model can be efficiently implemented by the use of the FFT on appropriate arrays. This makes the incorporation of the model in existing pseudospectral algorithms rather straightforward. We plan to apply the model in a parallel setting which will allow a better assessment of the properties of the flow field that is predicted by the model.

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