Supplementary Information for
A Statistical Dynamical Model to Predict Extreme Events and Anomalous Features in Shallow Water Waves with Abrupt Depth Change
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A. Details about the TKdV model non-dimensionalization

We provide more details about the derivation and properties for the truncated KdV equation with abrupt depth change used as the prediction model in the main text. A unified formulation with dependence on different parameter values is provided.

A.1. Mathematical formulation of the truncated KdV equation as a Hamiltonian system. The classic Korteweg-de Vries (KdV) equation (1) can be written in the standard form as

\[ u_t + uu_x + u_{xxx} = 0, \quad x \in [-\pi L_0, \pi L_0]. \]  

The state variable \( u(x, t) \) for the leading-order surface wave disturbance is defined on a periodic geometry of length \( 2\pi L_0 \). The KdV equation (1) can be written in the standard form as

\[ u_t + uu_x + u_{xxx} = 0, \quad x \in [-\pi L_0, \pi L_0]. \]

The corresponding Hamiltonian functional becomes the difference between the cubic term and the quadratic term

\[ H = \int_{-\pi L_0}^{\pi L_0} \left( \frac{1}{6} u^3 - \frac{1}{2} u_x^2 \right) dx. \]

Immediately, we have the conservation of the Hamiltonian in [2], \( H_t = \{ H, \mathcal{H} \} = 0 \). Besides the Hamiltonian \( H \), two other important conserved quantities in the KdV equation are the momentum \( M \) and energy \( E \) defined as

\[ M(u) = \int_{-\pi L_0}^{\pi L_0} u dx, \quad E(u) = \frac{1}{2} \int_{-\pi L_0}^{\pi L_0} u^2 dx. \]

In applications using the KdV equation for modeling water waves, it is useful to consider a normalized version of the equation. For convenience, the total momentum is normalized to zero \( M = 0 \) without loss of generality due to the Galilean invariance and the total energy is rescaled to unity \( E = 1 \). To achieve this, consider the following change of variables

\[ t = \tilde{t}, \quad x = L_0 \tilde{x} + M_0 \tilde{t}, \quad u = E_0^{1/2} L_0^{-1/2} \tilde{u} + M_0, \]

where \( M_0 = \int_{-\pi L_0}^{\pi L_0} u dx \) is the conserved total momentum and \( E_0 = \frac{1}{2} \int_{-\pi L_0}^{\pi L_0} u^2 dx - \pi L_0 M_0^2 \) is the conserved total energy from the original system [1], and \( L_0 \) defines the characteristic length scale of the system. The additional shift in time \( M_0 t \) in the new coordinate creates the Doppler shift from the non-zero mean momentum \( M_0 \). The normalized KdV equation with zero momentum and unit total energy for the scaled state variable \( \tilde{u}(\tilde{x}, \tilde{t}) \) becomes

\[ \tilde{u}_{\tilde{t}} + E_0^{1/2} L_0^{-3/2} \tilde{u}_{\tilde{x}x} + L_0^{-3} \tilde{u}_{\tilde{t}x} = 0, \quad \tilde{x} \in [-\pi, \pi], \]

where the system is defined on the normalized periodic domain, and we have the normalized conservatives

\[ \tilde{M} = \int_{-\pi}^{\pi} \tilde{u} \tilde{x} d\tilde{x} = 0, \quad \tilde{E} = \frac{1}{2} \int_{-\pi}^{\pi} \tilde{u}^2 d\tilde{x} = 1. \]

The corresponding Hamiltonian functional becomes the difference between the cubic term \( H_3 \) and the quadratic term \( H_2 \)

\[ \tilde{H}(\tilde{u}) = E_0^{1/2} L_0^{-3/2} H_3(\tilde{u}) - L_0^{-3} H_2(\tilde{u}), \quad H_3(\tilde{u}) = \frac{1}{6} \int_{0}^{2\pi} \tilde{u}^3 d\tilde{x}, \quad H_2(\tilde{u}) = \frac{1}{2} \int_{0}^{2\pi} \tilde{u}_x^2 d\tilde{x}. \]

Notice that the total momentum \( M_0 \) has no explicit contribution in the rescaled equation. \( E_0 \) is the total energy of the system depending on the domain size \( L_0 \), while \( E_0/(2\pi L_0) \) defines the energy density at unit scale length. The entire dynamics of the equation is determined by setting the two parameters \( (E_0, L_0) \). The choice of these parameter values will be discussed next based on a proper non-dimensionalization of the physical model. The advantage of adopting the normalized formulation [3] with Hamiltonian [4] is that it enables us to easily control the different cases with changing energy from a unified model setup.
Truncated KdV equation on the spectral domain. To investigate the turbulent dynamics generated from the KdV equation, usually a Galerkin projection $P_A$ is applied to the state variable $u$ with a high wavenumber truncation up to $\Lambda$

$$u_A(x,t) \equiv P_A u = \sum_{|k| \leq \Lambda} \hat{u}_k(t) e^{ikx}, \quad [5]$$

with in total $J = 2\Lambda + 1$ grid points. Accordingly, the truncated version of the KdV equation (TKdV) can be formulated by projecting the continuous equation [3] to the truncated subspace as

$$\frac{\partial}{\partial t} u_A + \frac{1}{2} E_0^{1/2} L^{-3/2}_0 \frac{\partial}{\partial x} P_A \left( u_A^2 \right) + L^{-3}_0 \frac{\partial^3}{\partial x^3} u_A = 0. \quad [6]$$

The additional projection in front of the quadratic term $u_A^2$ is used to remove the aliasing modes that go beyond the range $|k| > \Lambda$.

The TKdV equation [6] can be also written in the spectral domain for each Fourier mode $\hat{u}_k$ as

$$\frac{d\hat{u}_k}{dt} = -\frac{ik}{2} E_0^{1/2} L^{-3/2}_0 \sum_{|m| \leq \Lambda} \hat{u}_m^* \hat{u}_{m-k} + ik^3 L^{-3}_0 \hat{u}_k.$$

It can be shown that the three conserved quantities above are still conserved in this truncated system. With the spectral representation, the conserved momentum and energy set the constraints on the spectral coefficients

$$\mathcal{M}_A = \int_0^{2\pi} u_A dx = (2\pi) \hat{u}_0 = 0, \quad \mathcal{E}_A = \frac{1}{2} \int_0^{2\pi} P_A \left( u_A^2 \right) dx = 2\pi \sum_{k=1}^{\Lambda} |\hat{u}_k|^2 = 1.$$

The discretized Hamiltonian can be written accordingly as

$$\mathcal{H}_A = E_0^{1/2} L^{-3/2}_0 H_{3,\Lambda} - L^{-3}_0 H_{2,\Lambda}, \quad H_{3,\Lambda} = \frac{\pi}{3} \sum_{|m|,|n| \leq \Lambda} \hat{u}_m \hat{u}_n \hat{u}_{m+n}, \quad H_{2,\Lambda} = \pi \sum_{|k| \leq \Lambda} k^2 |\hat{u}_k|^2.$$

Especially, the Hamiltonian structure of the previous continuous equation is maintained in this semi-discrete TKdV equation.

The truncated equation [6] is still a Hamiltonian system with the corresponding discrete Hamiltonian $\mathcal{H}_A$. Furthermore, the truncated system [6] satisfies the Liouville property (2, 3), thus equilibrium statistical mechanics can be constructed based on the conserved quantities.

A.2. The rescaled TKdV model with non-dimensionalized parameters. Since the KdV equation is derived from the leading-order terms in the asymptotic expansion of the Euler equations, we find the sizes of the model parameters directly from the leading-order assumption for characterizing the physical problem from the experiments.

The KdV equation with depth dependence. The formulation for the KdV equation with a depth dependence $D_0$ can be introduced using the new coordinate system $(X, \xi)$ for the convenience of derivation as

$$X = \epsilon x, \quad \xi = D_0^{-1/2} x - t. \quad [7]$$

The equation is used to describe the far-field and long-time variability at $X = O(1), \xi = O(1)$ as $x \to \infty$ and $t \to \infty$. With the water depth dependence, the leading-order wave disturbance $\eta(\xi, X)$ obeys the generalized KdV equation (derivation can be found in (1))

$$D_0^{1/4} \eta_X + \frac{3}{2} D_0^{-5/4} \eta_{\xi\xi} + \frac{1}{6} D_0^{3/4} \eta_{\xi\xi\xi} = 0. \quad [8]$$

It needs to be emphasized that the spatial variable $X$ here plays the original role of the time variable $t$, and the new variable $\xi$ tracks the right-moving waves along the characteristics in the new equation.

Model parameters from non-dimensionalization. To derive the model parameter values, we first propose scales in the state variables according to the dimensional formulation. Then the parameter values in the rescaled model can be found by comparing the equation with its non-dimensionalized version. For the basic scales in the problem, we consider: i) the typical depth of the water tank $H_0$; ii) the surface wave disturbance amplitude $a$; and iii) the characteristic surface wavelength $\lambda_c = c_0/f_c$ with $f_c$ the characteristic forcing frequency and the gravity wave speed $c_0 = \sqrt{gH_0}$. The important characterizing parameters that can be measured from the experiments include

$$\epsilon = \frac{a}{H_0}, \quad \delta = \frac{H_0}{\lambda_c}, \quad D_0 = \frac{d}{H_0}. \quad [9]$$

Above, $\epsilon$ is the wave amplitude to water depth ratio; $\delta$ is the water depth to wavelength scale ratio; and $D_0$ defines the normalized wave depth ratio in the depth change from $d = H_0$ to $d < H_0$. The interpretations and reference values of these model parameters are based on the experimental setup (4).
The model parameters are determined by the non-dimensional measurements of wave location variable \( \lambda \), spatial variable \( x \), and temporal variable \( t \), based on the scales in (1):

\[
[\eta] = \epsilon H_0, \quad [x] = \epsilon^{-1/2} H_0, \quad [t] = \epsilon^{-1/2} H_0/c_0.
\]

We use \((\eta, X, \xi)\) to represent the variables in the dimensional formulation [8]; and \((\tilde{u}, \tilde{t}, \tilde{x})\) for the final non-dimensionalized variables in [3]. The above scales give the relations between the two sets of variables

\[
\tilde{u} = U^{-1} [\epsilon H_0] \eta, \quad \tilde{x} = L_d^{-1} \left[ \epsilon^{-\frac{2}{3}} D_0^\frac{2}{3} H_0 \right] \xi, \quad \tilde{t} = L_s^{-1} \left[ \epsilon^{-\frac{2}{3}} \frac{2}{3} H_0 \right] X,
\]

where \((U, L_d, L_s)\) are the scales in \((\eta, \xi, X)\) to normalize the original system to the standard form with unit energy and domain size \(2\pi\) as in [3]. We introduce the length scale \(L_d\) according to the characteristic wavenumber \(\lambda_c\) and the spatial scale \(L_s\) according to slow-varying far-field variables

\[
L_d = M\lambda_c = \frac{M\sqrt{d}}{f_0}, \quad L_s = \epsilon^{-3/2} H_0 = \frac{\sqrt{H_0^2}}{\sqrt{a^3}}, \quad \gamma = \frac{U}{a} = \frac{U}{\epsilon H_0}.
\]

where \(M\) is an integer representing a wave package of multiple wavelengths. Furthermore, \(\gamma\) defines the additional ratio that normalizes the total energy in \(u\) to unit. Substituting the above relations into the equation [8], we find the final non-dimensionalized equation about \(\tilde{u}(\tilde{t}, \tilde{x})\)

\[
\frac{\partial \tilde{u}}{\partial \tilde{t}} + \epsilon^{-\frac{1}{2}} \frac{3\delta}{4\gamma M} D_0^{-\frac{1}{3}} \frac{\partial}{\partial \tilde{x}} (\tilde{u}^2) + \epsilon^{-\frac{2}{3}} \frac{\delta^3}{6M^3} D_0^\frac{1}{3} \frac{\partial^3 \tilde{u}}{\partial \tilde{x}^3} = 0.
\]

Therefore, we formulate the non-dimensional equations [11] based on the non-dimensional parameters [9].

In the final step, we determine the parameter values in the computational models [3] by comparing with the corresponding non-dimensionalized form [11] with the measured quantities. After proper rearrangement and leaving the tildes in the normalized variables, we find the normalized KdV equation (as eqn. [1] in the main text)

\[
\partial_t u + E_0^{1/2} L_0^{-3/2} D_0^{-3/2} u u_x + L_0^{-3} D_0^{1/2} u_{xxx} = 0, \quad x \in [-\pi, \pi],
\]

The model parameters are determined by the non-dimensional measurements

\[
L_0 = 6 \frac{\delta}{\gamma} \left( M \frac{\epsilon^2}{\delta^2} \delta^{-1} \right), \quad E_0 = \frac{27}{2} \gamma^{-2} \left( M \frac{\epsilon^2}{\delta^2} \delta^{-1} \right).
\]

Notice that in the above model ‘\(t\)’ is used to represent either the far-field spatial variable ‘\(X\)’ and ‘\(x\)’ is used to represent the wave location variable ‘\(\lambda\)’. We also drop the ‘tildes’ in the variables for simplicity.

In addition, to estimate the parameter \(M\) for the computational domain size \(2\pi L_d = 2\pi M\lambda_c\) determined as \(M\)-multiple of the characteristic wavelength \(\lambda_c\), we consider the spatial discretization \(J = 2\lambda + 1 = 32\) so that the smallest resolved scale is comparable with the characteristic wavelength, that is, \(2\pi M\lambda_c/J = \lambda_c\). This estimation gives \(M = \frac{\lambda_c}{2\pi} \approx 5\), inferring a computational domain with 5 multiples of the characteristic wavelength. Using the reference experimental measurements, we can calculate the reference values for the model scales \(L_0\) and \(E_0\) used in the direct numerical simulations. The wave amplitude to depth ratio \(\epsilon\) changes in the range \([0.0024, 0.024]\), thus we can estimate the minimum and maximum values of the scales in Table S1. We adopt this basic guideline in choosing parameter values in the numerical confirmation of the theory in the main text.
B. Sampling strategy and statistical link for the invariant measures

Here we show the detailed strategy and algorithm used for computing the sampled Gibbs invariant measure and the statistical matching at the abrupt depth change shown in the main text. The invariant Gibbs measure is defined based on microcanonical ensemble in the quadratic energy \( E_{\Lambda} \) on the isosurface and canonical ensemble in the Hamiltonian \( H_{\Lambda} \) \((3, 5)\). The invariant Gibbs measure for the TKdV model \([12]\) about the normalized state variable \( u_{\Lambda} \) with unit energy can be written explicitly as

\[
G_{g_{\pm}}(u_{\Lambda}) = C_{g_{\pm}} \exp \left(-\theta \left\{ h_{\Lambda}^{\pm} \sum_{1 \leq |m| \leq \Lambda} \hat{u}_m \bar{u}_m^{*} u_{\Lambda}^{m+n} - h_{\Lambda}^{\pm} \sum_{1 \leq |m| \leq \Lambda} k^2 |\hat{u}_m|^2 \right\} \right) \delta \left( \pi \sum_{1 \leq |m| \leq \Lambda} |\hat{u}_m|^2 - 1 \right), \tag{14}
\]

with the coefficients \( h_{\Lambda}^{\pm} = \pi E_{0}^{1/2} L_0^{-3/2} D_{\pm}^{3/2} \) and \( h_{\Lambda}^{\pm} = \pi E_{0}^{-3} D_{\pm}^{1/2} \) depending on the model parameters. The proper statistical matching condition is proposed to connect the invariant measures in incoming and outgoing flow statistics with different inverse temperatures \( \theta^{\pm} \).

In predicting the statistical distribution of state variables after the abrupt depth change, we start with the assumption that we have full knowledge about the invariant distribution about the incoming flow field, \( G_{g_{-}} \) at equilibrium. Thus the parameters for \( \theta^{+} \) are determined with accuracy from the observations or statistical simulations in the incoming flow, and the total energy \( E_0 \) is conserved. The condition to get the outgoing flow statistics with parameter \( \theta^{+} \) is through the following matching condition

\[
\langle H_{\Lambda}^{+} \rangle_{G_{g_{-}}} = \langle H_{\Lambda}^{+} \rangle_{G_{g_{+}}}, \tag{15}
\]

with the Gibbs measures \( G_{g_{\pm}} \) before and after the abrupt depth change defined in \([14]\). Notice in the statistical matching condition \([15]\), we don’t require the statistical information for the incoming Hamiltonian \( H_{\Lambda}^{-} \). This is considering the near-Gaussian statistics in the incoming waves and its weak dependence on the Hamiltonian in the Gibbs measure. Therefore the Hamiltonian before the depth change is only used to construct the invariant measure, \( G_{g_{-}} \).

B.1. Sampling algorithm for the incoming Gibbs measure. First, the task is to compute the expectations \( \langle H^{+} \rangle_{g_{+}} \) under the proper invariant measures from \([14]\). The central question is how to find a method to properly sample the equilibrium distribution according to the chosen inverse temperature \( \theta \). One method to directly sample the Gibbs measure is described in \((2)\), where a Metropolis-Hasting Monte-Carlo sampling is applied. The following algorithm is adopted to sample the mixed canonical-microcanonical ensemble distribution \([14]\) for the numerical results in the main text:

\begin{algorithm}
1. Assign i.i.d standard Gaussian random variables to the \( \Lambda \)-dimensional complex valued samples, \( \vec{w} = \{ \hat{w}_1, \cdots, \hat{w}_{\Lambda} \} \in \mathbb{C}^{\Lambda} \); 
2. Normalize for each component \( \hat{u}_k = \sqrt{\hat{w}_k} \) with the total energy \( E = 2\pi \sum_{k=1}^{\Lambda} |\hat{w}_k|^2 \); 
3. Starting from \( \vec{u}_t = \vec{w} \), generate a new candidate \( \vec{w} \) according to a transition probability, \( g(\vec{w} | \vec{u}) \) (a simple choice is \( g \sim \exp \left(-\frac{\theta}{2} |\vec{w} - \vec{u}|^2 \right) \)), and normalize \( \vec{w} = \sqrt{\sum_{k=1}^{\Lambda} \hat{w}_k} \); 
4. Calculate the acceptance probability \( a = \min \left\{ 1, \exp \left(-\theta \left[ H(\vec{u}) - H(\vec{w}) \right] \right) \frac{g(\vec{w} | \vec{u})}{g(\vec{u} | \vec{w})} \right\} \); 
5. Generate a uniform random number \( r = \mathcal{U}[0, 1] \). Accept \( \vec{u}_{t+1} = \vec{w} \) if \( r \leq a \) and reject \( \vec{u}_{t+1} = \vec{u} \) if \( r > a \); 
6. Repeat the previous steps and sample at every \( n = 10 \) steps.
\end{algorithm}

We normalize the vector at each step to keep it on the fixed energy surface. It is also safer in practice to skip the first \( M = 10^{5} \) samples from the start for the system to reach stationary. Then the expectations can be calculated by averaging among all the samples from the above algorithms

\[
\langle H_{\Lambda}^{+} \rangle_{\mu} = \frac{1}{N} \sum_{t=1}^{N} H_{\Lambda}^{+}(\vec{u}_t) = E_{0}^{1/2} L_0^{-3/2} D_{+}^{3/2} \frac{1}{N} \sum_{t=1}^{N} H_{\Lambda}(\vec{u}_t) - L_0^{-3} D_{+}^{1/2} \frac{1}{N} \sum_{t=1}^{N} H_{\Lambda}(\vec{u}_t),
\]

where \( H_{\Lambda}(\vec{u}) \) computes the cubic term \( u_{\Lambda}^3 \) and \( H_{\Lambda}(\vec{u}) \) computes the wave slope energy \( u_{\Lambda}^2 \) according to the definition in \([2]\). The PDFs are achieved by “bin counting” from the histogram of the samples \( \{\vec{u}_t\}_{t=1}^{N} \). As one illustration, Figure S1 displays the sampled statistics, the PDFs for the Hamiltonian \( H_{\Lambda} \), the PDFs for the state variable \( u_{\Lambda} \), and the corresponding energy spectra for each Fourier mode. Typically we test the range of values of the inverse temperature from positive to negative \( \theta = 0.5, 0.25, 0, -0.25, -0.5 \). Supporting the claim in the main text, the negative inverse temperature \( \theta < 0 \) produces the proper physical regime for the incoming flow statistics, where the energy spectrum is decaying from larger scale to small scale. In contrast with \( \theta > 0 \), the small scale modes get more energetic and a flat top is generated in the PDF of \( u_{\Lambda} \) with two peaks. With \( \theta = 0 \), the Gibbs measure is the uniform distribution on the constant energy shell \( E_0 \), thus equipartition of energy in each mode is expected.
B.2. Computing downstream parameter from the statistical matching. In the second step, we need to enforce the statistical matching condition [15] to discover the corresponding downstream inverse temperature $\theta^+$. For the efficiency of the method in finding the optimal value of $\theta^+$, we use a secant method by solving a nonlinear equation following the algorithm:

**Algorithm 2** Recovering the downstream inverse temperature $\theta^+$ from the statistical matching condition

1. Pick up the upstream value of $\theta^-$ and compute the expectation under incoming flow Gibbs measure $\langle H^+_{\Lambda, \theta} \rangle_{\theta^=} -$;
2. Define a function of $\theta^+$ as the difference of the upstream and downstream expectations, $F(\theta^+) = \langle H^+_{\Lambda, \theta} \rangle_{\theta^=} - \langle H^+_{\Lambda, \theta} \rangle_{\theta^=} -$;
3. Find the solution of $F(\theta^+) = 0$ using the secant method, so that the optimal downstream value $\theta^+$ is found.

In addition, to remove stiffness with large values of $L_0$ and $E_0$, we introduce the normalized parameter $\tilde{\theta} = L_0^{-3/2} \theta$ so that $G_{\tilde{\theta}} \sim \exp \left(-\tilde{\theta} \left[ E_0^{1/2} D_0^{-3/2} H_3 - L_0^{-3/2} D_0^{1/2} H_2 \right] \right)$. In the numerical computations, we tune the parameter $\tilde{\theta}$ instead for the matching condition. Using this modified secant method, it is found that the optimal value can be usually reached after 2 to 5 iteration steps as long as the scheme is guaranteed non-stiff.

**Numerical comparison of the matching results.** The solid black lines in Figure S2 plot the downstream skewness $\kappa^+$ computed according to the matched downstream $\theta^+$ as shown in Fig. 1 of the main text. Here as a further comparison, the results with different total energy levels $E_0$ starting from various various values of incoming $\theta^-$ are compared. Further, Table S2 compares the incoming (with $\theta^-$) and outgoing (with $\theta^+$) wave statistics using the parameters achieved from the statistical matching condition. In the last two rows, the contributions of the cubic term $H_3$ and quadratic term $H_2$ in the Hamiltonian are compared in the incoming and outgoing flow statistics. It is found that in the upstream statistics $\theta^-$, the quadratic part $H_2$ is dominant, while in the downstream flow statistics, the cubic part becomes important in the total Hamiltonian, implying stronger non-Gaussian statistics.

As a further comparison, in the main text we also use the Gamma distribution

$$\rho(u; k, \alpha) = \frac{e^{-k} u^{-\alpha - 1}}{\Gamma(k)} \left( k + \alpha^{-1} u \right)^{k - 1} e^{-u^{-\alpha - 1}}, \quad \sigma^2 = k \alpha^2, \quad \kappa_3 = \frac{2}{\sqrt{k}}$$

to fit the downstream state $u^+$ according to the measured skewness $\kappa_3$ and variance $\sigma^2$ for the parameters $(k, \alpha)$. Table S2 also shows the fitting parameter values according to the downstream statistics with $\theta^+$. The Gamma fitting can also give an estimation for the kurtosis as $\kappa_4 = \frac{6}{k} + 3$, which is shown to have good agreement with the truth value $\kappa_4^+$ especially in the larger skewness cases.

**An explicit formula from the statistical matching condition.** In complementary to the main text, we offer a more detailed derivation about the statistical link between the upstream and downstream wave slope energy and the skewness. The statistical matching condition with the downstream Hamiltonian can be written as

$$\langle H^+_{\Lambda, \theta} \rangle_{\mu^=} = \langle H^+_{\Lambda, \theta} \rangle_{\mu^+}, \quad H^+_{\Lambda, \theta} = E_0^{1/2} L_0^{-3/2} D_0^{-3/2} H_3 - L_0^{-3} D_0^{1/2} H_2.$$

In general, we can have the expected value of $H_3$ and $H_2$ about any probability measure $\mu$ as

$$\langle H_3 \rangle_\mu = \frac{1}{\beta} \int_0^{2\pi} \langle u^3 \rangle_{\mu} \, dx = \frac{2\pi}{\beta} \sum_{j=1}^J \langle u^3 \rangle_{\mu^j}, \quad \langle H_2 \rangle_\mu = \frac{1}{2} \int_0^{2\pi} \langle u_x^2 \rangle_{\mu} \, dx.$$  

Above the quadratic term $H_2$ characterizes the slopes of the surface waves, $u_x$, in the incoming and outgoing flows. Beside, the skewness the state variable $u$ at one grid point is defined as the ratio between the third and second moments (note that in the
Table S2. Statistics computed from incoming and outgoing flow Gibbs measures with $E_0 = 100$. The skewness $\kappa_3$ and kurtosis $\kappa_4$ predicted from the Gibbs measures with different $\theta$ are compared. The outgoing statistics are compared with the Gamma function fitting with parameter $(k, \alpha)$ and the corresponding kurtosis $\kappa_4 = 6/k + 3$. The expected values of the cubic and quadratic terms in the Hamiltonian, $\bar{H}_3 = E_0^{1/2} L_0^{-3/2} D_v^{-3/2} \frac{1}{6} \int u^3$, and $\bar{H}_2 = L_0^{-3} D_v^{1/2} \frac{1}{2} \int u^2$ are also compared.

<table>
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<th>$\theta^-$</th>
<th>$E_0 = 100$</th>
<th>-0.5</th>
<th>-0.3</th>
<th>-0.1</th>
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<tbody>
<tr>
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<td>2.88</td>
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</tr>
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</table>

Above the total energy of the system is normalized to one in the computations. Now we use the two assumptions introduced in the main text:

- The upstream equilibrium measure $\mu_-$ has a near-Gaussian distribution, so that the third moments in the cubic term stay in a small value
  $$\langle H_3 \rangle_{\mu^-} = \frac{1}{6} \int_0^{2\pi} \langle u^3 \rangle_{\mu^-} dx = \epsilon;$$

- The downstream equilibrium measure $\mu_+$ is homogeneous at each physical grid point, so that the second and third moments are invariant at each grid point
  $$\langle u^3 \rangle_{\mu^+} = \sigma^3 = \pi^{-1}, \quad \langle u^3 \rangle_{\mu^+} = \sigma^3 \kappa_3 = \pi^{-\frac{3}{2}} \kappa_3.$$

With the homogeneous assumption above, the downstream cubic term can be simplified as
  $$\langle H_3 \rangle_{\mu^+} = \frac{1}{3} \sum_{j=1}^{j} \langle u^3 \rangle_{\mu^+} = \frac{1}{3} \pi^{-\frac{3}{2}} \kappa_3.$$

With the above assumptions, the above statistical matching condition can be simplified as

$$E_0^{1/2} L_0^{-3/2} D_v^{-3/2} \epsilon - L_0^{-3} D_v^{1/2} \langle H_2 \rangle_{\mu^-} = E_0^{1/2} L_0^{-3/2} D_v^{-3/2} \langle H_3 \rangle_{\mu^-} - L_0^{-3} D_v^{1/2} \langle H_2 \rangle_{\mu^-}$$

$$\Rightarrow E_0^{1/2} L_0^{-3/2} D_v^{1/2} \epsilon + \pi L_0^{-3} D_v^{1/2} \sum_{1 \leq |k| \leq \Lambda} k^2 (r_k^+ - r_k^-) = 1 \pi^{-1/2} E_0^{1/2} L_0^{-3/2} D_v^{3/2} \kappa_3,$$

with $r_k^- = \langle |\hat{u}_k|^2 \rangle_{\mu^-}$ the variance in upstream statistics and $r_k^+ = \langle |\hat{u}_k|^2 \rangle_{\mu^+}$ the statistics from downstream statistics. Therefore we can estimate the skewness of the state variable $u$ from the energy spectra in inflow and outflow statistics

$$\kappa_3^+ = 3 \pi^2 L_0^{-3} E_0^{-2} D_v^2 \sum_{1 \leq |k| \leq \Lambda} k^2 (r_k^+ - r_k^-) + 3 \pi^2 \epsilon.$$

With further assumption that the incoming flow gets zero third moment, then the skewness above can be calculated by

$$\kappa_3^+ = C \sum_{1 \leq |k| \leq \Lambda} k^2 (r_k^+ - r_k^-), \quad C = 3 \pi^2 L_0^{-3} E_0^{-\frac{1}{2}} D_v^2,$$
Above we proposed two approaches to predict the statistics in the downstream wave disturbance. The equilibrium statistical mechanics is an easy way to confirm the matching condition and get the equilibrium distribution in far-field long-time limit. Here we summarize the direct numerical methods used for the validation of the statistical theory in the main text, C. Symplectic scheme for direct numerical simulations of the TKdV equation

Here we summarize the direct numerical methods used for the validation of the statistical theory in the main text. In general, the strategy by using direct numerical simulations to confirm the anomalous statistics follows the steps:

1. Generate samples from the incoming equilibrium distribution $\mu_{\infty}$ according to Algorithm 1 for the incoming waves. The model parameters ($E_0, L_0, \Lambda$) are determined from the experimental measurements and statistics in the incoming flow;

2. Compute the expectation of the outflow Hamiltonian, $\langle H^+_{\Lambda} \rangle_{\mu_{\infty}}$ under the chosen incoming Gibbs measure $\mu_{\infty} = \mathcal{G}_{\theta^+};$

3. The outgoing flow statistics can be predicted by both the equilibrium statistical mechanics and the direct model simulations.

   (a) Equilibrium statistical mechanics: find the optimal outgoing flow parameter $\theta^+_{\ast}$ from matching the expectation in [15] with the inflow value according to Algorithm 2. Then the new equilibrium distribution $\mu_{\infty}^{\ast}$ after the abrupt depth change can be sampled using the same method;

   (b) Direct model simulations: run the dynamical model [12] with outflow depth $D_{\ast} < 1$ starting from the state sampled from the inflow distribution $\mathcal{G}_{\theta^+}$. The outgoing flow statistics can be captured by the solution from the direction simulations.

Above we proposed two approaches to predict the statistics in the downstream wave disturbance. The equilibrium statistical mechanics is an easy way to confirm the matching condition and get the equilibrium distribution in far-field long-time limit. On the other hand, the direct simulations of the dynamical model can offer estimates about the transient statistics before the far-field equilibrium state is reached. These two approaches can be used to confirm each other in the numerical results.

C.1. Conservation of energy and Hamiltonian form the symplectic scheme. For running the direct numerical simulations of the TKdV equations, a proper symplectic integrator is necessary to guarantee the conservation of the Hamiltonian and energy in time. As emphasized in the main text, the conservation property plays a central role in the statistical matching and the final equilibrium measure, otherwise the non-symplectic schemes introduce high numerical dissipations. The symplectic scheme used here for the time integration of equation [12] is the 4th-order midpoint method (6). Let $\mathbf{u}^n$ be the solution at time $t_n$ and $\mathbf{u}^{n+1}$
be the solution at next time step \( t_{n+1} = t_n + \Delta t \), and \( \mathbf{F} \) the discretized operator for the spectral modes. The midpoint method has two intermediate stages and the auxiliary vectors \( \mathbf{y}_1, \mathbf{y}_2 \)

\[
\begin{align*}
y_1 - u^n &= w_1 \Delta t \mathbf{F} \left( \frac{1}{2} [y_1 + u^n] \right), \\
y_2 - y_1 &= w_2 \Delta t \mathbf{F} \left( \frac{1}{2} [y_2 + y_1] \right), \\
u^{n+1} - y_2 &= w_3 \Delta t \mathbf{F} \left( \frac{1}{2} [u^{n+1} + y_2] \right),
\end{align*}
\]

with the time increments \( w_1 = \left( 2 + 2^{1/3} + 2^{-1/3} \right) / 3, w_2 = 1 - 2w_1, \) and \( w_3 = w_1 \). Then the PDFs can be counted from the solution \( u_{\Lambda} (t) \), and the statistical expectations can be calculated by averaging along the trajectory

\[
\langle F(u_{\Lambda}) \rangle = \frac{1}{T} \int_{t_0}^{t_0 + T} F(u_{\Lambda}(t)) \, dt,
\]

given ergodicity in the solution with \( u_{\Lambda}(t_0) \) already reaching the stationary state.

We compare the statistical spectra from the direct Monte Carlo model simulation with the prediction from the equilibrium statistical mechanics in Figure S3. The flow energy spectra computed both from the direct TKdV model simulations and from the Gibbs invariant measure in the equilibrium statistical mechanics are compared. Starting from two different \( \theta^- \) initial states, the direct model simulations offer the final equilibrium energy spectrum in a comparable decaying structure as predicted from the Gibbs measure and statistical matching condition. This first comparison offers the consistency in these two distinct approaches.

### C.2. Ergodicity and mixing properties in the TKdV solutions

Ergodicity is another crucial feature that needs to be confirmed in the TKdV equation solution. To achieve reasonable statistics from averaging the model time solutions, it is important to make sure that the solution is turbulent mixing. Following we display the direct numerical simulation solutions from the TKdV equation using initial states sampled from different inverse temperatures.

**Mixing solutions starting from negative inverse temperature \( \theta < 0 \).** In the main text, the results shown always use negative inverse temperature. As shown from \( (2, 5) \) and Figure S1 before, this is a reasonable choice to generate a decaying energy spectrum. Here as a further illustration, we show that the negative inverse temperature case also generates fast mixing solutions to calculate reasonable statistical averages.

In Figure S4 and S5, we display the single trajectory solutions for both downstream \( D_+ = 0.24 \) and upstream \( D_- = 1 \) flows. The initial state is taken from the samples in Algorithm 1 with a negative initial temperature \( \theta < 0 \). The other parameters are kept the same as in the main text. On the left panel, the contour plot for the time-series of \( u_{\Lambda} \) is shown for the downstream and upstream flows. Both cases display turbulent dynamics, while the downstream solution with \( D_+ < 0 \) generates more rare events and stronger turbulence.

For more detailed comparison, the right panels show one slice of the time-series of \( u_{\Lambda} (x_1) \) at one physical location as well as the first two Fourier modes, \( \hat{u}_1, \hat{u}_2 \), in the spectral domain. The rare events in the downstream solution in Figure S4 is obvious with high peaks representing the generation of skewed PDFs in the positive direction. In comparison, the upstream solution in Figure S5 displays almost symmetric values in positive and negative disturbances, implying near-Gaussian statistics in the incoming flow. The mixing properties are characterized by the autocorrelation and decorrelation time of the state variable. The decorrelation time \( T_k \) and absolute decorrelation time \( T^{abs}_k \) together with the autocorrelation function \( R_k \) in each spectral mode are defined as

\[
T_k = \left| \int R_k (t) \, dt \right|, \quad T^{abs}_k = \int \left| R_k (t) \right| \, dt, \quad R_k (t) = \langle \hat{u}_k (0) \hat{u}_k (t)^* \rangle / \langle |\hat{u}_k|^2 \rangle.
\]

Decaying autocorrelations in both the physical state and the spectral modes confirm the fast mixing and turbulent dynamics in both the tested regimes used in the main text (though weaker in upstream). We can observe the much faster time scales in...
the downstream flow in the autocorrelation functions in the corresponding time-series. The longer decorrelation times in the upstream flow indicate oscillating autocorrelations for a long time.

As an additional comparison, we show the power spectra of the upstream and downstream flows from the above test cases in Figure S6. The power spectrum can be calculated by the Fourier transform of the autocorrelation function

$$S(\lambda) = \int_{-\infty}^{\infty} R(\tau) e^{-i\lambda\tau} d\tau,$$

which offers the characterization of the decay in power at each time frequency. Again, consistent with the observations from experiments (4) and the numerics in the main text, the downstream power spectrum gets more energetic small time scales at low frequency modes and a slower decay rate in the high frequency modes compared with the upstream case. Also notice that the peak in the incoming flow power spectrum is taken place at a much lower frequency. This illustrates the slower mixing rate in the incoming flow state.

Non-mixing solutions starting from positive inverse temperature $\theta > 0$. In comparison, we also compute the solution starting from an initial state with a positive inverse temperature $\theta > 0$. In contrast to the previous case in Figure S5, the only difference in the present case is the initial state taken from the samples in an invariant measure with positive inverse temperature. Figure S7 displays the corresponding solution in this $\theta > 0$ case. In contrast to the previous mixing solution with negative inverse temperature, drastic difference performance is generated due to only the change in the initial state. The solution is recurrent in time with interacting small scales. As a result, the solution in stationary waves lacks the proper mixing property and the autocorrelation just oscillates for an extremely long time.

Accordingly in the autocorrelation functions and decorrelation times, the autocorrelation functions just keep oscillating and all the modes get long absolute decorrelation time (the decorrelation time is small due to the cancellation from the oscillations). This non-mixing feature with positive temperature supports the conclusion in the main text that this regime is not appropriate for correctly representing the physical statistics in the experimental setups. This simple test further confirms the unrealistic energy spectra in the positive temperature case with increasing energy in the smaller scales shown in (5) and Figure S1.

References

Fig. S5. Realization of the upstream flow solution with $D_0 = 1$ and initial state from samples with negative inverse temperature.

Fig. S6. Power spectra of the upstream and downstream states $u_X$.


Fig. S7. Realization of the upstream flow solution with $D_{u_0} = 1$ and initial state from samples with positive inverse temperature.