# HiSD Package: solscape V-1.0

A Python package for constructing solution landscapes using High-index Saddle Dynamics (HiSD). This toolkit enables numerical computation of saddle points and their hierarchical organization in dynamical systems.

# Installation & Import

Step 1: Download the Code First, download the code from GitHub:

GitHub Repository: https://github.com/HiSDpackage/solscape

**Step 2: Add the solscape-1.0 Directory to the System Path** After downloading the code, you need to add the path of the solscape-1.0 directory to the system path. This will allow you to access the package from anywhere on your system.

Use the following Python code to add the directory to the system path:

```
import sys
sys.path.append('/path/to/solscape-1.0')
```

Replace '/path/to/solscape-1.0' with the actual path where the solscape-1.0 directory is located.

**Step 3: Import the Main Class** Once the path is set, you can import the main class Landscape from the solscape package as follows:

```
from solscape import Landscape
```

This will allow you to use the Landscape class and other functionalities provided by the package.

# **Configuration Parameters**

The configuration parameters for the solscape package are divided into different categories, each focusing on a specific part of the algorithm. Below is an overview of the different categories and the associated parameters.

## System Parameters

These parameters are related to the system setup and its general properties: Dim, EnergyFunction, Grad, AutoDiff, NumericalGrad, DimerLength, SymmetryCheck and GradientSystem.

#### Dim (Optional)

## Description

System dimension specification.

## **Data Type**

int (positive)

#### **Behavior**

- Default: Inferred from initial point
- Manual override must match dimensions of:
  - initial point
  - InitialSearchDirection

### **EnergyFunction (Conditional Required) Description**

Specifies the energy function for gradient systems.

## **Data Type**

- Python function: Callable[[np.ndarray], float]
  - Input: np.ndarray with shape (d, 1) (column vector)
- Symbolic expression: str
  - **Specification**: Use x1, x2, ..., xd as variables & Supports full sympy syntax

## **Example**

```
EnergyFunction = '''
0.4*x1**2 - 0.2*x1*x2 + 0.25*x2**2
- 5*(atan(x1-5) + atan(x2-5))
'''
```

## **Grad (Conditional Required)**

## Description

Specifies the gradient function for gradient systems or vector field for non-gradient systems.

## Requirements

- For gradient systems:  $\nabla E$
- For non-gradient systems: Direct specification of F in  $\dot{x}$  = -F(x)

## **Data Types**

- Python function: Callable[[np.ndarray], np.ndarray]
  - Input/Output: np.ndarray with shape (d, 1) (column vector)
- Symbolic expression list: list[str] (e.g., ["2\*x1", "cos(x2)"])

### **Example**

```
# Gradient system
Grad = ["2*x1", "2*x2"]
```

```
# Non-gradient system

Grad = ["x2", "-x1 - 0.1*x2"]
```

## **AutoDiff (Optional)**

## Description

Enables automatic differentiation of the energy function.

### **Data Type**

bool

## **Options**

- True: Requires EnergyFunction parameter
- False: Requires manual specification of Grad parameter

#### **Behavior**

- Defaults to True if EnergyFunction is provided
- Defaults to False if Grad is provided

## NumericalGrad (Optional)

#### Description

Enables numerical gradient approximation.

## **Applicability**

Only active when Grad is unspecified

#### **Data Type**

bool

## **Options**

- True: Use finite difference approximation
- False: (Default) Use analytical gradient from EnergyFunction

## **DimerLength (Optional)**

## Description

Displacement length for numerical gradient approximations.

## **Data Type**

float (positive)

#### **Default**

1e-5

## **SymmetryCheck (Optional)**

## **Description**

Verifies gradient system properties via Hessian symmetry.

## **Data Type**

bool

#### **Behavior**

• True: (default) Auto-detect

• False: Bypass checks (improves performance)

## **GradientSystem (Optional)**

## Description

Explicit declaration of gradient system nature.

#### Usage

Override automatic detection when known a priori.

## **Data Type**

bool

#### **Behavior**

missing: Auto-detect

## **Hessian Parameters**

These parameters are related to the Hessian matrix: ExactHessian and HessianDimerLength.

## **ExactHessian (Optional)**

## Description

Controls Hessian matrix computation method.

#### **Data Type**

bool

## **Options**

- True: Analytical Hessian from symbolic gradient
- False: (Default) Dimer-based approximation

## Requirement

Requires **Grad** as symbolic expressions

## HessianDimerLength (Optional)

## Description

Displacement length for numerical Hessian-Vector product approximations.

## **Data Type**

float (positive)

### **Default**

1e-5

## **Eigen Parameters**

These parameters control how the eigen pairs (eigenvalues and eigenvectors) are computed: EigenMethod, EigenMaxIter, EigenStepSize and PrecisionTol.

## **EigenMethod (Optional)**

## Description

Eigensolver selection for stability analysis.

## **Data Type**

str

### **Options**

Method	System Type	Description	
lobpcg	Gradient (Default)	Locally Optimal Block PCG	
euler	Gradient & Non-gradient	Explicit Euler Discretization	
power	Non-gradient (Default)	Power Method	

## EigenMaxIter (Optional)

## Description

Maximum iterations for eigenpair computation.

## **Data Type**

int (positive)

#### Default

10

## **EigenStepSize (Optional)**

## Description

Discretization step for Euler/power methods.

## **Data Type**

float (positive)

## **Default**

1e-7

## PrecisionTol (Optional)

## Description

Precision tolerance for eigenvalues. (We treat eigenvalues as 0 if its absolute value less than tolerance.)

## **Data Type**

float (non-negative)

#### **Default**

1e-5

#### **Acceleration Parameters**

These parameters are related to improving the speed and efficiency of the algorithm: BBStep, Acceleration, NesterovChoice, NesterovRestart and Momentum.

## **BBStep (Optional)**

## Description

Enables Barzilai-Borwein adaptive step sizing.

## **Data Type**

bool

#### **Default**

False

#### Trade-off

- May accelerate convergence
- Can cause instability in stiff systems

## **Acceleration (Optional)**

## Description

Convergence acceleration technique.

## Data Type str

#### **Options**

- none: (Default) No acceleration
- heavyball: Momentum-based acceleration
- nesterov: Nesterov-accelerated dynamics

## **NesterovChoice (Optional)**

## **Description**

Specifies the acceleration parameter sequence for Nesterov's method.

## **Data Type**

int

#### **Options**

- 1 : (default)  $\gamma_n = rac{n}{n+3}$  .
- 2:  $\gamma_n=rac{ heta_n-1}{ heta_{n+1}}$  , with  $heta_{n+1}=rac{1+\sqrt{1+4 heta_n^2}}{2}$  ,  $heta_0=1$  .

#### Default

1

## **NesterovRestart (Optional)**

## Description

Iteration interval for Nesterov momentum reset.

## **Data Type**

int (positive) | None

#### **Behavior**

- None: Disables momentum restart
- Integer n: Resets momentum every n iterations

#### **Default**

None

#### Momentum (Optional)

## Description

Momentum coefficient for heavy ball acceleration.

## **Data Type**

float (non-negative)

#### **Default**

0.0

#### **Constraints**

- 0.0 ≤ Momentum < 1.0
- 0.0: Equivalent to no acceleration

## **Solver Parameters**

These parameters are related to the solver process and control the behavior of the HiSD process:

InitialPoint, Tolerance, SearchArea, TimeStep, MaxIter, SaveTrajectory, Verbose and ReportInterval.

## **InitialPoint (Required) Description**

The starting coordinates for saddle point search.

#### **Data Types**

list | numpy.ndarray (1D array)

#### Example

initial\_point = [0.5, -1.2]

## **Tolerance (Optional)**

## Description

Convergence threshold for saddle point iterations.

## **Data Type**

float (positive)

## **Default**

1e-6

## **Stopping Criterion**

 $\|gradient\ vector\|_2 < Tolerance$ 

## SearchArea (Optional)

## Description

Maximum exploration radius from initial point.

## **Data Type**

float (positive)

#### **Default**

**1e3** 

#### **Effect**

Terminates search if  $\|\mathbf{x} - \mathbf{x}_0\|_2 > \text{SearchArea}$ 

## TimeStep (Optional)

## Description

Temporal discretization interval for dynamics.

## **Data Type**

float (positive)

#### **Default**

1e-4

## MaxIter (Optional)

## Description

Maximum number of HiSD iterations permitted.

## **Data Type**

int (positive)

#### **Default**

1000

## SaveTrajectory (Optional)

## Description

Records full optimization path during computation.

## **Data Type**

bool

#### **Default**

True

## **Verbose (Optional)**

## Description

Controls real-time progress reporting.

#### Output

Prints iteration count and gradient norm.

## **Data Type**

bool

#### **Default**

False

## ReportInterval (Optional)

## Description

Iteration frequency for console output when Verbose=True.

## **Data Type**

int (positive)

#### **Default**

100

## Landscape Parameters

These parameters are related to constructing and navigating the solution landscape: MaxIndex, MaxIndexGap, SameJudgement, InitialEigenVectors, PerturbationMethod, PerturbationRadius, PerturbationNumber and EigenCombination.

## MaxIndex (Optional) Description

Maximum saddle index (k) to compute.

- Index 0: Uses standard SD (Steepest Descent) method
- Index ≥1: Uses HiSD (High-index Saddle Dynamics) method

## **Data Type**

int (non-negative)

### **Default**

6

#### **Constraints**

0 ≤ max\_index ≤ Dim

## MaxIndexGap (Optional) Description

Maximum allowed index difference between parent and child saddle points during hierarchical search.

#### **Data Type**

```
int (positive)
```

#### **Default**

1

## **Example**

If MaxIndexGap = 2:

- Parent (index 4) → Children (indices 2, 3)
- Parent (index 3) → Children (indices 1, 2)

## SameJudgement (Optional)

#### Description

Saddle point equivalence criterion.

### **Data Types**

- float (positive): Threshold for 2-norm distance (default: 1e-3)
- Callable[[np.ndarray, np.ndarray], bool]: Custom comparison function
  - Input: np.ndarray with shape (d, 1) (column vector)

#### **Example**

```
# Custom similarity check
def custom_judge(a, b):
   return np.linalg.norm(a - b) < 0.5 and abs(a[0] - b[0]) < 0.1</pre>
```

#### InitialEigenVectors (Optional)

### Description

Initial guess for Hessian eigenvectors.

#### **Data Types**

- None: (Default) Auto-initialize with k smallest eigenvectors
- Manual input: np.ndarray of shape (d, k)
  where d is the dimension and k equals MaxIndex.

## Requirement

Column vectors must be orthonormal

### PerturbationMethod (Optional)

## Description

Statistical distribution for saddle point exploration.

## **Data Type**

str

## **Options**

value uniform (Default) gaussian		Description	
		Uniform sampling	
		Normally distributed displacements	

## PerturbationRadius (Optional)

## Description

Displacement magnitude for saddle perturbations.

## **Data Type**

float (positive)

## **Default**

1e-4

## PerturbationNumber (Optional)

## Description

Number of directional probes per saddle point.

## Note

Actual probes = 2 × PerturbationNumber (bidirectional)

## **Data Type**

int (positive)

## **Default**

2

## **EigenCombination (Optional)**

## Description

Strategy for eigenvector utilization in perturbations.

## **Data Type**

str

## **Options**

Value	<b>Computational Cost</b>	nputational Cost Completeness	
all (Default)	High	Exhaustive	
min	Low	Partial	

## .Run()

### Description

Initiates the solution landscape computation process.

#### **Parameters**

None

#### Usage

```
landscape = Landscape(
    MaxIndex=3,
    AutoDiff=True,
    EnergyFunction="1*(x1**2-1)**2+2*(x2**2-1)**2+3*(x3**2-1)**2",
    InitialPoint=np.array([0, 0, 0])
)
landscape.Run() # Starts computation
```

## Output

Access search trajectory by instance.DetailRecord. Each row of it is data for one successful search. sequentially contains ID of end saddle point, ID of start saddle point, ndarray of each step position and ndarray of each step time position.

Access saddle point data by instance.SaddleList. Each row of it is data for one saddle point. sequentially contains ID of the saddle point, position of the saddle point, Morse Index of the saddle point, the eigenvectors of negative eigenvalues and set of father saddles.

See Output Objects for more details.

```
.RestartFromPoint(RestartPoint, MaxIndex)
```

#### **Description**

Restarts computation from specified coordinates.

## **Parameters**

Name	Туре	Constraints	Description
RestartPoint	list or numpy.ndarray (1D)	Must match system dimension	Initial position vector
MaxIndex	int	0 ≤ max_index ≤ dim	Maximum saddle index to compute

### **Usage**

```
landscape.RestartFromPoint(
    RestartPoint=np.array([[0.2], [-0.8]]),
```

```
MaxIndex=2
)
```

## .RestartFromSaddle(BeginID, Perturbation, MaxIndex)

## Description

Restarts computation from existing saddle point.

#### **Parameters**

Name	Туре	Constraints	Description
BeginID	int	0 ≤ begin_id < len(SaddleList)	Valid saddle point ID
Perturbation	numpy.ndarray (1D)	Must match system dimension	Initial perturbation vector
MaxIndex	int	0 ≤ max_index ≤ dim	Maximum saddle index to compute

## Usage

```
landscape.RestartFromSaddle(
    BeginID=3,
    Perturbation=1e-3 * np.array([[-1], [0.5]]),
    MaxIndex=2
)
```

# .DrawTrajectory(\*\*kwargs)

## Description

Visualizes search trajectories and energy landscapes for systems. Supports contour overlays, style customization, and high-D projections.

## **Parameters**

Name	Туре	Constraints	Description & Default Value
DetailedTraj	bool	Requires saved trajectory data	Show full iteration path (default: False)
Contour	bool	2D systems only	Enable contour lines (default: True)
Contourf	bool	Contour=True	Enable color-filled contours (default: True)

Name	Туре	Constraints	Description & Default Value
ContourGridNum	int	≥ 1	Main grid divisions per axis (default: 50)
ContourGridOut	int	≥ 0	Extended grid divisions per axis (default: 10)
Title	str		Figure title text (defalut: "The Search Trajectory")
TrajectorySet	dict		<pre>Set the trajectory style (default: {"linewidth": 0.4, "linestyle": "- ", "color": "blue", "label": "Search Trajectory"})</pre>
SaddlePointSet	list[dict]		<pre>Set the saddle point style (default: [{"marker": "o", "color":   colors[20 * i + 20], "label": f"Index {i} Saddle Point"} for i in   range(instance.MaxIndex + 1)])</pre>
GridSet	dict		Set the grid style  (default: {"visible": True, "linestyle": " ", "linewidth": 0.1, "color": "gray"})
1DSamples	int	≥ 1 (1D only)	Function sampling density (default: 1e3)
1DSamplesOut	int	≥ Ø (1D only)	Extended sampling range (default: 1e2)
1DFunctionDraw	dict	1D only	<pre>Set the 1D function style (default: {"linewidth": 2, "linestyle": "-",   "color": "red", "label": "Function Curve"})</pre>
WhetherSave	bool		Save to file (default: False)
SaveFigurePath	str	File extension determines format	Output path (default: "Landscape_figure.png")
Projection	None or callable	Required for dim > 2	Projection function (Signature: ((n*dim) array) → ((n*2) array))

# Usage

```
# 2D system without detailed trajectory
landscape.DrawTrajectory(
    Contour=True,
```

```
Contourf=True,
    WhetherSave=True,
    ContourGridNum=100,
    ContourGridOut=25,
    SaveFigurePath="landscape.png"
)
# high-D system with detailed trajectory
import numpy as np
def proj_func(input):
    output = np.hstack((1.0 * input[:, [0]]+ 1.5 * input[:, [1]], 1.0 * input[:,
[0]]+ 2.5 * input[:, [2]]))
    return output
landscape.DrawTrajectory(
    DetailedTraj=True,
    ContourGridNum=100,
    ContourGridOut=25,
    Projection=proj_func
)
```

#### **Visualization Rules**

## **Dimensionality Handling:**

- 1D: X-axis = iteration count, Y-axis = function value
- 2D: Natural coordinates
- ≥3D: Requires Projection to 2D plane

#### **Data Requirements**

• Detailed trajectories require SaveTrajectory=True in .Run() call

## **Output**

- Matplotlib figure (interactive display)
- Image file when WhetherSave=True (PNG/PDF/SVG based on path)

```
.DrawConnection(**kwargs)
```

## Description

Visualizes connectivity between saddle points in the solution landscape. Displays markers for saddle nodes and search paths connecting them. Supports custom styling of points and connection lines.

## **Parameters**

Name	Туре	Default Value	Description
Title	str	"The Connection of Saddle Points"	Figure title text

Name	Туре	Default Value	Description
SaddlePointSet	list[dict]	<pre>[{"node_shape": "o", "node_color": colors[i], "label": f"Index {i} Saddle Point"}for i in range(instance.MaxIndex + 1)]</pre>	Set the saddle point style
TrajectorySet	dict	<pre>{"width": 0.4, "style": "solid", "edge_color": "blue", "label": "Search Trajectory"}</pre>	Set the connection path style
WhetherSave	bool	False	Save to file
SaveFigurePath	str	"Landscape_Connection_figure.png"	Output path (file extension determines format)

## Usage

```
# Draw solution landscape
landscape.DrawConnection(
    Title="Solution Landscape",
    WhetherSave=True,
    SaveFigurePath="landscape.png"
)
```

## .Save(filepath, fileformat)

## Description

Persists calculation results to persistent storage with specified serialization format.

## **Parameters**

Name	Type	Constraints	Description
filepath	str	Valid filesystem path	Target path without extension
fileformat	str	"json" \  "pickle" \  "mat" (default: "json")	Serialization format

## Usage

```
landscape.Save(filepath="/output/results", fileformat="pickle")
```

## **File Output**

Generates:

- filepath.json (JSON text format)
- filepath.pickle (Python binary serialization)
- filepath.mat (MATLAB-compatible binary)

Contains dictionary with aligned indices for:

- 1. SaddleID: Unique identification numbers
- 2. Position: Spatial coordinates (N-dimensional array)
- 3. MorseIndex: Number of negative eigenvalues
- 4. FatherSet: Ancestry relationships (list-formatted references)

### **Index Consistency**

All data structures maintain identical ordering:

```
SaddleID[n] \leftrightarrow Position[n] \leftrightarrow MorseIndex[n] \leftrightarrow FatherSet[n]
```

# **Output Objects**

```
.SaddleList
```

## Description

Collection of discovered saddle points with topological metadata.

#### Type

list[list]

#### **Node Structure**

### **Access**

```
# Get all index-1 saddles
index1_saddles = [node for node in landscape.SaddleList if node[2] == 1]
```

```
# Get position of saddle ID 5
saddle5_pos = landscape.SaddleList[5][1]
```

#### .DetailRecord

### **Description**

Complete search trajectory data.

## **Type**

list[list]

#### **Record Structure**

#### Access

```
# Plot first search path
first_search = landscape.DetailRecord[0]
plt.plot(first_search[2][:,0], first_search[2][:,1])
```

# Package Requirements

#### Overview

This component requires the following Python packages to be installed. Verify installations before usage.

## **Built-in Packages**

The following standard library modules are required (no installation needed):

```
import copy
import sys
import warnings
import inspect
import json
import itertools
```

```
import math
import pickle
```

## Third-party Dependencies

Required external packages with version specifications:

Package	Required Version	Installation Command
NumPy	2.2.3	pip install numpy==2.2.3
SciPy	1.15.2	pip install scipy==1.15.2
SymPy	1.13.3	pip install sympy==1.13.3
Matplotlib	3.10.0	pip install matplotlib==3.10.0
NetworkX	3.4.2	pip install networkx==3.4.2

## **Full Environment Setup**

```
pip install "numpy>=2.2.3" \
    "scipy>=1.15.2" \
    "sympy>=1.13.3" \
    "matplotlib>=3.10.0" \
    "networkx>=3.4.2"
```

## Version Verification

Check installed versions using:

```
import numpy, scipy, sympy, matplotlib, networkx

print(f"NumPy: {numpy.__version__}")  # Should show 2.2.3

print(f"SciPy: {scipy.__version__}")  # Should show 1.15.2

print(f"SymPy: {sympy.__version__}")  # Should show 1.13.3

print(f"Matplotlib: {matplotlib.__version__}")  # Should show 3.10.0

print(f"NetworkX: {networkx.__version__}")  # Should show 3.4.2
```