# **ODDT Documentation**

Release 0.6-128-gefc1a7b

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2 Contents

Installation

# 1.1 Requirements

- Python 2.7+ or 3.4+
- OpenBabel (2.3.2+) or/and RDKit (2016.03)
- Numpy (1.8+)
- Scipy (0.14+)
- Sklearn (0.18+)
- joblib (0.8+)
- pandas (0.17.1+)
- Skimage (0.10+) (optional, only for surface generation)

**Note:** All installation methods assume that one of toolkits is installed. For detailed installation procedure visit toolkit's website (OpenBabel, RDKit)

Most convenient way of installing ODDT is using PIP. All required python modules will be installed automatically, although toolkits, either OpenBabel (pip install openbabel) or RDKit need to be installed manually

```
pip install oddt
```

If you want to install cutting edge version (master branch from GitHub) of ODDT also using PIP

```
pip install git+https://github.com/oddt/oddt.git@master
```

Finally you can install ODDT straight from the source

```
wget https://github.com/oddt/oddt/archive/0.5.tar.gz
tar zxvf 0.5.tar.gz
cd oddt-0.5/
python setup.py install
```

# 1.2 Common installation problems

# **Usage Instructions**

You can use any supported toolkit united under common API (for reference see Pybel or Cinfony). All methods and software which based on Pybel/Cinfony should be drop in compatible with ODDT toolkits. In contrast to its predecessors, which were aimed to have minimalistic API, ODDT introduces extended methods and additional handles. This extensions allow to use toolkits at all its grace and some features may be backported from others to introduce missing functionalities. To name a few:

- · coordinates are returned as Numpy Arrays
- atoms and residues methods of Molecule class are lazy, ie. not returning a list of pointers, rather an object which allows indexing and iterating through atoms/residues
- Bond object (similar to Atom)
- atom\_dict, ring\_dict, res\_dict comprehensive Numpy Arrays containing common information about given entity, particularly useful for high performance computing, ie. interactions, scoring etc.
- lazy Molecule (asynchronous), which is not converted to an object in reading phase, rather passed as a string and read in when underlying object is called
- pickling introduced for Pybel Molecule (internally saved to mol2 string)

# 2.1 Atom, residues, bonds iteration

One of the most common operation would be iterating through molecules atoms

```
mol = oddt.toolkit.readstring('smi', 'clccccl')
for atom in mol:
    print(atom.idx)
```

Note: mol.atoms, returns an object (AtomStack) which can be access via indexes or iterated

Iterating over residues is also very convenient, especially for proteins

```
for res in mol.residues:
    print(res.name)
```

Additionally residues can fetch atoms belonging to them:

```
for res in mol.residues:
    for atom in res:
        print(atom.idx)
```

Bonds are also iterable, similar to residues:

```
for bond in mol.bonds:
    print(bond.order)
    for atom in bond:
        print(atom.idx)
```

# 2.2 Reading molecules

Reading molecules is mostly identical to Pybel.

Reading from file

```
for mol in oddt.toolkit.readfile('smi', 'test.smi'):
    print(mol.title)
```

Reading from string

```
mol = oddt.toolkit.readstring('smi', 'clccccc1 benzene'):
    print(mol.title)
```

**Note:** You can force molecules to be read in asynchronously, aka "lazy molecules". Current default is not to produce lazy molecules due to OpenBabel's Memory Leaks in OBConverter. Main advantage of lazy molecules is using them in multiprocessing, then conversion is spreaded on all jobs.

Reading molecules from file in asynchronous manner

```
for mol in oddt.toolkit.readfile('smi', 'test.smi', lazy=True):
    pass
```

This example will execute instantaneously, since no molecules were evaluated.

# 2.3 Numpy Dictionaries - store your molecule as an uniform structure

Most important and handy property of Molecule in ODDT are Numpy dictionaries containing most properties of supplied molecule. Some of them are straightforward, other require some calculation, ie. atom features. Dictionaries are provided for major entities of molecule: atoms, bonds, residues and rings. It was primarily used for interactions calculations, although it is applicable for any other calculation. The main benefit is marvelous Numpy broadcasting and subsetting.

Each dictionary is defined as a format in Numpy.

# 2.3.1 atom dict

#### Atom basic information

- 'coords', type: float 32, shape: (3) atom coordinates
- 'charge', type: float32 atom's charge
- 'atomicnum', type: int8 atomic number
- 'atomtype', type: a4 Sybyl atom's type
- 'hybridization', type: int8 atoms hybrydization
- 'neighbors', type: float32, shape: (4,3) coordinates of non-H neighbors coordinates for angles (max of 4 neighbors should be enough)

#### Residue information for current atom

- 'resid', type: int16 residue ID
- 'resnumber', type: int16 residue number
- 'resname', type: a3 Residue name (3 letters)
- 'isbackbone', type: bool is atom part of backbone

# Atom properties

- 'isacceptor', type: bool is atom H-bond acceptor
- 'isdonor', type: bool is atom H-bond donor
- 'isdonorh', type: bool is atom H-bond donor Hydrogen
- 'ismetal', type: bool is atom a metal
- 'ishydrophobe', type: bool is atom hydrophobic
- 'isaromatic', type: bool is atom aromatic
- 'isminus', type: bool is atom negatively charged/chargable
- 'isplus', type: bool is atom positively charged/chargable
- 'ishalogen', type: bool is atom a halogen

# Secondary structure

- 'isalpha', type: bool is atom a part of alpha helix
- 'isbeta', type: bool' is atom a part of beta strand

# 2.3.2 ring dict

- 'centroid', type: float 32, shape: 3 coordinates of ring's centroid
- 'vector', type: float 32, shape: 3 normal vector for ring
- 'isalpha', type: bool is ring a part of alpha helix
- 'isbeta', type: bool' is ring a part of beta strand

# 2.3.3 res\_dict

- 'id', type: int16 residue ID
- 'resnumber', type: int16 residue number
- 'resname', type: a3 Residue name (3 letters)
- 'N', type: float 32, shape: 3 cordinates of backbone N atom
- 'CA', type: float 32, shape: 3 cordinates of backbone CA atom
- 'C', type: float 32, shape: 3 coordinates of backbone C atom
- 'isalpha', type: bool is residue a part of alpha helix
- 'isbeta', type: bool' is residue a part of beta strand

**Note:** All aforementioned dictionaries are generated "on demand", and are cached for molecule, thus can be shared between calculations. Caching of dictionaries brings incredible performance gain, since in some applications their generation is the major time consuming task.

Get all acceptor atoms:

```
mol.atom_dict['isacceptor']
```

# 2.4 Interaction Fingerprints

Module, where interactions between two molecules are calculated and stored in fingerprint.

# 2.4.1 The most common usage

Firstly, loading files

```
protein = next(oddt.toolkit.readfile('pdb', 'protein.pdb'))
protein.protein = True
ligand = next(oddt.toolkit.readfile('sdf', 'ligand.sdf'))
```

**Note:** You have to mark a variable with file as protein, otherwise You won't be able to get access to e.g. 'resname; , 'resid' etc. It can be done as above.

File with more than one molecule

```
mols = list(oddt.toolkit.readfile('sdf', 'ligands.sdf'))
```

When files are loaded, You can check interactions between molecules. Let's find out, which amino acids creates hydrogen bonds

```
protein_atoms, ligand_atoms, strict = hbonds(protein, ligand)
print(protein_atoms['resname'])
```

Or check hydrophobic contacts between molecules

```
protein_atoms, ligand_atoms = hydrophobic_contacts(protein, ligand)
print(protein_atoms, ligand_atoms)
```

But instead of checking interactions one by one, You can use fingerprints module.

```
IFP = InteractionFingerprint(ligand, protein)
SIFP = SimpleInteractionFingerprint(ligand, protein)
```

Very often we're looking for similar molecules. We can easily accomplish this by e.g.

```
results = []
reference = SimpleInteractionFingerprint(ligand, protein)
for el in query:
    fp_query = SimpleInteractionFingerprint(el, protein)
    # similarity score for current query
    cur_score = dice(reference, fp_query)
    # score is the lowest, required similarity
    if cur_score > score:
        results.append(el)
return results
```

# 2.5 Molecular shape comparison

Three methods for molecular shape comparison are supported: USR and its two derivatives: USRCAT and Electroshape.

- USR (Ultrafast Shape Recognition) function usr(molecule) Ballester PJ, Richards WG (2007). Ultrafast shape recognition to search compound databases for similar molecular shapes. Journal of computational chemistry, 28(10):1711-23. http://dx.doi.org/10.1002/jcc.20681
- USRCAT (USR with Credo Atom Types) function usr\_cat(molecule) Adrian M Schreyer, Tom Blundell (2012). USRCAT: real-time ultrafast shape recognition with pharmacophoric constraints. Journal of Cheminformatics, 2012 4:27. http://dx.doi.org/10.1186/1758-2946-4-27
- Electroshape function electroshape(molecule) Armstrong, M. S. et al. ElectroShape: fast molecular similarity calculations incorporating shape, chirality and electrostatics. J Comput Aided Mol Des 24, 789-801 (2010). http://dx.doi.org/doi:10.1007/s10822-010-9374-0

Aside from spatial coordinates, atoms' charges are also used as the fourth dimension to describe shape of the molecule.

To find most similar molecules from the given set, each of these methods can be used.

Loading files:

```
query = next(oddt.toolkit.readfile('sdf', 'query.sdf'))
database = list(oddt.toolkit.readfile('sdf', 'database.sdf'))
```

Example code to find similar molecules:

```
results = []
query_shape = usr(query)
for mol in database:
    mol_shape = usr(mol)
    similarity = usr_similarity(query_shape, mol_shape)
    if similarity > 0.7:
        results.append(mol)
```

To use another method, replace usr(mol) with usr_cat(mol) or electroshape(mol).		

# ODDT command line interface (CLI)

There is an *oddt* command to interface with Open Drug Discovery Toolkit from terminal, without any programming knowleadge. It simply reproduces *oddt.virtualscreening.virtualscreening*. One can filter, dock and score ligands using methods implemented or compatible with ODDT. All positional arguments are treated as input ligands, whereas output must be assigned using *-O* option (following *obabel* convention). Input and output formats are defined using *-i* and *-o* accordingly. If output format is present and no output file is assigned, then molecules are printed to STDOUT.

To list all the available options issue -h option:

```
oddt_cli -h
```

1. Docking ligand using Autodock Vina (construct box using ligand from crystal structure) with additional RFscore v2 rescoring:

```
oddt_cli input_ligands.sdf --dock autodock_vina --receptor rec.mol2 --auto_ligand_

→crystal_ligand.mol2 --score rfscore_v2 -0 output_ligands.sdf
```

2. Filtering ligands using Lipinski RO5 and PAINS. Afterwards dock with Autodock Vina:

```
oddt_cli input_ligands.sdf --filter ro5 --filter pains --dock autodock_vina --

→receptor rec.mol2 --auto_ligand crystal_ligand.mol2 -0 output_ligands.sdf
```

3. Dock with Autodock Vina, with precise box position and dimensions. Fix seed for reproducibility and increase exhaustiveness:

```
oddt_cli ampc/actives_final.mol2.gz --dock autodock_vina --receptor ampc/receptor.pdb_ \rightarrow--size '(8,8,8)' --center '(1,2,0.5)' --exhaustiveness 20 --seed 1 -O ampc_docked. \rightarrowsdf
```

4. Rescore ligands using 3 versions of RFscore and pre-trained scoring function (either pickle from ODDT or any other SF implementing oddt.scoring.scorer API):

```
oddt_cli docked_ligands.sdf --receptor rec.mol2 --score rfscore_v1 --score rfscore_v2_

---score rfscore_v3 --score TrainedNN.pickle -O docked_ligands_rescored.sdf
```

# Development and contributions guide

1. Indicies All indicies within toolkit are 0-based, but for backward compatibility with OpenBabel there is mol.idx property. If you develop using ODDT you are encouraged to use 0-based indicies and/or mol.idx0 and mol.idx1 properties to be exact which convention you adhere to. Otherwise you can run into bags which are hard to catch, when writing toolkit independent code.

# ODDT API documentation

# 5.1 oddt package

# 5.1.1 Subpackages

oddt.docking package

**Submodules** 

# oddt.docking.AutodockVina module

Bases: object

Autodock Vina docking engine, which extends it's capabilities: automatic box (auto-centering on ligand).

# **Parameters**

**protein: oddt.toolkit.Molecule object (default=None)** Protein object to be used while generating descriptors.

**auto\_ligand: oddt.toolkit.Molecule object or string (default=None)** Ligand use to center the docking box. Either ODDT molecule or a file (opened based on extesion and read to ODDT molecule). Box is centered on geometric center of molecule.

size: tuple, shape=[3] (default=(20, 20, 20)) Dimentions of docking box (in Angstroms) center: tuple, shape=[3] (default=(0,0,0)) The center of docking box in cartesian space.

exhaustiveness: int (default=8) Exhaustiveness parameter of Autodock Vina

**num\_modes:** int (default=9) Number of conformations generated by Autodock Vina. The maximum number of docked poses is 9 (due to Autodock Vina limitation).

energy\_range: int (default=3) Energy range cutoff for Autodock Vina

seed: int or None (default=None) Random seed for Autodock Vina

prefix\_dir: string (default=/tmp) Temporary directory for Autodock Vina files

**executable: string or None (default=None)** Autodock Vina executable location in the system. It's realy necessary if autodetection fails.

**autocleanup: bool (default=True)** Should the docking engine clean up after execution?

**skip\_bad\_mols:** bool (default=True) Should molecules that crash Autodock Vina be skipped.

## **Attributes**

tmp\_dir

#### **Methods**

dock(ligands[, protein])	Automated docking procedure.	
<pre>predict_ligand(ligand)</pre>	Local method to score one ligand and update it's	
	scores.	
predict_ligands(ligands)	Method to score ligands lazily	
score(ligands[, protein])	Automated scoring procedure.	
set_protein(protein)	Change protein to dock to.	

clean

#### clean()

dock (ligands, protein=None)

Automated docking procedure.

# **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to dock

**protein: oddt.toolkit.Molecule object or None** Protein object to be used. If None, then the default one is used, else the protein is new default.

# Returns

**ligands** [array of oddt.toolkit.Molecule objects] Array of ligands (scores are stored in mol.data method)

#### predict\_ligand(ligand)

Local method to score one ligand and update it's scores.

## **Parameters**

ligand: oddt.toolkit.Molecule object Ligand to be scored

## Returns

ligand: oddt.toolkit.Molecule object Scored ligand with updated scores

#### predict\_ligands (ligands)

Method to score ligands lazily

#### **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to be scored

#### **Returns**

ligand: iterator of oddt.toolkit.Molecule objects Scored ligands with updated scores

score (ligands, protein=None)

Automated scoring procedure.

#### **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to score

**protein: oddt.toolkit.Molecule object or None** Protein object to be used. If None, then the default one is used, else the protein is new default.

#### Returns

**ligands** [array of oddt.toolkit.Molecule objects] Array of ligands (scores are stored in mol.data method)

set\_protein (protein)

Change protein to dock to.

#### **Parameters**

protein: oddt.toolkit.Molecule object Protein object to be used.

# tmp\_dir

oddt.docking.AutodockVina.parse\_vina\_docking\_output(output)

Function parsing Autodock Vina docking output to a dictionary

#### **Parameters**

output [string] Autodock Vina standard ouptud (STDOUT).

#### Returns

out [dict] dicitionary containing scores computed by Autodock Vina

oddt.docking.AutodockVina.parse\_vina\_scoring\_output(output)

Function parsing Autodock Vina scoring output to a dictionary

## **Parameters**

**output** [string] Autodock Vina standard ouptud (STDOUT).

#### Returns

out [dict] dicitionary containing scores computed by Autodock Vina

oddt.docking.AutodockVina.write\_vina\_pdbqt(mol, directory, flexible=True,

name\_id=None)

Write single PDBQT molecule to a given directory. For proteins use *flexible=False* to avoid encoding torsions. Additionally an name ID can be appended to a name to avoid conflicts.

# oddt.docking.internal module

ODDT's internal docking/scoring engines

```
oddt.docking.internal.change_dihedral(coords, a1, a2, a3, a4, target_angle, rot_mask)
oddt.docking.internal.get_children(molecule, mother, restricted)
oddt.docking.internal.get_close_neighbors(molecule, a_idx, num_bonds=1)
oddt.docking.internal.num_rotors_pdbqt(lig)
class oddt.docking.internal.vina_docking(rec, lig=None, box=None, box_size=1.0, weights=None)
Bases: object
```

#### **Methods**

```
correct_radius (atom_dict)
score (coords=None)
score_inter (coords=None)
score_intra (coords=None)
score_total (coords=None)
set_box (box)
set_coords (coords)
set_ligand (lig)
set_protein (rec)
weighted_inter (coords=None)
weighted_intra (coords=None)
weighted_total (coords=None)
class oddt.docking.internal.vina_ligand (c0, num_rotors, engine, box_size=I)
Bases: object
```

#### **Methods**

mutate

mutate (x2, force=False)

#### **Module contents**

class oddt.docking.autodock\_vina (protein=None, auto\_ligand=None, size=(20, 20, 20), center=(0, 0, 0), exhaustiveness=8, num\_modes=9, energy\_range=3, seed=None, prefix\_dir='/tmp', n\_cpu=1, executable=None, autocleanup=True, skip\_bad\_mols=True)

Bases: object

Autodock Vina docking engine, which extends it's capabilities: automatic box (auto-centering on ligand).

#### **Parameters**

**protein: oddt.toolkit.Molecule object (default=None)** Protein object to be used while generating descriptors.

auto\_ligand: oddt.toolkit.Molecule object or string (default=None) Ligand use to center the docking box. Either ODDT molecule or a file (opened based on extesion and read to ODDT molecule). Box is centered on geometric center of molecule.

size: tuple, shape=[3] (default=(20, 20, 20)) Dimentions of docking box (in Angstroms)

center: tuple, shape=[3] (default=(0,0,0)) The center of docking box in cartesian space.

exhaustiveness: int (default=8) Exhaustiveness parameter of Autodock Vina

**num\_modes:** int (default=9) Number of conformations generated by Autodock Vina. The maximum number of docked poses is 9 (due to Autodock Vina limitation).

energy\_range: int (default=3) Energy range cutoff for Autodock Vina

seed: int or None (default=None) Random seed for Autodock Vina

prefix\_dir: string (default=/tmp) Temporary directory for Autodock Vina files

**executable: string or None** (**default=None**) Autodock Vina executable location in the system. It's realy necessary if autodetection fails.

autocleanup: bool (default=True) Should the docking engine clean up after execution?

skip\_bad\_mols: bool (default=True) Should molecules that crash Autodock Vina be skipped.

# **Attributes**

tmp\_dir

## **Methods**

dock(ligands[, protein])	Automated docking procedure.	
predict_ligand(ligand)	igand(ligand) Local method to score one ligand and update it	
	scores.	
<pre>predict_ligands(ligands)</pre>	Method to score ligands lazily	
score(ligands[, protein])	Automated scoring procedure.	
set_protein(protein)	Change protein to dock to.	

clean

#### clean()

dock (ligands, protein=None)

Automated docking procedure.

#### **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to dock

**protein: oddt.toolkit.Molecule object or None** Protein object to be used. If None, then the default one is used, else the protein is new default.

#### **Returns**

**ligands** [array of oddt.toolkit.Molecule objects] Array of ligands (scores are stored in mol.data method)

#### predict\_ligand(ligand)

Local method to score one ligand and update it's scores.

#### **Parameters**

ligand: oddt.toolkit.Molecule object Ligand to be scored

#### Returns

ligand: oddt.toolkit.Molecule object Scored ligand with updated scores

## predict\_ligands (ligands)

Method to score ligands lazily

## **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to be scored

# Returns

ligand: iterator of oddt.toolkit.Molecule objects Scored ligands with updated scores

score (ligands, protein=None)

Automated scoring procedure.

# **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to score

**protein: oddt.toolkit.Molecule object or None** Protein object to be used. If None, then the default one is used, else the protein is new default.

#### Returns

**ligands** [array of oddt.toolkit.Molecule objects] Array of ligands (scores are stored in mol.data method)

# set\_protein (protein)

Change protein to dock to.

#### **Parameters**

protein: oddt.toolkit.Molecule object Protein object to be used.

## tmp\_dir

# oddt.scoring package

# **Subpackages**

# oddt.scoring.descriptors package

#### **Submodules**

# oddt.scoring.descriptors.binana module

Internal implementation of binana software (http://nbcr.ucsd.edu/data/sw/hosted/binana/)

class oddt.scoring.descriptors.binana.binana\_descriptor(protein=None)
 Bases: object

Descriptor build from binana script (as used in NNScore 2.0

#### **Parameters**

**protein: oddt.toolkit.Molecule object (default=None)** Protein object to be used while generating descriptors.

#### **Methods**

build(ligands[, protein])	Descriptor building method
set_protein(protein)	One function to change all relevant proteins

**build** (*ligands*, *protein=None*)

Descriptor building method

#### **Parameters**

**ligands: array-like** An array of generator of oddt.toolkit.Molecule objects for which the descriptor is computed

**protein: oddt.toolkit.Molecule object (default=None)** Protein object to be used while generating descriptors. If none, then the default protein (from constructor) is used. Otherwise, protein becomes new global and default protein.

#### **Returns**

**descs:** numpy array, shape=[n\_samples, 351] An array of binana descriptors, aligned with input ligands

# set\_protein (protein)

One function to change all relevant proteins

#### **Parameters**

**protein: oddt.toolkit.Molecule object** Protein object to be used while generating descriptors. Protein becomes new global and default protein.

#### **Module contents**

Bases: object

Close contacts descriptor which tallies atoms of type X in certain cutoff from atoms of type Y.

#### **Parameters**

protein: oddt.toolkit.Molecule or None (default=None) Default protein to use as reference

**cutoff:** int or list, shape=[n,] or shape=[n,2] (default=4) Cutoff for atoms in Angstroms given as an integer or a list of ranges, eg. [0, 4, 8, 12] or [[0,4],[4,8],[8,12]]. Upper bound is always inclusive, lower exclusive.

mode: string (default='atomic\_nums') Method of atoms selection, as used in atoms\_by\_type

ligand\_types: array List of ligand atom types to use

protein\_types: array List of protein atom types to use

**aligned\_pairs: bool (default=False)** Flag indicating should permutation of types should be done, otherwise the atoms are treated as aligned pairs.

#### **Methods**

build(ligands[, protein])

Builds descriptors for series of ligands

build (ligands, protein=None)
Builds descriptors for series of ligands

# **Parameters**

**ligands: iterable of oddt.toolkit.Molecules or oddt.toolkit.Molecule** A list or iterable of ligands to build the descriptor or a single molecule.

protein: oddt.toolkit.Molecule or None (default=None) Default protein to use as reference

class oddt.scoring.descriptors.fingerprints(fp='fp2', toolkit='ob')
 Bases: object

# Methods

build

build(mols)

Bases: object

# **Methods**

build	
set_protein	

build (ligands, protein=None)

set\_protein (protein)

Bases: object

#### **Methods**

build	
set_protein	

build (ligands, protein=None)

set\_protein (protein)

## oddt.scoring.functions package

## **Submodules**

# oddt.scoring.functions.NNScore module

```
class oddt.scoring.functions.NNScore.nnscore(protein=None, n_jobs=-1)
    Bases: oddt.scoring.scorer
```

NNScore implementation [1]. Based on Binana descriptors [2] and an ensemble of 20 best scored nerual networks with a hidden layer of 5 nodes. The NNScore predicts binding affinity (pKi/d).

# **Parameters**

protein [oddt.toolkit.Molecule object] Receptor for the scored ligands

**n\_jobs:** int (default=-1) Number of cores to use for scoring and training. By default (-1) all cores are allocated.

#### References

[1], [2]

## **Methods**

fit(ligands, target, *args, **kwargs)	Trains model on supplied ligands and target values
<pre>predict(ligands, *args, **kwargs)</pre>	Predicts values (eg.
	Continued on next page

Table 5 – continued from previous page

predict_ligand(ligand) Local method to score one ligand and upda		
	scores.	
predict_ligands(ligands)	Method to score ligands in a lazy fashion.	
save(filename)	Saves scoring function to a pickle file.	
score(ligands, target, *args, **kwargs)	Methods estimates the quality of prediction using	
	model's default score (accuracy for classification or	
	R^2 for regression)	
set_protein(protein)	Proxy method to update protein in all relevant places.	

gen_training_data	
load	
train	

fit (ligands, target, \*args, \*\*kwargs)

Trains model on supplied ligands and target values

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs] Ground truth (correct) target values.

classmethod load(filename=None, pdbbind\_version=2016)

Loads scoring function from a pickle file.

# **Parameters**

filename: string Pickle filename

#### Returns

sf: scorer-like object Scoring function object loaded from a pickle

predict (ligands, \*args, \*\*kwargs)

Predicts values (eg. affinity) for supplied ligands.

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

#### Returns

predicted: np.array or array of np.arrays of shape = [n\_ligands] Predicted scores for ligands

# predict\_ligand(ligand)

Local method to score one ligand and update it's scores.

#### **Parameters**

ligand: oddt.toolkit.Molecule object Ligand to be scored

#### Returns

ligand: oddt.toolkit.Molecule object Scored ligand with updated scores

predict\_ligands (ligands)

Method to score ligands in a lazy fashion.

#### **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to be scored

#### Returns

ligand: iterator of oddt.toolkit.Molecule objects Scored ligands with updated scores

save (filename)

Saves scoring function to a pickle file.

#### **Parameters**

filename: string Pickle filename

```
score (ligands, target, *args, **kwargs)
```

Methods estimates the quality of prediction using model's default score (accuracy for classification or R^2 for regression)

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs] Ground truth (correct) target values.

#### **Returns**

s: float Quality score (accuracy or R^2) for prediction

set\_protein (protein)

Proxy method to update protein in all relevant places.

#### **Parameters**

protein: oddt.toolkit.Molecule object New default protein

train (home\_dir=None, sf\_pickle=None, pdbbind\_version=2016)

# oddt.scoring.functions.PLECscore module

Bases: oddt.scoring.scorer

PLECscore - a novel scoring function based on PLEC fingerprints. The underlying model can be one of:

- · linear regression
- neural network (dense, 200x200x200)
- random forest (100 trees)

The scoring function is trained on PDBbind v2016 database and even with linear model outperforms other machine-learning ones in terms of Pearson correlation coefficient on "core set". For details see PLEC publication. PLECscore predicts binding affinity (pKi/d).

New in version 0.6.

#### **Parameters**

protein [oddt.toolkit.Molecule object] Receptor for the scored ligands

n\_jobs: int (default=-1) Number of cores to use for scoring and training. By default (-1) all cores are allocated.

- **version: str** (**default='linear'**) A version of scoring function ('linear', 'nn' or 'rf') which model should be used for the scoring function.
- **depth\_protein:** int (default=5) The depth of ECFP environments generated on the protein side of interaction. By default 6 (0 to 5) environments are generated.
- **depth\_ligand:** int (default=1) The depth of ECFP environments generated on the ligand side of interaction. By default 2 (0 to 1) environments are generated.
- **size:** int (default=65536) The final size of a folded PLEC fingerprint. This setting is not used to limit the data encoded in PLEC fingerprint (for that tune the depths), but only the final length. Setting it to too low value will lead to many collisions.

#### Methods

fit(ligands, target, *args, **kwargs)	Trains model on supplied ligands and target values
predict(ligands, *args, **kwargs)	Predicts values (eg.
predict_ligand(ligand)	Local method to score one ligand and update it's
	scores.
predict_ligands(ligands)	Method to score ligands in a lazy fashion.
save(filename)	Saves scoring function to a pickle file.
score(ligands, target, *args, **kwargs)	Methods estimates the quality of prediction using
	model's default score (accuracy for classification or
	R^2 for regression)
set_protein(protein)	Proxy method to update protein in all relevant places.

gen_json	
gen_training_data	
load	
train	

fit (ligands, target, \*args, \*\*kwargs)

Trains model on supplied ligands and target values

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs] Ground truth (correct) target values.

gen\_json (home\_dir=None, pdbbind\_version=2016)

classmethod load (filename=None, version='linear', pdbbind\_version=2016, depth\_protein=5, depth\_ligand=1, size=65536)

Loads scoring function from a pickle file.

#### **Parameters**

filename: string Pickle filename

# Returns

sf: scorer-like object Scoring function object loaded from a pickle

```
predict (ligands, *args, **kwargs)
```

Predicts values (eg. affinity) for supplied ligands.

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

#### **Returns**

predicted: np.array or array of np.arrays of shape = [n\_ligands] Predicted scores for ligands

# predict\_ligand(ligand)

Local method to score one ligand and update it's scores.

#### **Parameters**

ligand: oddt.toolkit.Molecule object Ligand to be scored

#### **Returns**

ligand: oddt.toolkit.Molecule object Scored ligand with updated scores

## predict\_ligands (ligands)

Method to score ligands in a lazy fashion.

#### **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to be scored

#### Returns

ligand: iterator of oddt.toolkit.Molecule objects Scored ligands with updated scores

#### save (filename)

Saves scoring function to a pickle file.

#### **Parameters**

**filename: string** Pickle filename

```
score (ligands, target, *args, **kwargs)
```

Methods estimates the quality of prediction using model's default score (accuracy for classification or R^2 for regression)

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs] Ground truth (correct) target values.

#### Returns

s: float Quality score (accuracy or R^2) for prediction

# set\_protein (protein)

Proxy method to update protein in all relevant places.

#### **Parameters**

protein: oddt.toolkit.Molecule object New default protein

train (home\_dir=None, sf\_pickle=None, pdbbind\_version=2016, ignore\_json=False)

# oddt.scoring.functions.RFScore module

```
class oddt.scoring.functions.RFScore.rfscore(protein=None, n\_jobs=-1, version=1, spr=0, **kwargs)
```

Bases: oddt.scoring.scorer

Scoring function implementing RF-Score variants. It predicts the binding affinity (pKi/d) of ligand in a complex utilizing simple descriptors (close contacts of atoms <12A) with sophisticated machine-learning model (random forest). The third variand supplements those contacts with Vina partial scores. For futher details see RF-Score publications v1[Rd9e4db499696-1]\_, v2[Rd9e4db499696-2]\_, v3[Rd9e4db499696-3]\_.

#### **Parameters**

protein [oddt.toolkit.Molecule object] Receptor for the scored ligands

**n\_jobs:** int (default=-1) Number of cores to use for scoring and training. By default (-1) all cores are allocated.

**version:** int (default=1) Scoring function variant. The deault is the simplest one (v1).

spr: int (default=0) The minimum number of contacts in each pair of atom types in the training set for the column to be included in training. This is a way of removal of not frequent and empty contacts.

#### References

[1], [2], [3]

# Methods

fit(ligands, target, *args, **kwargs)	Trains model on supplied ligands and target values
predict(ligands, *args, **kwargs)	Predicts values (eg.
predict_ligand(ligand)	Local method to score one ligand and update it's
	scores.
predict_ligands(ligands)	Method to score ligands in a lazy fashion.
save(filename)	Saves scoring function to a pickle file.
score(ligands, target, *args, **kwargs)	Methods estimates the quality of prediction using
	model's default score (accuracy for classification or
	R^2 for regression)
set_protein(protein)	Proxy method to update protein in all relevant places.

gen_training_data	
load	
train	

fit (ligands, target, \*args, \*\*kwargs)

Trains model on supplied ligands and target values

# **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

**target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs]** Ground truth (correct) target values.

classmethod load (filename=None, version=1, pdbbind\_version=2016)

Loads scoring function from a pickle file.

#### **Parameters**

filename: string Pickle filename

#### Returns

sf: scorer-like object Scoring function object loaded from a pickle

predict (ligands, \*args, \*\*kwargs)

Predicts values (eg. affinity) for supplied ligands.

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

#### **Returns**

predicted: np.array or array of np.arrays of shape = [n\_ligands] Predicted scores for ligands

#### predict\_ligand(ligand)

Local method to score one ligand and update it's scores.

#### **Parameters**

ligand: oddt.toolkit.Molecule object Ligand to be scored

#### Returns

ligand: oddt.toolkit.Molecule object Scored ligand with updated scores

# predict\_ligands (ligands)

Method to score ligands in a lazy fashion.

#### **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to be scored

#### Returns

ligand: iterator of oddt.toolkit.Molecule objects Scored ligands with updated scores

## save (filename)

Saves scoring function to a pickle file.

#### **Parameters**

filename: string Pickle filename

```
score (ligands, target, *args, **kwargs)
```

Methods estimates the quality of prediction using model's default score (accuracy for classification or R^2 for regression)

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs] Ground truth (correct) target values.

# Returns

s: float Quality score (accuracy or R^2) for prediction

#### set\_protein(protein)

Proxy method to update protein in all relevant places.

#### **Parameters**

protein: oddt.toolkit.Molecule object New default protein

train (home\_dir=None, sf\_pickle=None, pdbbind\_version=2016)

## **Module contents**

```
class oddt.scoring.functions.rfscore(protein=None, n\_jobs=-1, version=1, spr=0, **kwargs)
```

Bases: oddt.scoring.scorer

Scoring function implementing RF-Score variants. It predicts the binding affinity (pKi/d) of ligand in a complex utilizing simple descriptors (close contacts of atoms <12A) with sophisticated machine-learning model (random forest). The third variand supplements those contacts with Vina partial scores. For futher details see RF-Score publications v1[R062ccc3ea4fa-1]\_, v2[R062ccc3ea4fa-2]\_, v3[R062ccc3ea4fa-3]\_.

#### **Parameters**

protein [oddt.toolkit.Molecule object] Receptor for the scored ligands

**n\_jobs:** int (default=-1) Number of cores to use for scoring and training. By default (-1) all cores are allocated.

version: int (default=1) Scoring function variant. The deault is the simplest one (v1).

spr: int (default=0) The minimum number of contacts in each pair of atom types in the training set for the column to be included in training. This is a way of removal of not frequent and empty contacts.

#### References

[1], [2], [3]

#### **Methods**

fit(ligands, target, *args, **kwargs)	Trains model on supplied ligands and target values
<pre>predict(ligands, *args, **kwargs)</pre>	Predicts values (eg.
predict_ligand(ligand)	Local method to score one ligand and update it's
	scores.
predict_ligands(ligands)	Method to score ligands in a lazy fashion.
save(filename)	Saves scoring function to a pickle file.
score(ligands, target, *args, **kwargs)	Methods estimates the quality of prediction using
	model's default score (accuracy for classification or
	R^2 for regression)
set_protein(protein)	Proxy method to update protein in all relevant places.

gen_training_data	
load	
train	

```
fit (ligands, target, *args, **kwargs)
```

Trains model on supplied ligands and target values

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs] Ground truth (correct) target values.

classmethod load (filename=None, version=1, pdbbind\_version=2016)

Loads scoring function from a pickle file.

#### **Parameters**

filename: string Pickle filename

#### Returns

sf: scorer-like object Scoring function object loaded from a pickle

```
predict (ligands, *args, **kwargs)
```

Predicts values (eg. affinity) for supplied ligands.

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

#### Returns

predicted: np.array or array of np.arrays of shape = [n\_ligands] Predicted scores for ligands

# predict\_ligand(ligand)

Local method to score one ligand and update it's scores.

#### **Parameters**

ligand: oddt.toolkit.Molecule object Ligand to be scored

#### Returns

ligand: oddt.toolkit.Molecule object Scored ligand with updated scores

# predict\_ligands (ligands)

Method to score ligands in a lazy fashion.

#### **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to be scored

#### Returns

ligand: iterator of oddt.toolkit.Molecule objects Scored ligands with updated scores

# save (filename)

Saves scoring function to a pickle file.

#### **Parameters**

filename: string Pickle filename

```
score (ligands, target, *args, **kwargs)
```

Methods estimates the quality of prediction using model's default score (accuracy for classification or R^2 for regression)

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

**target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs]** Ground truth (correct) target values.

#### **Returns**

s: float Quality score (accuracy or R^2) for prediction

# set\_protein(protein)

Proxy method to update protein in all relevant places.

## **Parameters**

protein: oddt.toolkit.Molecule object New default protein

train (home\_dir=None, sf\_pickle=None, pdbbind\_version=2016)

```
class oddt.scoring.functions.nnscore(protein=None, n_jobs=-1)
    Bases: oddt.scoring.scorer
```

NNScore implementation [1]. Based on Binana descriptors [2] and an ensemble of 20 best scored nerual networks with a hidden layer of 5 nodes. The NNScore predicts binding affinity (pKi/d).

#### **Parameters**

protein [oddt.toolkit.Molecule object] Receptor for the scored ligands

n\_jobs: int (default=-1) Number of cores to use for scoring and training. By default (-1) all cores are allocated.

#### References

[1],[2]

## **Methods**

<pre>fit(ligands, target, *args, **kwargs)</pre>	Trains model on supplied ligands and target values
predict(ligands, *args, **kwargs)	Predicts values (eg.
<pre>predict_ligand(ligand)</pre>	Local method to score one ligand and update it's
	scores.
<pre>predict_ligands(ligands)</pre>	Method to score ligands in a lazy fashion.
save(filename)	Saves scoring function to a pickle file.
score(ligands, target, *args, **kwargs)	Methods estimates the quality of prediction using
	model's default score (accuracy for classification or
	R^2 for regression)
set_protein(protein)	Proxy method to update protein in all relevant places.

Q	en_training_data	
$\vdash$	oad	
t	rain	

fit (ligands, target, \*args, \*\*kwargs)

Trains model on supplied ligands and target values

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs] Ground truth (correct) target values.

classmethod load (filename=None, pdbbind\_version=2016)

Loads scoring function from a pickle file.

#### **Parameters**

filename: string Pickle filename

#### Returns

sf: scorer-like object Scoring function object loaded from a pickle

predict (ligands, \*args, \*\*kwargs)

Predicts values (eg. affinity) for supplied ligands.

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

#### Returns

predicted: np.array or array of np.arrays of shape = [n\_ligands] Predicted scores for ligands

### predict\_ligand(ligand)

Local method to score one ligand and update it's scores.

# **Parameters**

ligand: oddt.toolkit.Molecule object Ligand to be scored

### Returns

ligand: oddt.toolkit.Molecule object Scored ligand with updated scores

# predict\_ligands (ligands)

Method to score ligands in a lazy fashion.

## **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to be scored

#### Returns

ligand: iterator of oddt.toolkit.Molecule objects Scored ligands with updated scores

#### save (filename)

Saves scoring function to a pickle file.

## **Parameters**

filename: string Pickle filename

# score (ligands, target, \*args, \*\*kwargs)

Methods estimates the quality of prediction using model's default score (accuracy for classification or R^2 for regression)

# **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs] Ground truth (correct) target values.

#### Returns

s: float Quality score (accuracy or R^2) for prediction

set\_protein (protein)

Proxy method to update protein in all relevant places.

#### **Parameters**

protein: oddt.toolkit.Molecule object New default protein

train (home\_dir=None, sf\_pickle=None, pdbbind\_version=2016)

bases. Oddt. Scoring. Scorer

PLECscore - a novel scoring function based on PLEC fingerprints. The underlying model can be one of:

- · linear regression
- neural network (dense, 200x200x200)
- random forest (100 trees)

The scoring function is trained on PDBbind v2016 database and even with linear model outperforms other machine-learning ones in terms of Pearson correlation coefficient on "core set". For details see PLEC publication. PLECscore predicts binding affinity (pKi/d).

New in version 0.6.

#### **Parameters**

protein [oddt.toolkit.Molecule object] Receptor for the scored ligands

- n\_jobs: int (default=-1) Number of cores to use for scoring and training. By default (-1) all cores are allocated.
- **version: str** (**default='linear'**) A version of scoring function ('linear', 'nn' or 'rf') which model should be used for the scoring function.
- **depth\_protein:** int (default=5) The depth of ECFP environments generated on the protein side of interaction. By default 6 (0 to 5) environments are generated.
- **depth\_ligand:** int (default=1) The depth of ECFP environments generated on the ligand side of interaction. By default 2 (0 to 1) environments are generated.
- **size:** int (default=65536) The final size of a folded PLEC fingerprint. This setting is not used to limit the data encoded in PLEC fingerprint (for that tune the depths), but only the final lenght. Setting it to too low value will lead to many collisions.

### **Methods**

fit		
predict		
predict_ligand		
predict_ligands		
save		

Continued on next page

Table 10 - continued from previous page

score
set\_protein

gen_json	
gen_training_data	
load	
train	

fit (ligands, target, \*args, \*\*kwargs)

Trains model on supplied ligands and target values

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs] Ground truth (correct) target values.

gen\_json (home\_dir=None, pdbbind\_version=2016)

Loads scoring function from a pickle file.

# **Parameters**

filename: string Pickle filename

## Returns

sf: scorer-like object Scoring function object loaded from a pickle

predict (ligands, \*args, \*\*kwargs)

Predicts values (eg. affinity) for supplied ligands.

### **Parameters**

**ligands:** array-like of ligands Molecules to featurize and feed into the model

## Returns

predicted: np.array or array of np.arrays of shape = [n\_ligands] Predicted scores for ligands

#### predict\_ligand(ligand)

Local method to score one ligand and update it's scores.

# **Parameters**

ligand: oddt.toolkit.Molecule object Ligand to be scored

#### Returns

ligand: oddt.toolkit.Molecule object Scored ligand with updated scores

# predict\_ligands (ligands)

Method to score ligands in a lazy fashion.

### **Parameters**

ligands: iterable of oddt.toolkit.Molecule objects Ligands to be scored

#### Returns

ligand: iterator of oddt.toolkit.Molecule objects Scored ligands with updated scores

save (filename)

Saves scoring function to a pickle file.

#### **Parameters**

filename: string Pickle filename

score (ligands, target, \*args, \*\*kwargs)

Methods estimates the quality of prediction using model's default score (accuracy for classification or R^2 for regression)

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs] Ground truth (correct) target values.

#### Returns

s: float Quality score (accuracy or R^2) for prediction

set\_protein (protein)

Proxy method to update protein in all relevant places.

#### **Parameters**

protein: oddt.toolkit.Molecule object New default protein

train (home\_dir=None, sf\_pickle=None, pdbbind\_version=2016, ignore\_json=False)

# oddt.scoring.models package

# **Submodules**

# oddt.scoring.models.classifiers module

```
oddt.scoring.models.classifiers. {\bf randomforest}\\ alias of {\tt sklearn.ensemble.forest.RandomForestClassifier}
```

class oddt.scoring.models.classifiers.svm(\*args, \*\*kwargs)
 Bases: oddt.scoring.models.classifiers.OddtClassifier

## Methods

fit	
get_params	
predict	
predict_log_proba	
predict_proba	
score	
set_params	

```
fit (descs, target_values, **kwargs)
     get_params (deep=True)
     predict (descs)
     predict_log_proba (descs)
     predict_proba (descs)
     score (descs, target values)
          Returns the mean accuracy on the given test data and labels.
          In multi-label classification, this is the subset accuracy which is a harsh metric since you require for each
          sample that each label set be correctly predicted.
              Parameters
                  X [array-like, shape = (n_samples, n_features)] Test samples.
                  y [array-like, shape = (n_samples) or (n_samples, n_outputs)] True labels for X.
                  sample_weight [array-like, shape = [n_samples], optional] Sample weights.
              Returns
                  score [float] Mean accuracy of self.predict(X) wrt. y.
     set_params (**kwargs)
class oddt.scoring.models.classifiers.neuralnetwork(*args, **kwargs)
     Bases: oddt.scoring.models.classifiers.OddtClassifier
```

#### **Methods**

fit	
get_params	
predict	
predict_log_proba	
predict_proba	
score	
set_params	

In multi-label classification, this is the subset accuracy which is a harsh metric since you require for each sample that each label set be correctly predicted.

#### **Parameters**

**X** [array-like, shape = (n\_samples, n\_features)] Test samples.

```
y [array-like, shape = (n_samples) or (n_samples, n_outputs)] True labels for X.
sample_weight [array-like, shape = [n_samples], optional] Sample weights.
Returns
```

```
score [float] Mean accuracy of self.predict(X) wrt. y.
set_params (**kwargs)
```

# oddt.scoring.models.regressors module

Collection of regressors models

```
oddt.scoring.models.regressors.randomforest
    alias of sklearn.ensemble.forest.RandomForestRegressor
class oddt.scoring.models.regressors.svm(*args, **kwargs)
    Bases: oddt.scoring.models.regressors.OddtRegressor
```

#### **Methods**

fit	
get_params	
predict	
score	
set_params	

```
fit (descs, target_values, **kwargs)
get_params (deep=True)
predict (descs)
score (descs, target_values)
    Returns the coefficient of determination R^2 of the prediction.
```

The coefficient R^2 is defined as (1 - u/v), where u is the residual sum of squares ((y\_true - y\_pred) \*\* 2).sum() and v is the total sum of squares ((y\_true - y\_true.mean()) \*\* 2).sum(). The best possible score is 1.0 and it can be negative (because the model can be arbitrarily worse). A constant model that always predicts the expected value of y, disregarding the input features, would get a R^2 score of 0.0.

### **Parameters**

**Returns** 

```
    X [array-like, shape = (n_samples, n_features)] Test samples.
    y [array-like, shape = (n_samples) or (n_samples, n_outputs)] True values for X.
    sample_weight [array-like, shape = [n_samples], optional] Sample weights.
```

```
score [float] R^2 of self.predict(X) wrt. y.
set_params (**kwargs)

oddt.scoring.models.regressors.pls
    alias of sklearn.cross_decomposition.pls_.PLSRegression

class oddt.scoring.models.regressors.neuralnetwork (*args, **kwargs)
    Bases: oddt.scoring.models.regressors.OddtRegressor
```

# **Methods**

fit	
get_params	
predict	
score	
set_params	

```
fit (descs, target_values, **kwargs)
get_params (deep=True)
predict (descs)
score (descs, target_values)
    Returns the coefficient of determination R^2 of the prediction.
```

The coefficient R^2 is defined as (1 - u/v), where u is the residual sum of squares ((y\_true - y\_pred) \*\* 2).sum() and v is the total sum of squares ((y\_true - y\_true.mean()) \*\* 2).sum(). The best possible score is 1.0 and it can be negative (because the model can be arbitrarily worse). A constant model that always predicts the expected value of y, disregarding the input features, would get a R^2 score of 0.0.

# **Parameters**

```
\label{eq:continuous} \begin{split} \textbf{X} \;\; & [\text{array-like, shape} = (\text{n\_samples}, \, \text{n\_features})] \; \text{Test samples}. \\ \textbf{y} \;\; & [\text{array-like, shape} = (\text{n\_samples}) \; \text{or} \; (\text{n\_samples}, \, \text{n\_outputs})] \; \text{True values for } X. \\ \textbf{sample\_weight} \;\; & [\text{array-like, shape} = [\text{n\_samples}], \; \text{optional}] \; \text{Sample weights}. \end{split}
```

# Returns

```
score [float] R^2 of self.predict(X) wrt. y.
```

```
set_params (**kwargs)
```

```
oddt.scoring.models.regressors.mlr
alias of sklearn.linear_model.base.LinearRegression
```

#### **Module contents**

#### Module contents

```
oddt.scoring.cross_validate (model, cv_set, cv_target, n=10, shuffle=True, n_jobs=1)
Perform cross validation of model using provided data
```

#### **Parameters**

```
model: object Model to be tested
cv_set: array-like of shape = [n_samples, n_features] Estimated target values.
cv_target: array-like of shape = [n_samples] or [n_samples, n_outputs] Estimated target values.
n: integer (default = 10) How many folds to be created from dataset
shuffle: bool (default = True) Should data be shuffled before folding.
n_jobs: integer (default = 1) How many CPUs to use during cross validation
```

### Returns

**r2:** array of shape = [n] R<sup>2</sup> score for each of generated folds

 $\verb"class" oddt.scoring.ensemble_descriptor" (\textit{descriptor}\_\textit{generators})$ 

Bases: object

Proxy class to build an ensemble of destriptors with an API as one

#### **Parameters**

models: array An array of models

#### **Methods**

build	
set_protein	

build (mols, \*args, \*\*kwargs)

set\_protein (protein)

class oddt.scoring.ensemble\_model(models)

Bases: object

Proxy class to build an ensemble of models with an API as one

#### **Parameters**

models: array An array of models

## **Methods**

fit	
predict	
score	

```
fit (X, y, *args, **kwargs)
```

predict (X, \*args, \*\*kwargs)

score (X, y, \*args, \*\*kwargs)

class oddt.scoring.scorer(model\_instance, descriptor\_generator\_instance, score\_title='score')

Bases: object

Scorer class is parent class for scoring functions.

### **Parameters**

model\_instance: model Medel compatible with sklearn API (fit, predict and score methods)

descriptor\_generator\_instance: array of descriptors Descriptor generator object

**score\_title: string** Title of score to be used.

# **Methods**

fit(ligands, target, *args, **kwargs)	Trains model on supplied ligands and target values
load(filename)	Loads scoring function from a pickle file.
predict(ligands, *args, **kwargs)	Predicts values (eg.
<pre>predict_ligand(ligand)</pre>	Local method to score one ligand and update it's
	scores.
predict_ligands(ligands)	Method to score ligands in a lazy fashion.
save(filename)	Saves scoring function to a pickle file.
score(ligands, target, *args, **kwargs)	Methods estimates the quality of prediction using
	model's default score (accuracy for classification or
	R^2 for regression)
set_protein(protein)	Proxy method to update protein in all relevant places.

# fit (ligands, target, \*args, \*\*kwargs)

Trains model on supplied ligands and target values

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

**target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs]** Ground truth (correct) target values.

# classmethod load(filename)

Loads scoring function from a pickle file.

# **Parameters**

filename: string Pickle filename

#### Returns

sf: scorer-like object Scoring function object loaded from a pickle

# predict (ligands, \*args, \*\*kwargs)

Predicts values (eg. affinity) for supplied ligands.

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

# Returns

predicted: np.array or array of np.arrays of shape = [n\_ligands] Predicted scores for ligands

# ${\tt predict\_ligand}\,(ligand)$

Local method to score one ligand and update it's scores.

# **Parameters**

ligand: oddt.toolkit.Molecule object Ligand to be scored

## Returns

ligand: oddt.toolkit.Molecule object Scored ligand with updated scores

### predict\_ligands (ligands)

Method to score ligands in a lazy fashion.

# **Parameters**

```
ligands: iterable of oddt.toolkit.Molecule objects Ligands to be scored
```

### Returns

ligand: iterator of oddt.toolkit.Molecule objects Scored ligands with updated scores

save (filename)

Saves scoring function to a pickle file.

#### **Parameters**

filename: string Pickle filename

score (ligands, target, \*args, \*\*kwargs)

Methods estimates the quality of prediction using model's default score (accuracy for classification or R^2 for regression)

#### **Parameters**

ligands: array-like of ligands Molecules to featurize and feed into the model

**target: array-like of shape = [n\_samples] or [n\_samples, n\_outputs]** Ground truth (correct) target values.

#### **Returns**

s: float Quality score (accuracy or R^2) for prediction

set\_protein (protein)

Proxy method to update protein in all relevant places.

#### **Parameters**

protein: oddt.toolkit.Molecule object New default protein

oddt.toolkits package

**Subpackages** 

oddt.toolkits.extras package

**Subpackages** 

oddt.toolkits.extras.rdkit package

**Submodules** 

oddt.toolkits.extras.rdkit.fixer module

```
exception oddt.toolkits.extras.rdkit.fixer.AddAtomsError
    Bases: exceptions.Exception
```

args

message

oddt.toolkits.extras.rdkit.fixer.**AddMissingAtoms** (*protein*, *residue*, *amap*, *template*)
Add missing atoms to protein molecule only at the residue according to template.

**Parameters** 

# protein: rdkit.Chem.rdchem.RWMol

Mol with whole protein. Note that it is modified in place.

residue: Mol with residue only

 $\textbf{amap: list} \ \, \text{List mapping atom IDs in residue to atom IDs in whole protein } (\text{amap[i]} = j$ 

means that i'th atom in residue corresponds to j'th atom in protein)

template: Residue template

#### Returns

\_\_\_

protein: rdkit.Chem.rdchem.RWMol Modified protein

visited\_bonds: list Bonds that match the template

is\_complete: bool Indicates whether all atoms in template were found in residue

oddt.toolkits.extras.rdkit.fixer.**ExtractPocketAndLigand** (mol, cutoff=12.0, expandResidues=True, ligand\_residue=None, ligand\_residue\_blacklist=None, append residues=None)

Function extracting a ligand (the largest HETATM residue) and the protein pocket within certain cutoff. The selection of pocket atoms can be expanded to contain whole residues. The single atom HETATM residues are attributed to pocket (metals and waters)

#### **Parameters**

#### mol: rdkit.Chem.rdchem.Mol

Molecule with a protein ligand complex

cutoff: float (default=12.) Distance cutoff for the pocket atoms

**expandResidues:** bool (default=True) Expand selection to whole residues within cutoff.

**ligand\_residue:** string (default None) Residue name which explicitly pint to a ligand(s).

**ligand\_residue\_blacklist:** array-like, optional (default None) List of residues to ignore during ligand lookup.

**append\_residues: array-like, optional (default None)** List of residues to append to pocket, even if they are HETATM, such as MSE, ATP, AMP, ADP, etc.

#### Returns

#### pocket: rdkit.Chem.rdchem.RWMol

Pocket constructed of protein residues/atoms around ligand

ligand: rdkit.Chem.rdchem.RWMol Largest HETATM residue contained in input molecule

oddt.toolkits.extras.rdkit.fixer.**FetchAffinityTable** (*pdbids*, *affinity\_types*) Fetch affinity data from RCSB PDB server.

# **Parameters**

pdbids: array-like

List of PDB IDs of structres with protein-ligand complexes.

**affinity\_types:** array-like List of types of affinity data to retrieve. Available types are: Ki, Kd, EC50, IC50, deltaG, deltaH, deltaS, Ka.

# Returns

**ligand\_affinity: pd.DataFrame** Table with protein-ligand binding affinities. Table contains following columns: structureId, ligandId, ligandFormula, ligandMolecularWeight + columns named after affinity types specified byt the user.

oddt.toolkits.extras.rdkit.fixer.**FetchStructure**(pdbid, sanitize=False, re-moveHs=True, cache\_dir=None)

Fetch the structure in PDB format from RCSB PDB server and read it with rdkit.

#### **Parameters**

#### pdbid: str

PDB IDs of the structre

sanitize: bool, optional (default False) Toggles molecule sanitationremoveHs: bool, optional (default False) Indicates wheter Hs should be removed during reading

#### Returns

mol: Chem.rdchem.Mol Retrieved molecule

exception oddt.toolkits.extras.rdkit.fixer.FixerError
Bases: exceptions.Exception

# args

# message

oddt.toolkits.extras.rdkit.fixer.**GetAtomResidueId**(atom)

Return (residue number, residue name, chain id) for a given atom

oddt.toolkits.extras.rdkit.fixer.**GetResidues** (mol, atom\_list=None)

Create dictrionary that maps residues to atom IDs: (res number, res name, chain id) -> [atom1 idx, atom2 idx, ...]

oddt.toolkits.extras.rdkit.fixer.**IsResidueConnected**(*mol*, *atom\_ids*)

Check if residue with given atom IDs is connected to other residues in the molecule.

oddt.toolkits.extras.rdkit.fixer.MolToTemplates (mol)

Prepare set of templates for a given PDB residue.

oddt.toolkits.extras.rdkit.fixer.**PrepareComplexes**(pdbids, pocket\_dist\_cutoff=12.0, affinity\_types=None,

cache dir=None)

Fetch structures and affinity data from RCSB PDB server and prepare ligand-pocket pairs for small molecules with known activites.

#### **Parameters**

### pdbids: array-like

List of PDB IDs of structres with protein-ligand complexes.

pocket\_dist\_cutoff: float, optional (default 12.) Distance cutoff for the pocket atoms

**affinity\_types: array-like, optional (default None)** List of types of affinity data to retrieve. Available types are: Ki, Kd, EC50, IC50, deltaG, deltaH, deltaS, Ka. If not specified Ki, Kd, EC50, and IC50 are used.

#### Returns

**complexes: dict** Dictionary with pocket-ligand paris, structured as follows: {'pdbid': {'ligid': (pocket\_mol, ligand\_mol)}. Ligands have binding affinity data stored as properties.

```
oddt.toolkits.extras.rdkit.fixer.PreparePDBMol (mol, removeHs=True, remove-
HOHs=True, residue_whitelist=None,
residue_blacklist=None, re-
move_incomplete=False,
add_missing_atoms=False,
custom_templates=None, re-
place_default_templates=False)
```

# Prepares protein molecule by:

- Removing Hs by hard using atomic number [default=True]
- Removes HOH [default=True]
- Assign bond orders from smiles of PDB residues (over 24k templates)
- · Removes bonds to metals

#### **Parameters**

#### mol: rdkit.Chem.rdchem.Mol

Mol with whole protein.

**removeHs: bool, optional (default True)** If True, hydrogens will be forcefully removed **removeHOHs: bool, optional (default True)** If True, remove waters using residue name

**residue\_whitelist: array-like, optional (default None)** List of residues to clean. If not specified, all residues present in the structure will be used.

**residue\_blacklist: array-like, optional (default None)** List of residues to ignore during cleaning. If not specified, all residues present in the structure will be cleaned.

**remove\_incomplete: bool, optional (default False)** If True, remove residues that do not fully match the template

add\_missing\_atoms: bool (default=False) Switch to add missing atoms accordingly to template SMILES structure.

custom\_templates: str or dict, optional (default None) Custom templates for residues.
Can be either path to SMILES file, or dictionary mapping names to SMILES or Mol objects

**replace\_default\_templates: bool, optional (default False)** Indicates whether default default templates should be replaced by cusom ones. If False, default templates will be updated with custom ones. This argument is ignored if custom\_templates is None.

#### Returns

new\_mol: rdkit.Chem.rdchem.RWMol Modified protein

oddt.toolkits.extras.rdkit.fixer.PreparePDBResidue (protein, residue, amap, template)

#### **Parameters**

# protein: rdkit.Chem.rdchem.RWMol

Mol with whole protein. Note that it is modified in place.

residue: Mol with residue only

**amap:** list List mapping atom IDs in residue to atom IDs in whole protein (amap[i] = j

means that i'th atom in residue corresponds to j'th atom in protein)

template: Residue template

#### **Returns**

\_\_\_\_

protein: rdkit.Chem.rdchem.RWMol Modified protein

visited\_bonds: list Bonds that match the template

is\_complete: bool Indicates whether all atoms in template were found in residue

oddt.toolkits.extras.rdkit.fixer.ReadTemplates (filename, resnames)

Load templates from file for specified residues

exception oddt.toolkits.extras.rdkit.fixer.SanitizeError

Bases: exceptions. Exception

args

message

oddt.toolkits.extras.rdkit.fixer.SimplifyMol(mol)

Change all bonds to single and discharge/dearomatize all atoms. The molecule is modified in-place (no copy is made).

exception oddt.toolkits.extras.rdkit.fixer.SubstructureMatchError

Bases: exceptions. Exception

args

message

oddt.toolkits.extras.rdkit.fixer.**UFFConstrainedOptimize**(*mol, moving\_atoms=None, fixed atoms=None, cut-*

off=5.0, verbose=False)

Minimize a molecule using UFF forcefield with a set of moving/fixed atoms. If both moving and fixed atoms are provided, fixed\_atoms parameter will be ignored. The minimization is done in-place (without copying molecule).

#### **Parameters**

## mol: rdkit.Chem.rdchem.Mol

Molecule to be minimized.

**moving\_atoms:** array-like (default=None) Indices of freely moving atoms. If None, fixed atoms are assigned based on *fixed\_atoms*. These two arguments are mutually exclusive.

**fixed\_atoms:** array-like (**default=None**) Indices of fixed atoms. If None, fixed atoms are assigned based on *moving\_atoms*. These two arguments are mutually exclusive.

cutoff: float (default=10.) Distance cutoff for the UFF minimization

#### Returns

mol: rdkit.Chem.rdchem.Mol Molecule with mimimized moving\_atoms

# **Module contents**

oddt.toolkits.extras.rdkit.**AtomListToSubMol** (mol, amap, includeConformer=False)

#### **Parameters**

#### mol: rdkit.Chem.rdchem.Mol

Molecule

amap: array-like List of atom indices (zero-based)

**includeConformer: bool (default=True)** Toogle to include atoms coordinates in submolecule.

#### Returns

submol: rdkit.Chem.rdchem.RWMol Submol determined by specified atom list

oddt.toolkits.extras.rdkit.MolFromPDBBlock (molBlock, sanitize=True, removeHs=True, flavor=0)

oddt.toolkits.extras.rdkit.**MolFromPDBQTBlock** (block, sanitize=True, removeHs=True) Read PDBQT block to a RDKit Molecule

#### **Parameters**

## block: string

Residue name which explicitly pint to a ligand(s).

sanitize: bool (default=True) Should the sanitization be performedremoveHs: bool (default=True) Should hydrogens be removed when reading molecule.

# Returns

mol: rdkit.Chem.rdchem.Mol Molecule read from PDBQT

oddt.toolkits.extras.rdkit.**MolToPDBQTBlock**(*mol*, *flexible=True*, *addHs=False*, *com- puteCharges=False*)

Write RDKit Molecule to a PDBQT block

#### **Parameters**

#### mol: rdkit.Chem.rdchem.Mol

Molecule with a protein ligand complex

**flexible: bool (default=True)** Should the molecule encode torsions. Ligands should be flexible, proteins in turn can be rigid.

addHs: bool (default=False) The PDBQT format requires at least polar Hs on donors. By default Hs are added.

**computeCharges: bool (default=False)** Should the partial charges be automatically computed. If the Hs are added the charges must and will be recomputed. If there are no partial charge information, they are set to 0.0.

#### Returns

block: str String wit PDBQT encoded molecule

oddt.toolkits.extras.rdkit.PDBQTAtomLines (mol, donors, acceptors)

Create a list with PDBQT atom lines for each atom in molecule. Donors and acceptors are given as a list of atom indices.

oddt.toolkits.extras.rdkit.PathFromAtomList(mol, amap)

#### **Module contents**

# **Submodules**

#### oddt.toolkits.common module

Code common to all toolkits

```
oddt.toolkits.common.canonize_ring_path(path)
```

Make a canonic path - list of consecutive atom IDXs bonded in a ring sorted in an uniform fasion.

- 1. Move the smallest index to position 0
- 2. Look for the smallest first step (delta IDX)
- 3. Ff -1 is smallest, inverse the path and move min IDX to position 0

#### **Parameters**

path [list of integers] A list of consecutive atom indices in a ring

#### Returns

canonic\_path [list of integers] Sorted list of atoms

```
oddt.toolkits.common.detect_secondary_structure(res_dict)
```

Detect alpha helices and beta sheets in res\_dict by phi and psi angles

# oddt.toolkits.ob module

# Attributes

atomicmass

atomicnum

bonds

cidx

coordidx

coords

exactmass

formalcharge

```
heterovalence
         hyb
         idx DEPRECATED: RDKit is 0-based and OpenBabel is 1-based.
         idx0 Note that this index is 0-based and OpenBabel's internal index in 1-based.
         idx1 Note that this index is 1-based as OpenBabel's internal index.
         implicitvalence
         isotope
         neighbors
         partialcharge
         residue
         spin
         type
         valence
         vector
atomicmass
atomicnum
bonds
cidx
coordidx
coords
exactmass
formalcharge
heavyvalence
heterovalence
hyb
idx
     DEPRECATED: RDKit is 0-based and OpenBabel is 1-based. State which convention you desire and use
     idx0 or idx1.
     Note that this index is 1-based as OpenBabel's internal index.
idx0
     Note that this index is 0-based and OpenBabel's internal index in 1-based. Changed to be compatible with
     RDKit
idx1
     Note that this index is 1-based as OpenBabel's internal index.
implicitvalence
isotope
neighbors
```

heavyvalence

```
partialcharge
     residue
     spin
    type
     valence
     vector
class oddt.toolkits.ob.AtomStack(OBMol)
     Bases: object
class oddt.toolkits.ob.Bond(OBBond)
     Bases: object
         Attributes
            atoms
            isrotor
            order
     atoms
     isrotor
     order
class oddt.toolkits.ob.BondStack(OBMol)
     Bases: object
class oddt.toolkits.ob.Fingerprint(fingerprint)
     Bases: pybel.Fingerprint
         Attributes
            bits
            raw
    bits
     raw
class oddt.toolkits.ob.Molecule(OBMol=None, source=None, protein=False)
     Bases: pybel.Molecule
         Attributes
            OBMol
            atom_dict
            atoms
            bonds
            canonic_order Returns np.array with canonic order of heavy atoms in the molecule
            charge
            charges
            clone
            conformers
```

coords

data

dim

energy

exactmass

formula

molwt

num\_rotors Number of strict rotatable

**protein** A flag for identifing the protein molecules, for which *atom\_dict* procedures may differ.

res\_dict

residues

ring\_dict

smiles

spin

sssr

title

unitcell

# **Methods**

addh([only_polar])	Add hydrogens
calccharges([model])	Calculate partial charges for a molecule.
calcdesc([descnames])	Calculate descriptor values.
calcfp([fptype])	Calculate a molecular fingerprint.
convertdbonds()	Convert Dative Bonds.
draw([show, filename, update, usecoords])	Create a 2D depiction of the molecule.
<pre>localopt([forcefield, steps])</pre>	Locally optimize the coordinates.
make2D()	Generate 2D coordinates for molecule
make3D([forcefield, steps])	Generate 3D coordinates
removeh()	Remove hydrogens

clone\_coords write

OBMol

addh (only\_polar=False)
 Add hydrogens

atom\_dict

atoms

bonds

```
calccharges (model='gasteiger')
     Calculate partial charges for a molecule. By default the Gasteiger charge model is used.
         Parameters
             model [str (default="gasteiger")] Method for generating partial charges. Supported models:
                * gasteiger * mmff94 * others supported by OpenBabel (obabel -L charges)
calcdesc(descnames=[])
     Calculate descriptor values.
     Optional parameter: descnames – a list of names of descriptors
     If descnames is not specified, all available descriptors are calculated. See the descs variable for a list of
     available descriptors.
calcfp (fptype='FP2')
     Calculate a molecular fingerprint.
     Optional parameters:
         fptype – the fingerprint type (default is "FP2"). See the fps variable for a list of of available fin-
             gerprint types.
canonic order
     Returns np.array with canonic order of heavy atoms in the molecule
charge
charges
clone
clone_coords (source)
conformers
convertdbonds()
     Convert Dative Bonds.
coords
data
dim
draw(show=True, filename=None, update=False, usecoords=False)
     Create a 2D depiction of the molecule.
     Optional parameters: show – display on screen (default is True) filename – write to file (default is None)
         update – update the coordinates of the atoms to those
             determined by the structure diagram generator (default is False)
         usecoords – don't calculate 2D coordinates, just use the current coordinates (default is False)
     Tkinter and Python Imaging Library are required for image display.
energy
exactmass
formula
localopt (forcefield='mmff94', steps=500)
     Locally optimize the coordinates.
```

# **Optional parameters:**

steps – default is 500

```
If the molecule does not have any coordinates, make 3D() is called before the optimization. Note that the
           molecule needs to have explicit hydrogens. If not, call addh().
     make2D()
           Generate 2D coordinates for molecule
     make3D (forcefield='mmff94', steps=50)
           Generate 3D coordinates
     molwt
     num rotors
           Number of strict rotatable
     protein
           A flag for identifing the protein molecules, for which atom_dict procedures may differ.
     removeh()
           Remove hydrogens
     res dict
     residues
     ring_dict
     smiles
     spin
     sssr
     title
     unitcell
     write (format='smi', filename=None, overwrite=False, opt=None, size=None)
           Write the molecule to a file or return a string.
           Optional parameters:
               format - see the informats variable for a list of available output formats (default is "smi")
               filename - default is None overwite - if the output file already exists, should it
                   be overwritten? (default is False)
               opt – a dictionary of format specific options For format options with no parameters, specify the
                   value as None.
           If a filename is specified, the result is written to a file. Otherwise, a string is returned containing the result.
           To write multiple molecules to the same file you should use the Outputfile class.
class oddt.toolkits.ob.MoleculeData(obmol)
     Bases: pybel.MoleculeData
```

forcefield – default is "mmff94". See the forcefields variable for a list of available forcefields.

# **Methods**

clear	
has_key	
items	
iteritems	
keys	
to_dict	
update	
values	

```
clear()
has_key(key)
items()
items()
keys()
to_dict()
update(dictionary)
values()
class oddt.toolkits.ob.Outputfile(format, filename, overwrite=False, opt=None)
Bases: pybel.Outputfile
```

# **Methods**

close()	Close the Outputfile to further writing.
write(molecule)	Write a molecule to the output file.

```
close()
```

Close the Outputfile to further writing.

write(molecule)

Write a molecule to the output file.

**Required parameters:** molecule

class oddt.toolkits.ob.Residue(OBResidue)

Bases: object

Represent a Pybel residue.

Required parameter: OBResidue – an Open Babel OBResidue

Attributes: atoms, idx, name.

(refer to the Open Babel library documentation for more info).

The original Open Babel atom can be accessed using the attribute: OBResidue

#### **Attributes**

atoms List of Atoms in the Residue

```
chain Resdiue chain ID
              idx DEPRECATED: Use idx0 instead.
              idx0 Internal index (0-based) of the Residue
              name Residue name
              number Residue number
     atoms
          List of Atoms in the Residue
     chain
          Resdiue chain ID
     idx
          DEPRECATED: Use idx0 instead.
          Internal index (0-based) of the Residue
     idx0
          Internal index (0-based) of the Residue
     name
          Residue name
     number
          Residue number
class oddt.toolkits.ob.ResidueStack(OBMol)
     Bases: object
class oddt.toolkits.ob.Smarts(smartspattern)
     Bases: pybel.Smarts
     Initialise with a SMARTS pattern.
```

# **Methods**

findall(molecule[, unique])	Find all matches of the SMARTS pattern to a partic-
	ular molecule
match(molecule)	Checks if there is any match.

```
findall (molecule, unique=True)
```

Find all matches of the SMARTS pattern to a particular molecule

```
match (molecule)
```

Checks if there is any match. Returns True or False

Produce diverse conformers using current conformer as starting point. Returns a generator. Each conformer is a copy of original molecule object.

New in version 0.6.

### **Parameters**

**mol** [oddt.toolkit.Molecule object] Molecule for which generating conformers

**n\_conf** [int (default=10)] Targer number of conformers

**method** [string (default='confab')] Method for generating conformers. Supported methods: \* confab \* ga

seed [None or int (default=None)] Random seed

**mutability** [int (default=5)] The inverse of probability of mutation. By default 5, which translates to 1/5 (20%) chance of mutation. This setting only works with genetic algorithm method ("ga").

**convergence** [int (default=5)] The number of generations with unchanged fitness, should the algorithm converge. This setting only works with genetic algorithm method ("ga").

**rmsd** [float (default=0.5)] The conformers are pruned unless their RMSD is higher than this cutoff. This setting only works with Confab method ("confab").

**nconf** [int (default=10000)] The number of initial conformers to generate before energy pruning. This setting only works with Confab method ("confab").

**energy\_gap** [float (default=5000.)] Energy gap from the lowest energy conformer to the highest possible. This setting only works with Confab method ("confab").

#### Returns

mols [list of oddt.toolkit.Molecule objects] Molecules with diverse conformers

oddt.toolkits.ob.readfile(format, filename, opt=None, lazy=False)

### oddt.toolkits.rdk module

rdkit - A Cinfony module for accessing the RDKit from CPython

**Global variables:** Chem and AllChem - the underlying RDKit Python bindings informats - a dictionary of supported input formats outformats - a dictionary of supported output formats descs - a list of supported descriptors fps - a list of supported fingerprint types forcefields - a list of supported forcefields

```
class oddt.toolkits.rdk.Atom(Atom)
```

Bases: object

Represent an rdkit Atom.

**Required parameters:** Atom – an RDKit Atom **Attributes:** atomicnum, coords, formalcharge

The original RDKit Atom can be accessed using the attribute: Atom

#### **Attributes**

atomicnum

bonds

coords

#### formalcharge

idx DEPRECATED: RDKit is 0-based and OpenBabel is 1-based.

idx0 Note that this index is 0-based as RDKit's

idx1 Note that this index is 1-based and RDKit's internal index in 0-based.

neighbors

```
atomicnum
     bonds
     coords
     formalcharge
     idx
          DEPRECATED: RDKit is 0-based and OpenBabel is 1-based. State which convention you desire and use
          idx0 or idx1.
          Note that this index is 1-based and RDKit's internal index in 0-based. Changed to be compatible with
              OpenBabel
     idx0
          Note that this index is 0-based as RDKit's
     idx1
          Note that this index is 1-based and RDKit's internal index in 0-based. Changed to be compatible with
          OpenBabel
     neighbors
     partialcharge
class oddt.toolkits.rdk.AtomStack(Mol)
     Bases: object
class oddt.toolkits.rdk.Bond(Bond)
     Bases: object
          Attributes
              atoms
              isrotor
              order
     atoms
     isrotor
     order
class oddt.toolkits.rdk.BondStack(Mol)
     Bases: object
class oddt.toolkits.rdk.Fingerprint(fingerprint)
     Bases: object
     A Molecular Fingerprint.
     Required parameters: fingerprint – a vector calculated by one of the fingerprint methods
     Attributes: fp – the underlying fingerprint object bits – a list of bits set in the Fingerprint
     Methods: The "|" operator can be used to calculate the Tanimoto coeff. For example, given two Fingerprints
          'a', and 'b', the Tanimoto coefficient is given by:
              tanimoto = a \mid b
          Attributes
```

partialcharge

#### raw

#### raw

class oddt.toolkits.rdk.Molecule(Mol=None, source=None, protein=False)
 Bases: object

Trap RDKit molecules which are 'None'

#### **Attributes**

Mol

atom\_dict

atoms

bonds

canonic\_order Returns np.array with canonic order of heavy atoms in the molecule

charges

clone

coords

data

formula

molwt

 $num\_rotors$ 

protein A flag for identifing the protein molecules, for which atom\_dict procedures may
differ.

res\_dict

residues

ring\_dict

smiles

sssr

title

# **Methods**

addh([only_polar])	Add hydrogens.
calccharges([model])	Calculate partial charges for a molecule.
calcdesc([descnames])	Calculate descriptor values.
calcfp([fptype, opt])	Calculate a molecular fingerprint.
<pre>localopt([forcefield, steps])</pre>	Locally optimize the coordinates.
make2D()	Generate 2D coordinates for molecule
make3D([forcefield, steps])	Generate 3D coordinates.
removeh(**kwargs)	Remove hydrogens.
write([format, filename, overwrite, size])	Write the molecule to a file or return a string.

```
clone_coords
Mol
addh (only_polar=False, **kwargs)
     Add hydrogens.
atom_dict
atoms
bonds
calccharges (model='gasteiger')
     Calculate partial charges for a molecule. By default the Gasteiger charge model is used.
         Parameters
             model [str (default="gasteiger")] Method for generating partial charges. Supported models:
                * gasteiger * mmff94
calcdesc(descnames=None)
     Calculate descriptor values.
     Optional parameter: descnames – a list of names of descriptors
     If descnames is not specified, all available descriptors are calculated. See the descs variable for a list of
     available descriptors.
calcfp (fptype='rdkit', opt=None)
     Calculate a molecular fingerprint.
     Optional parameters:
         fptype - the fingerprint type (default is "rdkit"). See the fps variable for a list of of available fin-
             gerprint types.
         opt – a dictionary of options for fingerprints. Currently only used for radius and bitInfo in Mor-
             gan fingerprints.
canonic order
     Returns np.array with canonic order of heavy atoms in the molecule
charges
clone
clone_coords (source)
coords
data
formula
localopt (forcefield='uff', steps=500)
     Locally optimize the coordinates.
```

**Optional parameters:** 

steps – default is 500

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If the molecule does not have any coordinates, make 3D() is called before the optimization.

forcefield – default is "uff". See the forcefields variable for a list of available forcefields.

make2D()

```
Generate 2D coordinates for molecule
      make3D (forcefield='mmff94', steps=50)
           Generate 3D coordinates.
           Optional parameters:
               forcefield – default is "uff". See the forcefields variable for a list of available forcefields.
               steps – default is 50
           Once coordinates are generated, a quick local optimization is carried out with 50 steps and the UFF force-
           field. Call localopt() if you want to improve the coordinates further.
      molwt
      num rotors
      protein
           A flag for identifing the protein molecules, for which atom_dict procedures may differ.
      removeh (**kwargs)
           Remove hydrogens.
      res_dict
      residues
      ring dict
      smiles
      sssr
      title
      write (format='smi', filename=None, overwrite=False, size=None, **kwargs)
           Write the molecule to a file or return a string.
           Optional parameters:
               format - see the informats variable for a list of available output formats (default is "smi")
               filename – default is None overwite – if the output file already exists, should it
                   be overwritten? (default is False)
           If a filename is specified, the result is written to a file. Otherwise, a string is returned containing the result.
           To write multiple molecules to the same file you should use the Outputfile class.
class oddt.toolkits.rdk.MoleculeData(Mol)
      Bases: object
      Store molecule data in a dictionary-type object
      Required parameters: Mol – an RDKit Mol
      Methods and accessor methods are like those of a dictionary except that the data is retrieved on-the-fly from the
      underlying Mol.
```

Example: >>> mol = next(readfile("sdf", 'head.sdf')) >>> data = mol.data >>> print(data) { 'Comment': 'CO-RINA 2.61 0041 25.10.2001', 'NSC': '1'} >>> print(len(data), data.keys(), data.has\_key("NSC")) 2 [ 'Comment', 'NSC'] True >>> print(data['Comment']) CORINA 2.61 0041 25.10.2001 >>> data['Comment'] = 'This is a new comment' >>> for k,v in data.items(): ... print(k, "->", v) Comment -> This is a new comment NSC -> 1 >>> del data['NSC'] >>> print(len(data), data.keys(), data.has\_key("NSC")) 1 ['Comment'] False

# **Methods**

clear()

clear	
has_key	
items	
iteritems	
keys	
to_dict	
update	
values	

```
has_key (key)
items()
iteritems()
keys()
to_dict()
update (dictionary)
values()
class oddt.toolkits.rdk.Outputfile (format, filename, overwrite=False)
Bases: object
```

Represent a file to which *output* is to be sent.

# **Required parameters:**

 $\textbf{format-see} \ \textbf{the outformats variable for a list of available} \ \ \textbf{output formats}$ 

filename

# **Optional parameters:**

**overwite – if the output file already exists, should it** be overwritten? (default is False)

**Methods:** write(molecule) close()

### **Methods**

close()	Close the Outputfile to further writing.
write(molecule)	Write a molecule to the output file.

```
close()
     Close the Outputfile to further writing.

write(molecule)
     Write a molecule to the output file.

Required parameters: molecule

class oddt.toolkits.rdk.Residue(ParentMol, atom_path, idx=0)
     Bases: object
```

Represent a RDKit residue.

Required parameter: ParentMol – Parent molecule (Mol) object path – atoms path of a residue

Attributes: atoms, idx, name.

(refer to the Open Babel library documentation for more info).

The Mol object constucted of residues' atoms can be accessed using the attribute: Residue

#### Attributes

```
atoms List of Atoms in the Residue
chain Resdiue chain ID
idx DEPRECATED: Use idx0 instead.
idx0 Internal index (0-based) of the Residue
name Residue name
```

#### atoms

List of Atoms in the Residue

number Residue number

### chain

Resdiue chain ID

#### idx

DEPRECATED: Use idx0 instead.

Internal index (0-based) of the Residue

# idx0

Internal index (0-based) of the Residue

#### name

Residue name

## number

Residue number

```
class oddt.toolkits.rdk.ResidueStack(Mol, paths)
```

Bases: object

class oddt.toolkits.rdk.Smarts(smartspattern)

Bases: object

Initialise with a SMARTS pattern.

## **Methods**

<pre>findal1(molecule[, unique])</pre>	Find all matches of the SMARTS pattern to a partic-
	ular molecule.
match(molecule)	Find all matches of the SMARTS pattern to a partic-
	ular molecule.

#### **findall** (*molecule*, *unique=True*)

Find all matches of the SMARTS pattern to a particular molecule.

# Required parameters: molecule match (molecule) Find all matches of the SMARTS pattern to a particular molecule. Required parameters: molecule oddt.toolkits.rdk.base feature factory = <rdkit.Chem.rdMolChemicalFeatures.MolChemicalFeatures. Global feature factory based on BaseFeatures.fdef oddt.toolkits.rdk.descs = ['fr\_C\_0\_noCOO', 'PEOE\_VSA3', 'Chi4v', 'fr\_Ar\_COO', 'fr\_SH', 'Chi A list of supported descriptors oddt.toolkits.rdk.diverse\_conformers\_generator(mol, $n\_conf=10$ , method='etkdg',seed=None, rmsd=0.5) Produce diverse conformers using current conformer as starting point. Each conformer is a copy of original molecule object. New in version 0.6. **Parameters** mol [oddt.toolkit.Molecule object] Molecule for which generating conformers **n conf** [int (default=10)] Targer number of conformers method [string (default='etkdg')] Method for generating conformers. Supported methods: "etkdg", "etdg", "kdg", "dg". seed [None or int (default=None)] Random seed rmsd [float (default=0.5)] The minimum RMSD that separates conformers to be ratained (otherwise, they will be pruned). Returns mols [list of oddt.toolkit.Molecule objects] Molecules with diverse conformers oddt.toolkits.rdk.forcefields = ['mmff94', 'uff'] A list of supported forcefields oddt.toolkits.rdk.fps = ['rdkit', 'layered', 'maccs', 'atompairs', 'torsions', 'morgan'] A list of supported fingerprint types oddt.toolkits.rdk.informats = {'inchi': 'InChI', 'mol': 'MDL MOL file', 'mol2': 'Tripos A dictionary of supported input formats oddt.toolkits.rdk.outformats = {'can': 'Canonical SMILES', 'inchi': 'InChI', 'inchikey': A dictionary of supported output formats oddt.toolkits.rdk.readfile(format, filename, lazy=False, opt=None, \*\*kwargs) Iterate over the molecules in a file.

### **Required parameters:**

format - see the informats variable for a list of available input formats

filename

You can access the first molecule in a file using the next() method of the iterator:

```
mol = next(readfile("smi", "myfile.smi"))
```

You can make a list of the molecules in a file using: mols = list(readfile("smi", "myfile.smi"))

You can iterate over the molecules in a file as shown in the following code snippet: >>> atomtotal = 0 >>> for mol in readfile("sdf", "head.sdf"): ... atomtotal += len(mol.atoms) ... >>> print(atomtotal) 43

oddt.toolkits.rdk.readstring(format, string, \*\*kwargs)

Read in a molecule from a string.

# **Required parameters:**

format - see the informats variable for a list of available input formats

string

Example: >>> input = "C1=CC=CS1" >>> mymol = readstring("smi", input) >>> len(mymol.atoms) 5

#### **Module contents**

# 5.1.2 Submodules

# 5.1.3 oddt.datasets module

Datasets wrapped in convenient models

class oddt.datasets.CASF(home)

Load CASF dataset as described in Li, Y. et al. Comparative Assessment of Scoring Functions on an Updated Benchmark: 2. Evaluation Methods and General Results. J. Chem. Inf. Model. 54, 1717-1736. (2014) http://dx.doi.org/10.1021/ci500081m

#### **Parameters**

**home: string** Path to CASF dataset main directory

### **Methods**

<pre>precomputed_score([scoring_function])</pre>	Load precomputed results of scoring power test for various scoring functions.
<pre>precomputed_screening([scoring_function,</pre>	Load precomputed results of screening power test for
])	various scoring functions

# precomputed\_score (scoring\_function=None)

Load precomputed results of scoring power test for various scoring functions.

### **Parameters**

**scoring\_function: string (default=None)** Name of the scoring function to get results If None, all results are returned.

precomputed\_screening(scoring\_function=None, cluster\_id=None)

Load precomputed results of screening power test for various scoring functions

### **Parameters**

**scoring\_function: string (default=None)** Name of the scoring function to get results If None, all results are returned

cluster\_id: int (default=None) Number of the protein cluster to get results If None, all results are returned

# 5.1.4 oddt.fingerprints module

Module checks interactions between two molecules and creates interacion fingerprints.

```
oddt.fingerprints.InteractionFingerprint(ligand, protein, strict=True)
```

Interaction fingerprint accomplished by converting the molecular interaction of ligand-protein into bit array according to the residue of choice and the interaction. For every residue (One row = one residue) there are eight bits which represent eight type of interactions:

- (Column 0) hydrophobic contacts
- (Column 1) aromatic face to face
- (Column 2) aromatic edge to face
- (Column 3) hydrogen bond (protein as hydrogen bond donor)
- (Column 4) hydrogen bond (protein as hydrogen bond acceptor)
- (Column 5) salt bridges (protein positively charged)
- (Column 6) salt bridges (protein negatively charged)
- (Column 7) salt bridges (ionic bond with metal ion)

## **Parameters**

**ligand, protein** [oddt.toolkit.Molecule object] Molecules, which are analysed in order to find interactions.

**strict** [bool (deafult = True)] If False, do not include condition, which informs whether atoms form 'strict' H-bond (pass all angular cutoffs).

# Returns

**InteractionFingerprint** [numpy array] Vector of calculated IFP (size = no residues \* 8 type of interaction)

oddt.fingerprints.SimpleInteractionFingerprint (ligand, protein, strict=True)

Based on http://dx.doi.org/10.1016/j.csbj.2014.05.004. Every IFP consists of 8 bits per amino acid (One row = one amino acid) and present eight type of interaction:

• (Column 0) hydrophobic contacts

- (Column 1) aromatic face to face
- (Column 2) aromatic edge to face
- (Column 3) hydrogen bond (protein as hydrogen bond donor)
- (Column 4) hydrogen bond (protein as hydrogen bond acceptor)
- (Column 5) salt bridges (protein positively charged)
- (Column 6) salt bridges (protein negatively charged)
- (Column 7) salt bridges (ionic bond with metal ion)

Returns matrix, which is sorted according to this pattern: 'ALA', 'ARG', 'ASN', 'ASP', 'CYS', 'GLN', 'GLU', 'GLY', 'HIS', 'ILE', 'LEU', 'LYS', 'MET', 'PHE', 'PRO', 'SER', 'THR', 'TRP', 'TYR', 'VAL', ''. The '' means cofactor. Index of amino acid in pattern coresponds to row in returned matrix.

### **Parameters**

**ligand, protein** [oddt.toolkit.Molecule object] Molecules, which are analysed in order to find interactions.

**strict** [bool (deafult = True)] If False, do not include condition, which informs whether atoms form 'strict' H-bond (pass all angular cutoffs).

#### Returns

**InteractionFingerprint** [numpy array] Vector of calculated IFP (size = 168)

oddt.fingerprints.**SPLIF** (*ligand*, *protein*, *depth=1*, *size=4096*, *distance\_cutoff=4.5*)

Calculates structural protein-ligand interaction fingerprint (SPLIF), based on http://pubs.acs.org/doi/abs/10. 1021/ci500319f.

# **Parameters**

**ligand, protein** [oddt.toolkit.Molecule object] Molecules, which are analysed in order to find interactions.

**depth** [int (deafult = 1)] The depth of the fingerprint, i.e. the number of bonds in Morgan algorithm. Note: For ECFP2: depth = 1, ECFP4: depth = 2, etc.

size: int (default = 4096) SPLIF is folded to given size.

**distance\_cutoff: float (default=4.5)** Cutoff distance for close contacts.

#### Returns

**SPLIF** [numpy array] Calculated SPLIF.shape = (no. of atoms, ). Every row consists of three elements:

row[0] = index of hashed atoms <math>row[1].shape = (7, 3) -> ligand's atom coords and 6 his neigbor's row[2].shape = (7, 3) -> protein's atom coords and 6 his neigbor's

oddt.fingerprints.similarity\_SPLIF(reference, query, rmsd\_cutoff=1.0)

Calculates similarity between structural interaction fingerprints, based on doi:http://pubs.acs.org/doi/abs/10.1021/ci500319f.

## **Parameters**

reference, query: numpy.array SPLIFs, which are compared in order to determine similarity.rmsd\_cutoff [int (default = 1)] Specific treshold for which, bits are considered as fully matching.

#### Returns

**SimilarityScore** [float] Similarity between given fingerprints.

oddt.fingerprints.**ECFP**(*mol*, *depth=2*, *size=4096*, *count\_bits=True*, *sparse=True*, *use pharm features=False*)

Extended connectivity fingerprints (ECFP) with an option to include atom features (FCPF). Depth of a fingerprint is counted as bond-steps, thus the depth for ECFP2 = 1, ECPF4 = 2, ECFP6 = 3, etc.

Reference: Rogers D, Hahn M. Extended-connectivity fingerprints. J Chem Inf Model. 2010;50: 742-754. http://dx.doi.org/10.1021/ci100050t

#### **Parameters**

mol [oddt.toolkit.Molecule object] Input molecule for the FP calculations

**depth** [int (deafult = 2)] The depth of the fingerprint, i.e. the number of bonds in Morgan algorithm. Note: For ECFP2: depth = 1, ECFP4: depth = 2, etc.

size [int (default = 4096)] Final size of fingerprint to which it is folded.

**count\_bits** [bool (default = True)] Should the bits be counted or unique. In dense representation it translates to integer array (count\_bits=True) or boolean array if False.

**sparse** [bool (default=True)] Should fingerprints be dense (contain all bits) or sparse (just the on bits).

**use\_pharm\_features** [bool (default=False)] Switch to use pharmacophoric features as atom representation instead of explicit atomic numbers etc.

#### Returns

**fingerprint** [numpy array] Calsulated FP of fixed size (dense) or on bits indices (sparse). Dtype is either integer or boolean.

Protein ligand extended connectivity fingerprint. For every pair of atoms in contact, compute ECFP and then hash every single, corresponding depth.

# **Parameters**

**ligand, protein** [oddt.toolkit.Molecule object] Molecules, which are analysed in order to find interactions.

**depth\_ligand, depth\_protein** [int (deafult = (2, 4))] The depth of the fingerprint, i.e. the number of bonds in Morgan algorithm. Note: For ECFP2: depth = 1, ECFP4: depth = 2, etc.

size: int (default = 16384) SPLIF is folded to given size.

distance cutoff: float (default=4.5) Cutoff distance for close contacts.

sparse [bool (default = True)] Should fingerprints be dense (contain all bits) or sparse (just the on bits).

**count\_bits** [bool (default = True)] Should the bits be counted or unique. In dense representation it translates to integer array (count\_bits=True) or boolean array if False.

**ignore\_hoh** [bool (default = True)] Should the water molecules be ignored. This is based on the name of the residue ('HOH').

### Returns

**PLEC** [numpy array] fp (size = atoms in contacts \* max(depth\_protein, depth\_ligand))

oddt.fingerprints.dice(a, b, sparse=False)

Calculates the Dice coefficient, the ratio of the bits in common to the arithmetic mean of the number of 'on' bits in the two fingerprints. Supports integer and boolean fingerprints.

#### **Parameters**

a, b [numpy array] Interaction fingerprints, which are compared in order to determine similarity.sparse [bool (default=False)] Type of FPs to use. Defaults to dense form.

#### Returns

**score** [float] Similarity between a, b.

oddt.fingerprints.tanimoto(a, b, sparse=False)

Tanimoto coefficient, supports boolean fingerprints. Integer fingerprints are casted to boolean.

#### **Parameters**

a, b [numpy array] Interaction fingerprints, which are compared in order to determine similarity.sparse [bool (default=False)] Type of FPs to use. Defaults to dense form.

### Returns

score [float] Similarity between a, b.

# 5.1.5 oddt.interactions module

Module calculates interactions between two molecules (proein-protein, protein-ligand, small-small). Currently following interacions are implemented:

- · hydrogen bonds
- · halogen bonds
- pi stacking (parallel and perpendicular)
- · salt bridges
- hydrophobic contacts
- · pi-cation
- · metal coordination
- pi-metal

oddt.interactions.close\_contacts(x, y, cutoff,  $x\_column='coords'$ ,  $y\_column='coords'$ ,  $cut-off\_low=0.0$ )

Returns pairs of atoms which are within close contac distance cutoff. The cutoff is semi-inclusive, i.e (cutoff low, cutoff].

# **Parameters**

**x, y** [atom\_dict-type numpy array] Atom dictionaries generated by oddt.toolkit.Molecule objects.

cutoff [float] Cutoff distance for close contacts

**x\_column, ycolumn** [string, (default='coords')] Column containing coordinates of atoms (or pseudo-atoms, i.e. ring centroids)

**cutoff\_low** [float (default=0.)] Lower bound of contacts to find (exclusive). Zero by default. .. versionadded:: 0.6

#### Returns

**x\_, y\_** [atom\_dict-type numpy array] Aligned pairs of atoms in close contact for further processing.

oddt.interactions.hbond\_acceptor\_donor(mol1, mol2, cutoff=3.5, base\_angle=120, toler-ance=30)

Returns pairs of acceptor-donor atoms, which meet H-bond criteria

#### **Parameters**

**mol1, mol2** [oddt.toolkit.Molecule object] Molecules to compute H-bond acceptor and H-bond donor pairs

cutoff [float, (default=3.5)] Distance cutoff for A-D pairs

**base\_angle** [int, (default=120)] Base angle determining allowed direction of hydrogen bond formation, which is devided by the number of neighbors of acceptor atom to establish final directional angle

**tolerance** [int, (default=30)] Range (+/- tolerance) from perfect direction (base\_angle/n\_neighbors) in which H-bonds are considered as strict.

#### Returns

**a, d** [atom\_dict-type numpy array] Aligned arrays of atoms forming H-bond, firstly acceptors, secondly donors.

**strict** [numpy array, dtype=bool] Boolean array align with atom pairs, informing whether atoms form 'strict' H-bond (pass all angular cutoffs). If false, only distance cutoff is met, therefore the bond is 'crude'.

oddt.interactions.hbonds (mol1, mol2, \*args, \*\*kwargs)
Calculates H-bonds between molecules

### **Parameters**

mol1, mol2 [oddt.toolkit.Molecule object] Molecules to compute H-bond acceptor and H-bond donor pairs

cutoff [float, (default=3.5)] Distance cutoff for A-D pairs

**base\_angle** [int, (default=120)] Base angle determining allowed direction of hydrogen bond formation, which is devided by the number of neighbors of acceptor atom to establish final directional angle

**tolerance** [int, (default=30)] Range (+/- tolerance) from perfect direction (base\_angle/n\_neighbors) in which H-bonds are considered as strict.

### Returns

mol1\_atoms, mol2\_atoms [atom\_dict-type numpy array] Aligned arrays of atoms forming H-bond

**strict** [numpy array, dtype=bool] Boolean array align with atom pairs, informing whether atoms form 'strict' H-bond (pass all angular cutoffs). If false, only distance cutoff is met, therefore the bond is 'crude'.

oddt.interactions.halogenbond\_acceptor\_halogen (mol1, mol2, base\_angle\_acceptor=120, base\_angle\_halogen=180, toler-ance=30, cutoff=4)

Returns pairs of acceptor-halogen atoms, which meet halogen bond criteria

#### **Parameters**

mol1, mol2 [oddt.toolkit.Molecule object] Molecules to compute halogen bond acceptor and halogen pairs

cutoff [float, (default=4)] Distance cutoff for A-H pairs

- **base\_angle\_acceptor** [int, (default=120)] Base angle determining allowed direction of halogen bond formation, which is devided by the number of neighbors of acceptor atom to establish final directional angle
- **base\_angle\_halogen** [int (default=180)] Ideal base angle between halogen bond and halogen-neighbor bond
- **tolerance** [int, (default=30)] Range (+/- tolerance) from perfect direction (base angle/n neighbors) in which halogen bonds are considered as strict.

### Returns

- **a, h** [atom\_dict-type numpy array] Aligned arrays of atoms forming halogen bond, firstly acceptors, secondly halogens
- **strict** [numpy array, dtype=bool] Boolean array align with atom pairs, informing whether atoms form 'strict' halogen bond (pass all angular cutoffs). If false, only distance cutoff is met, therefore the bond is 'crude'.

oddt.interactions.halogenbonds(mol1, mol2, \*\*kwargs)

Calculates halogen bonds between molecules

#### **Parameters**

- **mol1, mol2** [oddt.toolkit.Molecule object] Molecules to compute halogen bond acceptor and halogen pairs
- cutoff [float, (default=4)] Distance cutoff for A-H pairs
- **base\_angle\_acceptor** [int, (default=120)] Base angle determining allowed direction of halogen bond formation, which is devided by the number of neighbors of acceptor atom to establish final directional angle
- **base\_angle\_halogen** [int (default=180)] Ideal base angle between halogen bond and halogenneighbor bond
- **tolerance** [int, (default=30)] Range (+/- tolerance) from perfect direction (base\_angle/n\_neighbors) in which halogen bonds are considered as strict.

### Returns

- mol1\_atoms, mol2\_atoms [atom\_dict-type numpy array] Aligned arrays of atoms forming halogen bond
- **strict** [numpy array, dtype=bool] Boolean array align with atom pairs, informing whether atoms form 'strict' halogen bond (pass all angular cutoffs). If false, only distance cutoff is met, therefore the bond is 'crude'.

oddt.interactions.pi\_stacking (mol1, mol2, cutoff=5, tolerance=30)
Returns pairs of rings, which meet pi stacking criteria

### **Parameters**

- mol1, mol2 [oddt.toolkit.Molecule object] Molecules to compute ring pairs
- cutoff [float, (default=5)] Distance cutoff for Pi-stacking pairs
- **tolerance** [int, (default=30)] Range (+/- tolerance) from perfect direction (parallel or perpendicular) in which pi-stackings are considered as strict.

# Returns

r1, r2 [ring\_dict-type numpy array] Aligned arrays of rings forming pi-stacking

**strict\_parallel** [numpy array, dtype=bool] Boolean array align with ring pairs, informing whether rings form 'strict' parallel pi-stacking. If false, only distance cutoff is met, therefore the stacking is 'crude'.

**strict\_perpendicular** [numpy array, dtype=bool] Boolean array align with ring pairs, informing whether rings form 'strict' perpendicular pi-stacking (T-shaped, T-face, etc.). If false, only distance cutoff is met, therefore the stacking is 'crude'.

oddt.interactions.salt\_bridge\_plus\_minus(mol1, mol2, cutoff=4)

Returns pairs of plus-mins atoms, which meet salt bridge criteria

#### **Parameters**

mol1, mol2 [oddt.toolkit.Molecule object] Molecules to compute plus and minus pairs cutoff [float, (default=4)] Distance cutoff for A-H pairs

### **Returns**

**plus, minus** [atom\_dict-type numpy array] Aligned arrays of atoms forming salt bridge, firstly plus, secondly minus

oddt.interactions.salt\_bridges (mol1, mol2, \*args, \*\*kwargs)
Calculates salt bridges between molecules

#### **Parameters**

**mol1, mol2** [oddt.toolkit.Molecule object] Molecules to compute plus and minus pairs **cutoff** [float, (default=4)] Distance cutoff for plus-minus pairs

#### Returns

**mol1\_atoms**, **mol2\_atoms** [atom\_dict-type numpy array] Aligned arrays of atoms forming salt bridges

oddt.interactions.hydrophobic\_contacts(mol1, mol2, cutoff=4)

Calculates hydrophobic contacts between molecules

### **Parameters**

mol1, mol2 [oddt.toolkit.Molecule object] Molecules to compute hydrophobe pairs
cutoff [float, (default=4)] Distance cutoff for hydrophobe pairs

### Returns

**mol1\_atoms**, **mol2\_atoms** [atom\_dict-type numpy array] Aligned arrays of atoms forming hydrophobic contacts

oddt.interactions.pi cation (mol1, mol2, cutoff=5, tolerance=30)

Returns pairs of ring-cation atoms, which meet pi-cation criteria

### **Parameters**

**mol1, mol2** [oddt.toolkit.Molecule object] Molecules to compute ring-cation pairs **cutoff** [float, (default=5)] Distance cutoff for Pi-cation pairs

**tolerance** [int, (default=30)] Range (+/- tolerance) from perfect direction (perpendicular) in which pi-cation are considered as strict.

# Returns

r1 [ring\_dict-type numpy array] Aligned rings forming pi-stackingplus2 [atom\_dict-type numpy array] Aligned cations forming pi-cation

**strict\_parallel** [numpy array, dtype=bool] Boolean array align with ring-cation pairs, informing whether they form 'strict' pi-cation. If false, only distance cutoff is met, therefore the interaction is 'crude'.

oddt.interactions.acceptor\_metal (mol1, mol2, base\_angle=120, tolerance=30, cutoff=4)
Returns pairs of acceptor-metal atoms, which meet metal coordination criteria Note: This function is directional (mol1 holds acceptors, mol2 holds metals)

#### **Parameters**

mol1, mol2 [oddt.toolkit.Molecule object] Molecules to compute acceptor and metal pairs cutoff [float, (default=4)] Distance cutoff for A-M pairs

**base\_angle** [int, (default=120)] Base angle determining allowed direction of metal coordination, which is devided by the number of neighbors of acceptor atom to establish final directional angle

**tolerance** [int, (default=30)] Range (+/- tolerance) from perfect direction (base\_angle/n\_neighbors) in metal coordination are considered as strict.

#### Returns

**a, d** [atom\_dict-type numpy array] Aligned arrays of atoms forming metal coordination, firstly acceptors, secondly metals.

**strict** [numpy array, dtype=bool] Boolean array align with atom pairs, informing whether atoms form 'strict' metal coordination (pass all angular cutoffs). If false, only distance cutoff is met, therefore the interaction is 'crude'.

oddt.interactions.pi\_metal (mol1, mol2, cutoff=5, tolerance=30)
Returns pairs of ring-metal atoms, which meet pi-metal criteria

### **Parameters**

mol1, mol2 [oddt.toolkit.Molecule object] Molecules to compute ring-metal pairs

cutoff [float, (default=5)] Distance cutoff for Pi-metal pairs

**tolerance** [int, (default=30)] Range (+/- tolerance) from perfect direction (perpendicular) in which pi-metal are considered as strict.

# Returns

- **r1** [ring dict-type numpy array] Aligned rings forming pi-metal
- m [atom\_dict-type numpy array] Aligned metals forming pi-metal

**strict\_parallel** [numpy array, dtype=bool] Boolean array align with ring-metal pairs, informing whether they form 'strict' pi-metal. If false, only distance cutoff is met, therefore the interaction is 'crude'.

# 5.1.6 oddt.metrics module

Metrics for estimating performance of drug discovery methods implemented in ODDT

oddt.metrics.roc(y\_true, y\_score, pos\_label=None, sample\_weight=None, drop\_intermediate=True)
Compute Receiver operating characteristic (ROC)

Note: this implementation is restricted to the binary classification task.

Read more in the User Guide.

# Parameters

- **y\_true** [array, shape =  $[n_samples]$ ] True binary labels in range  $\{0, 1\}$  or  $\{-1, 1\}$ . If labels are not binary, pos\_label should be explicitly given.
- y\_score [array, shape = [n\_samples]] Target scores, can either be probability estimates of the positive class, confidence values, or non-thresholded measure of decisions (as returned by "decision\_function" on some classifiers).
- **pos\_label** [int or str, default=None] Label considered as positive and others are considered negative.
- **sample\_weight** [array-like of shape = [n\_samples], optional] Sample weights.
- **drop\_intermediate** [boolean, optional (default=True)] Whether to drop some suboptimal thresholds which would not appear on a plotted ROC curve. This is useful in order to create lighter ROC curves.

New in version 0.17: parameter *drop\_intermediate*.

### Returns

- **fpr** [array, shape = [>2]] Increasing false positive rates such that element i is the false positive rate of predictions with score >= thresholds[i].
- **tpr** [array, shape = [>2]] Increasing true positive rates such that element i is the true positive rate of predictions with score >= thresholds[i].
- **thresholds** [array, shape =  $[n_{thresholds}]$ ] Decreasing thresholds on the decision function used to compute fpr and tpr. thresholds[0] represents no instances being predicted and is arbitrarily set to  $max(y\_score) + 1$ .

#### See also:

roc\_auc\_score Compute the area under the ROC curve

# **Notes**

Since the thresholds are sorted from low to high values, they are reversed upon returning them to ensure they correspond to both fpr and tpr, which are sorted in reversed order during their calculation.

### References

[1]

# **Examples**

```
>>> import numpy as np
>>> from sklearn import metrics
>>> y = np.array([1, 1, 2, 2])
>>> scores = np.array([0.1, 0.4, 0.35, 0.8])
>>> fpr, tpr, thresholds = metrics.roc_curve(y, scores, pos_label=2)
>>> fpr
array([ 0. ,  0.5,  0.5,  1. ])
>>> tpr
array([ 0.5,  0.5,  1. ,  1. ])
>>> thresholds
array([ 0.8 ,  0.4 ,  0.35,  0.1 ])
```

```
oddt.metrics.auc(x, y, reorder=False)
```

Compute Area Under the Curve (AUC) using the trapezoidal rule

This is a general function, given points on a curve. For computing the area under the ROC-curve, see roc\_auc\_score(). For an alternative way to summarize a precision-recall curve, see average\_precision\_score().

#### **Parameters**

- $\mathbf{x}$  [array, shape = [n]] x coordinates.
- y [array, shape = [n]] y coordinates.

**reorder** [boolean, optional (default=False)] If True, assume that the curve is ascending in the case of ties, as for an ROC curve. If the curve is non-ascending, the result will be wrong.

#### Returns

auc [float]

See also:

roc\_auc\_score Compute the area under the ROC curve

average\_precision\_score Compute average precision from prediction scores

precision\_recall\_curve Compute precision-recall pairs for different probability thresholds

# **Examples**

```
>>> import numpy as np
>>> from sklearn import metrics
>>> y = np.array([1, 1, 2, 2])
>>> pred = np.array([0.1, 0.4, 0.35, 0.8])
>>> fpr, tpr, thresholds = metrics.roc_curve(y, pred, pos_label=2)
>>> metrics.auc(fpr, tpr)
0.75
```

oddt.metrics.roc\_auc(y\_true, y\_score, pos\_label=None, ascending\_score=True)
Computes ROC AUC score

### **Parameters**

- **y\_true** [array, shape=[n\_samples]] True binary labels, in range {0,1} or {-1,1}. If positive label is different than 1, it must be explicitly defined.
- **y\_score** [array, shape=[n\_samples]] Scores for tested series of samples

**pos\_label: int** Positive label of samples (if other than 1)

**ascending\_score: bool (default=True)** Indicates if your score is ascendig. Ascending score icreases with deacreasing activity. In other words it ascends on ranking list (where actives are on top).

# Returns

```
roc_auc [float] ROC AUC in range 0:1
```

```
oddt.metrics.roc_log_auc(y_true, y_score, pos_label=None, ascending_score=True, log_min=0.001, log_max=1.0)

Computes area under semi-log ROC.
```

### **Parameters**

- **y\_true** [array, shape=[n\_samples]] True binary labels, in range {0,1} or {-1,1}. If positive label is different than 1, it must be explicitly defined.
- **y\_score** [array, shape=[n\_samples]] Scores for tested series of samples
- pos\_label: int Positive label of samples (if other than 1)
- **ascending\_score: bool (default=True)** Indicates if your score is ascendig. Ascending score icreases with deacreasing activity. In other words it ascends on ranking list (where actives are on top).
- **log\_min** [float (default=0.001)] Minimum value for estimating AUC. Lower values will be clipped for numerical stability.
- log\_max [float (default=1.)] Maximum value for estimating AUC. Higher values will be ignored.

### Returns

auc [float] semi-log ROC AUC

oddt.metrics.enrichment\_factor (y\_true, y\_score, percentage=1, pos\_label=None, kind='fold')
Computes enrichment factor for given percentage, i.e. EF\_1% is enrichment factor for first percent of given samples. This function assumes that results are already sorted and samples with best predictions are first.

#### **Parameters**

- **y\_true** [array, shape=[n\_samples]] True binary labels, in range {0,1} or {-1,1}. If positive label is different than 1, it must be explicitly defined.
- **y\_score** [array, shape=[n\_samples]] Scores for tested series of samples
- percentage [int or float] The percentage for which EF is being calculated
- **pos\_label: int** Positive label of samples (if other than 1)
- **kind:** 'fold' or 'percentage' (default='fold') Two kinds of enrichment factor: fold and percentage. Fold shows the increase over random distribution (1 is random, the higher EF the better enrichment). Percentage returns the fraction of positive labels within the top x% of dataset.

### Returns

ef [float] Enrichment Factor for given percenage in range 0:1

oddt.metrics.random\_roc\_log\_auc (log\_min=0.001, log\_max=1.0) Computes area under semi-log ROC for random distribution.

#### **Parameters**

**log\_min** [float (default=0.001)] Minimum logarithm value for estimating AUC

log\_max [float (default=1.)] Maximum logarithm value for estimating AUC.

### Returns

auc [float] semi-log ROC AUC for random distribution

oddt.metrics.**rmse** $(y\_true, y\_pred)$ 

Compute Root Mean Squared Error (RMSE)

# **Parameters**

- **y\_true** [array-like of shape = [n\_samples] or [n\_samples, n\_outputs]] Ground truth (correct) target values.
- y pred [array-like of shape = [n samples] or [n samples, n outputs]] Estimated target values.

### Returns

rmse [float] A positive floating point value (the best value is 0.0).

```
oddt.metrics.rie(y_true, y_score, alpha=20, pos_label=None)
```

Computes Robust Initial Enhancement [1]. This function assumes that results are already sorted and samples with best predictions are first.

#### **Parameters**

```
y_true [array, shape=[n_samples]] True binary labels, in range {0,1} or {-1,1}. If positive label is different than 1, it must be explicitly defined.
```

**y\_score** [array, shape=[n\_samples]] Scores for tested series of samples

alpha: float Alpha. 1/Alpha should be proportional to the percentage in EF.

**pos\_label: int** Positive label of samples (if other than 1)

#### Returns

rie\_score [float] Robust Initial Enhancement

### References

[1]

```
oddt.metrics.bedroc(y_true, y_score, alpha=20.0, pos_label=None)
```

Computes Boltzmann-Enhanced Discrimination of Receiver Operating Characteristic [1]. This function assumes that results are already sorted and samples with best predictions are first.

#### **Parameters**

**y\_true** [array, shape=[n\_samples]] True binary labels, in range {0,1} or {-1,1}. If positive label is different than 1, it must be explicitly defined.

y\_score [array, shape=[n\_samples]] Scores for tested series of samples

alpha: float Alpha. 1/Alpha should be proportional to the percentage in EF.

**pos\_label: int** Positive label of samples (if other than 1)

## Returns

**bedroc\_score** [float] Boltzmann-Enhanced Discrimination of Receiver Operating Characteristic

# References

[1]

# 5.1.7 oddt.pandas module

Pandas extension for chemical analysis

```
 \begin{array}{c} \textbf{class} \text{ oddt.pandas.ChemDataFrame} \ (\textit{data=None}, \quad \textit{index=None}, \quad \textit{columns=None}, \quad \textit{dtype=None}, \\ \textit{copy=False}) \\ \text{Bases: pandas.core.frame.DataFrame} \end{array}
```

Chemical DataFrame object, which contains molecules column of *oddt.toolkit.Molecule* objects. Rich display of molecules (2D) is available in iPython Notebook. Additional *to\_sdf* and *to\_mol2* methods make writing to molecular formats easy.

New in version 0.3.

### **Notes**

Thanks to: http://blog.snapdragon.cc/2015/05/05/subclass-pandas-dataframe-to-save-custom-attributes/

### **Attributes**

- **T** Transpose index and columns.
- at Access a single value for a row/column label pair.
- **axes** Return a list representing the axes of the DataFrame.
- **blocks** Internal property, property synonym for as\_blocks()
- columns The column labels of the DataFrame.
- **dtypes** Return the dtypes in the DataFrame.
- empty Indicator whether DataFrame is empty.
- ftypes Return the ftypes (indication of sparse/dense and dtype) in DataFrame.
- iat Access a single value for a row/column pair by integer position.
- iloc Purely integer-location based indexing for selection by position.
- index The index (row labels) of the DataFrame.

# is\_copy

- ix A primarily label-location based indexer, with integer position fallback.
- **loc** Access a group of rows and columns by label(s) or a boolean array.
- ndim Return an int representing the number of axes / array dimensions.
- **shape** Return a tuple representing the dimensionality of the DataFrame.
- **size** Return an int representing the number of elements in this object.
- **style** Property returning a Styler object containing methods for building a styled HTML representation fo the DataFrame.
- **values** Return a Numpy representation of the DataFrame.

### **Methods**

abs()	Return a Series/DataFrame with absolute numeric value of each element.
add(other[, axis, level, fill_value])	Addition of dataframe and other, element-wise (binary operator <i>add</i> ).
add_prefix(prefix)	Prefix labels with string <i>prefix</i> .
add_suffix(suffix)	Suffix labels with string <i>suffix</i> .

Continued on next page

Table 19 – continue	d from previous page
agg(func[, axis])	Aggregate using one or more operations over the specified axis.
aggregate(func[, axis])	Aggregate using one or more operations over the specified axis.
align(other[, join, axis, level, copy,])	Align two objects on their axes with the specified join method for each axis Index
all([axis, bool_only, skipna, level])	Return whether all elements are True, potentially over an axis.
any([axis, bool_only, skipna, level])	Return whether any element is True over requested axis.
append(other[, ignore_index,])	Append rows of <i>other</i> to the end of this frame, returning a new object.
apply(func[, axis, broadcast, raw, reduce,])	Apply a function along an axis of the DataFrame.
applymap(func)	Apply a function to a Dataframe elementwise.
as_blocks([copy])	Convert the frame to a dict of dtype -> Constructor
_ (173)	Types that each has a homogeneous dtype.
as_matrix([columns])	Convert the frame to its Numpy-array representation.
asfreq(freq[, method, how, normalize,])	Convert TimeSeries to specified frequency.
asof(where[, subset])	The last row without any NaN is taken (or the last
	row without NaN considering only the subset of
	columns in the case of a DataFrame)
assign(**kwargs)	Assign new columns to a DataFrame, returning a
	new object (a copy) with the new columns added to
	the original ones.
astype(**kwargs)	Cast a pandas object to a specified dtype dtype.
<pre>at_time(time[, asof])</pre>	Select values at particular time of day (e.g.
between_time(start_time, end_time[,])	Select values between particular times of the day
	(e.g., 9:00-9:30 AM).
bfill([axis, inplace, limit, downcast])	Synonym for DataFrame.
	fillna(method='bfill')
bool()	Return the bool of a single element PandasObject.
boxplot([column, by, ax, fontsize, rot,])	Make a box plot from DataFrame columns.
clip([lower, upper, axis, inplace])	Trim values at input threshold(s).
<pre>clip_lower(threshold[, axis, inplace])</pre>	Return copy of the input with values below a thresh-
	old truncated.
<pre>clip_upper(threshold[, axis, inplace])</pre>	Return copy of input with values above given value(s) truncated.
combine(other, func[, fill_value, overwrite])	Add two DataFrame objects and do not propagate
	NaN values, so if for a (column, time) one frame is
	missing a value, it will default to the other frame's
	value (which might be NaN as well)
combine_first(other)	Combine two DataFrame objects and default to non-
	null values in frame calling the method.
compound([axis, skipna, level])	Return the compound percentage of the values for
	the requested axis
consolidate([inplace])	Compute NDFrame with "consolidated" internals
	(data of each dtype grouped together in a single ndar-
	ray).
<pre>convert_objects([convert_dates,])</pre>	Attempt to infer better dtype for object columns.
copy([deep])	Make a copy of this object's indices and data.
	Continued on next page

	ed from previous page
corr([method, min_periods])	Compute pairwise correlation of columns, excluding NA/null values
corrwith(other[, axis, drop])	Compute pairwise correlation between rows or
	columns of two DataFrame objects.
count([axis, level, numeric_only])	Count non-NA cells for each column or row.
cov([min_periods])	Compute pairwise covariance of columns, excluding
\L _1 3/	NA/null values.
cummax([axis, skipna])	Return cumulative maximum over a DataFrame or
1 1/	Series axis.
cummin([axis, skipna])	Return cumulative minimum over a DataFrame or Series axis.
cumprod([axis, skipna])	Return cumulative product over a DataFrame or Series axis.
cumsum([axis, skipna])	Return cumulative sum over a DataFrame or Series axis.
describe([percentiles, include, exclude])	Generates descriptive statistics that summarize the central tendency, dispersion and shape of a dataset's distribution, excluding NaN values.
diff([periods, axis])	First discrete difference of element.
div(other[, axis, level, fill_value])	Floating division of dataframe and other, elementwise (binary operator <i>truediv</i> ).
divide(other[, axis, level, fill_value])	Floating division of dataframe and other, elementwise (binary operator <i>truediv</i> ).
dot(other)	Matrix multiplication with DataFrame or Series objects.
drop([labels, axis, index, columns, level,])	Drop specified labels from rows or columns.
drop_duplicates([subset, keep, inplace])	Return DataFrame with duplicate rows removed, op-
arep_aapricases([sueses, meep, mpmee])	tionally only considering certain columns
dropna([axis, how, thresh, subset, inplace])	Remove missing values.
duplicated([subset, keep])	Return boolean Series denoting duplicate rows, op-
1	tionally only considering certain columns
eq(other[, axis, level])	Wrapper for flexible comparison methods eq
equals(other)	Determines if two NDFrame objects contain the
	same elements.
eval(expr[, inplace])	Evaluate a string describing operations on
( 1 1 1 3)	DataFrame columns.
ewm([com, span, halflife, alpha,])	Provides exponential weighted functions
expanding([min_periods, center, axis])	Provides expanding transformations.
ffill([axis, inplace, limit, downcast])	Synonym for DataFrame. fillna(method='ffill')
fillna([value, method, axis, inplace,])	Fill NA/NaN values using the specified method
filter([items, like, regex, axis])	Subset rows or columns of dataframe according to
	labels in the specified index.
first(offset)	Convenience method for subsetting initial periods of
()	time series data based on a date offset.
first_valid_index()	Return index for first non-NA/null value.
floordiv(other[, axis, level, fill_value])	Integer division of dataframe and other, element-
	wise (binary operator <i>floordiv</i> ).
from_csv(path[, header, sep, index_col,])	Read CSV file.
from_dict(data[, orient, dtype, columns])	Construct DataFrame from dict of array-like or dicts.
from_items(items[, columns, orient])	Construct a dataframe from a list of tuples
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Table 19 - continued from previous page

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<pre>from_records(data[, index, exclude,])</pre>	Convert structured or record ndarray to DataFrame
ge(other[, axis, level])	Wrapper for flexible comparison methods ge
get(key[, default])	Get item from object for given key (DataFrame col-
	umn, Panel slice, etc.).
get_dtype_counts()	Return counts of unique dtypes in this object.
get_ftype_counts()	Return counts of unique ftypes in this object.
<pre>get_value(index, col[, takeable])</pre>	Quickly retrieve single value at passed column and
	index
get_values()	Return an ndarray after converting sparse values to
	dense.
groupby([by, axis, level, as_index, sort,])	Group series using mapper (dict or key function, ap-
	ply given function to group, return result as series) or
	by a series of columns.
gt(other[, axis, level])	Wrapper for flexible comparison methods gt
head([n])	Return the first <i>n</i> rows.
hist([column, by, grid, xlabelsize, xrot,])	Make a histogram of the DataFrame's.
idxmax([axis, skipna])	Return index of first occurrence of maximum over
• •	requested axis.
idxmin([axis, skipna])	Return index of first occurrence of minimum over re-
•	quested axis.
infer_objects()	Attempt to infer better dtypes for object columns.
<pre>info([verbose, buf, max_cols, memory_usage,</pre>	Print a concise summary of a DataFrame.
])	•
<pre>insert(loc, column, value[, allow_duplicates])</pre>	Insert column into DataFrame at specified location.
<pre>interpolate([method, axis, limit, inplace,])</pre>	Interpolate values according to different methods.
isin(values)	Return boolean DataFrame showing whether each el-
	ement in the DataFrame is contained in values.
isna()	Detect missing values.
isnull()	Detect missing values.
items()	Iterator over (column name, Series) pairs.
iteritems()	Iterator over (column name, Series) pairs.
iterrows()	Iterate over DataFrame rows as (index, Series) pairs.
<pre>itertuples([index, name])</pre>	Iterate over DataFrame rows as namedtuples, with
	index value as first element of the tuple.
<pre>join(other[, on, how, lsuffix, rsuffix, sort])</pre>	Join columns with other DataFrame either on index
J = ( <u>-</u>	or on a key column.
keys()	Get the 'info axis' (see Indexing for more)
kurt([axis, skipna, level, numeric_only])	Return unbiased kurtosis over requested axis using
(fr 1) - 1 - 1) - 1 - 1 - 1 - 1 - 1	Fisher's definition of kurtosis (kurtosis of normal ==
	0.0).
kurtosis([axis, skipna, level, numeric_only])	Return unbiased kurtosis over requested axis using
	Fisher's definition of kurtosis (kurtosis of normal ==
	0.0).
last(offset)	Convenience method for subsetting final periods of
	time series data based on a date offset.
last_valid_index()	Return index for last non-NA/null value.
le(other[, axis, level])	Wrapper for flexible comparison methods le
lookup(row_labels, col_labels)	Label-based "fancy indexing" function for
	DataFrame.
lt(other[, axis, level])	Wrapper for flexible comparison methods lt
C L) ** ** ** 1/	Continued on next page
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the requested axis  Return an object of same shape as self and whose corresponding entries are from self where cond is False and otherwise are from other.  Max([axis, skipna, level, numeric_only])  Median([axis, skipna, level, numeric_only])  Merge(fight[, how, on, left_on, right_on,])  Merge DataFrame objects by performing a database style join operation by columns or indexes.  Min([axis, skipna, level, numeric_only])  Modulo of dataframe and other, element-wise (binary operator mod).  Modulo of dataframe and other, element-wise (binary operator mod).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise (binary operator mul).  Multiplication of dataframe and other, element-wise operator mul).  Multiplication of dataframe and other, element-wise operator mul).  Multiplication of dataframe and other, element-wise operator mul.  Multiplication of dataframe and other, element-wise		d from previous page
mask(cond[, other, inplace, axis, level,])         Return an object of same shape as self where cond is False and otherwise are from other.           max([axis, skipna, level, numeric_only])         This method returns the maximum of the values in the object.           mean([axis, skipna, level, numeric_only])         Return the mean of the values for the requested axis median([axis, skipna, level, numeric_only])           melta([id_vars, value_vars, var_name,])         "Unpivots" a DataFrame from wide format to long format, optionally leaving identifier variables set.           memory_usage([index, deep])         Return the median of the values for the requested axis           merge(rightl, how, on, left_on, right_on,])         Return the memory usage of each column in bytes.           merge(rightl, how, on, left_on, right_on,])         Merge DataFrame objects by performing a database-style join operation by columns or indexes.           min([axis, skipna, level, numeric_only])         This method returns the minimum of the values in the object.           mod(otherf, axis, level, fill_value])         Modulo of dataframe and other, element-wise (binary operator mod).           mode([axis, numeric_only])         Gets the mode(s) of each element along the axis selected.           mul(otherf, axis, level, fill_value])         Multiplication of dataframe and other, element-wise (binary operator mul).           multiply (otherf, axis, level, fill_value])         Wultiplication of dataframe and other, element-wise (binary operator mul).           netotherf, axis, lev	mad([axis, skipna, level])	Return the mean absolute deviation of the values for
corresponding entries are from self where cond is False and otherwise are from other.  max([axis, skipna, level, numeric_only])  mean([axis, skipna, level, numeric_only])  mean([axis, skipna, level, numeric_only])  median([axis, skipna, level, numeric_only])  median([axis, skipna, level, numeric_only])  metit([id_vars, value_vars, var_name,])  metit([id_vars, value_vars, var_name,])  memory_usage([index, deep])  metit([id_vars, value_vars, var_name,])  metit([id_vars, value_vars, value_var_name,])  metit([id_vars, value_var_name,])  metit([id_var_name,])  metit([id_var_name,])  metit([id_var_name,])  metit([id_var_name,])  metit([id_var_name,		1
### False and otherwise are from other.  #### This method returns the maximum of the values in the object.  #### This method returns the maximum of the values in the object.  ##### This method returns the maximum of the values for the requested axis median([axis, skipna, level, numeric_only])  ##### Return the mean of the values for the requested axis median([axis, skipna, level, numeric_only])  ########## This method returns the maximum of the values for the requested axis median of the values for the requested axis memory_usage([index, deep])  ###################################	<pre>mask(cond[, other, inplace, axis, level,])</pre>	
max([axis, skipna, level, numeric_only])         This method returns the maximum of the values in the object.           mean([axis, skipna, level, numeric_only])         Return the mean of the values for the requested axis           median([axis, skipna, level, numeric_only])         Return the mean of the values for the requested axis           melt([id_vars, value_vars, var_name,])         "Unpivots" a DataFrame from wide format to long format, optionally leaving identifier variables set.           memory_usage([index, deep])         Return the memory usage of each column in bytes.           merge(right[, how, on, left_on, right_on,])         Merge DataFrame objects by performing a database-style join operation by columns or indexes.           min([axis, skipna, level, numeric_only])         Merge DataFrame objects by performing a database-style join operation by columns or indexes.           min([axis, skipna, level, nill_value])         Modulo of dataframe and other, element-wise (binary operator mod).           mode([axis, numeric_only])         Gets the mode(s) of each element along the axis selected.           mull(other[, axis, level, fill_value])         Multiplication of dataframe and other, element-wise (binary operator mul).           multiply(other[, axis, level, fill_value])         Multiplication of dataframe and other, element-wise (binary operator mul).           ne(other[, axis, level], fill_value])         Wrapper for flexible comparison methods ne           Return the first n rows ordered by columns in descending order.		corresponding entries are from self where cond is
the object.  Return the mean of the values for the requested axis  median([axis, skipna, level, numeric_only]) Return the median of the values for the requested axis  melt([id_vars, value_vars, var_name,])  "Unpivots" a DataFrame from wide format to long format, optionally leaving identifier variables set.  memory_usage([index, deep]) Return the memory usage of each column in bytes.  merge(right[, how, on, left_on, right_on,])  merge(right[, how, on, left_on, right_on,])  mode([axis, skipna, level, numeric_only])  mode([axis, skipna, level, numeric_only])  mode([axis, skipna, level, fill_value])  mode([axis, numeric_only])  Modulo of dataframe and other, element-wise (binary operator mod).  mode([axis, numeric_only])  Multiplication of dataframe and other, element-wise (binary operator mul).  multiply (other[, axis, level, fill_value])  Multiplication of dataframe and other, element-wise (binary operator mul).  me(other[, axis, level, fill_value])  Ne(other[, axis, level, fill_value])  Ne(other[, axis, level, fill_value])  Ne(other[, axis, level, fill_value])  Neturn the median of the values for the requested axis  multiply (other[, axis, level, fill_value])  Multiplication of dataframe and other, element-wise (binary operator mul).  Neturn the median of the values for the requested axis  not null()  Detect existing (non-missing) values.  Detect existing (non-missing) values.  Detect existing (non-missing) values.  Detect existing (non-missing) values.  Percentage change between the current and a prior element.  pipe(func, *args, **kwargs)  pivot([index, columns, values])  Return reshaped DataFrame organized by given index / column values.  pivot_table([values, index, columns,])  Precentage change between the current and a prior element.  pow(other[, axis, level, fill_value])  Return tern and drop from frame.  Exponential power of dataframe and other, element-wise (binary operator pow).  Procd([axis, skipna, level, numeric_only,])  Return the median of the values for the requested axis		False and otherwise are from <i>other</i> .
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Pivot([index, columns, values])       Return reshaped DataFrame organized by given index / column values.         Pivot_table([values, index, columns,])       Create a spreadsheet-style pivot table as a DataFrame.         Plot       alias of pandas.plottingcore. FramePlotMethods         Pop(item)       Return item and drop from frame.         Exponential power of dataframe and other, elementwise (binary operator pow).         Prod([axis, skipna, level, numeric_only,])       Return the product of the values for the requested axis	-	
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DataFrame.       plot     alias of pandas.plottingcore. FramePlotMethods       pop(item)     Return item and drop from frame.       pow(other[, axis, level, fill_value])     Exponential power of dataframe and other, elementwise (binary operator pow).       prod([axis, skipna, level, numeric_only,])     Return the product of the values for the requested axis	<pre>pivot_table([values, index, columns,])</pre>	
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axis	prod([axis, skipna, level, numeric_only,])	
product([axis, skipna, level, numeric only]) Return the product of the values for the requested		
return the product of the values for the requested	<pre>product([axis, skipna, level, numeric_only,])</pre>	Return the product of the values for the requested
axis	-	
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Table 19 – continue	d from previous page
quantile([q, axis, numeric_only, interpolation])	Return values at the given quantile over requested
	axis, a la numpy.percentile.
query(expr[, inplace])	Query the columns of a frame with a boolean expres-
1 · · · · · · · · · · · · · · · · · · ·	sion.
radd(other[, axis, level, fill_value])	Addition of dataframe and other, element-wise (bi-
rada(onier[, axis, ievei, im_varae])	nary operator <i>radd</i> ).
rank([axis, method, numeric_only,])	Compute numerical data ranks (1 through n) along
Tank([axis, method, numeric_omy,])	axis.
7' (-d. f. '. 1. 1.011 .1.1)	
rdiv(other[, axis, level, fill_value])	Floating division of dataframe and other, element-
, a (dald )	wise (binary operator <i>rtruediv</i> ).
reindex(**kwargs)	Conform DataFrame to new index with optional fill-
	ing logic, placing NA/NaN in locations having no
	value in the previous index.
reindex_axis(labels[, axis, method, level,])	Conform input object to new index with optional fill-
	ing logic, placing NA/NaN in locations having no
	value in the previous index.
<pre>reindex_like(other[, method, copy, limit,])</pre>	Return an object with matching indices to myself.
rename(**kwargs)	Alter axes labels.
rename_axis(mapper[, axis, copy, inplace])	Alter the name of the index or columns.
reorder_levels(order[, axis])	Rearrange index levels using input order.
replace([to_replace, value, inplace, limit,])	Replace values given in to_replace with value.
resample(rule[, how, axis, fill_method,])	Convenience method for frequency conversion and
resumprectulet, now, axis, mi_method,])	resampling of time series.
reset_index([level, drop, inplace,])	For DataFrame with multi-level index, return new
reset_index([level, drop, inplace,])	DataFrame with labeling information in the columns
	under the index names, defaulting to 'level_0',
	_
of the state of th	'level_1', etc.
rfloordiv(other[, axis, level, fill_value])	Integer division of dataframe and other, element-
*/ ·1	wise (binary operator <i>rfloordiv</i> ).
rmod(other[, axis, level, fill_value])	Modulo of dataframe and other, element-wise (bi-
	nary operator <i>rmod</i> ).
rmul(other[, axis, level, fill_value])	Multiplication of dataframe and other, element-wise
	(binary operator <i>rmul</i> ).
<pre>rolling(window[, min_periods, center,])</pre>	Provides rolling window calculations.
round([decimals])	Round a DataFrame to a variable number of decimal
	places.
rpow(other[, axis, level, fill_value])	Exponential power of dataframe and other, element-
	wise (binary operator <i>rpow</i> ).
rsub(other[, axis, level, fill_value])	Subtraction of dataframe and other, element-wise
· · · · · · · · · · · · · · · · · · ·	(binary operator <i>rsub</i> ).
rtruediv(other[, axis, level, fill_value])	Floating division of dataframe and other, element-
refacely (others, axis, rever, mi_varae))	wise (binary operator <i>rtruediv</i> ).
sample([n, frac, replace, weights,])	Return a random sample of items from an axis of
camp 10([ii, iiac, icpiace, weights,])	object.
select(crit[, axis])	Return data corresponding to axis labels matching
SELECCION, axis])	
7 (['1 1 1 1 1 ])	criteria
<pre>select_dtypes([include, exclude])</pre>	Return a subset of the DataFrame's columns based
	on the column dtypes.
<pre>sem([axis, skipna, level, ddof, numeric_only])</pre>	Return unbiased standard error of the mean over re-
	quested axis.
<pre>set_axis(labels[, axis, inplace])</pre>	Assign desired index to given axis.
	Continued on next page

	d from previous page
<pre>set_index(keys[, drop, append, inplace,])</pre>	Set the DataFrame index (row labels) using one or
	more existing columns.
set_value(index, col, value[, takeable])	Put single value at passed column and index
shift([periods, freq, axis])	Shift index by desired number of periods with an op-
• -	tional time freq
skew([axis, skipna, level, numeric_only])	Return unbiased skew over requested axis Normal-
(En -) - 1	ized by N-1
slice_shift([periods, axis])	Equivalent to <i>shift</i> without copying data.
sort_index([axis, level, ascending,])	Sort object by labels (along an axis)
sort_values(by[, axis, ascending, inplace,])	Sort by the values along either axis
sortlevel([level, axis, ascending, inplace,])	Sort multilevel index by chosen axis and primary
borere ver (lievel, axis, ascending, inplace,)	level.
squeeze([axis])	Squeeze length 1 dimensions.
stack([level, dropna])	Stack the prescribed level(s) from columns to index.
std([axis, skipna, level, ddof, numeric_only])	Return sample standard deviation over requested
Sca([axis, skipha, level, ddol, humeric_omy])	
	axis.
<pre>sub(other[, axis, level, fill_value])</pre>	Subtraction of dataframe and other, element-wise
	(binary operator <i>sub</i> ).
<pre>subtract(other[, axis, level, fill_value])</pre>	Subtraction of dataframe and other, element-wise
	(binary operator <i>sub</i> ).
sum([axis, skipna, level, numeric_only,])	Return the sum of the values for the requested axis
swapaxes(axis1, axis2[, copy])	Interchange axes and swap values axes appropriately
swaplevel([i, j, axis])	Swap levels i and j in a MultiIndex on a particular
	axis
tail([n])	Return the last <i>n</i> rows.
<pre>take(indices[, axis, convert, is_copy])</pre>	Return the elements in the given <i>positional</i> indices
	along an axis.
to_clipboard([excel, sep])	Copy object to the system clipboard.
to_csv(*args, **kwargs)	Write DataFrame to a comma-separated values (csv)
	file
to_dense()	Return dense representation of NDFrame (as op-
	posed to sparse)
to_dict([orient, into])	Convert the DataFrame to a dictionary.
to_excel(*args, **kwargs)	Write DataFrame to an excel sheet
to_feather(fname)	write out the binary feather-format for DataFrames
to_gbq(destination_table, project_id[,])	Write a DataFrame to a Google BigQuery table.
to_hdf(path_or_buf, key, **kwargs)	Write the contained data to an HDF5 file using HDF-
	Store.
to_html(*args, **kwargs)	Render a DataFrame as an HTML table.
to_json([path_or_buf, orient, date_format,])	Convert the object to a JSON string.
to_latex([buf, columns, col_space, header,])	Render an object to a tabular environment table.
to_mol2([filepath_or_buffer,])	Write DataFrame to Mol2 file.
to_msgpack([path_or_buf, encoding])	msgpack (serialize) object to input file path
to_panel()	Transform long (stacked) format (DataFrame) into
to_panet()	— · · · · · · · · · · · · · · · · · · ·
+ + (fnomo[ or -::::::::::	wide (3D, Panel) format.
to_parquet(fname[, engine, compression])	Write a DataFrame to the binary parquet format.
<pre>to_period([freq, axis, copy])</pre>	Convert DataFrame from DatetimeIndex to Peri-
	odIndex with desired frequency (inferred from index
	if not passed)
to_pickle(path[, compression, protocol])	Pickle (serialize) object to file.
to_records([index, convert_datetime64])	Convert DataFrame to a NumPy record array.
	Continued on next page

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to_sdf([filepath_or_buffer,])	Write DataFrame to SDF file.
to_sparse([fill_value, kind])	Convert to SparseDataFrame
to_sql(name, con[, schema, if_exists,])	Write records stored in a DataFrame to a SQL
	database.
to_stata(fname[, convert_dates,])	Export Stata binary dta files.
to_string([buf, columns, col_space, header,	Render a DataFrame to a console-friendly tabular
])	output.
to_timestamp([freq, how, axis, copy])	Cast to DatetimeIndex of timestamps, at beginning
	of period
to_xarray()	Return an xarray object from the pandas object.
transform(func, *args, **kwargs)	Call function producing a like-indexed NDFrame
	and return a NDFrame with the transformed values
transpose(*args, **kwargs)	Transpose index and columns.
truediv(other[, axis, level, fill_value])	Floating division of dataframe and other, element-
	wise (binary operator <i>truediv</i> ).
<pre>truncate([before, after, axis, copy])</pre>	Truncate a Series or DataFrame before and after
	some index value.
tshift([periods, freq, axis])	Shift the time index, using the index's frequency if
	available.
tz_convert(tz[, axis, level, copy])	Convert tz-aware axis to target time zone.
tz_localize(tz[, axis, level, copy, ambiguous])	Localize tz-naive TimeSeries to target time zone.
unstack([level, fill_value])	Pivot a level of the (necessarily hierarchical) index
	labels, returning a DataFrame having a new level of
	column labels whose inner-most level consists of the
	pivoted index labels.
update(other[, join, overwrite,])	Modify in place using non-NA values from another
	DataFrame.
<pre>var([axis, skipna, level, ddof, numeric_only])</pre>	Return unbiased variance over requested axis.
where(cond[, other, inplace, axis, level,])	Return an object of same shape as self and whose
	corresponding entries are from self where cond is
	True and otherwise are from <i>other</i> .
xs(key[, axis, level, drop_level])	Returns a cross-section (row(s) or column(s)) from
	the Series/DataFrame.

Т

Transpose index and columns.

Reflect the DataFrame over its main diagonal by writing rows as columns and vice-versa. The property T is an accessor to the method transpose().

### **Parameters**

**copy** [bool, default False] If True, the underlying data is copied. Otherwise (default), no copy is made if possible.

\*args, \*\*kwargs Additional keywords have no effect but might be accepted for compatibility with numpy.

# Returns

**DataFrame** The transposed DataFrame.

# See also:

numpy.transpose Permute the dimensions of a given array.

### **Notes**

Transposing a DataFrame with mixed dtypes will result in a homogeneous DataFrame with the *object* dtype. In such a case, a copy of the data is always made.

### **Examples**

# Square DataFrame with homogeneous dtype

```
>>> d1 = {'col1': [1, 2], 'col2': [3, 4]}
>>> df1 = pd.DataFrame(data=d1)
>>> df1
    col1 col2
0    1    3
1    2    4
```

When the dtype is homogeneous in the original DataFrame, we get a transposed DataFrame with the same dtype:

```
>>> df1.dtypes
col1   int64
col2   int64
dtype: object
>>> df1_transposed.dtypes
0   int64
1   int64
dtype: object
```

# Non-square DataFrame with mixed dtypes

```
>>> d2 = {'name': ['Alice', 'Bob'],
... 'score': [9.5, 8],
... 'employed': [False, True],
... 'kids': [0, 0]}
>>> df2 = pd.DataFrame(data=d2)
>>> df2
name score employed kids
0 Alice 9.5 False 0
1 Bob 8.0 True 0
```

When the DataFrame has mixed dtypes, we get a transposed DataFrame with the *object* dtype:

```
>>> df2.dtypes
name object
score float64
employed bool
kids int64
dtype: object
>>> df2_transposed.dtypes
0 object
1 object
dtype: object
```

### abs()

Return a Series/DataFrame with absolute numeric value of each element.

This function only applies to elements that are all numeric.

#### Returns

**abs** Series/DataFrame containing the absolute value of each element.

### See also:

numpy.absolute calculate the absolute value element-wise.

# **Notes**

For complex inputs, 1.2 + 1 j, the absolute value is  $\sqrt{a^2 + b^2}$ .

# **Examples**

Absolute numeric values in a Series.

```
>>> s = pd.Series([-1.10, 2, -3.33, 4])

>>> s.abs()

0    1.10

1    2.00

2    3.33

3    4.00

dtype: float64
```

Absolute numeric values in a Series with complex numbers.

```
>>> s = pd.Series([1.2 + 1j])
>>> s.abs()
0 1.56205
dtype: float64
```

Absolute numeric values in a Series with a Timedelta element.

```
>>> s = pd.Series([pd.Timedelta('1 days')])
>>> s.abs()
0  1 days
dtype: timedelta64[ns]
```

Select rows with data closest to certain value using argsort (from StackOverflow).

```
>>> df = pd.DataFrame({
        'a': [4, 5, 6, 7],
        'b': [10, 20, 30, 40],
        'c': [100, 50, -30, -50]
. . . })
>>> df
          b
               С
0
         10
             100
1
         20
              50
2
     6
         30
             -30
3
     7
         40 -50
>>> df.loc[(df.c - 43).abs().argsort()]
         b
         20
     5
              50
0
     4
         10
             100
         30
             -30
2
     6
             -50
3
     7
         40
```

add (other, axis='columns', level=None, fill\_value=None)

Addition of dataframe and other, element-wise (binary operator add).

Equivalent to dataframe + other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

```
other [Series, DataFrame, or constant]
```

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

### Returns

result [DataFrame]

### See also:

DataFrame.radd

### **Notes**

Mismatched indices will be unioned together

# **Examples**

```
>>> a = pd.DataFrame([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'],
... columns=['one'])
>>> a
one
a 1.0
```

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```
1.0
c 1.0
d NaN
>>> b = pd.DataFrame(dict(one=[1, np.nan, 1, np.nan],
                       two=[np.nan, 2, np.nan, 2]),
                   index=['a', 'b', 'd', 'e'])
. . .
>>> b
  one two
  1.0 NaN
b NaN 2.0
d 1.0 NaN
e NaN 2.0
>>> a.add(b, fill_value=0)
  one two
a 2.0 NaN
b 1.0 2.0
  1.0 NaN
  1.0 NaN
  NaN 2.0
```

# add\_prefix (prefix)

Prefix labels with string prefix.

For Series, the row labels are prefixed. For DataFrame, the column labels are prefixed.

### **Parameters**

**prefix** [str] The string to add before each label.

## Returns

**Series or DataFrame** New Series or DataFrame with updated labels.

See also:

Series.add\_suffix Suffix row labels with string suffix.

DataFrame.add\_suffix Suffix column labels with string suffix.

# **Examples**

```
>>> s = pd.Series([1, 2, 3, 4])

>>> s

0    1

1    2

2    3

3    4

dtype: int64
```

```
>>> s.add_prefix('item_')
item_0   1
item_1   2
item_2   3
item_3   4
dtype: int64
```

```
>>> df = pd.DataFrame({'A': [1, 2, 3, 4], 'B': [3, 4, 5, 6]})
>>> df

A B
0 1 3
1 2 4
2 3 5
3 4 6
```

### add\_suffix (suffix)

Suffix labels with string *suffix*.

For Series, the row labels are suffixed. For DataFrame, the column labels are suffixed.

### **Parameters**

**suffix** [str] The string to add after each label.

### **Returns**

**Series or DataFrame** New Series or DataFrame with updated labels.

See also:

Series.add\_prefix Prefix row labels with string prefix.

DataFrame.add\_prefix Prefix column labels with string prefix.

# **Examples**

```
>>> df = pd.DataFrame({'A': [1, 2, 3, 4], 'B': [3, 4, 5, 6]})
>>> df
A B
0 1 3
1 2 4
```

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```
2 3 5
3 4 6
```

agg (func, axis=0, \*args, \*\*kwargs)

Aggregate using one or more operations over the specified axis.

New in version 0.20.0.

#### **Parameters**

**func** [function, string, dictionary, or list of string/functions] Function to use for aggregating the data. If a function, must either work when passed a DataFrame or when passed to DataFrame.apply. For a DataFrame, can pass a dict, if the keys are DataFrame column names.

Accepted combinations are:

- string function name.
- function.
- · list of functions.
- dict of column names -> functions (or list of functions).

axis [{0 or 'index', 1 or 'columns'}, default 0]

- 0 or 'index': apply function to each column.
- 1 or 'columns': apply function to each row.

\*args Positional arguments to pass to func.

\*\*kwargs Keyword arguments to pass to func.

### **Returns**

aggregated [DataFrame]

See also:

**DataFrame.apply** Perform any type of operations.

**DataFrame.transform** Perform transformation type operations.

pandas.core.groupby.GroupBy Perform operations over groups.

pandas.core.resample.Resampler Perform operations over resampled bins.

pandas.core.window.Rolling Perform operations over rolling window.

pandas.core.window.Expanding Perform operations over expanding window.

pandas.core.window.EWM Perform operation over exponential weighted window.

### **Notes**

agg is an alias for aggregate. Use the alias.

A passed user-defined-function will be passed a Series for evaluation.

The aggregation operations are always performed over an axis, either the index (default) or the column axis. This behavior is different from *numpy* aggregation functions (*mean*, *median*, *prod*, *sum*, *std*, *var*), where the default is to compute the aggregation of the flattened array, e.g., numpy.mean(arr\_2d) as opposed to numpy.mean(arr\_2d, axis=0).

agg is an alias for aggregate. Use the alias.

# **Examples**

```
>>> df = pd.DataFrame([[1, 2, 3],
...
[4, 5, 6],
...
[7, 8, 9],
...
[np.nan, np.nan, np.nan]],
...
columns=['A', 'B', 'C'])
```

Aggregate these functions over the rows.

```
>>> df.agg(['sum', 'min'])

A B C

sum 12.0 15.0 18.0

min 1.0 2.0 3.0
```

Different aggregations per column.

```
>>> df.agg({'A' : ['sum', 'min'], 'B' : ['min', 'max']})

A B

max NaN 8.0

min 1.0 2.0

sum 12.0 NaN
```

Aggregate over the columns.

aggregate (func, axis=0, \*args, \*\*kwargs)

Aggregate using one or more operations over the specified axis.

New in version 0.20.0.

# **Parameters**

**func** [function, string, dictionary, or list of string/functions] Function to use for aggregating the data. If a function, must either work when passed a DataFrame or when passed to DataFrame.apply. For a DataFrame, can pass a dict, if the keys are DataFrame column names.

Accepted combinations are:

- string function name.
- function.
- · list of functions.
- dict of column names -> functions (or list of functions).

axis [{0 or 'index', 1 or 'columns'}, default 0]

- 0 or 'index': apply function to each column.
- 1 or 'columns': apply function to each row.

\*args Positional arguments to pass to func.

\*\*kwargs Keyword arguments to pass to func.

#### Returns

```
aggregated [DataFrame]
```

See also:

**DataFrame.apply** Perform any type of operations.

**DataFrame.transform** Perform transformation type operations.

pandas.core.groupby.GroupBy Perform operations over groups.

pandas.core.resample.Resampler Perform operations over resampled bins.

pandas.core.window.Rolling Perform operations over rolling window.

pandas.core.window.Expanding Perform operations over expanding window.

pandas.core.window.EWM Perform operation over exponential weighted window.

# **Notes**

agg is an alias for aggregate. Use the alias.

A passed user-defined-function will be passed a Series for evaluation.

The aggregation operations are always performed over an axis, either the index (default) or the column axis. This behavior is different from *numpy* aggregation functions (*mean*, *median*, *prod*, *sum*, *std*, *var*), where the default is to compute the aggregation of the flattened array, e.g., numpy.mean(arr\_2d) as opposed to numpy.mean(arr\_2d, axis=0).

agg is an alias for aggregate. Use the alias.

# **Examples**

Aggregate these functions over the rows.

```
>>> df.agg(['sum', 'min'])

A B C

sum 12.0 15.0 18.0

min 1.0 2.0 3.0
```

Different aggregations per column.

```
>>> df.agg({'A' : ['sum', 'min'], 'B' : ['min', 'max']})

A B

max NaN 8.0

min 1.0 2.0

sum 12.0 NaN
```

Aggregate over the columns.

align(other, join='outer', axis=None, level=None, copy=True, fill\_value=None, method=None,
limit=None, fill\_axis=0, broadcast\_axis=None)

Align two objects on their axes with the specified join method for each axis Index

#### **Parameters**

```
other [DataFrame or Series]
join [{'outer', 'inner', 'left', 'right'}, default 'outer']
```

**axis** [allowed axis of the other object, default None] Align on index (0), columns (1), or both (None)

**level** [int or level name, default None] Broadcast across a level, matching Index values on the passed MultiIndex level

**copy** [boolean, default True] Always returns new objects. If copy=False and no reindexing is required then original objects are returned.

**fill\_value** [scalar, default np.NaN] Value to use for missing values. Defaults to NaN, but can be any "compatible" value

```
method [str, default None]
```

```
limit [int, default None]
```

fill\_axis [{0 or 'index', 1 or 'columns'}, default 0] Filling axis, method and limit

**broadcast\_axis** [{0 or 'index', 1 or 'columns'}, default None] Broadcast values along this axis, if aligning two objects of different dimensions

#### Returns

```
(left, right) [(DataFrame, type of other)] Aligned objects
```

**all** (axis=0, bool\_only=None, skipna=True, level=None, \*\*kwargs)
Return whether all elements are True, potentially over an axis.

Returns True if all elements within a series or along a Dataframe axis are non-zero, not-empty or not-False.

### **Parameters**

axis [{0 or 'index', 1 or 'columns', None}, default 0] Indicate which axis or axes should be reduced.

- 0 / 'index' : reduce the index, return a Series whose index is the original column labels.
- 1 / 'columns' : reduce the columns, return a Series whose index is the original index.
- None: reduce all axes, return a scalar.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series.

**bool\_only** [boolean, default None] Include only boolean columns. If None, will attempt to use everything, then use only boolean data. Not implemented for Series.

\*\*kwargs [any, default None] Additional keywords have no effect but might be accepted for compatibility with NumPy.

#### Returns

**all** [Series or DataFrame (if level specified)]

#### See also:

```
pandas.Series.all Return True if all elements are True
pandas.DataFrame.any Return True if one (or more) elements are True
```

# **Examples**

## Series

```
>>> pd.Series([True, True]).all()
True
>>> pd.Series([True, False]).all()
False
```

## DataFrames

Create a dataframe from a dictionary.

```
>>> df = pd.DataFrame({'col1': [True, True], 'col2': [True, False]})
>>> df
    col1    col2
0    True    True
1    True    False
```

Default behaviour checks if column-wise values all return True.

```
>>> df.all()
col1 True
col2 False
dtype: bool
```

Specify axis='columns' to check if row-wise values all return True.

```
>>> df.all(axis='columns')
0 True
1 False
dtype: bool
```

Or axis=None for whether every value is True.

```
>>> df.all(axis=None)
False
```

any (axis=0, bool\_only=None, skipna=True, level=None, \*\*kwargs)

Return whether any element is True over requested axis.

Unlike DataFrame.all(), this performs an *or* operation. If any of the values along the specified axis is True, this will return True.

#### **Parameters**

axis [{0 or 'index', 1 or 'columns', None}, default 0] Indicate which axis or axes should be reduced.

- 0 / 'index' : reduce the index, return a Series whose index is the original column labels.
- 1 / 'columns' : reduce the columns, return a Series whose index is the original index.
- None: reduce all axes, return a scalar.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series.

**bool\_only** [boolean, default None] Include only boolean columns. If None, will attempt to use everything, then use only boolean data. Not implemented for Series.

\*\*kwargs [any, default None] Additional keywords have no effect but might be accepted for compatibility with NumPy.

### Returns

any [Series or DataFrame (if level specified)]

See also:

pandas.DataFrame.all Return whether all elements are True.

# **Examples**

#### Series

For Series input, the output is a scalar indicating whether any element is True.

```
>>> pd.Series([True, False]).any()
True
```

## **DataFrame**

Whether each column contains at least one True element (the default).

```
>>> df = pd.DataFrame({"A": [1, 2], "B": [0, 2], "C": [0, 0]})
>>> df

A B C
0 1 0 0
1 2 2 0
```

```
>>> df.any()
A True
B True
C False
dtype: bool
```

Aggregating over the columns.

```
>>> df.any(axis='columns')
0 True
1 True
dtype: bool
```

```
>>> df = pd.DataFrame({"A": [True, False], "B": [1, 0]})
>>> df

A B

0 True 1

1 False 0
```

```
>>> df.any(axis='columns')
0    True
1    False
dtype: bool
```

Aggregating over the entire DataFrame with axis=None.

```
>>> df.any(axis=None)
True
```

any for an empty DataFrame is an empty Series.

```
>>> pd.DataFrame([]).any()
Series([], dtype: bool)
```

append (other, ignore\_index=False, verify\_integrity=False, sort=None)

Append rows of *other* to the end of this frame, returning a new object. Columns not in this frame are added as new columns.

### **Parameters**

**other** [DataFrame or Series/dict-like object, or list of these] The data to append.

ignore\_index [boolean, default False] If True, do not use the index labels.

**verify\_integrity** [boolean, default False] If True, raise ValueError on creating index with duplicates.

sort [boolean, default None] Sort columns if the columns of *self* and *other* are not aligned. The default sorting is deprecated and will change to not-sorting in a future version of pandas. Explicitly pass sort=True to silence the warning and sort. Explicitly pass sort=False to silence the warning and not sort.

New in version 0.23.0.

#### Returns

**appended** [DataFrame]

See also:

pandas.concat General function to concatenate DataFrame, Series or Panel objects

# **Notes**

If a list of dict/series is passed and the keys are all contained in the DataFrame's index, the order of the columns in the resulting DataFrame will be unchanged.

Iteratively appending rows to a DataFrame can be more computationally intensive than a single concatenate. A better solution is to append those rows to a list and then concatenate the list with the original DataFrame all at once.

# **Examples**

With *ignore\_index* set to True:

```
>>> df.append(df2, ignore_index=True)

A B
0 1 2
1 3 4
2 5 6
3 7 8
```

The following, while not recommended methods for generating DataFrames, show two ways to generate a DataFrame from multiple data sources.

Less efficient:

```
>>> df = pd.DataFrame(columns=['A'])
>>> for i in range(5):
... df = df.append({'A': i}, ignore_index=True)
```

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```
>>> df
A
0 0
1 1
2 2
3 3
4 4
```

### More efficient:

**apply** (func, axis=0, broadcast=None, raw=False, reduce=None, result\_type=None, args=(), \*\*kwds) Apply a function along an axis of the DataFrame.

Objects passed to the function are Series objects whose index is either the DataFrame's index (axis=0) or the DataFrame's columns (axis=1). By default (result\_type=None), the final return type is inferred from the return type of the applied function. Otherwise, it depends on the *result\_type* argument.

#### **Parameters**

**func** [function] Function to apply to each column or row.

axis [{0 or 'index', 1 or 'columns'}, default 0] Axis along which the function is applied:

- 0 or 'index': apply function to each column.
- 1 or 'columns': apply function to each row.

**broadcast** [bool, optional] Only relevant for aggregation functions:

- False or None: returns a Series whose length is the length of the index or the number of columns (based on the *axis* parameter)
- True: results will be broadcast to the original shape of the frame, the original index and columns will be retained.

Deprecated since version 0.23.0: This argument will be removed in a future version, replaced by result\_type='broadcast'.

raw [bool, default False]

- False: passes each row or column as a Series to the function.
- True: the passed function will receive ndarray objects instead. If you are just applying a NumPy reduction function this will achieve much better performance.

**reduce** [bool or None, default None] Try to apply reduction procedures. If the DataFrame is empty, *apply* will use *reduce* to determine whether the result should be a Series or a DataFrame. If reduce=None (the default), *apply*'s return value will be guessed by calling *func* on an empty Series (note: while guessing, exceptions raised by *func* will be ignored). If reduce=True a Series will always be returned, and if reduce=False a DataFrame will always be returned.

Deprecated since version 0.23.0: This argument will be removed in a future version, replaced by result\_type='reduce'.

**result\_type** [{'expand', 'reduce', 'broadcast', None}, default None] These only act when axis=1 (columns):

- 'expand': list-like results will be turned into columns.
- 'reduce': returns a Series if possible rather than expanding list-like results. This is the opposite of 'expand'.
- 'broadcast': results will be broadcast to the original shape of the DataFrame, the original index and columns will be retained.

The default behaviour (None) depends on the return value of the applied function: list-like results will be returned as a Series of those. However if the apply function returns a Series these are expanded to columns.

New in version 0.23.0.

**args** [tuple] Positional arguments to pass to *func* in addition to the array/series.

\*\*kwds Additional keyword arguments to pass as keywords arguments to func.

#### Returns

```
applied [Series or DataFrame]
```

See also:

DataFrame.applymap For elementwise operations

DataFrame.aggregate only perform aggregating type operations

DataFrame.transform only perform transformating type operations

# **Notes**

In the current implementation apply calls *func* twice on the first column/row to decide whether it can take a fast or slow code path. This can lead to unexpected behavior if *func* has side-effects, as they will take effect twice for the first column/row.

# **Examples**

```
>>> df = pd.DataFrame([[4, 9],] * 3, columns=['A', 'B'])
>>> df

A B
0 4 9
1 4 9
2 4 9
```

Using a numpy universal function (in this case the same as np.sqrt(df)):

```
>>> df.apply(np.sqrt)

A B
0 2.0 3.0
1 2.0 3.0
2 2.0 3.0
```

Using a reducing function on either axis

```
>>> df.apply(np.sum, axis=0)
A 12
B 27
dtype: int64
```

```
>>> df.apply(np.sum, axis=1)
0 13
1 13
2 13
dtype: int64
```

Retuning a list-like will result in a Series

```
>>> df.apply(lambda x: [1, 2], axis=1)
0    [1, 2]
1    [1, 2]
2    [1, 2]
dtype: object
```

Passing result\_type='expand' will expand list-like results to columns of a Dataframe

Returning a Series inside the function is similar to passing result\_type='expand'. The resulting column names will be the Series index.

```
>>> df.apply(lambda x: pd.Series([1, 2], index=['foo', 'bar']), axis=1)
foo bar
0 1 2
1 1 2
2 1 2
```

Passing result\_type='broadcast' will ensure the same shape result, whether list-like or scalar is returned by the function, and broadcast it along the axis. The resulting column names will be the originals.

```
>>> df.apply(lambda x: [1, 2], axis=1, result_type='broadcast')

A B
0 1 2
1 1 2
2 1 2
```

# applymap (func)

Apply a function to a Dataframe elementwise.

This method applies a function that accepts and returns a scalar to every element of a DataFrame.

# **Parameters**

func [callable] Python function, returns a single value from a single value.

# Returns

**DataFrame** Transformed DataFrame.

### See also:

**DataFrame.apply** Apply a function along input axis of DataFrame

# **Examples**

Note that a vectorized version of *func* often exists, which will be much faster. You could square each number elementwise.

But it's better to avoid applymap in that case.

# as\_blocks (copy=True)

Convert the frame to a dict of dtype -> Constructor Types that each has a homogeneous dtype.

Deprecated since version 0.21.0.

NOTE: the dtypes of the blocks WILL BE PRESERVED HERE (unlike in as\_matrix)

# **Parameters**

```
copy [boolean, default True]
```

### **Returns**

```
values [a dict of dtype -> Constructor Types]
```

# as\_matrix(columns=None)

Convert the frame to its Numpy-array representation.

Deprecated since version 0.23.0: Use <code>DataFrame.values()</code> instead.

#### **Parameters**

**columns: list, optional, default:None** If None, return all columns, otherwise, returns specified columns.

# Returns

**values** [ndarray] If the caller is heterogeneous and contains booleans or objects, the result will be of dtype=object. See Notes.

#### See also:

pandas.DataFrame.values

### **Notes**

Return is NOT a Numpy-matrix, rather, a Numpy-array.

The dtype will be a lower-common-denominator dtype (implicit upcasting); that is to say if the dtypes (even of numeric types) are mixed, the one that accommodates all will be chosen. Use this with care if you are not dealing with the blocks.

e.g. If the dtypes are float16 and float32, dtype will be upcast to float32. If dtypes are int32 and uint8, dtype will be upcase to int32. By numpy.find\_common\_type convention, mixing int64 and uint64 will result in a flot64 dtype.

This method is provided for backwards compatibility. Generally, it is recommended to use '.values'.

**asfreq** (freq, method=None, how=None, normalize=False, fill\_value=None) Convert TimeSeries to specified frequency.

Optionally provide filling method to pad/backfill missing values.

Returns the original data conformed to a new index with the specified frequency. resample is more appropriate if an operation, such as summarization, is necessary to represent the data at the new frequency.

#### **Parameters**

freq [DateOffset object, or string]

**method** [{'backfill'/'bfill', 'pad'/'ffill'}, default None] Method to use for filling holes in reindexed Series (note this does not fill NaNs that already were present):

- 'pad' / 'ffill': propagate last valid observation forward to next valid
- 'backfill' / 'bfill': use NEXT valid observation to fill

**how** [{'start', 'end'}, default end] For PeriodIndex only, see PeriodIndex.asfreq

normalize [bool, default False] Whether to reset output index to midnight

**fill\_value: scalar, optional** Value to use for missing values, applied during upsampling (note this does not fill NaNs that already were present).

New in version 0.20.0.

#### Returns

**converted** [type of caller]

### See also:

reindex

### **Notes**

To learn more about the frequency strings, please see this link.

# **Examples**

Start by creating a series with 4 one minute timestamps.

Upsample the series into 30 second bins.

```
>>> df.asfreq(freq='30S')

s
2000-01-01 00:00:00 0.0
2000-01-01 00:00:30 NaN
2000-01-01 00:01:00 NaN
2000-01-01 00:01:30 NaN
2000-01-01 00:02:00 2.0
2000-01-01 00:02:30 NaN
2000-01-01 00:02:30 NaN
2000-01-01 00:03:00 3.0
```

Upsample again, providing a fill value.

```
>>> df.asfreq(freq='30S', fill_value=9.0)

s
2000-01-01 00:00:00 0.0
2000-01-01 00:00:30 9.0
2000-01-01 00:01:00 NaN
2000-01-01 00:01:30 9.0
2000-01-01 00:02:00 2.0
2000-01-01 00:02:30 9.0
2000-01-01 00:02:30 9.0
2000-01-01 00:03:00 3.0
```

Upsample again, providing a method.

```
>>> df.asfreq(freq='30S', method='bfill')
s
2000-01-01 00:00:00 0.0
2000-01-01 00:00:30 NaN
2000-01-01 00:01:00 NaN
2000-01-01 00:01:30 2.0
2000-01-01 00:02:00 2.0
2000-01-01 00:02:30 3.0
2000-01-01 00:03:00 3.0
```

asof (where, subset=None)

The last row without any NaN is taken (or the last row without NaN considering only the subset of columns in the case of a DataFrame)

New in version 0.19.0: For DataFrame

If there is no good value, NaN is returned for a Series a Series of NaN values for a DataFrame

**Parameters** 

where [date or array of dates]

**subset** [string or list of strings, default None] if not None use these columns for NaN propagation

### **Returns**

## where is scalar

- value or NaN if input is Series
- · Series if input is DataFrame

where is Index: same shape object as input

### See also:

merge\_asof

#### **Notes**

Dates are assumed to be sorted Raises if this is not the case

```
assign(**kwargs)
```

Assign new columns to a DataFrame, returning a new object (a copy) with the new columns added to the original ones. Existing columns that are re-assigned will be overwritten.

#### **Parameters**

**kwargs** [keyword, value pairs] keywords are the column names. If the values are callable, they are computed on the DataFrame and assigned to the new columns. The callable must not change input DataFrame (though pandas doesn't check it). If the values are not callable, (e.g. a Series, scalar, or array), they are simply assigned.

### Returns

**df** [DataFrame] A new DataFrame with the new columns in addition to all the existing columns.

#### **Notes**

Assigning multiple columns within the same assign is possible. For Python 3.6 and above, later items in '\*\*kwargs' may refer to newly created or modified columns in 'df'; items are computed and assigned into 'df' in order. For Python 3.5 and below, the order of keyword arguments is not specified, you cannot refer to newly created or modified columns. All items are computed first, and then assigned in alphabetical order.

Changed in version 0.23.0: Keyword argument order is maintained for Python 3.6 and later.

### **Examples**

```
>>> df = pd.DataFrame({'A': range(1, 11), 'B': np.random.randn(10)})
```

Where the value is a callable, evaluated on df:

```
>>> df.assign(ln_A = lambda x: np.log(x.A))

A B ln_A

0 1 0.426905 0.000000
1 2 -0.780949 0.693147
2 3 -0.418711 1.098612
3 4 -0.269708 1.386294
4 5 -0.274002 1.609438
5 6 -0.500792 1.791759
6 7 1.649697 1.945910
7 8 -1.495604 2.079442
8 9 0.549296 2.197225
9 10 -0.758542 2.302585
```

Where the value already exists and is inserted:

```
>>> newcol = np.log(df['A'])
>>> df.assign(ln_A=newcol)
   Α
            В
                   ln_A
  1 0.426905 0.000000
  2 -0.780949 0.693147
   3 -0.418711 1.098612
3
   4 -0.269708 1.386294
  5 -0.274002 1.609438
   6 -0.500792 1.791759
      1.649697
               1.945910
   8 -1.495604 2.079442
      0.549296 2.197225
  10 -0.758542 2.302585
```

Where the keyword arguments depend on each other

```
>>> df = pd.DataFrame({'A': [1, 2, 3]})
```

```
>>> df.assign(B=df.A, C=lambda x:x['A']+ x['B'])

A B C
0 1 1 2
1 2 2 4
2 3 3 6
```

#### astype (\*\*kwargs)

Cast a pandas object to a specified dtype dtype.

## **Parameters**

**dtype** [data type, or dict of column name -> data type] Use a numpy.dtype or Python type to cast entire pandas object to the same type. Alternatively, use {col: dtype, ...}, where col is a column label and dtype is a numpy.dtype or Python type to cast one or more of the DataFrame's columns to column-specific types.

**copy** [bool, default True.] Return a copy when copy=True (be very careful setting copy=False as changes to values then may propagate to other pandas objects).

**errors** [{'raise', 'ignore'}, default 'raise'.] Control raising of exceptions on invalid data for provided dtype.

- raise: allow exceptions to be raised
- ignore: suppress exceptions. On error return original object

New in version 0.20.0.

raise\_on\_error [raise on invalid input] Deprecated since version 0.20.0: Use errors instead

**kwargs** [keyword arguments to pass on to the constructor]

### **Returns**

**casted** [type of caller]

See also:

pandas.to\_datetime Convert argument to datetime.

pandas.to\_timedelta Convert argument to timedelta.

pandas.to\_numeric Convert argument to a numeric type.

numpy.ndarray.astype Cast a numpy array to a specified type.

# **Examples**

```
>>> ser = pd.Series([1, 2], dtype='int32')
>>> ser
0    1
1    2
dtype: int32
>>> ser.astype('int64')
0    1
1    2
dtype: int64
```

### Convert to categorical type:

```
>>> ser.astype('category')
0 1
1 2
dtype: category
Categories (2, int64): [1, 2]
```

Convert to ordered categorical type with custom ordering:

Note that using copy=False and changing data on a new pandas object may propagate changes:

```
>>> s1 = pd.Series([1,2])
>>> s2 = s1.astype('int64', copy=False)
>>> s2[0] = 10
>>> s1 # note that s1[0] has changed too
0    10
1    2
dtype: int64
```

at

Access a single value for a row/column label pair.

Similar to loc, in that both provide label-based lookups. Use at if you only need to get or set a single value in a DataFrame or Series.

#### Raises

KeyError When label does not exist in DataFrame

See also:

DataFrame.iat Access a single value for a row/column pair by integer position

DataFrame.loc Access a group of rows and columns by label(s)

Series.at Access a single value using a label

# **Examples**

```
>>> df = pd.DataFrame([[0, 2, 3], [0, 4, 1], [10, 20, 30]],
... index=[4, 5, 6], columns=['A', 'B', 'C'])
>>> df
    A B C
4 0 2 3
5 0 4 1
6 10 20 30
```

Get value at specified row/column pair

```
>>> df.at[4, 'B']
2
```

Set value at specified row/column pair

```
>>> df.at[4, 'B'] = 10
>>> df.at[4, 'B']
10
```

Get value within a Series

```
>>> df.loc[5].at['B']
4
```

at\_time (time, asof=False)

Select values at particular time of day (e.g. 9:30AM).

# **Parameters**

time [datetime.time or string]

Returns

values\_at\_time [type of caller]

Raises

TypeError If the index is not a DatetimeIndex

See also:

between\_time Select values between particular times of the day

first Select initial periods of time series based on a date offset

last Select final periods of time series based on a date offset

**DatetimeIndex.indexer\_at\_time** Get just the index locations for values at particular time of the day

# **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='12H')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 00:00:00 1
2018-04-09 12:00:00 2
2018-04-10 00:00:00 3
2018-04-10 12:00:00 4
```

#### axes

Return a list representing the axes of the DataFrame.

It has the row axis labels and column axis labels as the only members. They are returned in that order.

# **Examples**

```
>>> df = pd.DataFrame({'col1': [1, 2], 'col2': [3, 4]})
>>> df.axes
[RangeIndex(start=0, stop=2, step=1), Index(['col1', 'col2'],
dtype='object')]
```

 $\textbf{between\_time} \ (\textit{start\_time}, \textit{end\_time}, \textit{include\_start=True}, \textit{include\_end=True})$ 

Select values between particular times of the day (e.g., 9:00-9:30 AM).

By setting start\_time to be later than end\_time, you can get the times that are *not* between the two times.

# **Parameters**

```
start_time [datetime.time or string]
end_time [datetime.time or string]
include_start [boolean, default True]
include_end [boolean, default True]

Returns
values_between_time [type of caller]

Raises
TypeError If the index is not a DatetimeIndex
```

See also:

at\_time Select values at a particular time of the day

first Select initial periods of time series based on a date offset

last Select final periods of time series based on a date offset

**DatetimeIndex.indexer\_between\_time** Get just the index locations for values between particular times of the day

# **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='1D20min')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 00:00:00 1
2018-04-10 00:20:00 2
2018-04-11 00:40:00 3
2018-04-12 01:00:00 4
```

```
>>> ts.between_time('0:15', '0:45')

A
2018-04-10 00:20:00 2
2018-04-11 00:40:00 3
```

You get the times that are *not* between two times by setting start\_time later than end\_time:

```
>>> ts.between_time('0:45', '0:15')

A
2018-04-09 00:00:00 1
2018-04-12 01:00:00 4
```

```
bfill (axis=None, inplace=False, limit=None, downcast=None)
Synonym for DataFrame.fillna (method='bfill')
```

## blocks

Internal property, property synonym for as\_blocks()

Deprecated since version 0.21.0.

#### bool()

Return the bool of a single element PandasObject.

This must be a boolean scalar value, either True or False. Raise a ValueError if the PandasObject does not have exactly 1 element, or that element is not boolean

```
boxplot (column=None, by=None, ax=None, fontsize=None, rot=0, grid=True, figsize=None, layout=None, return_type=None, **kwds)

Make a box plot from DataFrame columns.
```

Make a box-and-whisker plot from DataFrame columns, optionally grouped by some other columns. A box plot is a method for graphically depicting groups of numerical data through their quartiles. The box extends from the Q1 to Q3 quartile values of the data, with a line at the median (Q2). The whiskers extend from the edges of box to show the range of the data. The position of the whiskers is set by default to I.5 \* IQR (IQR = Q3 - Q1) from the edges of the box. Outlier points are those past the end of the whiskers.

For further details see Wikipedia's entry for boxplot.

### **Parameters**

**column** [str or list of str, optional] Column name or list of names, or vector. Can be any valid input to pandas.DataFrame.groupby().

**by** [str or array-like, optional] Column in the DataFrame to pandas.DataFrame. groupby(). One box-plot will be done per value of columns in by.

**ax** [object of class matplotlib.axes.Axes, optional] The matplotlib axes to be used by boxplot.

**fontsize** [float or str] Tick label font size in points or as a string (e.g., *large*).

**rot** [int or float, default 0] The rotation angle of labels (in degrees) with respect to the screen coordinate sytem.

**grid** [boolean, default True] Setting this to True will show the grid.

figsize [A tuple (width, height) in inches] The size of the figure to create in matplotlib.

**layout** [tuple (rows, columns), optional] For example, (3, 5) will display the subplots using 3 columns and 5 rows, starting from the top-left.

return\_type [{'axes', 'dict', 'both'} or None, default 'axes'] The kind of object to return. The default is axes.

- 'axes' returns the matplotlib axes the boxplot is drawn on.
- 'dict' returns a dictionary whose values are the matplotlib Lines of the boxplot.
- 'both' returns a namedtuple with the axes and dict.
- when grouping with by, a Series mapping columns to return\_type is returned.

If return\_type is *None*, a NumPy array of axes with the same shape as layout is returned.

\*\*kwds All other plotting keyword arguments to be passed to matplotlib.pyplot. boxplot().

#### **Returns**

**result :** The return type depends on the *return\_type* parameter:

- 'axes': object of class matplotlib.axes.Axes
- 'dict' : dict of matplotlib.lines.Line2D objects
- 'both': a nametuple with strucure (ax, lines)

For data grouped with by:

- Series
- array (for return\_type = None)

### See also:

Series.plot.hist Make a histogram.

matplotlib.pyplot.boxplot Matplotlib equivalent plot.

### **Notes**

Use return\_type='dict' when you want to tweak the appearance of the lines after plotting. In this case a dict containing the Lines making up the boxes, caps, fliers, medians, and whiskers is returned.

# **Examples**

Boxplots can be created for every column in the dataframe by df.boxplot() or indicating the columns to be used:

Boxplots of variables distributions grouped by the values of a third variable can be created using the option by. For instance:

A list of strings (i.e. ['X', 'Y']) can be passed to boxplot in order to group the data by combination of the variables in the x-axis:

The layout of boxplot can be adjusted giving a tuple to layout:

Additional formatting can be done to the boxplot, like suppressing the grid (grid=False), rotating the labels in the x-axis (i.e. rot=45) or changing the fontsize (i.e. fontsize=15):

The parameter return\_type can be used to select the type of element returned by *boxplot*. When return\_type='axes' is selected, the matplotlib axes on which the boxplot is drawn are returned:

```
>>> boxplot = df.boxplot(column=['Col1','Col2'], return_type='axes')
>>> type(boxplot)
<class 'matplotlib.axes._subplots.AxesSubplot'>
```

When grouping with by, a Series mapping columns to return\_type is returned:

```
>>> boxplot = df.boxplot(column=['Col1', 'Col2'], by='X',
... return_type='axes')
>>> type(boxplot)
<class 'pandas.core.series.Series'>
```

If return\_type is None, a NumPy array of axes with the same shape as layout is returned:

```
>>> boxplot = df.boxplot(column=['Col1', 'Col2'], by='X',
... return_type=None)
>>> type(boxplot)
<class 'numpy.ndarray'>
```

clip (lower=None, upper=None, axis=None, inplace=False, \*args, \*\*kwargs)
Trim values at input threshold(s).

Assigns values outside boundary to boundary values. Thresholds can be singular values or array like, and in the latter case the clipping is performed element-wise in the specified axis.

# **Parameters**

**lower** [float or array\_like, default None] Minimum threshold value. All values below this threshold will be set to it.

**upper** [float or array\_like, default None] Maximum threshold value. All values above this threshold will be set to it.

**axis** [int or string axis name, optional] Align object with lower and upper along the given axis.

**inplace** [boolean, default False] Whether to perform the operation in place on the data.

New in version 0.21.0.

\*args, \*\*kwargs Additional keywords have no effect but might be accepted for compatibility with numpy.

# Returns

**Series or DataFrame** Same type as calling object with the values outside the clip boundaries replaced

#### See also:

```
clip_lower Clip values below specified threshold(s).clip_upper Clip values above specified threshold(s).
```

# **Examples**

```
>>> data = {'col_0': [9, -3, 0, -1, 5], 'col_1': [-2, -7, 6, 8, -5]}
>>> df = pd.DataFrame(data)
>>> df
   col_0 col_1
      9
0
             -2
1
      -3
             -7
2
      0
3
      -1
             8
4
       5
             -5
```

Clips per column using lower and upper thresholds:

```
>>> df.clip(-4, 6)
   col_0 col_1
0
       6
1
      -3
             -4
2
       0
              6
3
      -1
              6
4
       5
             -4
```

Clips using specific lower and upper thresholds per column element:

```
>>> df.clip(t, t + 4, axis=0)
   col_0 col_1
0
       6
1
      -3
              -4
2
       0
               3
3
               8
       6
       5
               3
4
```

clip\_lower (threshold, axis=None, inplace=False)

Return copy of the input with values below a threshold truncated.

#### **Parameters**

**threshold** [numeric or array-like] Minimum value allowed. All values below threshold will be set to this value.

- float : every value is compared to threshold.
- array-like: The shape of *threshold* should match the object it's compared to. When *self* is a Series, *threshold* should be the length. When *self* is a DataFrame, *threshold* should 2-D and the same shape as *self* for axis=None, or 1-D and the same length as the axis being compared.

axis [{0 or 'index', 1 or 'columns'}, default 0] Align self with threshold along the given axis

inplace [boolean, default False] Whether to perform the operation in place on the data.

New in version 0.21.0.

#### Returns

**clipped** [same type as input]

See also:

**Series.clip** Return copy of input with values below and above thresholds truncated.

Series.clip\_upper Return copy of input with values above threshold truncated.

# **Examples**

Series single threshold clipping:

Series clipping element-wise using an array of thresholds. *threshold* should be the same length as the Series.

```
>>> elemwise_thresholds = [4, 8, 7, 2, 5]
>>> s.clip_lower(elemwise_thresholds)
0    5
1    8
2    7
3    8
4    9
dtype: int64
```

DataFrames can be compared to a scalar.

```
>>> df = pd.DataFrame({"A": [1, 3, 5], "B": [2, 4, 6]})
>>> df
    A B
0 1 2
1 3 4
2 5 6
```

```
>>> df.clip_lower(3)

A B

0 3 3

1 3 4

2 5 6
```

Or to an array of values. By default, threshold should be the same shape as the DataFrame.

```
>>> df.clip_lower(np.array([[3, 4], [2, 2], [6, 2]]))

A B
0 3 4
1 3 4
2 6 6
```

Control how *threshold* is broadcast with *axis*. In this case *threshold* should be the same length as the axis specified by *axis*.

```
>>> df.clip_lower(np.array([3, 3, 5]), axis='index')

A B
0 3 3
1 3 4
2 5 6
```

```
>>> df.clip_lower(np.array([4, 5]), axis='columns')

A B
0 4 5
1 4 5
2 5 6
```

### clip upper (threshold, axis=None, inplace=False)

Return copy of input with values above given value(s) truncated.

# **Parameters**

```
threshold [float or array_like]
```

axis [int or string axis name, optional] Align object with threshold along the given axis.

inplace [boolean, default False] Whether to perform the operation in place on the data

New in version 0.21.0.

### Returns

clipped [same type as input]

# See also:

clip

### columns

The column labels of the DataFrame.

```
combine (other, func, fill_value=None, overwrite=True)
```

Add two DataFrame objects and do not propagate NaN values, so if for a (column, time) one frame is missing a value, it will default to the other frame's value (which might be NaN as well)

### **Parameters**

```
other [DataFrame]
```

func [function] Function that takes two series as inputs and return a Series or a scalar

```
fill_value [scalar value]
```

**overwrite** [boolean, default True] If True then overwrite values for common keys in the calling frame

### **Returns**

result [DataFrame]

See also:

DataFrame.combine\_first Combine two DataFrame objects and default to non-null values in frame calling the method

# **Examples**

```
>>> df1 = DataFrame({'A': [0, 0], 'B': [4, 4]})
>>> df2 = DataFrame({'A': [1, 1], 'B': [3, 3]})
>>> df1.combine(df2, lambda s1, s2: s1 if s1.sum() < s2.sum() else s2)

A B
0 0 3
1 0 3
```

# combine\_first (other)

Combine two DataFrame objects and default to non-null values in frame calling the method. Result index columns will be the union of the respective indexes and columns

#### **Parameters**

other [DataFrame]

# Returns

combined [DataFrame]

See also:

DataFrame.combine Perform series-wise operation on two DataFrames using a given function

# **Examples**

df1's values prioritized, use values from df2 to fill holes:

compound (axis=None, skipna=None, level=None)

Return the compound percentage of the values for the requested axis

# **Parameters**

```
axis [{index (0), columns (1)}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**numeric\_only** [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

## Returns

**compounded** [Series or DataFrame (if level specified)]

### consolidate(inplace=False)

Compute NDFrame with "consolidated" internals (data of each dtype grouped together in a single ndarray).

Deprecated since version 0.20.0: Consolidate will be an internal implementation only.

Attempt to infer better dtype for object columns.

Deprecated since version 0.21.0.

### **Parameters**

**convert\_dates** [boolean, default True] If True, convert to date where possible. If 'coerce', force conversion, with unconvertible values becoming NaT.

**convert\_numeric** [boolean, default False] If True, attempt to coerce to numbers (including strings), with unconvertible values becoming NaN.

**convert\_timedeltas** [boolean, default True] If True, convert to timedelta where possible. If 'coerce', force conversion, with unconvertible values becoming NaT.

**copy** [boolean, default True] If True, return a copy even if no copy is necessary (e.g. no conversion was done). Note: This is meant for internal use, and should not be confused with inplace.

# Returns

converted [same as input object]

#### See also:

pandas.to\_datetime Convert argument to datetime.

pandas.to\_timedelta Convert argument to timedelta.

pandas.to\_numeric Return a fixed frequency timedelta index, with day as the default.

# copy (deep=True)

Make a copy of this object's indices and data.

When deep=True (default), a new object will be created with a copy of the calling object's data and indices. Modifications to the data or indices of the copy will not be reflected in the original object (see notes below).

When deep=False, a new object will be created without copying the calling object's data or index (only references to the data and index are copied). Any changes to the data of the original will be reflected in the shallow copy (and vice versa).

#### **Parameters**

**deep** [bool, default True] Make a deep copy, including a copy of the data and the indices. With deep=False neither the indices nor the data are copied.

### Returns

**copy** [Series, DataFrame or Panel] Object type matches caller.

## **Notes**

When deep=True, data is copied but actual Python objects will not be copied recursively, only the reference to the object. This is in contrast to *copy.deepcopy* in the Standard Library, which recursively copies object data (see examples below).

While Index objects are copied when deep=True, the underlying numpy array is not copied for performance reasons. Since Index is immutable, the underlying data can be safely shared and a copy is not needed.

# **Examples**

```
>>> s = pd.Series([1, 2], index=["a", "b"])
>>> s
a    1
b    2
dtype: int64
```

# Shallow copy versus default (deep) copy:

```
>>> s = pd.Series([1, 2], index=["a", "b"])
>>> deep = s.copy()
>>> shallow = s.copy(deep=False)
```

Shallow copy shares data and index with original.

```
>>> s is shallow
False
>>> s.values is shallow.values and s.index is shallow.index
True
```

Deep copy has own copy of data and index.

```
>>> s is deep
False
>>> s.values is deep.values or s.index is deep.index
False
```

Updates to the data shared by shallow copy and original is reflected in both; deep copy remains unchanged.

```
>>> s[0] = 3
>>> shallow[1] = 4
>>> s
a 3
b 4
dtype: int64
```

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Note that when copying an object containing Python objects, a deep copy will copy the data, but will not do so recursively. Updating a nested data object will be reflected in the deep copy.

corr (method='pearson', min\_periods=1)

Compute pairwise correlation of columns, excluding NA/null values

### **Parameters**

```
method [{'pearson', 'kendall', 'spearman'}]
```

- pearson: standard correlation coefficient
- · kendall: Kendall Tau correlation coefficient
- spearman : Spearman rank correlation

**min\_periods** [int, optional] Minimum number of observations required per pair of columns to have a valid result. Currently only available for pearson and spearman correlation

### Returns

y [DataFrame]

```
corrwith (other, axis=0, drop=False)
```

Compute pairwise correlation between rows or columns of two DataFrame objects.

# **Parameters**

```
other [DataFrame, Series]
```

**axis** [{0 or 'index', 1 or 'columns'}, default 0] 0 or 'index' to compute column-wise, 1 or 'columns' for row-wise

drop [boolean, default False] Drop missing indices from result, default returns union of all

# Returns

```
correls [Series]
```

count (axis=0, level=None, numeric only=False)

Count non-NA cells for each column or row.

The values *None*, *NaN*, *NaT*, and optionally *numpy.inf* (depending on *pan-das.options.mode.use\_inf\_as\_na*) are considered NA.

#### **Parameters**

axis [{0 or 'index', 1 or 'columns'}, default 0] If 0 or 'index' counts are generated for each column. If 1 or 'columns' counts are generated for each row.

**level** [int or str, optional] If the axis is a *MultiIndex* (hierarchical), count along a particular *level*, collapsing into a *DataFrame*. A *str* specifies the level name.

numeric\_only [boolean, default False] Include only float, int or boolean data.

### **Returns**

**Series or DataFrame** For each column/row the number of non-NA/null entries. If *level* is specified returns a *DataFrame*.

### See also:

Series.count number of non-NA elements in a Series

DataFrame.shape number of DataFrame rows and columns (including NA elements)

DataFrame.isna boolean same-sized DataFrame showing places of NA elements

# **Examples**

Constructing DataFrame from a dictionary:

```
>>> df = pd.DataFrame({"Person":
                       ["John", "Myla", None, "John", "Myla"],
. . .
                       "Age": [24., np.nan, 21., 33, 26],
. . .
                       "Single": [False, True, True, True, False] })
. . .
>>> df
  Person Age Single
     John 24.0
                 False
0
                  True
1
     Myla
           NaN
    None 21.0
2
                   True
3
     John 33.0
                   True
     Myla 26.0
                  False
```

Notice the uncounted NA values:

```
>>> df.count()
Person 4
Age 4
Single 5
dtype: int64
```

# Counts for each row:

Counts for one level of a MultiIndex:

# cov (min periods=None)

Compute pairwise covariance of columns, excluding NA/null values.

Compute the pairwise covariance among the series of a DataFrame. The returned data frame is the covariance matrix of the columns of the DataFrame.

Both NA and null values are automatically excluded from the calculation. (See the note below about bias from missing values.) A threshold can be set for the minimum number of observations for each value created. Comparisons with observations below this threshold will be returned as NaN.

This method is generally used for the analysis of time series data to understand the relationship between different measures across time.

### **Parameters**

**min\_periods** [int, optional] Minimum number of observations required per pair of columns to have a valid result.

#### Returns

**DataFrame** The covariance matrix of the series of the DataFrame.

See also:

```
pandas.Series.cov compute covariance with another Series
pandas.core.window.EWM.cov expoential weighted sample covariance
pandas.core.window.Expanding.cov expanding sample covariance
pandas.core.window.Rolling.cov rolling sample covariance
```

## **Notes**

Returns the covariance matrix of the DataFrame's time series. The covariance is normalized by N-1.

For DataFrames that have Series that are missing data (assuming that data is missing at random) the returned covariance matrix will be an unbiased estimate of the variance and covariance between the member Series.

However, for many applications this estimate may not be acceptable because the estimate covariance matrix is not guaranteed to be positive semi-definite. This could lead to estimate correlations having absolute values which are greater than one, and/or a non-invertible covariance matrix. See Estimation of covariance matrices for more details.

# **Examples**

```
>>> df = pd.DataFrame([(1, 2), (0, 3), (2, 0), (1, 1)],
...
columns=['dogs', 'cats'])
>>> df.cov()
```

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```
dogs cats
dogs 0.666667 -1.000000
cats -1.000000 1.666667
```

## Minimum number of periods

This method also supports an optional min\_periods keyword that specifies the required minimum number of non-NA observations for each column pair in order to have a valid result:

```
>>> np.random.seed(42)
>>> df = pd.DataFrame(np.random.randn(20, 3),
... columns=['a', 'b', 'c'])
>>> df.loc[df.index[:5], 'a'] = np.nan
>>> df.loc[df.index[5:10], 'b'] = np.nan
>>> df.cov(min_periods=12)

a b c
a 0.316741 NaN -0.150812
b NaN 1.248003 0.191417
c -0.150812 0.191417 0.895202
```

**cummax** (axis=None, skipna=True, \*args, \*\*kwargs)

Return cumulative maximum over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative maximum.

# **Parameters**

axis [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

#### **Returns**

**cummax** [Series or DataFrame]

See also:

pandas.core.window.Expanding.max Similar functionality but ignores NaN values.

DataFrame.max Return the maximum over DataFrame axis.

DataFrame.cummax Return cumulative maximum over DataFrame axis.

DataFrame.cummin Return cumulative minimum over DataFrame axis.

DataFrame.cumsum Return cumulative sum over DataFrame axis.

**DataFrame.cumprod** Return cumulative product over DataFrame axis.

# **Examples**

### **Series**

By default, NA values are ignored.

```
>>> s.cummax()
0 2.0
1 NaN
2 5.0
3 5.0
4 5.0
dtype: float64
```

To include NA values in the operation, use skipna=False

```
>>> s.cummax(skipna=False)
0 2.0
1 NaN
2 NaN
3 NaN
4 NaN
dtype: float64
```

### **DataFrame**

By default, iterates over rows and finds the maximum in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cummax()

A B

0 2.0 1.0

1 3.0 NaN

2 3.0 1.0
```

To iterate over columns and find the maximum in each row, use axis=1

```
>>> df.cummax(axis=1)

A B
0 2.0 2.0
1 3.0 NaN
2 1.0 1.0
```

```
cummin (axis=None, skipna=True, *args, **kwargs)
```

Return cumulative minimum over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative minimum.

### **Parameters**

axis [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

#### Returns

**cummin** [Series or DataFrame]

See also:

pandas.core.window.Expanding.min Similar functionality but ignores NaN values.

DataFrame.min Return the minimum over DataFrame axis.

DataFrame.cummax Return cumulative maximum over DataFrame axis.

DataFrame.cummin Return cumulative minimum over DataFrame axis.

DataFrame.cumsum Return cumulative sum over DataFrame axis.

DataFrame.cumprod Return cumulative product over DataFrame axis.

# **Examples**

### **Series**

```
>>> s = pd.Series([2, np.nan, 5, -1, 0])

>>> s

0    2.0

1    NaN

2    5.0

3    -1.0

4    0.0

dtype: float64
```

By default, NA values are ignored.

```
>>> s.cummin()
0 2.0
1 NaN
2 2.0
```

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```
3 -1.0
4 -1.0
dtype: float64
```

To include NA values in the operation, use skipna=False

```
>>> s.cummin(skipna=False)
0 2.0
1 NaN
2 NaN
3 NaN
4 NaN
dtype: float64
```

#### DataFrame

```
>>> df = pd.DataFrame([[2.0, 1.0],
... [3.0, np.nan],
... [1.0, 0.0]],
... columns=list('AB'))
>>> df
A B
0 2.0 1.0
1 3.0 NaN
2 1.0 0.0
```

By default, iterates over rows and finds the minimum in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cummin()

A B

0 2.0 1.0

1 2.0 NaN

2 1.0 0.0
```

To iterate over columns and find the minimum in each row, use axis=1

cumprod (axis=None, skipna=True, \*args, \*\*kwargs)

Return cumulative product over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative product.

# **Parameters**

axis [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

### **Returns**

**cumprod** [Series or DataFrame]

See also:

pandas.core.window.Expanding.prod Similar functionality but ignores NaN values.

**DataFrame.prod** Return the product over DataFrame axis.

**DataFrame.cummax** Return cumulative maximum over DataFrame axis.

DataFrame.cummin Return cumulative minimum over DataFrame axis.

DataFrame.cumsum Return cumulative sum over DataFrame axis.

**DataFrame.cumprod** Return cumulative product over DataFrame axis.

# **Examples**

#### **Series**

By default, NA values are ignored.

```
>>> s.cumprod()
0 2.0
1 NaN
2 10.0
3 -10.0
4 -0.0
dtype: float64
```

To include NA values in the operation, use skipna=False

```
>>> s.cumprod(skipna=False)

0 2.0

1 NaN

2 NaN

3 NaN

4 NaN

dtype: float64
```

### **DataFrame**

```
>>> df = pd.DataFrame([[2.0, 1.0],
... [3.0, np.nan],
... [1.0, 0.0]],
... columns=list('AB'))
>>> df
A B
```

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```
0 2.0 1.0
1 3.0 NaN
2 1.0 0.0
```

By default, iterates over rows and finds the product in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cumprod()

A B

0 2.0 1.0

1 6.0 NaN

2 6.0 0.0
```

To iterate over columns and find the product in each row, use axis=1

```
>>> df.cumprod(axis=1)

A B
0 2.0 2.0
1 3.0 NaN
2 1.0 0.0
```

cumsum(axis=None, skipna=True, \*args, \*\*kwargs)

Return cumulative sum over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative sum.

### **Parameters**

axis [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

### **Returns**

**cumsum** [Series or DataFrame]

See also:

pandas.core.window.Expanding.sum Similar functionality but ignores NaN values.

DataFrame.sum Return the sum over DataFrame axis.

DataFrame.cummax Return cumulative maximum over DataFrame axis.

DataFrame.cummin Return cumulative minimum over DataFrame axis.

DataFrame.cumsum Return cumulative sum over DataFrame axis.

DataFrame.cumprod Return cumulative product over DataFrame axis.

# **Examples**

Series

By default, NA values are ignored.

```
>>> s.cumsum()
0 2.0
1 NaN
2 7.0
3 6.0
4 6.0
dtype: float64
```

To include NA values in the operation, use skipna=False

```
>>> s.cumsum(skipna=False)
0 2.0
1 NaN
2 NaN
3 NaN
4 NaN
dtype: float64
```

# **DataFrame**

By default, iterates over rows and finds the sum in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cumsum()

A B

0 2.0 1.0
1 5.0 NaN
2 6.0 1.0
```

To iterate over columns and find the sum in each row, use axis=1

```
>>> df.cumsum(axis=1)

A B
0 2.0 3.0
1 3.0 NaN
2 1.0 1.0
```

**describe** (percentiles=None, include=None, exclude=None)

Generates descriptive statistics that summarize the central tendency, dispersion and shape of a dataset's distribution, excluding NaN values.

Analyzes both numeric and object series, as well as DataFrame column sets of mixed data types. The output will vary depending on what is provided. Refer to the notes below for more detail.

#### **Parameters**

**percentiles** [list-like of numbers, optional] The percentiles to include in the output. All should fall between 0 and 1. The default is [.25, .5, .75], which returns the 25th, 50th, and 75th percentiles.

**include** ['all', list-like of dtypes or None (default), optional] A white list of data types to include in the result. Ignored for Series. Here are the options:

- 'all' : All columns of the input will be included in the output.
- A list-like of dtypes: Limits the results to the provided data types. To limit the result to numeric types submit numpy.number. To limit it instead to object columns submit the numpy.object data type. Strings can also be used in the style of select\_dtypes (e.g. df.describe(include=['O'])). To select pandas categorical columns, use 'category'
- None (default): The result will include all numeric columns.

**exclude** [list-like of dtypes or None (default), optional,] A black list of data types to omit from the result. Ignored for Series. Here are the options:

- A list-like of dtypes: Excludes the provided data types from the result. To exclude numeric types submit numpy.number. To exclude object columns submit the data type numpy.object. Strings can also be used in the style of select\_dtypes (e.g. df.describe(include=['O'])). To exclude pandas categorical columns, use 'category'
- None (default): The result will exclude nothing.

### Returns

# summary: Series/DataFrame of summary statistics

# See also:

DataFrame.count, DataFrame.max, DataFrame.min, DataFrame.mean, DataFrame.std, DataFrame.select\_dtypes

#### Notes

For numeric data, the result's index will include count, mean, std, min, max as well as lower, 50 and upper percentiles. By default the lower percentile is 25 and the upper percentile is 75. The 50 percentile is the same as the median.

For object data (e.g. strings or timestamps), the result's index will include count, unique, top, and freq. The top is the most common value. The freq is the most common value's frequency. Timestamps also include the first and last items.

If multiple object values have the highest count, then the count and top results will be arbitrarily chosen from among those with the highest count.

For mixed data types provided via a DataFrame, the default is to return only an analysis of numeric columns. If the dataframe consists only of object and categorical data without any numeric columns,

the default is to return an analysis of both the object and categorical columns. If include='all' is provided as an option, the result will include a union of attributes of each type.

The *include* and *exclude* parameters can be used to limit which columns in a DataFrame are analyzed for the output. The parameters are ignored when analyzing a Series.

# **Examples**

Describing a numeric Series.

```
>>> s = pd.Series([1, 2, 3])
>>> s.describe()
count 3.0
        2.0
mean
std
        1.0
min
        1.0
25%
        1.5
50%
        2.0
75%
        2.5
        3.0
max
```

Describing a categorical Series.

```
>>> s = pd.Series(['a', 'a', 'b', 'c'])
>>> s.describe()
count    4
unique    3
top          a
freq    2
dtype: object
```

Describing a timestamp Series.

```
>>> s = pd.Series([
... np.datetime64("2000-01-01"),
    np.datetime64("2010-01-01"),
     np.datetime64("2010-01-01")
...])
>>> s.describe()
count
                           3
unique
top
        2010-01-01 00:00:00
freq
        2000-01-01 00:00:00
first
        2010-01-01 00:00:00
last
dtype: object
```

Describing a DataFrame. By default only numeric fields are returned.

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```
    std
    1.0

    min
    1.0

    25%
    1.5

    50%
    2.0

    75%
    2.5

    max
    3.0
```

Describing all columns of a DataFrame regardless of data type.

```
>>> df.describe(include='all')
    categorical numeric object
      3 3.0 3
count
            3
                 NaN
                         3
unique
top
            f
                 NaN
                         C
            1
                         1
freq
                  NaN
          NaN
NaN
                       NaN
                  2.0
mean
                  1.0
           NaN
std
                        NaN
min
           NaN
                  1.0
25%
           NaN
                  1.5
                        NaN
50%
           NaN
                   2.0
                        NaN
75%
           NaN
                   2.5
                        NaN
           NaN
                   3.0
                        NaN
max
```

Describing a column from a DataFrame by accessing it as an attribute.

```
>>> df.numeric.describe()
count 3.0
        2.0
mean
std
        1.0
min
        1.0
25%
        1.5
50%
        2.0
75%
        2.5
       3.0
max
Name: numeric, dtype: float64
```

Including only numeric columns in a DataFrame description.

```
>>> df.describe(include=[np.number])
      numeric
count
      3.0
mean
          2.0
std
          1.0
         1.0
min
25%
         1.5
50%
          2.0
75%
          2.5
          3.0
max
```

Including only string columns in a DataFrame description.

Including only categorical columns from a DataFrame description.

Excluding numeric columns from a DataFrame description.

Excluding object columns from a DataFrame description.

```
>>> df.describe(exclude=[np.object])
        categorical numeric
                 3
                        3.0
count
                 3
                        NaN
unique
                f
top
                        NaN
                1
freq
                        NaN
               NaN
                        2.0
mean
               NaN
                        1.0
std
min
               NaN
                        1.0
25%
               NaN
                        1.5
50%
                        2.0
               NaN
75%
                        2.5
               NaN
max
               NaN
                        3.0
```

# **diff** (*periods*=1, *axis*=0)

First discrete difference of element.

Calculates the difference of a DataFrame element compared with another element in the DataFrame (default is the element in the same column of the previous row).

### **Parameters**

```
periods [int, default 1] Periods to shift for calculating difference, accepts negative values.axis [{0 or 'index', 1 or 'columns'}, default 0] Take difference over rows (0) or columns (1).New in version 0.16.1..
```

#### Returns

diffed [DataFrame]

See also:

Series.diff First discrete difference for a Series.

DataFrame.pct\_change Percent change over given number of periods.

DataFrame.shift Shift index by desired number of periods with an optional time freq.

# **Examples**

# Difference with previous row

```
>>> df = pd.DataFrame({'a': [1, 2, 3, 4, 5, 6],
... 'b': [1, 1, 2, 3, 5, 8],
... 'c': [1, 4, 9, 16, 25, 36]})
>>> df

a b c
0 1 1 1
1 2 1 4
2 3 2 9
3 4 3 16
4 5 5 25
5 6 8 36
```

```
>>> df.diff()
    a    b    c
0    NaN    NaN    NaN
1    1.0    0.0    3.0
2    1.0    1.0    5.0
3    1.0    1.0    7.0
4    1.0    2.0    9.0
5    1.0    3.0    11.0
```

# Difference with previous column

```
>>> df.diff(axis=1)
    a    b    c
0 NaN 0.0 0.0
1 NaN -1.0 3.0
2 NaN -1.0 7.0
3 NaN -1.0 13.0
4 NaN 0.0 20.0
5 NaN 2.0 28.0
```

# Difference with 3rd previous row

```
>>> df.diff(periods=3)
    a    b    c
0 NaN NaN NaN NaN
1 NaN NaN NaN
2 NaN NaN NaN
3 3.0 2.0 15.0
4 3.0 4.0 21.0
5 3.0 6.0 27.0
```

### Difference with following row

```
>>> df.diff(periods=-1)
    a     b     c
0 -1.0     0.0     -3.0
1 -1.0 -1.0     -5.0
2 -1.0 -1.0     -7.0
3 -1.0 -2.0     -9.0
4 -1.0 -3.0 -11.0
5 NaN NaN NaN
```

```
div (other, axis='columns', level=None, fill value=None)
```

Floating division of dataframe and other, element-wise (binary operator truediv).

Equivalent to dataframe / other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

other [Series, DataFrame, or constant]

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

fill\_value [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

# Returns

result [DataFrame]

#### See also:

DataFrame.rtruediv

## **Notes**

Mismatched indices will be unioned together

# **Examples**

None

divide (other, axis='columns', level=None, fill\_value=None)

Floating division of dataframe and other, element-wise (binary operator truediv).

Equivalent to dataframe / other, but with support to substitute a fill\_value for missing data in one of the inputs.

# **Parameters**

```
other [Series, DataFrame, or constant]
```

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

#### **Returns**

result [DataFrame]

# See also:

DataFrame.rtruediv

### **Notes**

Mismatched indices will be unioned together

# **Examples**

None

#### dot (other)

Matrix multiplication with DataFrame or Series objects. Can also be called using *self* @ *other* in Python >= 3.5.

### **Parameters**

other [DataFrame or Series]

#### Returns

dot\_product [DataFrame or Series]

**drop** (*labels=None*, *axis=0*, *index=None*, *columns=None*, *level=None*, *inplace=False*, *errors='raise'*) Drop specified labels from rows or columns.

Remove rows or columns by specifying label names and corresponding axis, or by specifying directly index or column names. When using a multi-index, labels on different levels can be removed by specifying the level.

### **Parameters**

labels [single label or list-like] Index or column labels to drop.

**axis** [{0 or 'index', 1 or 'columns'}, default 0] Whether to drop labels from the index (0 or 'index') or columns (1 or 'columns').

index, columns [single label or list-like] Alternative to specifying axis (labels, axis=1 is equivalent to columns=labels).

New in version 0.21.0.

**level** [int or level name, optional] For MultiIndex, level from which the labels will be removed.

inplace [bool, default False] If True, do operation inplace and return None.

**errors** [{'ignore', 'raise'}, default 'raise'] If 'ignore', suppress error and only existing labels are dropped.

### **Returns**

**dropped** [pandas.DataFrame]

## Raises

**KeyError** If none of the labels are found in the selected axis

# See also:

DataFrame.loc Label-location based indexer for selection by label.

**DataFrame.dropna** Return DataFrame with labels on given axis omitted where (all or any) data are missing

DataFrame.drop\_duplicates Return DataFrame with duplicate rows removed, optionally only considering certain columns

**Series.drop** Return Series with specified index labels removed.

# **Examples**

```
>>> df = pd.DataFrame(np.arange(12).reshape(3,4),
... columns=['A', 'B', 'C', 'D'])
>>> df

A B C D

0 0 1 2 3

1 4 5 6 7

2 8 9 10 11
```

# Drop columns

```
>>> df.drop(['B', 'C'], axis=1)

A D
0 0 3
1 4 7
2 8 11
```

```
>>> df.drop(columns=['B', 'C'])

A D
0 0 3
1 4 7
2 8 11
```

# Drop a row by index

```
>>> df.drop([0, 1])
    A B C D
2 8 9 10 11
```

# Drop columns and/or rows of MultiIndex DataFrame

```
>>> midx = pd.MultiIndex(levels=[['lama', 'cow', 'falcon'],
                                 ['speed', 'weight', 'length']],
. . .
                        labels=[[0, 0, 0, 1, 1, 1, 2, 2, 2],
. . .
                                [0, 1, 2, 0, 1, 2, 0, 1, 2]])
>>> df = pd.DataFrame(index=midx, columns=['big', 'small'],
                      data=[[45, 30], [200, 100], [1.5, 1], [30, 20],
                           [250, 150], [1.5, 0.8], [320, 250],
. . .
                            [1, 0.8], [0.3, 0.2]])
. . .
>>> df
               big
                       small
       speed 45.0
                       30.0
lama
       weight 200.0 100.0
       length 1.5
                       1.0
COW
       speed 30.0
                       20.0
       weight 250.0 150.0
       length 1.5
                       0.8
falcon
       speed 320.0 250.0
       weight 1.0
                       0.8
       length 0.3
                       0.2
```

```
>>> df.drop(index='length', level=1)
                big
                        small
                45.0
                        30.0
lama
        speed
        weight 200.0
                       100.0
COW
        speed
                30.0
                        20.0
        weight 250.0
                        150.0
                320.0
                        250.0
falcon
        speed
        weight
               1.0
```

drop\_duplicates (subset=None, keep='first', inplace=False)

Return DataFrame with duplicate rows removed, optionally only considering certain columns

#### **Parameters**

**subset** [column label or sequence of labels, optional] Only consider certain columns for identifying duplicates, by default use all of the columns

**keep** [{'first', 'last', False}, default 'first']

- first: Drop duplicates except for the first occurrence.
- last: Drop duplicates except for the last occurrence.
- False: Drop all duplicates.

**inplace** [boolean, default False] Whether to drop duplicates in place or to return a copy

### Returns

# deduplicated [DataFrame]

**dropna** (axis=0, how='any', thresh=None, subset=None, inplace=False) Remove missing values.

See the User Guide for more on which values are considered missing, and how to work with missing data.

#### **Parameters**

**axis** [{0 or 'index', 1 or 'columns'}, default 0] Determine if rows or columns which contain missing values are removed.

- 0, or 'index': Drop rows which contain missing values.
- 1, or 'columns': Drop columns which contain missing value.

Deprecated since version 0.23.0:: Pass tuple or list to drop on multiple

axes.

**how** [{'any', 'all'}, default 'any'] Determine if row or column is removed from DataFrame, when we have at least one NA or all NA.

- 'any': If any NA values are present, drop that row or column.
- 'all' : If all values are NA, drop that row or column.

**thresh** [int, optional] Require that many non-NA values.

**subset** [array-like, optional] Labels along other axis to consider, e.g. if you are dropping rows these would be a list of columns to include.

inplace [bool, default False] If True, do operation inplace and return None.

### **Returns**

DataFrame DataFrame with NA entries dropped from it.

See also:

DataFrame.isna Indicate missing values.

DataFrame.notna Indicate existing (non-missing) values.

DataFrame.fillna Replace missing values.

Series.dropna Drop missing values.

Index.dropna Drop missing indices.

# **Examples**

```
>>> df = pd.DataFrame({"name": ['Alfred', 'Batman', 'Catwoman'],
                      "toy": [np.nan, 'Batmobile', 'Bullwhip'],
                      "born": [pd.NaT, pd.Timestamp("1940-04-25"),
. . .
                               pd.NaT] })
. . .
>>> df
                           born
      name
                  toy
0
    Alfred
                 NaN
                            NaT
1
    Batman Batmobile 1940-04-25
  Catwoman Bullwhip
                             NaT
```

Drop the rows where at least one element is missing.

```
>>> df.dropna()
name toy born
1 Batman Batmobile 1940-04-25
```

Drop the columns where at least one element is missing.

Drop the rows where all elements are missing.

```
>>> df.dropna(how='all')
name toy born

O Alfred NaN NaT

Batman Batmobile 1940-04-25

Catwoman Bullwhip NaT
```

Keep only the rows with at least 2 non-NA values.

```
>>> df.dropna(thresh=2)
name toy born
1 Batman Batmobile 1940-04-25
2 Catwoman Bullwhip NaT
```

Define in which columns to look for missing values.

Keep the DataFrame with valid entries in the same variable.

```
>>> df.dropna(inplace=True)
>>> df
name toy born
1 Batman Batmobile 1940-04-25
```

# dtypes

Return the dtypes in the DataFrame.

This returns a Series with the data type of each column. The result's index is the original DataFrame's columns. Columns with mixed types are stored with the object dtype. See the User Guide for more.

### **Returns**

**pandas.Series** The data type of each column.

See also:

pandas.DataFrame.ftypes dtype and sparsity information.

# **Examples**

## duplicated (subset=None, keep='first')

Return boolean Series denoting duplicate rows, optionally only considering certain columns

## **Parameters**

**subset** [column label or sequence of labels, optional] Only consider certain columns for identifying duplicates, by default use all of the columns

```
keep [{'first', 'last', False}, default 'first']
```

- first: Mark duplicates as True except for the first occurrence.
- last: Mark duplicates as True except for the last occurrence.

• False: Mark all duplicates as True.

### **Returns**

```
duplicated [Series]
```

# empty

Indicator whether DataFrame is empty.

True if DataFrame is entirely empty (no items), meaning any of the axes are of length 0.

# Returns

bool If DataFrame is empty, return True, if not return False.

# See also:

```
pandas.Series.dropna, pandas.DataFrame.dropna
```

### **Notes**

If DataFrame contains only NaNs, it is still not considered empty. See the example below.

# **Examples**

An example of an actual empty DataFrame. Notice the index is empty:

```
>>> df_empty = pd.DataFrame({'A' : []})
>>> df_empty
Empty DataFrame
Columns: [A]
Index: []
>>> df_empty.empty
True
```

If we only have NaNs in our DataFrame, it is not considered empty! We will need to drop the NaNs to make the DataFrame empty:

```
eq (other, axis='columns', level=None)
```

Wrapper for flexible comparison methods eq

```
equals (other)
```

Determines if two NDFrame objects contain the same elements. NaNs in the same location are considered equal.

```
eval (expr, inplace=False, **kwargs)
```

Evaluate a string describing operations on DataFrame columns.

Operates on columns only, not specific rows or elements. This allows *eval* to run arbitrary code, which can make you vulnerable to code injection if you pass user input to this function.

### **Parameters**

**expr** [str] The expression string to evaluate.

**inplace** [bool, default False] If the expression contains an assignment, whether to perform the operation inplace and mutate the existing DataFrame. Otherwise, a new DataFrame is returned.

New in version 0.18.0..

**kwargs** [dict] See the documentation for eval() for complete details on the keyword arguments accepted by query().

# Returns

ndarray, scalar, or pandas object The result of the evaluation.

### See also:

**DataFrame.query** Evaluates a boolean expression to query the columns of a frame.

**DataFrame.assign** Can evaluate an expression or function to create new values for a column.

pandas.eval Evaluate a Python expression as a string using various backends.

# **Notes**

For more details see the API documentation for eval (). For detailed examples see enhancing performance with eval.

# **Examples**

```
>>> df = pd.DataFrame({'A': range(1, 6), 'B': range(10, 0, -2)})
>>> df
  Α
      В
0
  1
     10
  2
       8
  3
       6
3
  4
       4
  5
       2
>>> df.eval('A + B')
    11
     10
      9
3
      8
dtype: int64
```

Assignment is allowed though by default the original DataFrame is not modified.

```
>>> df.eval('C = A + B')
   Α
       В
           C
      10
          11
  1
  2
       8
          10
  3
       6
            9
3
            8
   4
       4
4
   5
       2
            7
```

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```
>>> df
   Α
       В
0
      10
  1
  2
       8
  3
        6
3
   4
        4
4
   5
        2
```

Use inplace=True to modify the original DataFrame.

```
>>> df.eval('C = A + B', inplace=True)
>>> df
  Α
     В
          С
0
  1
     10 11
       8
          10
2
           9
       6
3
           8
  4
       4
4
   5
       2
           7
```

**ewm** (com=None, span=None, halflife=None, alpha=None, min\_periods=0, adjust=True, ignore\_na=False, axis=0)

Provides exponential weighted functions

New in version 0.18.0.

#### **Parameters**

```
 \begin{tabular}{ll} \textbf{com} & [float, optional] Specify decay in terms of center of mass, $\alpha=1/(1+com)$, for $com \geq 0$ \\ \hline \\ 0 & \\ \hline \\ 0 & \\ \hline \\ \end{array}
```

```
span [float, optional] Specify decay in terms of span, \alpha = 2/(span + 1), for span > 1
```

**halflife** [float, optional] Specify decay in terms of half-life,  $\alpha = 1 - exp(log(0.5)/halflife)$ , for halflife > 0

**alpha** [float, optional] Specify smoothing factor  $\alpha$  directly,  $0 < \alpha \le 1$ 

New in version 0.18.0.

**min\_periods** [int, default 0] Minimum number of observations in window required to have a value (otherwise result is NA).

**adjust** [boolean, default True] Divide by decaying adjustment factor in beginning periods to account for imbalance in relative weightings (viewing EWMA as a moving average)

**ignore\_na** [boolean, default False] Ignore missing values when calculating weights; specify True to reproduce pre-0.15.0 behavior

### **Returns**

a Window sub-classed for the particular operation

## See also:

```
rolling Provides rolling window calculations
```

**expanding** Provides expanding transformations.

## **Notes**

Exactly one of center of mass, span, half-life, and alpha must be provided. Allowed values and relationship between the parameters are specified in the parameter descriptions above; see the link at the end of this section for a detailed explanation.

When adjust is True (default), weighted averages are calculated using weights (1-alpha)\*\*(n-1), (1-alpha)\*\*(n-2), ..., 1-alpha, 1.

When adjust is False, weighted averages are calculated recursively as: weighted\_average[0] = arg[0]; weighted\_average[i] = (1-alpha)\*weighted\_average[i-1] + alpha\*arg[i].

When ignore\_na is False (default), weights are based on absolute positions. For example, the weights of x and y used in calculating the final weighted average of [x, None, y] are (1-alpha)\*\*2 and 1 (if adjust is True), and (1-alpha)\*\*2 and alpha (if adjust is False).

When ignore\_na is True (reproducing pre-0.15.0 behavior), weights are based on relative positions. For example, the weights of x and y used in calculating the final weighted average of [x, None, y] are 1-alpha and 1 (if adjust is True), and 1-alpha and alpha (if adjust is False).

More details can be found at http://pandas.pydata.org/pandas-docs/stable/computation.html# exponentially-weighted-windows

# **Examples**

```
>>> df = DataFrame({'B': [0, 1, 2, np.nan, 4]})
B
0 0.0
1 1.0
2 2.0
3 NaN
4 4.0
```

```
>>> df.ewm(com=0.5).mean()

B
0 0.000000
1 0.750000
2 1.615385
3 1.615385
4 3.670213
```

**expanding** (*min\_periods*=1, *center*=*False*, *axis*=0)

Provides expanding transformations.

New in version 0.18.0.

## **Parameters**

**min\_periods** [int, default 1] Minimum number of observations in window required to have a value (otherwise result is NA).

**center** [boolean, default False] Set the labels at the center of the window.

axis [int or string, default 0]

## Returns

a Window sub-classed for the particular operation

See also:

rolling Provides rolling window calculations

ewm Provides exponential weighted functions

### **Notes**

By default, the result is set to the right edge of the window. This can be changed to the center of the window by setting center=True.

### **Examples**

```
>>> df = DataFrame({'B': [0, 1, 2, np.nan, 4]})
B
0 0.0
1 1.0
2 2.0
3 NaN
4 4.0
```

```
>>> df.expanding(2).sum()

B

NaN

1 1.0

2 3.0

3 3.0

4 7.0
```

**ffill** (axis=None, inplace=False, limit=None, downcast=None)
Synonym for DataFrame.fillna (method='ffill')

Fill NA/NaN values using the specified method

### **Parameters**

value [scalar, dict, Series, or DataFrame] Value to use to fill holes (e.g. 0), alternately a dict/Series/DataFrame of values specifying which value to use for each index (for a Series) or column (for a DataFrame). (values not in the dict/Series/DataFrame will not be filled). This value cannot be a list.

method [{'backfill', 'bfill', 'pad', 'ffill', None}, default None] Method to use for filling holes in reindexed Series pad / ffill: propagate last valid observation forward to next valid backfill / bfill: use NEXT valid observation to fill gap

```
axis [{0 or 'index', 1 or 'columns'}]
```

**inplace** [boolean, default False] If True, fill in place. Note: this will modify any other views on this object, (e.g. a no-copy slice for a column in a DataFrame).

**limit** [int, default None] If method is specified, this is the maximum number of consecutive NaN values to forward/backward fill. In other words, if there is a gap with more than this number of consecutive NaNs, it will only be partially filled. If method is not specified, this is the maximum number of entries along the entire axis where NaNs will be filled. Must be greater than 0 if not None.

**downcast** [dict, default is None] a dict of item->dtype of what to downcast if possible, or the string 'infer' which will try to downcast to an appropriate equal type (e.g. float64 to int64 if possible)

### **Returns**

filled [DataFrame]

### See also:

interpolate Fill NaN values using interpolation.

```
reindex, asfreq
```

## **Examples**

```
>>> df = pd.DataFrame([[np.nan, 2, np.nan, 0],
                       [3, 4, np.nan, 1],
. . .
                       [np.nan, np.nan, np.nan, 5],
. . .
                       [np.nan, 3, np.nan, 4]],
. . .
                       columns=list('ABCD'))
. . .
>>> df
        B C D
    Α
  NaN 2.0 NaN 0
  3.0 4.0 NaN
  NaN NaN NaN 5
  NaN 3.0 NaN 4
```

Replace all NaN elements with 0s.

```
>>> df.fillna(0)

A B C D

0 0.0 2.0 0.0 0

1 3.0 4.0 0.0 1

2 0.0 0.0 0.0 5

3 0.0 3.0 0.0 4
```

We can also propagate non-null values forward or backward.

```
>>> df.fillna(method='ffill')

A B C D

0 NaN 2.0 NaN 0

1 3.0 4.0 NaN 1

2 3.0 4.0 NaN 5

3 3.0 3.0 NaN 4
```

Replace all NaN elements in column 'A', 'B', 'C', and 'D', with 0, 1, 2, and 3 respectively.

Only replace the first NaN element.

```
>>> df.fillna(value=values, limit=1)

A B C D

0 0.0 2.0 2.0 0

1 3.0 4.0 NaN 1

2 NaN 1.0 NaN 5

3 NaN 3.0 NaN 4
```

## **filter** (*items=None*, *like=None*, *regex=None*, *axis=None*)

Subset rows or columns of dataframe according to labels in the specified index.

Note that this routine does not filter a dataframe on its contents. The filter is applied to the labels of the index.

### **Parameters**

```
items [list-like] List of info axis to restrict to (must not all be present)
```

**like** [string] Keep info axis where "arg in col == True"

**regex** [string (regular expression)] Keep info axis with re.search(regex, col) == True

**axis** [int or string axis name] The axis to filter on. By default this is the info axis, 'index' for Series, 'columns' for DataFrame

### **Returns**

## same type as input object

### See also:

```
pandas.DataFrame.loc
```

### **Notes**

The items, like, and regex parameters are enforced to be mutually exclusive.

 ${\tt axis}$  defaults to the info axis that is used when indexing with  $[\,]\,.$ 

## **Examples**

```
>>> df
one two three
mouse 1 2 3
rabbit 4 5 6
```

```
>>> # select columns by name
>>> df.filter(items=['one', 'three'])
one three
mouse 1 3
rabbit 4 6
```

```
>>> # select columns by regular expression
>>> df.filter(regex='e$', axis=1)
one three
mouse 1 3
rabbit 4 6
```

```
>>> # select rows containing 'bbi'
>>> df.filter(like='bbi', axis=0)
one two three
rabbit 4 5 6
```

### first (offset)

Convenience method for subsetting initial periods of time series data based on a date offset.

### **Parameters**

offset [string, DateOffset, dateutil.relativedelta]

#### Returns

subset [type of caller]

### **Raises**

TypeError If the index is not a DatetimeIndex

### See also:

last Select final periods of time series based on a date offset

at\_time Select values at a particular time of the day

between\_time Select values between particular times of the day

## **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='2D')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 1
2018-04-11 2
2018-04-13 3
2018-04-15 4
```

Get the rows for the first 3 days:

```
>>> ts.first('3D')

A
2018-04-09 1
2018-04-11 2
```

Notice the data for 3 first calender days were returned, not the first 3 days observed in the dataset, and therefore data for 2018-04-13 was not returned.

## first\_valid\_index()

Return index for first non-NA/null value.

# Returns

scalar [type of index]

## **Notes**

If all elements are non-NA/null, returns None. Also returns None for empty NDFrame.

**floordiv** (other, axis='columns', level=None, fill\_value=None)

Integer division of dataframe and other, element-wise (binary operator floordiv).

Equivalent to dataframe // other, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

other [Series, DataFrame, or constant]

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

fill\_value [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

## Returns

result [DataFrame]

#### See also:

DataFrame.rfloordiv

### **Notes**

Mismatched indices will be unioned together

### **Examples**

None

classmethod from\_csv (path, header=0, sep=', ', index\_col=0, parse\_dates=True, encoding=None, tupleize\_cols=None, infer\_datetime\_format=False)

Read CSV file.

Deprecated since version 0.21.0: Use pandas.read\_csv() instead.

It is preferable to use the more powerful pandas.read\_csv() for most general purposes, but from\_csv makes for an easy roundtrip to and from a file (the exact counterpart of to\_csv), especially with a DataFrame of time series data.

This method only differs from the preferred pandas.read\_csv() in some defaults:

- index col is 0 instead of None (take first column as index by default)
- parse\_dates is True instead of False (try parsing the index as datetime by default)

So a pd.DataFrame.from\_csv(path) can be replaced by pd.read\_csv(path, index\_col=0, parse\_dates=True).

### **Parameters**

```
path [string file path or file handle / StringIO]header [int, default 0] Row to use as header (skip prior rows)sep [string, default ','] Field delimiter
```

index\_col [int or sequence, default 0] Column to use for index. If a sequence is given, a MultiIndex is used. Different default from read table

parse\_dates [boolean, default True] Parse dates. Different default from read\_table

**tupleize\_cols** [boolean, default False] write multi\_index columns as a list of tuples (if True) or new (expanded format) if False)

**infer\_datetime\_format: boolean, default False** If True and *parse\_dates* is True for a column, try to infer the datetime format based on the first datetime string. If the format can be inferred, there often will be a large parsing speed-up.

### **Returns**

y [DataFrame]

## See also:

```
pandas.read_csv
```

classmethod from\_dict (data, orient='columns', dtype=None, columns=None)

Construct DataFrame from dict of array-like or dicts.

Creates DataFrame object from dictionary by columns or by index allowing dtype specification.

#### **Parameters**

```
data [dict] Of the form {field : array-like} or {field : dict}.
```

orient [{'columns', 'index'}, default 'columns'] The "orientation" of the data. If the keys of the passed dict should be the columns of the resulting DataFrame, pass 'columns' (default). Otherwise if the keys should be rows, pass 'index'.

dtype [dtype, default None] Data type to force, otherwise infer.

columns [list, default None] Column labels to use when orient='index'. Raises a
 ValueError if used with orient='columns'.

New in version 0.23.0.

#### Returns

### pandas.DataFrame

See also:

DataFrame.from\_records DataFrame from ndarray (structured dtype), list of tuples, dict, or DataFrame

**DataFrame** DataFrame object creation using constructor

## **Examples**

By default the keys of the dict become the DataFrame columns:

Specify orient='index' to create the DataFrame using dictionary keys as rows:

When using the 'index' orientation, the column names can be specified manually:

```
>>> pd.DataFrame.from_dict(data, orient='index',
... columns=['A', 'B', 'C', 'D'])

A B C D

row_1 3 2 1 0

row_2 a b c d
```

#### classmethod from items (items, columns=None, orient='columns')

Construct a dataframe from a list of tuples

Deprecated since version 0.23.0: from\_items is deprecated and will be removed in a future version. Use DataFrame.from\_dict(dict(items)) instead. DataFrame.from\_dict(OrderedDict(items)) may be used to preserve the key order.

Convert (key, value) pairs to DataFrame. The keys will be the axis index (usually the columns, but depends on the specified orientation). The values should be arrays or Series.

#### **Parameters**

items [sequence of (key, value) pairs] Values should be arrays or Series.

columns [sequence of column labels, optional] Must be passed if orient='index'.

**orient** [{'columns', 'index'}, default 'columns'] The "orientation" of the data. If the keys of the input correspond to column labels, pass 'columns' (default). Otherwise if the keys correspond to the index, pass 'index'.

#### Returns

**frame** [DataFrame]

Convert structured or record ndarray to DataFrame

## **Parameters**

data [ndarray (structured dtype), list of tuples, dict, or DataFrame]

**index** [string, list of fields, array-like] Field of array to use as the index, alternately a specific set of input labels to use

exclude [sequence, default None] Columns or fields to exclude

**columns** [sequence, default None] Column names to use. If the passed data do not have names associated with them, this argument provides names for the columns. Otherwise this argument indicates the order of the columns in the result (any names not found in the data will become all-NA columns)

**coerce\_float** [boolean, default False] Attempt to convert values of non-string, non-numeric objects (like decimal.Decimal) to floating point, useful for SQL result sets

#### Returns

df [DataFrame]

## ftypes

Return the ftypes (indication of sparse/dense and dtype) in DataFrame.

This returns a Series with the data type of each column. The result's index is the original DataFrame's columns. Columns with mixed types are stored with the object dtype. See the User Guide for more.

### Returns

pandas. Series The data type and indication of sparse/dense of each column.

See also:

```
pandas.DataFrame.dtypes Series with just dtype information.
```

pandas. SparseDataFrame Container for sparse tabular data.

### **Notes**

Sparse data should have the same dtypes as its dense representation.

## **Examples**

```
>>> import numpy as np
>>> arr = np.random.RandomState(0).randn(100, 4)
>>> arr[arr < .8] = np.nan
>>> pd.DataFrame(arr).ftypes
0    float64:dense
1    float64:dense
2    float64:dense
3    float64:dense
dtype: object
```

```
>>> pd.SparseDataFrame(arr).ftypes
0    float64:sparse
1    float64:sparse
2    float64:sparse
3    float64:sparse
dtype: object
```

```
ge (other, axis='columns', level=None)
```

Wrapper for flexible comparison methods ge

```
get (key, default=None)
```

Get item from object for given key (DataFrame column, Panel slice, etc.). Returns default value if not found.

### **Parameters**

key [object]

## Returns

value [type of items contained in object]

### get\_dtype\_counts()

Return counts of unique dtypes in this object.

## Returns

**dtype** [Series] Series with the count of columns with each dtype.

### See also:

dtypes Return the dtypes in this object.

## **Examples**

```
>>> a = [['a', 1, 1.0], ['b', 2, 2.0], ['c', 3, 3.0]]
>>> df = pd.DataFrame(a, columns=['str', 'int', 'float'])
>>> df
str int float
0 a 1 1.0
1 b 2 2.0
2 c 3 3.0
```

```
>>> df.get_dtype_counts()
float64    1
int64    1
object    1
dtype: int64
```

## get\_ftype\_counts()

Return counts of unique ftypes in this object.

Deprecated since version 0.23.0.

This is useful for SparseDataFrame or for DataFrames containing sparse arrays.

### Returns

**dtype** [Series] Series with the count of columns with each type and sparsity (dense/sparse)

# See also:

ftypes Return ftypes (indication of sparse/dense and dtype) in this object.

## **Examples**

```
>>> a = [['a', 1, 1.0], ['b', 2, 2.0], ['c', 3, 3.0]]
>>> df = pd.DataFrame(a, columns=['str', 'int', 'float'])
>>> df
    str int float
0 a 1 1.0
1 b 2 2.0
2 c 3 3.0
```

```
>>> df.get_ftype_counts()
float64:dense    1
int64:dense    1
object:dense    1
dtype: int64
```

# get\_value (index, col, takeable=False)

Quickly retrieve single value at passed column and index

Deprecated since version 0.21.0: Use .at[] or .iat[] accessors instead.

### **Parameters**

```
index [row label]col [column label]
```

**takeable** [interpret the index/col as indexers, default False]

#### Returns

value [scalar value]

### get\_values()

Return an ndarray after converting sparse values to dense.

This is the same as .values for non-sparse data. For sparse data contained in a *pandas.SparseArray*, the data are first converted to a dense representation.

### Returns

numpy.ndarray Numpy representation of DataFrame

### See also:

values Numpy representation of DataFrame.

pandas. SparseArray Container for sparse data.

# **Examples**

```
>>> df = pd.DataFrame({'a': [1, 2], 'b': [True, False],
... 'c': [1.0, 2.0]})
>>> df
a b c
0 1 True 1.0
1 2 False 2.0
```

```
>>> df.get_values() array([[1, True, 1.0], [2, False, 2.0]], dtype=object)
```

Group series using mapper (dict or key function, apply given function to group, return result as series) or by a series of columns.

### **Parameters**

by [mapping, function, label, or list of labels] Used to determine the groups for the groupby. If by is a function, it's called on each value of the object's index. If a dict or Series is passed, the Series or dict VALUES will be used to determine the groups (the Series' values are first aligned; see <code>.align()</code> method). If an ndarray is passed, the values are used as-is determine the groups. A label or list of labels may be passed to group by the columns in <code>self</code>. Notice that a tuple is interpreted a (single) key.

axis [int, default 0]

**level** [int, level name, or sequence of such, default None] If the axis is a MultiIndex (hierarchical), group by a particular level or levels

**as\_index** [boolean, default True] For aggregated output, return object with group labels as the index. Only relevant for DataFrame input. as\_index=False is effectively "SQL-style" grouped output

**sort** [boolean, default True] Sort group keys. Get better performance by turning this off. Note this does not influence the order of observations within each group. groupby preserves the order of rows within each group.

group\_keys [boolean, default True] When calling apply, add group keys to index to identify
pieces

**squeeze** [boolean, default False] reduce the dimensionality of the return type if possible, otherwise return a consistent type

**observed** [boolean, default False] This only applies if any of the groupers are Categoricals If True: only show observed values for categorical groupers. If False: show all values for categorical groupers.

New in version 0.23.0.

### Returns

## GroupBy object

### See also:

**resample** Convenience method for frequency conversion and resampling of time series.

### **Notes**

See the user guide for more.

## **Examples**

DataFrame results

```
>>> data.groupby(func, axis=0).mean()
>>> data.groupby(['col1', 'col2'])['col3'].mean()
```

DataFrame with hierarchical index

```
>>> data.groupby(['col1', 'col2']).mean()
```

gt (other, axis='columns', level=None)

Wrapper for flexible comparison methods gt

#### head(n=5)

Return the first *n* rows.

This function returns the first *n* rows for the object based on position. It is useful for quickly testing if your object has the right type of data in it.

## **Parameters**

n [int, default 5] Number of rows to select.

#### Returns

**obj\_head** [type of caller] The first *n* rows of the caller object.

See also:

pandas.DataFrame.tail Returns the last n rows.

## **Examples**

```
>>> df = pd.DataFrame({'animal':['alligator', 'bee', 'falcon', 'lion',
                        'monkey', 'parrot', 'shark', 'whale', 'zebra']})
>>> df
      animal
0
  alligator
1
         bee
2
      falcon
3
        lion
4
      monkey
5
      parrot
6
       shark
       whale
8
       zebra
```

### Viewing the first 5 lines

```
>>> df.head()
    animal
0 alligator
1 bee
2 falcon
3 lion
4 monkey
```

Viewing the first *n* lines (three in this case)

```
>>> df.head(3)
        animal
0 alligator
1      bee
2     falcon
```

hist (column=None, by=None, grid=True, xlabelsize=None, xrot=None, ylabelsize=None, yrot=None, ax=None, sharex=False, sharey=False, figsize=None, layout=None, bins=10, \*\*kwds)

Make a histogram of the DataFrame's.

A histogram is a representation of the distribution of data. This function calls matplotlib.pyplot. hist(), on each series in the DataFrame, resulting in one histogram per column.

#### **Parameters**

data [DataFrame] The pandas object holding the data.

column [string or sequence] If passed, will be used to limit data to a subset of columns.

by [object, optional] If passed, then used to form histograms for separate groups.

grid [boolean, default True] Whether to show axis grid lines.

**xlabelsize** [int, default None] If specified changes the x-axis label size.

**xrot** [float, default None] Rotation of x axis labels. For example, a value of 90 displays the x labels rotated 90 degrees clockwise.

ylabelsize [int, default None] If specified changes the y-axis label size.

**yrot** [float, default None] Rotation of y axis labels. For example, a value of 90 displays the y labels rotated 90 degrees clockwise.

**ax** [Matplotlib axes object, default None] The axes to plot the histogram on.

**sharex** [boolean, default True if ax is None else False] In case subplots=True, share x axis and set some x axis labels to invisible; defaults to True if ax is None otherwise False if an ax is passed in. Note that passing in both an ax and sharex=True will alter all x axis labels for all subplots in a figure.

**sharey** [boolean, default False] In case subplots=True, share y axis and set some y axis labels to invisible.

**figsize** [tuple] The size in inches of the figure to create. Uses the value in *mat-plotlib.rcParams* by default.

**layout** [tuple, optional] Tuple of (rows, columns) for the layout of the histograms.

**bins** [integer or sequence, default 10] Number of histogram bins to be used. If an integer is given, bins + 1 bin edges are calculated and returned. If bins is a sequence, gives bin edges, including left edge of first bin and right edge of last bin. In this case, bins is returned unmodified.

\*\*kwds All other plotting keyword arguments to be passed to matplotlib.pyplot. hist().

## Returns

axes [matplotlib.AxesSubplot or numpy.ndarray of them]

See also:

matplotlib.pyplot.hist Plot a histogram using matplotlib.

## **Examples**

## iat

Access a single value for a row/column pair by integer position.

Similar to iloc, in that both provide integer-based lookups. Use iat if you only need to get or set a single value in a DataFrame or Series.

## Raises

**IndexError** When integer position is out of bounds

See also:

DataFrame.at Access a single value for a row/column label pair

DataFrame.loc Access a group of rows and columns by label(s)

DataFrame.iloc Access a group of rows and columns by integer position(s)

## **Examples**

```
>>> df = pd.DataFrame([[0, 2, 3], [0, 4, 1], [10, 20, 30]],
                      columns=['A', 'B', 'C'])
>>> df
           С
       В
   Α
0
   0
       2
           3
   0
        4
            1
  10
      20
           30
```

Get value at specified row/column pair

```
>>> df.iat[1, 2]
1
```

Set value at specified row/column pair

```
>>> df.iat[1, 2] = 10
>>> df.iat[1, 2]
10
```

Get value within a series

```
>>> df.loc[0].iat[1]
2
```

idxmax (axis=0, skipna=True)

Return index of first occurrence of maximum over requested axis. NA/null values are excluded.

### **Parameters**

axis [{0 or 'index', 1 or 'columns'}, default 0] 0 or 'index' for row-wise, 1 or 'columns' for column-wise

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

#### Returns

idxmax [Series]

### **Raises**

# ValueError

• If the row/column is empty

#### See also:

Series.idxmax

### **Notes**

This method is the DataFrame version of ndarray.argmax.

### idxmin (axis=0, skipna=True)

Return index of first occurrence of minimum over requested axis. NA/null values are excluded.

#### **Parameters**

axis [{0 or 'index', 1 or 'columns'}, default 0] 0 or 'index' for row-wise, 1 or 'columns' for column-wise

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

### Returns

idxmin [Series]

#### Raises

### ValueError

• If the row/column is empty

### See also:

Series.idxmin

#### **Notes**

This method is the DataFrame version of ndarray.argmin.

### iloc

Purely integer-location based indexing for selection by position.

.iloc[] is primarily integer position based (from 0 to length-1 of the axis), but may also be used with a boolean array.

Allowed inputs are:

- An integer, e.g. 5.
- A list or array of integers, e.g. [4, 3, 0].
- A slice object with ints, e.g. 1:7.
- · A boolean array.
- A callable function with one argument (the calling Series, DataFrame or Panel) and that returns valid output for indexing (one of the above)
- .iloc will raise IndexError if a requested indexer is out-of-bounds, except *slice* indexers which allow out-of-bounds indexing (this conforms with python/numpy *slice* semantics).

See more at Selection by Position

#### index

The index (row labels) of the DataFrame.

## infer\_objects()

Attempt to infer better dtypes for object columns.

Attempts soft conversion of object-dtyped columns, leaving non-object and unconvertible columns unchanged. The inference rules are the same as during normal Series/DataFrame construction.

New in version 0.21.0.

## Returns

converted [same type as input object]

See also:

pandas.to\_datetime Convert argument to datetime.

pandas.to\_timedelta Convert argument to timedelta.

pandas.to\_numeric Convert argument to numeric typeR

## **Examples**

```
>>> df = pd.DataFrame({"A": ["a", 1, 2, 3]})
>>> df = df.iloc[1:]
>>> df
    A
1    1
2    2
3    3
```

```
>>> df.dtypes
A object
dtype: object
```

```
>>> df.infer_objects().dtypes
A int64
dtype: object
```

**info** (*verbose=None*, *buf=None*, *max\_cols=None*, *memory\_usage=None*, *null\_counts=None*) Print a concise summary of a DataFrame.

This method prints information about a DataFrame including the index dtype and column dtypes, non-null values and memory usage.

### **Parameters**

**verbose** [bool, optional] Whether to print the full summary. By default, the setting in pandas.options.display.max\_info\_columns is followed.

**buf** [writable buffer, defaults to sys.stdout] Where to send the output. By default, the output is printed to sys.stdout. Pass a writable buffer if you need to further process the output.

max\_cols [int, optional] When to switch from the verbose to the truncated output. If the DataFrame has more than *max\_cols* columns, the truncated output is used. By default, the setting in pandas.options.display.max\_info\_columns is used.

memory\_usage [bool, str, optional] Specifies whether total memory usage of the DataFrame
elements (including the index) should be displayed. By default, this follows the pandas.
options.display.memory\_usage setting.

True always show memory usage. False never shows memory usage. A value of 'deep' is equivalent to "True with deep introspection". Memory usage is shown in human-readable units (base-2 representation). Without deep introspection a memory estimation is made based in column dtype and number of rows assuming values consume the same memory

amount for corresponding dtypes. With deep memory introspection, a real memory usage calculation is performed at the cost of computational resources.

null\_counts [bool, optional] Whether to show the non-null counts. By default, this is shown only if the frame is smaller than pandas.options.display.max\_info\_rows and pandas.options.display.max\_info\_columns. A value of True always shows the counts, and False never shows the counts.

#### **Returns**

None This method prints a summary of a DataFrame and returns None.

See also:

DataFrame.describe Generate descriptive statistics of DataFrame columns.

DataFrame.memory\_usage Memory usage of DataFrame columns.

## **Examples**

```
>>> int_values = [1, 2, 3, 4, 5]
>>> text_values = ['alpha', 'beta', 'gamma', 'delta', 'epsilon']
>>> float_values = [0.0, 0.25, 0.5, 0.75, 1.0]
>>> df = pd.DataFrame({"int_col": int_values, "text_col": text_values,
                     "float_col": float_values})
>>> df
  int_col text_col float_col
                    0.00
0
       1 alpha
        2
             beta
                        0.25
2
        3
           gamma
                        0.50
3
        4
            delta
                         0.75
4
        5 epsilon
                         1.00
```

## Prints information of all columns:

```
>>> df.info(verbose=True)
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 5 entries, 0 to 4
Data columns (total 3 columns):
int_col 5 non-null int64
text_col 5 non-null object
float_col 5 non-null float64
dtypes: float64(1), int64(1), object(1)
memory usage: 200.0+ bytes
```

Prints a summary of columns count and its dtypes but not per column information:

```
>>> df.info(verbose=False)
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 5 entries, 0 to 4
Columns: 3 entries, int_col to float_col
dtypes: float64(1), int64(1), object(1)
memory usage: 200.0+ bytes
```

Pipe output of DataFrame.info to buffer instead of sys.stdout, get buffer content and writes to a text file:

```
>>> import io
>>> buffer = io.StringIO()
>>> df.info(buf=buffer)
>>> s = buffer.getvalue()
>>> with open("df_info.txt", "w", encoding="utf-8") as f:
... f.write(s)
260
```

The *memory\_usage* parameter allows deep introspection mode, specially useful for big DataFrames and fine-tune memory optimization:

```
>>> random_strings_array = np.random.choice(['a', 'b', 'c'], 10 ** 6)
>>> df = pd.DataFrame({
       'column_1': np.random.choice(['a', 'b', 'c'], 10 ** 6),
       'column_2': np.random.choice(['a', 'b', 'c'], 10 ** 6),
. . .
       'column_3': np.random.choice(['a', 'b', 'c'], 10 ** 6)
. . .
. . . })
>>> df.info()
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 1000000 entries, 0 to 999999
Data columns (total 3 columns):
column_1 1000000 non-null object
dtypes: object(3)
memory usage: 22.9+ MB
```

```
>>> df.info(memory_usage='deep')
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 1000000 entries, 0 to 999999
Data columns (total 3 columns):
column_1 1000000 non-null object
column_2 1000000 non-null object
column_3 1000000 non-null object
dtypes: object(3)
memory usage: 188.8 MB
```

insert (loc, column, value, allow duplicates=False)

Insert column into DataFrame at specified location.

Raises a ValueError if *column* is already contained in the DataFrame, unless *allow\_duplicates* is set to True.

### **Parameters**

```
loc [int] Insertion index. Must verify 0 <= loc <= len(columns)
column [string, number, or hashable object] label of the inserted column
value [int, Series, or array-like]
allow_duplicates [bool, optional]</pre>
```

Please note that only method='linear' is supported for DataFrames/Series with a MultiIndex.

### **Parameters**

```
method [{'linear', 'time', 'index', 'values', 'nearest', 'zero',]
```

'slinear', 'quadratic', 'cubic', 'barycentric', 'krogh', 'polynomial', 'spline', 'piecewise\_polynomial', 'from\_derivatives', 'pchip', 'akima'}

- 'linear': ignore the index and treat the values as equally spaced. This is the only method supported on MultiIndexes. default
- 'time': interpolation works on daily and higher resolution data to interpolate given length of interval
- 'index', 'values': use the actual numerical values of the index
- 'nearest', 'zero', 'slinear', 'quadratic', 'cubic', 'barycentric', 'polynomial' is passed to scipy.interpolate.interpld. Both 'polynomial' and 'spline' require that you also specify an *order* (int), e.g. df.interpolate(method='polynomial', order=4). These use the actual numerical values of the index.
- 'krogh', 'piecewise\_polynomial', 'spline', 'pchip' and 'akima' are all wrappers around the scipy interpolation methods of similar names. These use the actual numerical values of the index. For more information on their behavior, see the scipy documentation and tutorial documentation
- 'from\_derivatives' refers to BPoly.from\_derivatives which replaces 'piecewise\_polynomial' interpolation method in scipy 0.18

New in version 0.18.1: Added support for the 'akima' method Added interpolate method 'from\_derivatives' which replaces 'piecewise\_polynomial' in scipy 0.18; backwards-compatible with scipy < 0.18

**axis** [{0, 1}, default 0]

- 0: fill column-by-column
- 1: fill row-by-row

**limit** [int, default None.] Maximum number of consecutive NaNs to fill. Must be greater than 0.

limit\_direction [{'forward', 'backward', 'both'}, default 'forward']

limit\_area [{'inside', 'outside'}, default None]

- None: (default) no fill restriction
- 'inside' Only fill NaNs surrounded by valid values (interpolate).
- 'outside' Only fill NaNs outside valid values (extrapolate).

If limit is specified, consecutive NaNs will be filled in this direction.

New in version 0.21.0.

inplace [bool, default False] Update the NDFrame in place if possible.

downcast [optional, 'infer' or None, defaults to None] Downcast dtypes if possible.

**kwargs** [keyword arguments to pass on to the interpolating function.]

# Returns

Series or DataFrame of same shape interpolated at the NaNs

## See also:

```
reindex, replace, fillna
```

## **Examples**

## Filling in NaNs

## is\_copy

#### isin(values)

Return boolean DataFrame showing whether each element in the DataFrame is contained in values.

### **Parameters**

**values** [iterable, Series, DataFrame or dictionary] The result will only be true at a location if all the labels match. If *values* is a Series, that's the index. If *values* is a dictionary, the keys must be the column names, which must match. If *values* is a DataFrame, then both the index and column labels must match.

### **Returns**

#### DataFrame of booleans

# **Examples**

When values is a list:

```
>>> df = pd.DataFrame({'A': [1, 2, 3], 'B': ['a', 'b', 'f']})
>>> df.isin([1, 3, 12, 'a'])

A B
0 True True
1 False False
2 True False
```

When values is a dict:

When values is a Series or DataFrame:

### isna()

Detect missing values.

Return a boolean same-sized object indicating if the values are NA. NA values, such as None or numpy. NaN, gets mapped to True values. Everything else gets mapped to False values. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode. use inf as na = True).

#### Returns

**DataFrame** Mask of bool values for each element in DataFrame that indicates whether an element is not an NA value.

See also:

DataFrame.isnull alias of isna

DataFrame.notna boolean inverse of isna

DataFrame.dropna omit axes labels with missing values

isna top-level isna

## **Examples**

Show which entries in a DataFrame are NA.

```
>>> df = pd.DataFrame({ 'age': [5, 6, np.NaN],
                      'born': [pd.NaT, pd.Timestamp('1939-05-27'),
                               pd.Timestamp('1940-04-25')],
. . .
                      'name': ['Alfred', 'Batman', ''],
. . .
                       'toy': [None, 'Batmobile', 'Joker']})
. . .
>>> df
        born name
                               toy
  age
0 5.0
            NaT Alfred
                              None
1 6.0 1939-05-27 Batman Batmobile
2 NaN 1940-04-25
                              Joker
```

```
>>> df.isna()
age born name toy
0 False True False True
1 False False False
2 True False False
```

Show which entries in a Series are NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.isna()
0 False
1 False
2 True
dtype: bool
```

#### isnull()

Detect missing values.

Return a boolean same-sized object indicating if the values are NA. NA values, such as None or numpy. NaN, gets mapped to True values. Everything else gets mapped to False values. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True).

#### Returns

**DataFrame** Mask of bool values for each element in DataFrame that indicates whether an element is not an NA value.

See also:

DataFrame.isnull alias of isna

DataFrame.notna boolean inverse of isna

DataFrame.dropna omit axes labels with missing values

isna top-level isna

## **Examples**

Show which entries in a DataFrame are NA.

```
>>> df = pd.DataFrame({ 'age': [5, 6, np.NaN],
                       'born': [pd.NaT, pd.Timestamp('1939-05-27'),
                                pd.Timestamp('1940-04-25')],
. . .
                       'name': ['Alfred', 'Batman', ''],
. . .
                       'toy': [None, 'Batmobile', 'Joker']})
. . .
>>> df
           born name
                                toy
  age
0 5.0
            NaT Alfred
                               None
  6.0 1939-05-27 Batman Batmobile
2 NaN 1940-04-25
                               Joker
```

```
>>> df.isna()
age born name toy
0 False True False True
1 False False False
2 True False False
```

Show which entries in a Series are NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.isna()
0 False
1 False
2 True
dtype: bool
```

#### items()

Iterator over (column name, Series) pairs.

#### See also:

iterrows Iterate over DataFrame rows as (index, Series) pairs.

itertuples Iterate over DataFrame rows as namedtuples of the values.

#### iteritems()

Iterator over (column name, Series) pairs.

### See also:

iterrows Iterate over DataFrame rows as (index, Series) pairs.

itertuples Iterate over DataFrame rows as namedtuples of the values.

### iterrows()

Iterate over DataFrame rows as (index, Series) pairs.

#### Returns

it [generator] A generator that iterates over the rows of the frame.

### See also:

itertuples Iterate over DataFrame rows as namedtuples of the values.

iteritems Iterate over (column name, Series) pairs.

#### **Notes**

1. Because iterrows returns a Series for each row, it does **not** preserve dtypes across the rows (dtypes are preserved across columns for DataFrames). For example,

To preserve dtypes while iterating over the rows, it is better to use *itertuples()* which returns namedtuples of the values and which is generally faster than iterrows.

2. You should **never modify** something you are iterating over. This is not guaranteed to work in all cases. Depending on the data types, the iterator returns a copy and not a view, and writing to it will have no effect.

## itertuples (index=True, name='Pandas')

Iterate over DataFrame rows as namedtuples, with index value as first element of the tuple.

#### **Parameters**

**index** [boolean, default True] If True, return the index as the first element of the tuple.

**name** [string, default "Pandas"] The name of the returned namedtuples or None to return regular tuples.

#### See also:

iterrows Iterate over DataFrame rows as (index, Series) pairs.

iteritems Iterate over (column name, Series) pairs.

#### **Notes**

The column names will be renamed to positional names if they are invalid Python identifiers, repeated, or start with an underscore. With a large number of columns (>255), regular tuples are returned.

## **Examples**

## ix

A primarily label-location based indexer, with integer position fallback.

Warning: Starting in 0.20.0, the .ix indexer is deprecated, in favor of the more strict .iloc and .loc indexers.

- .ix[] supports mixed integer and label based access. It is primarily label based, but will fall back to integer positional access unless the corresponding axis is of integer type.
- .ix is the most general indexer and will support any of the inputs in .loc and .iloc. .ix also supports floating point label schemes. .ix is exceptionally useful when dealing with mixed positional and label based hierarchical indexes.

However, when an axis is integer based, ONLY label based access and not positional access is supported. Thus, in such cases, it's usually better to be explicit and use .iloc or .loc.

See more at Advanced Indexing.

```
join (other, on=None, how='left', lsuffix=", rsuffix=", sort=False)
```

Join columns with other DataFrame either on index or on a key column. Efficiently Join multiple DataFrame objects by index at once by passing a list.

## **Parameters**

**other** [DataFrame, Series with name field set, or list of DataFrame] Index should be similar to one of the columns in this one. If a Series is passed, its name attribute must be set, and that will be used as the column name in the resulting joined DataFrame

on [name, tuple/list of names, or array-like] Column or index level name(s) in the caller to join on the index in *other*, otherwise joins index-on-index. If multiple values given, the *other* DataFrame must have a MultiIndex. Can pass an array as the join key if it is not already contained in the calling DataFrame. Like an Excel VLOOKUP operation

**how** [{'left', 'right', 'outer', 'inner'}, default: 'left'] How to handle the operation of the two objects.

- left: use calling frame's index (or column if on is specified)
- right: use other frame's index
- outer: form union of calling frame's index (or column if on is specified) with other frame's index, and sort it lexicographically
- inner: form intersection of calling frame's index (or column if on is specified) with other frame's index, preserving the order of the calling's one

Isuffix [string] Suffix to use from left frame's overlapping columns

rsuffix [string] Suffix to use from right frame's overlapping columns

**sort** [boolean, default False] Order result DataFrame lexicographically by the join key. If False, the order of the join key depends on the join type (how keyword)

### **Returns**

joined [DataFrame]

See also:

**DataFrame.merge** For column(s)-on-columns(s) operations

#### **Notes**

on, lsuffix, and rsuffix options are not supported when passing a list of DataFrame objects Support for specifying index levels as the *on* parameter was added in version 0.23.0

## **Examples**

```
>>> caller = pd.DataFrame({'key': ['K0', 'K1', 'K2', 'K3', 'K4', 'K5'], 'A': ['A0', 'A1', 'A2', 'A3', 'A4', 'A5']})
```

```
>>> caller
    A key
  A0 K0
  A1
       К1
       K2
  A 2.
3
   A 3
       K3
4
   A4
       K4
5
   Α5
       K5
```

```
>>> other = pd.DataFrame({'key': ['K0', 'K1', 'K2'],
... 'B': ['B0', 'B1', 'B2']})
```

Join DataFrames using their indexes.

```
>>> caller.join(other, lsuffix='_caller', rsuffix='_other')
```

```
>>>
      A key_caller
                   B key_other
                  В0
   0 A0
          K0
                          K0
   1
     A1
              K1
                   В1
                           K1
     A2
              K2
                  В2
                           K2
   3 A3
               K3 NaN
                           NaN
   4 A4
               K4 NaN
                           NaN
              K5 NaN
   5 A5
                           NaN
```

If we want to join using the key columns, we need to set key to be the index in both caller and other. The joined DataFrame will have key as its index.

```
>>> caller.set_index('key').join(other.set_index('key'))
```

```
>>>
         Α
   key
   ΚO
        Α0
              В0
   K1
        A1
              В1
   K2.
        A 2.
             B2
   KЗ
        A3 NaN
   K4
        A4 NaN
   K5
        A5 NaN
```

Another option to join using the key columns is to use the on parameter. DataFrame.join always uses other's index but we can use any column in the caller. This method preserves the original caller's index in the result.

```
>>> caller.join(other.set_index('key'), on='key')
```

```
>>> A key B
0 A0 K0 B0
1 A1 K1 B1
2 A2 K2 B2
3 A3 K3 NaN
4 A4 K4 NaN
5 A5 K5 NaN
```

#### keys()

Get the 'info axis' (see Indexing for more)

This is index for Series, columns for DataFrame and major\_axis for Panel.

kurt (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0). Normalized by N-1

# **Parameters**

**axis** [{index (0), columns (1)}]

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**numeric\_only** [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

#### Returns

kurt [Series or DataFrame (if level specified)]

kurtosis (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0). Normalized by N-1

#### **Parameters**

```
axis [{index (0), columns (1)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**numeric\_only** [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### **Returns**

**kurt** [Series or DataFrame (if level specified)]

## last (offset)

Convenience method for subsetting final periods of time series data based on a date offset.

# **Parameters**

offset [string, DateOffset, dateutil.relativedelta]

#### Returns

**subset** [type of caller]

### Raises

 $\textbf{TypeError} \ \ \textbf{If the index is not a} \ \textbf{DatetimeIndex}$ 

#### See also:

first Select initial periods of time series based on a date offset

at\_time Select values at a particular time of the day

between\_time Select values between particular times of the day

### **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='2D')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 1
2018-04-11 2
```

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```
2018-04-13 3
2018-04-15 4
```

Get the rows for the last 3 days:

```
>>> ts.last('3D')

A
2018-04-13 3
2018-04-15 4
```

Notice the data for 3 last calender days were returned, not the last 3 observed days in the dataset, and therefore data for 2018-04-11 was not returned.

## last\_valid\_index()

Return index for last non-NA/null value.

### **Returns**

scalar [type of index]

### **Notes**

If all elements are non-NA/null, returns None. Also returns None for empty NDFrame.

**le** (*other*, *axis='columns'*, *level=None*)

Wrapper for flexible comparison methods le

#### loc

Access a group of rows and columns by label(s) or a boolean array.

.loc[] is primarily label based, but may also be used with a boolean array.

Allowed inputs are:

- A single label, e.g. 5 or 'a', (note that 5 is interpreted as a *label* of the index, and **never** as an integer position along the index).
- A list or array of labels, e.g. ['a', 'b', 'c'].
- A slice object with labels, e.g. 'a': 'f'.

Warning: Note that contrary to usual python slices, **both** the start and the stop are included

- A boolean array of the same length as the axis being sliced, e.g. [True, False, True].
- A callable function with one argument (the calling Series, DataFrame or Panel) and that returns valid output for indexing (one of the above)

See more at Selection by Label

## Raises

**KeyError:** when any items are not found

See also:

DataFrame.at Access a single value for a row/column label pair

**DataFrame.iloc** Access group of rows and columns by integer position(s)

**DataFrame.xs** Returns a cross-section (row(s) or column(s)) from the Series/DataFrame.

Series.loc Access group of values using labels

## **Examples**

### **Getting values**

Single label. Note this returns the row as a Series.

```
>>> df.loc['viper']
max_speed 4
shield 5
Name: viper, dtype: int64
```

List of labels. Note using [ [ ] ] returns a DataFrame.

```
>>> df.loc[['viper', 'sidewinder']]

max_speed shield
viper 4 5
sidewinder 7 8
```

Single label for row and column

```
>>> df.loc['cobra', 'shield']
2
```

Slice with labels for row and single label for column. As mentioned above, note that both the start and stop of the slice are included.

```
>>> df.loc['cobra':'viper', 'max_speed']
cobra 1
viper 4
Name: max_speed, dtype: int64
```

Boolean list with the same length as the row axis

```
>>> df.loc[[False, False, True]]

max_speed shield
sidewinder 7 8
```

Conditional that returns a boolean Series

Conditional that returns a boolean Series with column labels specified

```
>>> df.loc[df['shield'] > 6, ['max_speed']]

max_speed
sidewinder 7
```

#### Callable that returns a boolean Series

```
>>> df.loc[lambda df: df['shield'] == 8]

max_speed shield
sidewinder 7 8
```

## **Setting values**

Set value for all items matching the list of labels

### Set value for an entire row

# Set value for an entire column

### Set value for rows matching callable condition

## Getting values on a DataFrame with an index that has integer labels

Another example using integers for the index

```
>>> df = pd.DataFrame([[1, 2], [4, 5], [7, 8]],
... index=[7, 8, 9], columns=['max_speed', 'shield'])
>>> df
max_speed shield
1 1 2
```

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```
    8
    4
    5

    9
    7
    8
```

Slice with integer labels for rows. As mentioned above, note that both the start and stop of the slice are included.

### Getting values with a MultiIndex

A number of examples using a DataFrame with a MultiIndex

```
>>> tuples = [
      ('cobra', 'mark i'), ('cobra', 'mark ii'),
       ('sidewinder', 'mark i'), ('sidewinder', 'mark ii'),
. . .
       ('viper', 'mark ii'), ('viper', 'mark iii')
. . .
...]
>>> index = pd.MultiIndex.from_tuples(tuples)
>>> values = [[12, 2], [0, 4], [10, 20],
...
          [1, 4], [7, 1], [16, 36]]
>>> df = pd.DataFrame(values, columns=['max_speed', 'shield'], index=index)
>>> df
                    max_speed shield
        mark i
                           12
cobra
                           Ω
                                    4
          mark ii
sidewinder mark i
                          10
                                   20
                           1
          mark ii
                                    4
          mark ii
                            7
                                   1
viper
          mark iii
                           16
                                   36
```

Single label. Note this returns a DataFrame with a single index.

```
>>> df.loc['cobra']

max_speed shield

mark i 12 2

mark ii 0 4
```

Single index tuple. Note this returns a Series.

```
>>> df.loc[('cobra', 'mark ii')]

max_speed 0

shield 4

Name: (cobra, mark ii), dtype: int64
```

Single label for row and column. Similar to passing in a tuple, this returns a Series.

```
>>> df.loc['cobra', 'mark i']
max_speed 12
shield 2
Name: (cobra, mark i), dtype: int64
```

Single tuple. Note using [ [ ] ] returns a DataFrame.

```
>>> df.loc[[('cobra', 'mark ii')]]

max_speed shield

cobra mark ii 0 4
```

Single tuple for the index with a single label for the column

```
>>> df.loc[('cobra', 'mark i'), 'shield']
2
```

Slice from index tuple to single label

```
>>> df.loc[('cobra', 'mark i'):'viper']
                 max_speed shield
                      12
cobra
        mark i
        mark ii
                       0
                               4
sidewinder mark i
                      10
                              20
        mark ii
                       1
                       7
viper
        mark ii
                              1
        mark iii
                     16
                              36
```

Slice from index tuple to index tuple

```
>>> df.loc[('cobra', 'mark i'):('viper', 'mark ii')]
               max_speed shield
                 12
cobra
         mark i
        mark ii
                      0
                              4
                      10
sidewinder mark i
                              2.0
       mark ii
                      1
                               4
viper mark ii
                               1
```

## lookup (row\_labels, col\_labels)

Label-based "fancy indexing" function for DataFrame. Given equal-length arrays of row and column labels, return an array of the values corresponding to each (row, col) pair.

### **Parameters**

```
row_labels [sequence] The row labels to use for lookupcol_labels [sequence] The column labels to use for lookup
```

## **Notes**

Akin to:

```
result = []
for row, col in zip(row_labels, col_labels):
    result.append(df.get_value(row, col))
```

# **Examples**

values [ndarray] The found values

```
1t (other, axis='columns', level=None)
```

Wrapper for flexible comparison methods lt

mad (axis=None, skipna=None, level=None)

Return the mean absolute deviation of the values for the requested axis

#### **Parameters**

```
axis [\{index (0), columns (1)\}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**numeric\_only** [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

#### Returns

mad [Series or DataFrame (if level specified)]

mask (cond, other=nan, inplace=False, axis=None, level=None, errors='raise', try\_cast=False, raise on error=None)

Return an object of same shape as self and whose corresponding entries are from self where *cond* is False and otherwise are from *other*.

#### **Parameters**

**cond** [boolean NDFrame, array-like, or callable] Where *cond* is False, keep the original value. Where True, replace with corresponding value from *other*. If *cond* is callable, it is computed on the NDFrame and should return boolean NDFrame or array. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as cond.

**other** [scalar, NDFrame, or callable] Entries where *cond* is True are replaced with corresponding value from *other*. If other is callable, it is computed on the NDFrame and should return scalar or NDFrame. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as other.

**inplace** [boolean, default False] Whether to perform the operation in place on the data

axis [alignment axis if needed, default None]

level [alignment level if needed, default None]

errors [str, {'raise', 'ignore'}, default 'raise']

- raise: allow exceptions to be raised
- ignore: suppress exceptions. On error return original object

Note that currently this parameter won't affect the results and will always coerce to a suitable dtype.

try\_cast [boolean, default False] try to cast the result back to the input type (if possible),

raise\_on\_error [boolean, default True] Whether to raise on invalid data types (e.g. trying
to where on strings)

Deprecated since version 0.21.0.

## Returns

wh [same type as caller]

### See also:

```
DataFrame.where()
```

## **Notes**

The mask method is an application of the if-then idiom. For each element in the calling DataFrame, if cond is False the element is used; otherwise the corresponding element from the DataFrame other is used.

The signature for DataFrame.where() differs from numpy.where(). Roughly dfl.where(m, df2) is equivalent to np.where(m, df1, df2).

For further details and examples see the mask documentation in indexing.

# **Examples**

```
>>> s = pd.Series(range(5))

>>> s.where(s > 0)

0  NaN

1  1.0

2  2.0

3  3.0

4  4.0
```

(continues on next page)

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max (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

This method returns the maximum of the values in the object. If you want the *index* of the maximum, use idxmax. This is the equivalent of the numpy.ndarray method argmax.

#### **Parameters**

```
axis [\{index (0), columns (1)\}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**numeric\_only** [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

#### Returns

max [Series or DataFrame (if level specified)]

mean (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)
Return the mean of the values for the requested axis

# **Parameters**

```
axis [{index (0), columns (1)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**numeric\_only** [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

## Returns

```
mean [Series or DataFrame (if level specified)]
```

median (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)
Return the median of the values for the requested axis

# **Parameters**

```
axis [{index (0), columns (1)}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**numeric\_only** [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

median [Series or DataFrame (if level specified)]

melt (id\_vars=None, value\_vars=None, var\_name=None, value\_name='value', col\_level=None)
"Unpivots" a DataFrame from wide format to long format, optionally leaving identifier variables set.

This function is useful to massage a DataFrame into a format where one or more columns are identifier variables (*id\_vars*), while all other columns, considered measured variables (*value\_vars*), are "unpivoted" to the row axis, leaving just two non-identifier columns, 'variable' and 'value'.

New in version 0.20.0.

### **Parameters**

frame [DataFrame]

**id\_vars** [tuple, list, or ndarray, optional] Column(s) to use as identifier variables.

**value\_vars** [tuple, list, or ndarray, optional] Column(s) to unpivot. If not specified, uses all columns that are not set as *id\_vars*.

var\_name [scalar] Name to use for the 'variable' column. If None it uses frame.
columns.name or 'variable'.

value\_name [scalar, default 'value'] Name to use for the 'value' column.

col\_level [int or string, optional] If columns are a MultiIndex then use this level to melt.

#### See also:

```
melt, pivot_table, DataFrame.pivot
```

## **Examples**

```
>>> import pandas as pd
>>> df = pd.DataFrame({'A': {0: 'a', 1: 'b', 2: 'c'},
...
'B': {0: 1, 1: 3, 2: 5},
...
'C': {0: 2, 1: 4, 2: 6}})
>>> df
ABC
0 a 1 2
1 b 3 4
2 c 5 6
```

```
>>> df.melt(id_vars=['A'], value_vars=['B'])

A variable value

0 a B 1

1 b B 3

2 c B 5
```

```
>>> df.melt(id_vars=['A'], value_vars=['B', 'C'])
   A variable value
0
  а
            В
                   1
1 b
            В
                    3
2
  С
            В
3
  а
            С
                    2
4
            C
  b
                    4
5
            С
                    6
  C
```

The names of 'variable' and 'value' columns can be customized:

# If you have multi-index columns:

```
>>> df.columns = [list('ABC'), list('DEF')]
>>> df

A B C

D E F

0 a 1 2

1 b 3 4

2 c 5 6
```

```
>>> df.melt(id_vars=[('A', 'D')], value_vars=[('B', 'E')])
  (A, D) variable_0 variable_1 value
0
                   В
                               E
                                        1
       а
1
                   В
                                Ε
                                        3
       b
2
                   В
                                \mathbf{E}
                                        5
       С
```

### memory\_usage (index=True, deep=False)

Return the memory usage of each column in bytes.

The memory usage can optionally include the contribution of the index and elements of *object* dtype.

This value is displayed in *DataFrame.info* by default. This can be suppressed by setting pandas. options.display.memory\_usage to False.

# **Parameters**

index [bool, default True] Specifies whether to include the memory usage of the DataFrame's index in returned Series. If index=True the memory usage of the index the first item in the output.

**deep** [bool, default False] If True, introspect the data deeply by interrogating *object* dtypes for system-level memory consumption, and include it in the returned values.

## Returns

**sizes** [Series] A Series whose index is the original column names and whose values is the memory usage of each column in bytes.

#### See also:

numpy.ndarray.nbytes Total bytes consumed by the elements of an ndarray.

Series.memory\_usage Bytes consumed by a Series.

pandas.Categorical Memory-efficient array for string values with many repeated values.

DataFrame.info Concise summary of a DataFrame.

# **Examples**

```
>>> dtypes = ['int64', 'float64', 'complex128', 'object', 'bool']
>>> data = dict([(t, np.ones(shape=5000).astype(t))
               for t in dtypes])
. . .
>>> df = pd.DataFrame(data)
>>> df.head()
  int64 float64 complex128 object bool
    1 1.0 (1+0j) 1 True
\cap
           1.0
                    (1+0j)
1
      1
                               1 True
2
      1
           1.0
                    (1+0j)
                              1 True
3
      1
            1.0
                    (1+0j)
                              1 True
4
           1.0
                    (1+0j)
                              1 True
```

```
>>> df.memory_usage()
Index 80
int64 40000
float64 40000
complex128 80000
object 40000
bool 5000
dtype: int64
```

The memory footprint of *object* dtype columns is ignored by default:

```
>>> df.memory_usage(deep=True)
Index 80
int64 40000
float64 40000
complex128 80000
object 160000
bool 5000
dtype: int64
```

Use a Categorical for efficient storage of an object-dtype column with many repeated values.

```
>>> df['object'].astype('category').memory_usage(deep=True)
5168
```

```
merge (right, how='inner', on=None, left_on=None, right_on=None, left_index=False, right_index=False, sort=False, suffixes=('_x', '_y'), copy=True, indicator=False, validate=None)
```

Merge DataFrame objects by performing a database-style join operation by columns or indexes.

If joining columns on columns, the DataFrame indexes will be ignored. Otherwise if joining indexes on indexes or indexes on a column or columns, the index will be passed on.

#### **Parameters**

right [DataFrame]

how [{'left', 'right', 'outer', 'inner'}, default 'inner']

- left: use only keys from left frame, similar to a SQL left outer join; preserve key order
- right: use only keys from right frame, similar to a SQL right outer join; preserve key order
- outer: use union of keys from both frames, similar to a SQL full outer join; sort keys lexicographically
- inner: use intersection of keys from both frames, similar to a SQL inner join; preserve the order of the left keys
- **on** [label or list] Column or index level names to join on. These must be found in both DataFrames. If *on* is None and not merging on indexes then this defaults to the intersection of the columns in both DataFrames.
- **left\_on** [label or list, or array-like] Column or index level names to join on in the left DataFrame. Can also be an array or list of arrays of the length of the left DataFrame. These arrays are treated as if they are columns.
- right\_on [label or list, or array-like] Column or index level names to join on in the right DataFrame. Can also be an array or list of arrays of the length of the right DataFrame. These arrays are treated as if they are columns.
- **left\_index** [boolean, default False] Use the index from the left DataFrame as the join key(s). If it is a MultiIndex, the number of keys in the other DataFrame (either the index or a number of columns) must match the number of levels
- **right\_index** [boolean, default False] Use the index from the right DataFrame as the join key. Same caveats as left index
- **sort** [boolean, default False] Sort the join keys lexicographically in the result DataFrame. If False, the order of the join keys depends on the join type (how keyword)
- **suffixes** [2-length sequence (tuple, list, ...)] Suffix to apply to overlapping column names in the left and right side, respectively
- copy [boolean, default True] If False, do not copy data unnecessarily
- indicator [boolean or string, default False] If True, adds a column to output DataFrame called "\_merge" with information on the source of each row. If string, column with information on source of each row will be added to output DataFrame, and column will be named value of string. Information column is Categorical-type and takes on a value of "left\_only" for observations whose merge key only appears in 'left' DataFrame, "right\_only" for observations whose merge key only appears in 'right' DataFrame, and "both" if the observation's merge key is found in both.

validate [string, default None] If specified, checks if merge is of specified type.

- "one\_to\_one" or "1:1": check if merge keys are unique in both left and right datasets.
- "one\_to\_many" or "1:m": check if merge keys are unique in left dataset.
- "many\_to\_one" or "m:1": check if merge keys are unique in right dataset.
- "many\_to\_many" or "m:m": allowed, but does not result in checks.

New in version 0.21.0.

#### Returns

**merged** [DataFrame] The output type will the be same as 'left', if it is a subclass of DataFrame.

#### See also:

```
merge_ordered, merge_asof, DataFrame.join
```

#### **Notes**

Support for specifying index levels as the on, left\_on, and right\_on parameters was added in version 0.23.0

# **Examples**

```
>>> A
                  >>> B
   lkey value
                      rkey value
   foo 1
                  0
                      foo 5
        2
                  1
                      bar
                           6
   har
                           7
2
        3
                  2
                      qux
   baz
3
                  3
                      bar 8
    foo
```

```
>>> A.merge(B, left_on='lkey', right_on='rkey', how='outer')
   lkey value_x rkey value_y
        1
                 foo
                       5
                       5
  foo
        4
                 foo
2
        2
                       6
  bar
                 bar
3
        2
                       8
  bar
                 bar
        3
4
  baz
                 NaN
                       NaN
5
                       7
  NaN
        NaN
                 qux
```

min (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

This method returns the minimum of the values in the object. If you want the *index* of the minimum, use idxmin. This is the equivalent of the numpy.ndarray method argmin.

#### **Parameters**

```
axis [{index (0), columns (1)}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**numeric\_only** [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

## Returns

min [Series or DataFrame (if level specified)]

```
mod (other, axis='columns', level=None, fill_value=None)
```

Modulo of dataframe and other, element-wise (binary operator mod).

Equivalent to dataframe % other, but with support to substitute a fill\_value for missing data in one of the inputs.

# **Parameters**

other [Series, DataFrame, or constant]

```
axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on
```

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

# Returns

```
result [DataFrame]
```

### See also:

DataFrame.rmod

### **Notes**

Mismatched indices will be unioned together

# **Examples**

None

```
mode (axis=0, numeric only=False)
```

Gets the mode(s) of each element along the axis selected. Adds a row for each mode per label, fills in gaps with nan.

Note that there could be multiple values returned for the selected axis (when more than one item share the maximum frequency), which is the reason why a dataframe is returned. If you want to impute missing values with the mode in a dataframe df, you can just do this: df.fillna(df.mode().iloc[0])

### **Parameters**

```
axis [{0 or 'index', 1 or 'columns'}, default 0]
```

- 0 or 'index': get mode of each column
- 1 or 'columns': get mode of each row

numeric\_only [boolean, default False] if True, only apply to numeric columns

### **Returns**

```
modes [DataFrame (sorted)]
```

# **Examples**

```
>>> df = pd.DataFrame({'A': [1, 2, 1, 2, 1, 2, 3]})
>>> df.mode()

A
0 1
1 2
```

```
mul (other, axis='columns', level=None, fill_value=None)
```

Multiplication of dataframe and other, element-wise (binary operator mul).

Equivalent to dataframe \* other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

other [Series, DataFrame, or constant]

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

fill\_value [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

# Returns

result [DataFrame]

#### See also:

DataFrame.rmul

## **Notes**

Mismatched indices will be unioned together

# **Examples**

None

multiply (other, axis='columns', level=None, fill\_value=None)

Multiplication of dataframe and other, element-wise (binary operator mul).

Equivalent to dataframe \* other, but with support to substitute a fill\_value for missing data in one of the inputs.

# **Parameters**

other [Series, DataFrame, or constant]

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

### Returns

result [DataFrame]

# See also:

DataFrame.rmul

### **Notes**

Mismatched indices will be unioned together

# **Examples**

None

#### ndim

Return an int representing the number of axes / array dimensions.

Return 1 if Series. Otherwise return 2 if DataFrame.

### See also:

ndarray.ndim

# **Examples**

```
>>> s = pd.Series({'a': 1, 'b': 2, 'c': 3})
>>> s.ndim
1
```

```
>>> df = pd.DataFrame({'coll': [1, 2], 'col2': [3, 4]})
>>> df.ndim
2
```

ne (other, axis='columns', level=None)

Wrapper for flexible comparison methods ne

```
nlargest (n, columns, keep='first')
```

Return the first *n* rows ordered by *columns* in descending order.

Return the first *n* rows with the largest values in *columns*, in descending order. The columns that are not specified are returned as well, but not used for ordering.

This method is equivalent to df.sort\_values(columns, ascending=False).head(n), but more performant.

### **Parameters**

**n** [int] Number of rows to return.

**columns** [label or list of labels] Column label(s) to order by.

**keep** [{'first', 'last'}, default 'first'] Where there are duplicate values:

- first : prioritize the first occurrence(s)
- *last* : prioritize the last occurrence(s)

# Returns

**DataFrame** The first *n* rows ordered by the given columns in descending order.

See also:

**DataFrame.nsmallest** Return the first *n* rows ordered by *columns* in ascending order.

DataFrame.sort\_values Sort DataFrame by the values

**DataFrame.head** Return the first *n* rows without re-ordering.

# **Notes**

This function cannot be used with all column types. For example, when specifying columns with *object* or *category* dtypes, TypeError is raised.

# **Examples**

```
>>> df = pd.DataFrame({'a': [1, 10, 8, 10, -1],
. . .
                       'b': list('abdce'),
. . .
                       'c': [1.0, 2.0, np.nan, 3.0, 4.0]})
>>> df
   a b
      a 1.0
\cap
  1
  10 b 2.0
   8
      d NaN
  10
      C
         3.0
  -1 e 4.0
```

In the following example, we will use nlargest to select the three rows having the largest values in column "a".

```
>>> df.nlargest(3, 'a')
    a b c
1 10 b 2.0
3 10 c 3.0
2 8 d NaN
```

When using keep='last', ties are resolved in reverse order:

```
>>> df.nlargest(3, 'a', keep='last')
    a b c
3 10 c 3.0
1 10 b 2.0
2 8 d NaN
```

To order by the largest values in column "a" and then "c", we can specify multiple columns like in the next example.

```
>>> df.nlargest(3, ['a', 'c'])
    a b c
3 10 c 3.0
1 10 b 2.0
2 8 d NaN
```

Attempting to use nlargest on non-numeric dtypes will raise a TypeError:

```
>>> df.nlargest(3, 'b')
Traceback (most recent call last):
TypeError: Column 'b' has dtype object, cannot use method 'nlargest'
```

## notna()

Detect existing (non-missing) values.

Return a boolean same-sized object indicating if the values are not NA. Non-missing values get mapped to True. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True). NA values, such as None or numpy.NaN, get mapped to False values.

## Returns

**DataFrame** Mask of bool values for each element in DataFrame that indicates whether an element is not an NA value.

See also:

DataFrame.notnull alias of notna

DataFrame.isna boolean inverse of notna

DataFrame.dropna omit axes labels with missing values

notna top-level notna

# **Examples**

Show which entries in a DataFrame are not NA.

```
>>> df = pd.DataFrame({'age': [5, 6, np.NaN],
                      'born': [pd.NaT, pd.Timestamp('1939-05-27'),
                              pd.Timestamp('1940-04-25')],
                      'name': ['Alfred', 'Batman', ''],
. . .
                      'toy': [None, 'Batmobile', 'Joker']})
. . .
>>> df
           born name
                               tov
  age
           NaT Alfred
0 5.0
                             None
  6.0 1939-05-27 Batman Batmobile
2 NaN 1940-04-25
                             Joker
```

```
>>> df.notna()
   age born name toy
0 True False True False
1 True True True True
2 False True True True
```

Show which entries in a Series are not NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.notna()
0 True
1 True
2 False
dtype: bool
```

### notnull()

Detect existing (non-missing) values.

Return a boolean same-sized object indicating if the values are not NA. Non-missing values get mapped to True. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True). NA values, such as None or numpy.NaN, get mapped to False values.

## Returns

**DataFrame** Mask of bool values for each element in DataFrame that indicates whether an element is not an NA value.

See also:

```
DataFrame.notnull alias of notna
```

DataFrame.isna boolean inverse of notna

DataFrame.dropna omit axes labels with missing values

notna top-level notna

## **Examples**

Show which entries in a DataFrame are not NA.

```
>>> df = pd.DataFrame({'age': [5, 6, np.NaN],
                      'born': [pd.NaT, pd.Timestamp('1939-05-27'),
                              pd.Timestamp('1940-04-25')],
                      'name': ['Alfred', 'Batman', ''],
. . .
                      'toy': [None, 'Batmobile', 'Joker']})
. . .
>>> df
           born name
                               tov
  age
           NaT Alfred
0 5.0
                             None
  6.0 1939-05-27 Batman Batmobile
2 NaN 1940-04-25
                             Joker
```

```
>>> df.notna()
age born name toy
0 True False True False
1 True True True
2 False True True
```

Show which entries in a Series are not NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0     5.0
1     6.0
2     NaN
dtype: float64
```

```
>>> ser.notna()
0 True
1 True
2 False
dtype: bool
```

nsmallest (n, columns, keep='first')

Get the rows of a DataFrame sorted by the n smallest values of columns.

### **Parameters**

```
n [int] Number of items to retrieve
```

columns [list or str] Column name or names to order by

**keep** [{'first', 'last'}, default 'first'] Where there are duplicate values: - first : take the first occurrence. - last : take the last occurrence.

#### Returns

**DataFrame** 

# **Examples**

```
>>> df = pd.DataFrame({'a': [1, 10, 8, 11, -1],
... 'b': list('abdce'),
... 'c': [1.0, 2.0, np.nan, 3.0, 4.0]})
>>> df.nsmallest(3, 'a')
a b c
4 -1 e 4
0 1 a 1
2 8 d NaN
```

# nunique (axis=0, dropna=True)

Return Series with number of distinct observations over requested axis.

New in version 0.20.0.

### **Parameters**

```
axis [{0 or 'index', 1 or 'columns'}, default 0]
```

dropna [boolean, default True] Don't include NaN in the counts.

## Returns

nunique [Series]

# **Examples**

# pct\_change (periods=1, fill\_method='pad', limit=None, freq=None, \*\*kwargs)

Percentage change between the current and a prior element.

Computes the percentage change from the immediately previous row by default. This is useful in comparing the percentage of change in a time series of elements.

# **Parameters**

**periods** [int, default 1] Periods to shift for forming percent change.

fill\_method [str, default 'pad'] How to handle NAs before computing percent changes.

limit [int, default None] The number of consecutive NAs to fill before stopping.

**freq** [DateOffset, timedelta, or offset alias string, optional] Increment to use from time series API (e.g. 'M' or BDay()).

\*\*kwargs Additional keyword arguments are passed into DataFrame.shift or Series.shift.

# Returns

chg [Series or DataFrame] The same type as the calling object.

### See also:

**Series.diff** Compute the difference of two elements in a Series.

DataFrame.diff Compute the difference of two elements in a DataFrame.

**Series.shift** Shift the index by some number of periods.

DataFrame.shift Shift the index by some number of periods.

# **Examples**

#### Series

```
>>> s = pd.Series([90, 91, 85])
>>> s
0 90
1 91
2 85
dtype: int64
```

See the percentage change in a Series where filling NAs with last valid observation forward to next valid.

```
>>> s = pd.Series([90, 91, None, 85])
>>> s
0 90.0
1 91.0
2 NaN
3 85.0
dtype: float64
```

# **DataFrame**

Percentage change in French franc, Deutsche Mark, and Italian lira from 1980-01-01 to 1980-03-01.

```
>>> df = pd.DataFrame({
... 'FR': [4.0405, 4.0963, 4.3149],
... 'GR': [1.7246, 1.7482, 1.8519],
... 'IT': [804.74, 810.01, 860.13]},
... index=['1980-01-01', '1980-02-01', '1980-03-01'])
>>> df

FR GR IT

1980-01-01 4.0405 1.7246 804.74

1980-02-01 4.0963 1.7482 810.01

1980-03-01 4.3149 1.8519 860.13
```

Percentage of change in GOOG and APPL stock volume. Shows computing the percentage change between columns.

```
>>> df = pd.DataFrame({
... '2016': [1769950, 30586265],
... '2015': [1500923, 40912316],
... '2014': [1371819, 41403351]},
... index=['GOOG', 'APPL'])
>>> df

2016 2015 2014

GOOG 1769950 1500923 1371819
APPL 30586265 40912316 41403351
```

### **Parameters**

**func** [function] function to apply to the NDFrame. args, and kwargs are passed into func. Alternatively a (callable, data\_keyword) tuple where data\_keyword is a string indicating the keyword of callable that expects the NDFrame.

args [iterable, optional] positional arguments passed into func.

kwargs [mapping, optional] a dictionary of keyword arguments passed into func.

object [the return type of func.]

## See also:

```
pandas.DataFrame.apply, pandas.DataFrame.applymap, pandas.Series.map
```

#### **Notes**

Use .pipe when chaining together functions that expect Series, DataFrames or GroupBy objects. Instead of writing

```
>>> f(g(h(df), arg1=a), arg2=b, arg3=c)
```

You can write

```
>>> (df.pipe(h)
... .pipe(g, arg1=a)
... .pipe(f, arg2=b, arg3=c)
... )
```

If you have a function that takes the data as (say) the second argument, pass a tuple indicating which keyword expects the data. For example, suppose f takes its data as arg2:

```
>>> (df.pipe(h)
... .pipe(g, arg1=a)
... .pipe((f, 'arg2'), arg1=a, arg3=c)
... )
```

pivot (index=None, columns=None, values=None)

Return reshaped DataFrame organized by given index / column values.

Reshape data (produce a "pivot" table) based on column values. Uses unique values from specified *index* / *columns* to form axes of the resulting DataFrame. This function does not support data aggregation, multiple values will result in a MultiIndex in the columns. See the User Guide for more on reshaping.

# **Parameters**

**index** [string or object, optional] Column to use to make new frame's index. If None, uses existing index.

**columns** [string or object] Column to use to make new frame's columns.

**values** [string, object or a list of the previous, optional] Column(s) to use for populating new frame's values. If not specified, all remaining columns will be used and the result will have hierarchically indexed columns.

Changed in version 0.23.0: Also accept list of column names.

#### Returns

**DataFrame** Returns reshaped DataFrame.

# Raises

**ValueError:** When there are any *index*, *columns* combinations with multiple values. *DataFrame.pivot\_table* when you need to aggregate.

See also:

**DataFrame.pivot\_table** generalization of pivot that can handle duplicate values for one index/column pair.

DataFrame.unstack pivot based on the index values instead of a column.

#### **Notes**

For finer-tuned control, see hierarchical indexing documentation along with the related stack/unstack methods

## **Examples**

```
>>> df = pd.DataFrame({'foo': ['one', 'one', 'one', 'two', 'two',
                             'two'],
                     'bar': ['A', 'B', 'C', 'A', 'B', 'C'],
. . .
                     'baz': [1, 2, 3, 4, 5, 6],
. . .
                     'zoo': ['x', 'y', 'z', 'q', 'w', 't']})
. . .
>>> df
   foo
        bar baz zoo
0
        A 1
   one
                  Х
        В
1
   one
             2
2
   one
        С
             3
3
             4
   two
        A
                  q
             5
   two
        В
5
        С
             6
   t.wo
```

```
>>> df.pivot(index='foo', columns='bar', values='baz')
bar A B C
foo
one 1 2 3
two 4 5 6
```

```
bar A B C
foo
one 1 2 3
two 4 5 6
```

```
baz zoo
bar A B C A B C
foo
one 1 2 3 x y z
two 4 5 6 q w t
```

A ValueError is raised if there are any duplicates.

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```
2 two B 3
3 two C 4
```

Notice that the first two rows are the same for our *index* and *columns* arguments.

```
>>> df.pivot(index='foo', columns='bar', values='baz')
Traceback (most recent call last):
...
ValueError: Index contains duplicate entries, cannot reshape
```

pivot\_table (values=None, index=None, columns=None, aggfunc='mean', fill\_value=None, margins=False, dropna=True, margins name='All')

Create a spreadsheet-style pivot table as a DataFrame. The levels in the pivot table will be stored in MultiIndex objects (hierarchical indexes) on the index and columns of the result DataFrame

#### **Parameters**

values [column to aggregate, optional]

**index** [column, Grouper, array, or list of the previous] If an array is passed, it must be the same length as the data. The list can contain any of the other types (except list). Keys to group by on the pivot table index. If an array is passed, it is being used as the same manner as column values.

**columns** [column, Grouper, array, or list of the previous] If an array is passed, it must be the same length as the data. The list can contain any of the other types (except list). Keys to group by on the pivot table column. If an array is passed, it is being used as the same manner as column values.

**aggfunc** [function, list of functions, dict, default numpy.mean] If list of functions passed, the resulting pivot table will have hierarchical columns whose top level are the function names (inferred from the function objects themselves) If dict is passed, the key is column to aggregate and value is function or list of functions

fill\_value [scalar, default None] Value to replace missing values with

margins [boolean, default False] Add all row / columns (e.g. for subtotal / grand totals)

dropna [boolean, default True] Do not include columns whose entries are all NaN

margins\_name [string, default 'All'] Name of the row / column that will contain the totals when margins is True.

# Returns

table [DataFrame]

See also:

DataFrame.pivot pivot without aggregation that can handle non-numeric data

## **Examples**

```
>>> df = pd.DataFrame({"A": ["foo", "foo", "foo", "foo", "foo", "bar", "bar", "bar", "bar"],
...
"B": ["one", "one", "one", "two", "two",
"one", "one", "two", "two"],
...
"C": ["small", "large", "large", "small",
```

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```
"small", "large", "small", "small",
. . .
                           "large"],
. . .
                     "D": [1, 2, 2, 3, 3, 4, 5, 6, 7]})
. . .
>>> df
       В
             C D
    Α
  foo one small
  foo
      one large 2
  foo one large 2
3
  foo two small 3
  foo two small 3
  bar one large 4
  bar one small 5
  bar two small 6
 bar two large 7
```

```
>>> table = pivot_table(df, values='D', index=['A', 'B'],
                      columns=['C'], aggfunc=np.sum)
>>> table
С
       large small
  В
Α
          4.0
              5.0
bar one
          7.0
               6.0
   two
foo one
          4.0
                 1.0
          NaN
                6.0
```

```
>>> table = pivot_table(df, values='D', index=['A', 'B'],
                     columns=['C'], aggfunc=np.sum)
. . .
>>> table
С
       large small
  В
Α
         4.0
              5.0
bar one
        7.0 6.0
   two
foo one
       4.0 1.0
   two
       NaN 6.0
```

```
>>> table = pivot_table(df, values=['D', 'E'], index=['A', 'C'],
. . .
                       aggfunc={'D': np.mean,
                                'E': [min, max, np.mean] })
. . .
>>> table
                D
                    Ε
             mean max median min
A C
bar large 5.500000 16
                       14.5 13
   small 5.500000 15
                         14.5 14
foo large 2.000000 10
                         9.5
   small 2.333333 12
                         11.0
```

# plot

alias of pandas.plotting.\_core.FramePlotMethods

#### pop (item

Return item and drop from frame. Raise KeyError if not found.

### **Parameters**

item [str] Column label to be popped

## Returns

## popped [Series]

# **Examples**

```
>>> df = pd.DataFrame([('falcon', 'bird',
                                            389.0),
                       ('parrot', 'bird',
                                            24.0),
. . .
                                            80.5),
                      ('lion', 'mammal',
. . .
                      ('monkey', 'mammal', np.nan)],
. . .
                     columns=('name', 'class', 'max_speed'))
. . .
>>> df
    name class max_speed
  falcon bird 389.0
0
1
  parrot
           bird
                      24.0
2
    lion mammal
                      80.5
3
  monkey mammal
                       NaN
```

```
>>> df.pop('class')
0 bird
1 bird
2 mammal
3 mammal
Name: class, dtype: object
```

```
>>> df
    name max_speed
0 falcon 389.0
1 parrot 24.0
2 lion 80.5
3 monkey NaN
```

pow (other, axis='columns', level=None, fill\_value=None)

Exponential power of dataframe and other, element-wise (binary operator pow).

Equivalent to dataframe \*\* other, but with support to substitute a fill\_value for missing data in one of the inputs.

# **Parameters**

```
other [Series, DataFrame, or constant]
```

```
axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on
```

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

#### Returns

result [DataFrame]

## See also:

DataFrame.rpow

### **Notes**

Mismatched indices will be unioned together

# **Examples**

None

prod (axis=None, skipna=None, level=None, numeric\_only=None, min\_count=0, \*\*kwargs)
Return the product of the values for the requested axis

#### **Parameters**

```
axis [{index (0), columns (1)}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**numeric\_only** [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

min\_count [int, default 0] The required number of valid values to perform the operation. If fewer than min\_count non-NA values are present the result will be NA.

New in version 0.22.0: Added with the default being 0. This means the sum of an all-NA or empty Series is 0, and the product of an all-NA or empty Series is 1.

#### Returns

prod [Series or DataFrame (if level specified)]

### **Examples**

By default, the product of an empty or all-NA Series is 1

```
>>> pd.Series([]).prod()
1.0
```

This can be controlled with the min\_count parameter

```
>>> pd.Series([]).prod(min_count=1)
nan
```

Thanks to the skipna parameter, min\_count handles all-NA and empty series identically.

```
>>> pd.Series([np.nan]).prod()
1.0
```

```
>>> pd.Series([np.nan]).prod(min_count=1)
nan
```

product (axis=None, skipna=None, level=None, numeric\_only=None, min\_count=0, \*\*kwargs)
Return the product of the values for the requested axis

### **Parameters**

```
axis [{index (0), columns (1)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**numeric\_only** [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

min\_count [int, default 0] The required number of valid values to perform the operation. If fewer than min\_count non-NA values are present the result will be NA.

New in version 0.22.0: Added with the default being 0. This means the sum of an all-NA or empty Series is 0, and the product of an all-NA or empty Series is 1.

### Returns

prod [Series or DataFrame (if level specified)]

# **Examples**

By default, the product of an empty or all-NA Series is 1

```
>>> pd.Series([]).prod()
1.0
```

This can be controlled with the min\_count parameter

```
>>> pd.Series([]).prod(min_count=1)
nan
```

Thanks to the skipna parameter, min\_count handles all-NA and empty series identically.

```
>>> pd.Series([np.nan]).prod()
1.0
```

```
>>> pd.Series([np.nan]).prod(min_count=1)
nan
```

**quantile** (q=0.5, axis=0, numeric only=True, interpolation='linear')

Return values at the given quantile over requested axis, a la numpy.percentile.

# **Parameters**

**q** [float or array-like, default 0.5 (50% quantile)] 0 <= q <= 1, the quantile(s) to compute

axis [{0, 1, 'index', 'columns'} (default 0)] 0 or 'index' for row-wise, 1 or 'columns' for column-wise

**numeric\_only** [boolean, default True] If False, the quantile of datetime and timedelta data will be computed as well

**interpolation** [{'linear', 'lower', 'higher', 'midpoint', 'nearest'}] New in version 0.18.0.

This optional parameter specifies the interpolation method to use, when the desired quantile lies between two data points i and j:

- linear: i + (j i) \* fraction, where fraction is the fractional part of the index surrounded by i and j.
- lower: *i*.

- higher: *j*.
- nearest: *i* or *j* whichever is nearest.
- midpoint: (i + j) / 2.

**quantiles** [Series or DataFrame]

- If q is an array, a DataFrame will be returned where the index is q, the columns are the columns of self, and the values are the quantiles.
- If q is a float, a Series will be returned where the index is the columns of self and the values are the quantiles.

#### See also:

```
pandas.core.window.Rolling.quantile
```

# **Examples**

Specifying numeric\_only=False will also compute the quantile of datetime and timedelta data.

query (expr, inplace=False, \*\*kwargs)

Query the columns of a frame with a boolean expression.

## **Parameters**

**expr** [string] The query string to evaluate. You can refer to variables in the environment by prefixing them with an '@' character like @a + b.

**inplace** [bool] Whether the query should modify the data in place or return a modified copy New in version 0.18.0.

**kwargs** [dict] See the documentation for pandas.eval() for complete details on the keyword arguments accepted by DataFrame.query().

**q** [DataFrame]

### See also:

```
pandas.eval, DataFrame.eval
```

## Notes

The result of the evaluation of this expression is first passed to DataFrame.loc and if that fails because of a multidimensional key (e.g., a DataFrame) then the result will be passed to DataFrame.

\_\_getitem\_\_().

This method uses the top-level pandas.eval() function to evaluate the passed query.

The query () method uses a slightly modified Python syntax by default. For example, the & and | (bitwise) operators have the precedence of their boolean cousins, and and or. This is syntactically valid Python, however the semantics are different.

You can change the semantics of the expression by passing the keyword argument parser='python'. This enforces the same semantics as evaluation in Python space. Likewise, you can pass engine='python' to evaluate an expression using Python itself as a backend. This is not recommended as it is inefficient compared to using numexpr as the engine.

The DataFrame index and DataFrame columns attributes of the DataFrame instance are placed in the query namespace by default, which allows you to treat both the index and columns of the frame as a column in the frame. The identifier index is used for the frame index; you can also use the name of the index to identify it in a query. Please note that Python keywords may not be used as identifiers.

For further details and examples see the query documentation in indexing.

### **Examples**

```
>>> from numpy.random import randn
>>> from pandas import DataFrame
>>> df = pd.DataFrame(randn(10, 2), columns=list('ab'))
>>> df.query('a > b')
>>> df[df.a > df.b] # same result as the previous expression
```

radd (other, axis='columns', level=None, fill\_value=None)

Addition of dataframe and other, element-wise (binary operator radd).

Equivalent to other + dataframe, but with support to substitute a fill\_value for missing data in one of the inputs.

#### **Parameters**

```
other [Series, DataFrame, or constant]
```

```
axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on
```

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

fill\_value [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

result [DataFrame]

## See also:

DataFrame.add

### **Notes**

Mismatched indices will be unioned together

# **Examples**

```
>>> a = pd.DataFrame([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'],
                    columns=['one'])
>>> a
  one
a 1.0
b 1.0
c 1.0
d NaN
>>> b = pd.DataFrame(dict(one=[1, np.nan, 1, np.nan],
                          two=[np.nan, 2, np.nan, 2]),
. . .
                     index=['a', 'b', 'd', 'e'])
. . .
>>> b
  one
       two
  1.0 NaN
  NaN 2.0
  1.0 NaN
  NaN 2.0
>>> a.add(b, fill_value=0)
  one two
  2.0 NaN
  1.0
       2.0
  1.0 NaN
  1.0
       NaN
  NaN
        2.0
```

rank (axis=0, method='average', numeric\_only=None, na\_option='keep', ascending=True, pct=False)

Compute numerical data ranks (1 through n) along axis. Equal values are assigned a rank that is the average of the ranks of those values

### **Parameters**

```
axis [{0 or 'index', 1 or 'columns'}, default 0] index to direct ranking
method [{'average', 'min', 'max', 'first', 'dense'}]
```

- · average: average rank of group
- min: lowest rank in group
- max: highest rank in group
- first: ranks assigned in order they appear in the array
- dense: like 'min', but rank always increases by 1 between groups

numeric\_only [boolean, default None] Include only float, int, boolean data. Valid only for DataFrame or Panel objects

na\_option [{'keep', 'top', 'bottom'}]

- keep: leave NA values where they are
- top: smallest rank if ascending
- bottom: smallest rank if descending

ascending [boolean, default True] False for ranks by high (1) to low (N)

pct [boolean, default False] Computes percentage rank of data

#### Returns

ranks [same type as caller]

rdiv (other, axis='columns', level=None, fill\_value=None)

Floating division of dataframe and other, element-wise (binary operator rtruediv).

Equivalent to other / dataframe, but with support to substitute a fill\_value for missing data in one of the inputs.

#### **Parameters**

other [Series, DataFrame, or constant]

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

fill\_value [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

#### **Returns**

result [DataFrame]

# See also:

DataFrame.truediv

# **Notes**

Mismatched indices will be unioned together

# **Examples**

None

#### reindex (\*\*kwargs)

Conform DataFrame to new index with optional filling logic, placing NA/NaN in locations having no value in the previous index. A new object is produced unless the new index is equivalent to the current one and copy=False

## **Parameters**

labels [array-like, optional] New labels / index to conform the axis specified by 'axis' to.

**index, columns** [array-like, optional (should be specified using keywords)] New labels / index to conform to. Preferably an Index object to avoid duplicating data

**axis** [int or str, optional] Axis to target. Can be either the axis name ('index', 'columns') or number (0, 1).

**method** [{None, 'backfill', 'pad'/'ffill', 'nearest'}, optional] method to use for filling holes in reindexed DataFrame. Please note: this is only applicable to DataFrames/Series with a monotonically increasing/decreasing index.

- default: don't fill gaps
- pad / ffill: propagate last valid observation forward to next valid
- backfill / bfill: use next valid observation to fill gap
- nearest: use nearest valid observations to fill gap

copy [boolean, default True] Return a new object, even if the passed indexes are the same

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

**fill\_value** [scalar, default np.NaN] Value to use for missing values. Defaults to NaN, but can be any "compatible" value

limit [int, default None] Maximum number of consecutive elements to forward or backward fill

tolerance [optional] Maximum distance between original and new labels for inexact
matches. The values of the index at the matching locations most satisfy the equation
abs(index[indexer] - target) <= tolerance.</pre>

Tolerance may be a scalar value, which applies the same tolerance to all values, or list-like, which applies variable tolerance per element. List-like includes list, tuple, array, Series, and must be the same size as the index and its dtype must exactly match the index's type.

New in version 0.21.0: (list-like tolerance)

#### Returns

reindexed [DataFrame]

# **Examples**

DataFrame.reindex supports two calling conventions

```
• (index=index labels, columns=column labels, ...)
```

```
• (labels, axis={'index', 'columns'}, ...)
```

We highly recommend using keyword arguments to clarify your intent.

Create a dataframe with some fictional data.

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Firefox	200	0.04
Chrome	200	0.02
Safari	404	0.07
IE10	404	0.08
Konqueror	301	1.00

Create a new index and reindex the dataframe. By default values in the new index that do not have corresponding records in the dataframe are assigned NaN.

```
>>> new_index= ['Safari', 'Iceweasel', 'Comodo Dragon', 'IE10',
                'Chrome']
>>> df.reindex(new_index)
              http_status response_time
Safari
                     404.0
                                      0.07
Iceweasel
                                      NaN
                      NaN
Comodo Dragon
                       NaN
                                      NaN
IE10
                     404.0
                                      0.08
Chrome
                     200.0
                                      0.02
```

We can fill in the missing values by passing a value to the keyword fill\_value. Because the index is not monotonically increasing or decreasing, we cannot use arguments to the keyword method to fill the NaN values.

```
>>> df.reindex(new_index, fill_value=0)
              http_status response_time
Safari
                     404
                                     0.07
                       0
                                     0.00
Iceweasel
                        0
                                     0.00
Comodo Dragon
                                     0.08
IE10
                       404
Chrome
                       200
                                     0.02
```

```
>>> df.reindex(new_index, fill_value='missing')
            http_status response_time
                     404
                                  0.07
Safari
Iceweasel
                 missing
                               missing
Comodo Dragon
                               missing
                 missing
IE10
                                   0.08
                      404
Chrome
                      200
                                   0.02
```

We can also reindex the columns.

```
>>> df.reindex(columns=['http_status', 'user_agent'])
         http_status user_agent
Firefox
                   200
                               NaN
Chrome
                   200
                               NaN
Safari
                   404
                               NaN
IE10
                   404
                               NaN
Konqueror
                   301
                               NaN
```

Or we can use "axis-style" keyword arguments

```
>>> df.reindex(['http_status', 'user_agent'], axis="columns")
    http_status user_agent
Firefox 200 NaN
Chrome 200 NaN
Safari 404 NaN
```

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```
IE10 404 NaN
Konqueror 301 NaN
```

To further illustrate the filling functionality in reindex, we will create a dataframe with a monotonically increasing index (for example, a sequence of dates).

```
>>> date_index = pd.date_range('1/1/2010', periods=6, freq='D')
>>> df2 = pd.DataFrame({"prices": [100, 101, np.nan, 100, 89, 88]},
                       index=date_index)
. . .
>>> df2
           prices
2010-01-01
             100
2010-01-02
              101
2010-01-03
              NaN
2010-01-04
              100
2010-01-05
               89
2010-01-06
```

Suppose we decide to expand the dataframe to cover a wider date range.

```
>>> date_index2 = pd.date_range('12/29/2009', periods=10, freq='D')
>>> df2.reindex(date_index2)
           prices
2009-12-29
              NaN
2009-12-30
              NaN
2009-12-31
              NaN
2010-01-01
               100
2010-01-02
               101
2010-01-03
              NaN
2010-01-04
              100
               89
2010-01-05
2010-01-06
               88
2010-01-07
               NaN
```

The index entries that did not have a value in the original data frame (for example, '2009-12-29') are by default filled with NaN. If desired, we can fill in the missing values using one of several options.

For example, to backpropagate the last valid value to fill the NaN values, pass bfill as an argument to the method keyword.

```
>>> df2.reindex(date_index2, method='bfill')
            prices
2009-12-29
               100
2009-12-30
               100
2009-12-31
               100
2010-01-01
               100
2010-01-02
               101
2010-01-03
              NaN
2010-01-04
              100
2010-01-05
               89
2010-01-06
                88
2010-01-07
               NaN
```

Please note that the NaN value present in the original dataframe (at index value 2010-01-03) will not be filled by any of the value propagation schemes. This is because filling while reindexing does not look at dataframe values, but only compares the original and desired indexes. If you do want to fill in the NaN values present in the original dataframe, use the fillna() method.

See the user guide for more.

reindex\_axis (labels, axis=0, method=None, level=None, copy=True, limit=None, fill\_value=nan)

Conform input object to new index with optional filling logic, placing NA/NaN in locations having no value in the previous index. A new object is produced unless the new index is equivalent to the current one and copy=False

#### **Parameters**

labels [array-like] New labels / index to conform to. Preferably an Index object to avoid duplicating data

```
axis [{0 or 'index', 1 or 'columns'}]
```

**method** [{None, 'backfill', 'pad'/'ffill', 'nearest'}, optional] Method to use for filling holes in reindexed DataFrame:

- default: don't fill gaps
- pad / ffill: propagate last valid observation forward to next valid
- backfill / bfill: use next valid observation to fill gap
- nearest: use nearest valid observations to fill gap

copy [boolean, default True] Return a new object, even if the passed indexes are the same

**level** [int or name] Broadcast across a level, matching Index values on the passed MultiIndex level

**limit** [int, default None] Maximum number of consecutive elements to forward or backward fill

tolerance [optional] Maximum distance between original and new labels for inexact
matches. The values of the index at the matching locations most satisfy the equation
abs(index[indexer] - target) <= tolerance.</pre>

Tolerance may be a scalar value, which applies the same tolerance to all values, or list-like, which applies variable tolerance per element. List-like includes list, tuple, array, Series, and must be the same size as the index and its dtype must exactly match the index's type.

New in version 0.21.0: (list-like tolerance)

# Returns

reindexed [DataFrame]

#### See also:

reindex, reindex like

# **Examples**

```
>>> df.reindex_axis(['A', 'B', 'C'], axis=1)
```

**reindex\_like** (*other*, *method=None*, *copy=True*, *limit=None*, *tolerance=None*) Return an object with matching indices to myself.

## **Parameters**

```
other [Object]
method [string or None]
```

```
copy [boolean, default True]
```

**limit** [int, default None] Maximum number of consecutive labels to fill for inexact matches.

**tolerance** [optional] Maximum distance between labels of the other object and this object for inexact matches. Can be list-like.

New in version 0.21.0: (list-like tolerance)

#### Returns

reindexed [same as input]

#### **Notes**

**Like calling s.reindex(index=other.index, columns=other.columns,** method=...)

```
rename (**kwargs)
```

Alter axes labels.

Function / dict values must be unique (1-to-1). Labels not contained in a dict / Series will be left as-is. Extra labels listed don't throw an error.

See the user guide for more.

#### **Parameters**

mapper, index, columns [dict-like or function, optional] dict-like or functions transformations to apply to that axis' values. Use either mapper and axis to specify the axis to target with mapper, or index and columns.

**axis** [int or str, optional] Axis to target with mapper. Can be either the axis name ('index', 'columns') or number (0, 1). The default is 'index'.

copy [boolean, default True] Also copy underlying data

**inplace** [boolean, default False] Whether to return a new DataFrame. If True then value of copy is ignored.

**level** [int or level name, default None] In case of a MultiIndex, only rename labels in the specified level.

### Returns

renamed [DataFrame]

### See also:

```
pandas.DataFrame.rename axis
```

# **Examples**

DataFrame.rename supports two calling conventions

- (index=index\_mapper, columns=columns\_mapper, ...)
- (mapper, axis={'index', 'columns'}, ...)

We highly recommend using keyword arguments to clarify your intent.

```
>>> df.rename(index=str, columns={"A": "a", "C": "c"})

a B
0 1 4
1 2 5
2 3 6
```

## Using axis-style parameters

```
>>> df.rename(str.lower, axis='columns')
    a    b
0    1    4
1    2    5
2    3    6
```

```
>>> df.rename({1: 2, 2: 4}, axis='index')

A B
0 1 4
2 2 5
4 3 6
```

#### rename\_axis (mapper, axis=0, copy=True, inplace=False)

Alter the name of the index or columns.

# **Parameters**

```
mapper [scalar, list-like, optional] Value to set as the axis name attribute.
```

axis [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis.

copy [boolean, default True] Also copy underlying data.

**inplace** [boolean, default False] Modifies the object directly, instead of creating a new Series or DataFrame.

## Returns

**renamed** [Series, DataFrame, or None] The same type as the caller or None if *inplace* is True.

# See also:

```
pandas. Series. rename Alter Series index labels or name
```

pandas.DataFrame.rename Alter DataFrame index labels or name

pandas.Index.rename Set new names on index

### **Notes**

Prior to version 0.21.0, rename\_axis could also be used to change the axis *labels* by passing a mapping or scalar. This behavior is deprecated and will be removed in a future version. Use rename instead.

# **Examples**

#### Series

```
>>> s = pd.Series([1, 2, 3])
>>> s.rename_axis("foo")
foo
0    1
1    2
2    3
dtype: int64
```

### **DataFrame**

```
>>> df.rename_axis("bar", axis="columns")
bar A B
0 1 4
1 2 5
2 3 6
```

## reorder\_levels (order, axis=0)

Rearrange index levels using input order. May not drop or duplicate levels

### **Parameters**

**order** [list of int or list of str] List representing new level order. Reference level by number (position) or by key (label).

axis [int] Where to reorder levels.

## Returns

# type of caller (new object)

**replace** (to\_replace=None, value=None, inplace=False, limit=None, regex=False, method='pad') Replace values given in to\_replace with value.

Values of the DataFrame are replaced with other values dynamically. This differs from updating with .loc or .iloc, which require you to specify a location to update with some value.

### **Parameters**

**to\_replace** [str, regex, list, dict, Series, int, float, or None] How to find the values that will be replaced.

- numeric, str or regex:
  - numeric: numeric values equal to to\_replace will be replaced with value
  - str: string exactly matching to\_replace will be replaced with value
  - regex: regexs matching to\_replace will be replaced with value
- list of str, regex, or numeric:

- First, if *to\_replace* and *value* are both lists, they **must** be the same length.
- Second, if regex=True then all of the strings in both lists will be interpreted as
  regexs otherwise they will match directly. This doesn't matter much for *value* since
  there are only a few possible substitution regexes you can use.
- str, regex and numeric rules apply as above.

#### · dict:

- Dicts can be used to specify different replacement values for different existing values.
   For example, { 'a': 'b', 'y': 'z'} replaces the value 'a' with 'b' and 'y' with 'z'. To use a dict in this way the *value* parameter should be *None*.
- For a DataFrame a dict can specify that different values should be replaced in different columns. For example, { 'a': 1, 'b': 'z'} looks for the value 1 in column 'a' and the value 'z' in column 'b' and replaces these values with whatever is specified in *value*. The *value* parameter should not be None in this case. You can treat this as a special case of passing two lists except that you are specifying the column to search in.
- For a DataFrame nested dictionaries, e.g., { 'a': {'b': np.nan}}, are read as follows: look in column 'a' for the value 'b' and replace it with NaN. The *value* parameter should be None to use a nested dict in this way. You can nest regular expressions as well. Note that column names (the top-level dictionary keys in a nested dictionary) **cannot** be regular expressions.

#### • None:

This means that the *regex* argument must be a string, compiled regular expression, or
list, dict, ndarray or Series of such elements. If *value* is also None then this **must** be
a nested dictionary or Series.

See the examples section for examples of each of these.

value [scalar, dict, list, str, regex, default None] Value to replace any values matching to\_replace with. For a DataFrame a dict of values can be used to specify which value to use for each column (columns not in the dict will not be filled). Regular expressions, strings and lists or dicts of such objects are also allowed.

**inplace** [boolean, default False] If True, in place. Note: this will modify any other views on this object (e.g. a column from a DataFrame). Returns the caller if this is True.

limit [int, default None] Maximum size gap to forward or backward fill.

**regex** [bool or same types as *to\_replace*, default False] Whether to interpret *to\_replace* and/or *value* as regular expressions. If this is True then *to\_replace must* be a string. Alternatively, this could be a regular expression or a list, dict, or array of regular expressions in which case *to replace* must be None.

**method** [{'pad', 'ffill', 'bfill', *None*}] The method to use when for replacement, when *to\_replace* is a scalar, list or tuple and *value* is None.

Changed in version 0.23.0: Added to DataFrame.

#### Returns

DataFrame Object after replacement.

# Raises

#### AssertionError

If regex is not a bool and to\_replace is not None.

# **TypeError**

- If to\_replace is a dict and value is not a list, dict, ndarray, or Series
- If *to\_replace* is None and *regex* is not compilable into a regular expression or is a list, dict, ndarray, or Series.
- When replacing multiple bool or datetime64 objects and the arguments to to\_replace does not match the type of the value being replaced

#### ValueError

• If a list or an ndarray is passed to *to\_replace* and *value* but they are not the same length.

#### See also:

DataFrame.fillna Fill NA values

DataFrame.where Replace values based on boolean condition

Series.str.replace Simple string replacement.

### **Notes**

- Regex substitution is performed under the hood with re.sub. The rules for substitution for re.sub are the same.
- Regular expressions will only substitute on strings, meaning you cannot provide, for example, a regular expression matching floating point numbers and expect the columns in your frame that have a numeric dtype to be matched. However, if those floating point numbers *are* strings, then you can do this.
- This method has *a lot* of options. You are encouraged to experiment and play with this method to gain intuition about how it works.
- When dict is used as the *to\_replace* value, it is like key(s) in the dict are the to\_replace part and value(s) in the dict are the value parameter.

### **Examples**

# Scalar 'to\_replace' and 'value'

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```
1 1 6 b
2 2 7 c
3 3 8 d
4 4 9 e
```

# List-like 'to\_replace'

```
>>> df.replace([0, 1, 2, 3], 4)

A B C
0 4 5 a
1 4 6 b
2 4 7 c
3 4 8 d
4 4 9 e
```

```
>>> df.replace([0, 1, 2, 3], [4, 3, 2, 1])

A B C
0 4 5 a
1 3 6 b
2 2 7 c
3 1 8 d
4 4 9 e
```

# dict-like 'to\_replace'

```
>>> df.replace({0: 10, 1: 100})

A B C

0 10 5 a

1 100 6 b

2 2 7 c

3 3 8 d

4 4 9 e
```

```
>>> df.replace({'A': 0, 'B': 5}, 100)

A B C

0 100 100 a

1 1 6 b

2 2 7 c

3 3 8 d

4 4 9 e
```

```
>>> df.replace({'A': {0: 100, 4: 400}})

A B C

0 100 5 a

1 1 6 b

2 2 7 c

3 3 8 d
```

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```
4 400 9 e
```

# Regular expression 'to\_replace'

```
>>> df.replace({'A': r'^ba.$'}, {'A': 'new'}, regex=True)

A B

0 new abc

1 foo bar

2 bait xyz
```

```
>>> df.replace(regex=r'^ba.$', value='new')

A B

0 new abc

1 foo new

2 bait xyz
```

```
>>> df.replace(regex={r'^ba.$':'new', 'foo':'xyz'})
        A      B
0        new abc
1        xyz        new
2        bait       xyz
```

Note that when replacing multiple bool or datetime64 objects, the data types in the *to\_replace* parameter must match the data type of the value being replaced:

```
>>> df = pd.DataFrame({'A': [True, False, True],
... 'B': [False, True, False]})
>>> df.replace({'a string': 'new value', True: False}) # raises
Traceback (most recent call last):
...
TypeError: Cannot compare types 'ndarray(dtype=bool)' and 'str'
```

This raises a TypeError because one of the dict keys is not of the correct type for replacement.

Compare the behavior of s.replace ( $\{'a': None\}$ ) and s.replace ('a', None) to understand the pecularities of the *to\_replace* parameter:

```
>>> s = pd.Series([10, 'a', 'a', 'b', 'a'])
```

When one uses a dict as the *to\_replace* value, it is like the value(s) in the dict are equal to the *value* parameter. s.replace({'a': None}) is equivalent to s.replace(to\_replace={'a': None}, value=None, method=None):

```
>>> s.replace({'a': None})
0 10
1 None
2 None
3 b
4 None
dtype: object
```

When value=None and to\_replace is a scalar, list or tuple, replace uses the method parameter (default 'pad') to do the replacement. So this is why the 'a' values are being replaced by 10 in rows 1 and 2 and 'b' in row 4 in this case. The command s.replace('a', None) is actually equivalent to s. replace(to\_replace='a', value=None, method='pad'):

```
>>> s.replace('a', None)
0    10
1    10
2    10
3         b
4         b
dtype: object
```

resample (rule, how=None, axis=0, fill\_method=None, closed=None, label=None, convention='start', kind=None, loffset=None, limit=None, base=0, on=None, level=None)

Convenience method for frequency conversion and resampling of time series. Object must have a datetime-like index (DatetimeIndex, PeriodIndex, or TimedeltaIndex), or pass datetime-like values to the on or level keyword.

## **Parameters**

rule [string] the offset string or object representing target conversion

axis [int, optional, default 0]

**closed** [{'right', 'left'}] Which side of bin interval is closed. The default is 'left' for all frequency offsets except for 'M', 'A', 'Q', 'BM', 'BA', 'BQ', and 'W' which all have a default of 'right'.

label [{'right', 'left'}] Which bin edge label to label bucket with. The default is 'left' for all frequency offsets except for 'M', 'A', 'Q', 'BM', 'BA', 'BQ', and 'W' which all have a default of 'right'.

**convention** [{'start', 'end', 's', 'e'}] For PeriodIndex only, controls whether to use the start or end of *rule* 

kind: {'timestamp', 'period'}, optional Pass 'timestamp' to convert the resulting index to a DateTimeIndex or 'period' to convert it to a PeriodIndex. By default the input representation is retained.

**loffset** [timedelta] Adjust the resampled time labels

**base** [int, default 0] For frequencies that evenly subdivide 1 day, the "origin" of the aggregated intervals. For example, for '5min' frequency, base could range from 0 through 4. Defaults to 0

**on** [string, optional] For a DataFrame, column to use instead of index for resampling. Column must be datetime-like.

New in version 0.19.0.

**level** [string or int, optional] For a MultiIndex, level (name or number) to use for resampling. Level must be datetime-like.

New in version 0.19.0.

## Returns

# Resampler object

## See also:

**groupby** Group by mapping, function, label, or list of labels.

### **Notes**

See the user guide for more.

To learn more about the offset strings, please see this link.

# **Examples**

Start by creating a series with 9 one minute timestamps.

```
>>> index = pd.date_range('1/1/2000', periods=9, freq='T')
>>> series = pd.Series(range(9), index=index)
>>> series
2000-01-01 00:00:00
                       0
2000-01-01 00:01:00
                       1
2000-01-01 00:02:00
                       2
2000-01-01 00:03:00
2000-01-01 00:04:00
                       4
2000-01-01 00:05:00
                       5
2000-01-01 00:06:00
                       6
2000-01-01 00:07:00
                       7
2000-01-01 00:08:00
                       8
Freq: T, dtype: int64
```

Downsample the series into 3 minute bins and sum the values of the timestamps falling into a bin.

Downsample the series into 3 minute bins as above, but label each bin using the right edge instead of the left. Please note that the value in the bucket used as the label is not included in the bucket, which it labels. For example, in the original series the bucket 2000-01-01 00:03:00 contains the value 3, but the summed value in the resampled bucket with the label 2000-01-01 00:03:00 does not include 3 (if it did, the summed value would be 6, not 3). To include this value close the right side of the bin interval as illustrated in the example below this one.

Downsample the series into 3 minute bins as above, but close the right side of the bin interval.

Upsample the series into 30 second bins.

```
>>> series.resample('30S').asfreq()[0:5] #select first 5 rows
2000-01-01 00:00:00 0.0
2000-01-01 00:00:30 NaN
2000-01-01 00:01:00 1.0
2000-01-01 00:01:30 NaN
2000-01-01 00:02:00 2.0
Freq: 30S, dtype: float64
```

Upsample the series into 30 second bins and fill the NaN values using the pad method.

Upsample the series into 30 second bins and fill the NaN values using the bfill method.

Pass a custom function via apply

```
>>> def custom_resampler(array_like):
... return np.sum(array_like)+5
```

For a Series with a PeriodIndex, the keyword *convention* can be used to control whether to use the start or end of *rule*.

```
>>> s = pd.Series([1, 2], index=pd.period_range('2012-01-01', freq='A', periods=2))
>>> s
2012 1
```

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```
2013 2
Freq: A-DEC, dtype: int64
```

Resample by month using 'start' convention. Values are assigned to the first month of the period.

Resample by month using 'end' convention. Values are assigned to the last month of the period.

```
>>> s.resample('M', convention='end').asfreq()
2012-12
          1.0
2013-01
          NaN
2013-02
          NaN
2013-03
          NaN
2013-04
          NaN
2013-05
          NaN
2013-06
          NaN
2013-07
          NaN
2013-08
          NaN
2013-09
          NaN
2013-10
          NaN
2013-11
          NaN
2013-12
           2.0
Freq: M, dtype: float64
```

For DataFrame objects, the keyword on can be used to specify the column instead of the index for resampling.

For a DataFrame with MultiIndex, the keyword level can be used to specify on level the resampling needs to take place.

reset\_index (level=None, drop=False, inplace=False, col\_level=0, col\_fill=")

For DataFrame with multi-level index, return new DataFrame with labeling information in the columns under the index names, defaulting to 'level\_0', 'level\_1', etc. if any are None. For a standard index, the index name will be used (if set), otherwise a default 'index' or 'level\_0' (if 'index' is already taken) will be used.

## **Parameters**

**level** [int, str, tuple, or list, default None] Only remove the given levels from the index. Removes all levels by default

**drop** [boolean, default False] Do not try to insert index into dataframe columns. This resets the index to the default integer index.

**inplace** [boolean, default False] Modify the DataFrame in place (do not create a new object)

**col\_level** [int or str, default 0] If the columns have multiple levels, determines which level the labels are inserted into. By default it is inserted into the first level.

**col\_fill** [object, default ''] If the columns have multiple levels, determines how the other levels are named. If None then the index name is repeated.

### Returns

resetted [DataFrame]

# **Examples**

```
>>> df = pd.DataFrame([('bird',
                               389.0),
                      ('bird',
                                 24.0),
                      ('mammal', 80.5),
                      ('mammal', np.nan)],
                     index=['falcon', 'parrot', 'lion', 'monkey'],
. . .
                     columns=('class', 'max_speed'))
. . .
>>> df
        class max_speed
falcon
       bird
               389.0
                   24.0
parrot
        bird
lion
                    80.5
       mammal
monkey mammal
                     NaN
```

When we reset the index, the old index is added as a column, and a new sequential index is used:

```
>>> df.reset_index()
   index class max_speed
0 falcon bird 389.0
1 parrot bird 24.0
2 lion mammal 80.5
3 monkey mammal NaN
```

We can use the *drop* parameter to avoid the old index being added as a column:

```
>>> df.reset_index(drop=True)
class max_speed
0 bird 389.0
1 bird 24.0
2 mammal 80.5
3 mammal NaN
```

You can also use reset index with MultiIndex.

```
>>> index = pd.MultiIndex.from_tuples([('bird', 'falcon'),
                                         ('bird', 'parrot'),
. . .
                                         ('mammal', 'lion'),
. . .
                                         ('mammal', 'monkey')],
. . .
                                       names=['class', 'name'])
. . .
>>> columns = pd.MultiIndex.from_tuples([('speed', 'max'),
                                           ('species', 'type')])
>>> df = pd.DataFrame([(389.0, 'fly'),
                        ( 24.0, 'fly'),
                        ( 80.5, 'run'),
. . .
                       (np.nan, 'jump')],
. . .
                      index=index,
. . .
                      columns=columns)
>>> df
               speed species
                max
                      type
class name
      falcon 389.0
bird
                        fly
       parrot 24.0
                        fly
               80.5
mammal lion
                         run
      monkey NaN
                        jump
```

If the index has multiple levels, we can reset a subset of them:

```
>>> df.reset_index(level='class')
        class speed species
                max
                     type
name
falcon
        bird 389.0
                       fly
        bird
               24.0
parrot
                        fly
               80.5
lion
      mammal
                        run
monkey mammal
              NaN
                       jump
```

If we are not dropping the index, by default, it is placed in the top level. We can place it in another level:

```
>>> df.reset_index(level='class', col_level=1)
               speed species
        class
               max type
name
        bird 389.0
                        fly
falcon
parrot
         bird
               24.0
                        fly
                80.5
lion
       mammal
                        run
monkey mammal
                NaN
                        jump
```

When the index is inserted under another level, we can specify under which one with the parameter *col\_fill*:

```
>>> df.reset_index(level='class', col_level=1, col_fill='species')
             species speed species
              class
                     max
name
               bird 389.0
falcon
                              fly
                     24.0
parrot
              bird
                              fly
lion
             mammal
                     80.5
                              run
monkey
            mammal NaN
                              jump
```

If we specify a nonexistent level for *col\_fill*, it is created:

```
>>> df.reset_index(level='class', col_level=1, col_fill='genus')
               genus speed species
               class
                      max
                              type
name
                bird 389.0
falcon
                               fly
                      24.0
                               fly
parrot
                bird
lion
              mammal
                      80.5
                                run
monkey
              mammal
                       NaN
                               jump
```

**rfloordiv** (other, axis='columns', level=None, fill\_value=None)

Integer division of dataframe and other, element-wise (binary operator rfloordiv).

Equivalent to other // dataframe, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

other [Series, DataFrame, or constant]

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

#### Returns

result [DataFrame]

## See also:

DataFrame.floordiv

## **Notes**

Mismatched indices will be unioned together

# **Examples**

None

rmod (other, axis='columns', level=None, fill\_value=None)

Modulo of dataframe and other, element-wise (binary operator *rmod*).

Equivalent to other % dataframe, but with support to substitute a fill\_value for missing data in one of the inputs.

# **Parameters**

**other** [Series, DataFrame, or constant]

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

## Returns

result [DataFrame]

### See also:

DataFrame.mod

## **Notes**

Mismatched indices will be unioned together

# **Examples**

None

rmul (other, axis='columns', level=None, fill\_value=None)

Multiplication of dataframe and other, element-wise (binary operator rmul).

Equivalent to other \* dataframe, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

```
other [Series, DataFrame, or constant]
```

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

## **Returns**

result [DataFrame]

## See also:

DataFrame.mul

## **Notes**

Mismatched indices will be unioned together

# **Examples**

None

 $\begin{tabular}{ll} \textbf{rolling} (window, & min\_periods=None, & center=False, & win\_type=None, & on=None, & axis=0, \\ & & closed=None) \end{tabular}$ 

Provides rolling window calculations.

New in version 0.18.0.

#### **Parameters**

**window** [int, or offset] Size of the moving window. This is the number of observations used for calculating the statistic. Each window will be a fixed size.

If its an offset then this will be the time period of each window. Each window will be a variable sized based on the observations included in the time-period. This is only valid for datetimelike indexes. This is new in 0.19.0

**min\_periods** [int, default None] Minimum number of observations in window required to have a value (otherwise result is NA). For a window that is specified by an offset, this will default to 1.

**center** [boolean, default False] Set the labels at the center of the window.

win\_type [string, default None] Provide a window type. If None, all points are evenly weighted. See the notes below for further information.

**on** [string, optional] For a DataFrame, column on which to calculate the rolling window, rather than the index

**closed** [string, default None] Make the interval closed on the 'right', 'left', 'both' or 'neither' endpoints. For offset-based windows, it defaults to 'right'. For fixed windows, defaults to 'both'. Remaining cases not implemented for fixed windows.

New in version 0.20.0.

axis [int or string, default 0]

## Returns

a Window or Rolling sub-classed for the particular operation

## See also:

expanding Provides expanding transformations.

ewm Provides exponential weighted functions

# Notes

By default, the result is set to the right edge of the window. This can be changed to the center of the window by setting center=True.

To learn more about the offsets & frequency strings, please see this link.

The recognized win\_types are:

- boxcar
- triang
- blackman
- hamming
- bartlett
- parzen

- bohman
- blackmanharris
- nuttall
- barthann
- kaiser (needs beta)
- gaussian (needs std)
- general\_gaussian (needs power, width)
- slepian (needs width).

If win\_type=None all points are evenly weighted. To learn more about different window types see scipy.signal window functions.

# **Examples**

```
>>> df = pd.DataFrame({'B': [0, 1, 2, np.nan, 4]})
>>> df
B
0 0.0
1 1.0
2 2.0
3 NaN
4 4.0
```

Rolling sum with a window length of 2, using the 'triang' window type.

Rolling sum with a window length of 2, min\_periods defaults to the window length.

```
>>> df.rolling(2).sum()
B
0 NaN
1 1.0
2 3.0
3 NaN
4 NaN
```

Same as above, but explicitly set the min\_periods

```
>>> df.rolling(2, min_periods=1).sum()

B
0 0.0
1 1.0
2 3.0
3 2.0
4 4.0
```

A ragged (meaning not-a-regular frequency), time-indexed DataFrame

```
>>> df

B

2013-01-01 09:00:00 0.0
2013-01-01 09:00:02 1.0
2013-01-01 09:00:03 2.0
2013-01-01 09:00:05 NaN
2013-01-01 09:00:06 4.0
```

Contrasting to an integer rolling window, this will roll a variable length window corresponding to the time period. The default for min\_periods is 1.

```
>>> df.rolling('2s').sum()

B

2013-01-01 09:00:00 0.0

2013-01-01 09:00:02 1.0

2013-01-01 09:00:03 3.0

2013-01-01 09:00:05 NaN

2013-01-01 09:00:06 4.0
```

round (decimals=0, \*args, \*\*kwargs)

Round a DataFrame to a variable number of decimal places.

# **Parameters**

**decimals** [int, dict, Series] Number of decimal places to round each column to. If an int is given, round each column to the same number of places. Otherwise dict and Series round to variable numbers of places. Column names should be in the keys if *decimals* is a dict-like, or in the index if *decimals* is a Series. Any columns not included in *decimals* will be left as is. Elements of *decimals* which are not columns of the input will be ignored.

## Returns

## DataFrame object

## See also:

numpy.around, Series.round

## **Examples**

```
>>> df = pd.DataFrame(np.random.random([3, 3]),
... columns=['A', 'B', 'C'], index=['first', 'second', 'third'])
>>> df

A B C
first 0.028208 0.992815 0.173891
second 0.038683 0.645646 0.577595
third 0.877076 0.149370 0.491027
```

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```
>>> df.round(2)
               В
                     С
         Α
       0.03 0.99 0.17
first
second 0.04 0.65 0.58
third 0.88 0.15 0.49
>>> df.round({'A': 1, 'C': 2})
                  В
        Α
first
       0.0
           0.992815 0.17
second 0.0
           0.645646 0.58
       0.9 0.149370 0.49
third
>>> decimals = pd.Series([1, 0, 2], index=['A', 'B', 'C'])
>>> df.round(decimals)
        A B
       0.0 1 0.17
first
second 0.0 1 0.58
third
       0.9 0 0.49
```

**rpow** (other, axis='columns', level=None, fill\_value=None)

Exponential power of dataframe and other, element-wise (binary operator rpow).

Equivalent to other \*\* dataframe, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

```
other [Series, DataFrame, or constant]
```

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

## Returns

result [DataFrame]

## See also:

DataFrame.pow

## **Notes**

Mismatched indices will be unioned together

# **Examples**

None

```
rsub (other, axis='columns', level=None, fill_value=None)
```

Subtraction of dataframe and other, element-wise (binary operator rsub).

Equivalent to other - dataframe, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

```
other [Series, DataFrame, or constant]
```

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

fill\_value [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

### Returns

result [DataFrame]

## See also:

DataFrame.sub

## **Notes**

Mismatched indices will be unioned together

# **Examples**

```
>>> a = pd.DataFrame([2, 1, 1, np.nan], index=['a', 'b', 'c', 'd'],
                    columns=['one'])
>>> a
  one
  2.0
а
b 1.0
 1.0
C
d NaN
>>> b = pd.DataFrame(dict(one=[1, np.nan, 1, np.nan],
                          two=[3, 2, np.nan, 2]),
                     index=['a', 'b', 'd', 'e'])
. . .
>>> b
  one
       two
       3.0
  1.0
 NaN 2.0
b
d
  1.0
       NaN
  NaN 2.0
>>> a.sub(b, fill_value=0)
  one two
       -3.0
  1.0
       -2.0
  1.0
  1.0
       NaN
  -1.0
        NaN
е
  NaN
       -2.0
```

rtruediv (other, axis='columns', level=None, fill\_value=None)

Floating division of dataframe and other, element-wise (binary operator *rtruediv*).

Equivalent to other / dataframe, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

**other** [Series, DataFrame, or constant]

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

fill\_value [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

## Returns

result [DataFrame]

## See also:

DataFrame.truediv

## **Notes**

Mismatched indices will be unioned together

# **Examples**

None

**sample** (*n*=*None*, *frac*=*None*, *replace*=*False*, *weights*=*None*, *random\_state*=*None*, *axis*=*None*) Return a random sample of items from an axis of object.

You can use *random\_state* for reproducibility.

## **Parameters**

**n** [int, optional] Number of items from axis to return. Cannot be used with frac. Default = 1 if frac = None.

**frac** [float, optional] Fraction of axis items to return. Cannot be used with n.

**replace** [boolean, optional] Sample with or without replacement. Default = False.

weights [str or ndarray-like, optional] Default 'None' results in equal probability weighting. If passed a Series, will align with target object on index. Index values in weights not found in sampled object will be ignored and index values in sampled object not in weights will be assigned weights of zero. If called on a DataFrame, will accept the name of a column when axis = 0. Unless weights are a Series, weights must be same length as axis being sampled. If weights do not sum to 1, they will be normalized to sum to 1. Missing values in the weights column will be treated as zero. inf and -inf values not allowed.

**random\_state** [int or numpy.random.RandomState, optional] Seed for the random number generator (if int), or numpy RandomState object.

**axis** [int or string, optional] Axis to sample. Accepts axis number or name. Default is stat axis for given data type (0 for Series and DataFrames, 1 for Panels).

### Returns

A new object of same type as caller.

# **Examples**

Generate an example Series and DataFrame:

```
>>> s = pd.Series(np.random.randn(50))
>>> s.head()
  -0.038497
0
1
    1.820773
2
  -0.972766
3
  -1.598270
  -1.095526
dtype: float64
>>> df = pd.DataFrame(np.random.randn(50, 4), columns=list('ABCD'))
>>> df.head()
                   В
0 0.016443 -2.318952 -0.566372 -1.028078
1 -1.051921 0.438836 0.658280 -0.175797
2 -1.243569 -0.364626 -0.215065 0.057736
  1.768216 0.404512 -0.385604 -1.457834
4 1.072446 -1.137172 0.314194 -0.046661
```

Next extract a random sample from both of these objects...

3 random elements from the Series:

```
>>> s.sample(n=3)

27 -0.994689

55 -1.049016

67 -0.224565

dtype: float64
```

And a random 10% of the DataFrame with replacement:

```
>>> df.sample(frac=0.1, replace=True)

A B C D

35 1.981780 0.142106 1.817165 -0.290805
49 -1.336199 -0.448634 -0.789640 0.217116
40 0.823173 -0.078816 1.009536 1.015108
15 1.421154 -0.055301 -1.922594 -0.019696
6 -0.148339 0.832938 1.787600 -1.383767
```

You can use random state for reproducibility:

```
>>> df.sample(random_state=1)
A B C D
37 -2.027662 0.103611 0.237496 -0.165867
43 -0.259323 -0.583426 1.516140 -0.479118
12 -1.686325 -0.579510 0.985195 -0.460286
8 1.167946 0.429082 1.215742 -1.636041
9 1.197475 -0.864188 1.554031 -1.505264
```

select (crit, axis=0)

Return data corresponding to axis labels matching criteria

Deprecated since version 0.21.0: Use df.loc[df.index.map(crit)] to select via labels

## **Parameters**

```
crit [function] To be called on each index (label). Should return True or False
axis [int]
```

## Returns

**selection** [type of caller]

```
select dtypes (include=None, exclude=None)
```

Return a subset of the DataFrame's columns based on the column dtypes.

## **Parameters**

**include, exclude** [scalar or list-like] A selection of dtypes or strings to be included/excluded. At least one of these parameters must be supplied.

## **Returns**

**subset** [DataFrame] The subset of the frame including the dtypes in include and excluding the dtypes in exclude.

### Raises

### ValueError

- If both of include and exclude are empty
- If include and exclude have overlapping elements
- If any kind of string dtype is passed in.

## **Notes**

- To select all numeric types, use np.number or 'number'
- To select strings you must use the object dtype, but note that this will return all object dtype columns
- See the numpy dtype hierarchy
- To select datetimes, use np.datetime64, 'datetime' or 'datetime64'
- To select timedeltas, use np.timedelta64, 'timedelta' or 'timedelta64'
- To select Pandas categorical dtypes, use 'category'
- To select Pandas datetimetz dtypes, use 'datetimetz' (new in 0.20.0) or 'datetime64 [ns, tz]'

# **Examples**

```
>>> df = pd.DataFrame({'a': [1, 2] * 3,
                        'b': [True, False] * 3,
. . .
                        'c': [1.0, 2.0] * 3)
. . .
>>> df
               h
        а
            True 1.0
0
        1
        2 False 2.0
1
2
        1
           True
                 1.0
3
        2 False
                 2.0
```

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```
4 1 True 1.0
5 2 False 2.0
```

```
>>> df.select_dtypes(exclude=['int'])
0
    True
          1.0
          2.0
1
  False
2
          1.0
   True
3
          2.0
  False
4
    True
          1.0
5
  False
          2.0
```

**sem** (axis=None, skipna=None, level=None, ddof=1, numeric\_only=None, \*\*kwargs)
Return unbiased standard error of the mean over requested axis.

Normalized by N-1 by default. This can be changed using the ddof argument

## **Parameters**

```
axis [{index (0), columns (1)}]
```

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**ddof** [int, default 1] Delta Degrees of Freedom. The divisor used in calculations is N - ddof, where N represents the number of elements.

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

## Returns

sem [Series or DataFrame (if level specified)]

set\_axis (labels, axis=0, inplace=None)

Assign desired index to given axis.

Indexes for column or row labels can be changed by assigning a list-like or Index.

Changed in version 0.21.0: The signature is now *labels* and *axis*, consistent with the rest of pandas API. Previously, the *axis* and *labels* arguments were respectively the first and second positional arguments.

### **Parameters**

labels [list-like, Index] The values for the new index.

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The axis to update. The value 0 identifies the rows, and 1 identifies the columns.

**inplace** [boolean, default None] Whether to return a new %(klass)s instance.

**Warning:** inplace=None currently falls back to to True, but in a future version, will default to False. Use inplace=True explicitly rather than relying on the default.

## **Returns**

**renamed** [%(klass)s or None] An object of same type as caller if inplace=False, None otherwise.

See also:

pandas.DataFrame.rename axis Alter the name of the index or columns.

## **Examples**

## **Series**

The original object is not modified.

```
>>> s
0    1
1    2
2    3
dtype: int64
```

## **DataFrame**

```
>>> df = pd.DataFrame({"A": [1, 2, 3], "B": [4, 5, 6]})
```

Change the row labels.

```
>>> df.set_axis(['a', 'b', 'c'], axis='index', inplace=False)

A B
a 1 4
b 2 5
c 3 6
```

Change the column labels.

Now, update the labels inplace.

set\_index (keys, drop=True, append=False, inplace=False, verify\_integrity=False)

Set the DataFrame index (row labels) using one or more existing columns. By default yields a new object.

## **Parameters**

**keys** [column label or list of column labels / arrays]

**drop** [boolean, default True] Delete columns to be used as the new index

append [boolean, default False] Whether to append columns to existing index

**inplace** [boolean, default False] Modify the DataFrame in place (do not create a new object)

verify\_integrity [boolean, default False] Check the new index for duplicates. Otherwise defer the check until necessary. Setting to False will improve the performance of this method

## Returns

dataframe [DataFrame]

## **Examples**

```
>>> df = pd.DataFrame({'month': [1, 4, 7, 10],
                        'year': [2012, 2014, 2013, 2014],
                       'sale':[55, 40, 84, 31]})
  month sale year
          55
                2012
  1
                2014
  4
          40
2
  7
          84
                2013
  10
          31
                2014
```

Set the index to become the 'month' column:

Create a multi-index using columns 'year' and 'month':

Create a multi-index using a set of values and a column:

# set\_value (index, col, value, takeable=False)

Put single value at passed column and index

Deprecated since version 0.21.0: Use .at[] or .iat[] accessors instead.

# **Parameters**

```
index [row label]col [column label]value [scalar value]takeable [interpret the index/col as indexers, default False]
```

# Returns

**frame** [DataFrame] If label pair is contained, will be reference to calling DataFrame, otherwise a new object

## shape

Return a tuple representing the dimensionality of the DataFrame.

# See also:

```
ndarray.shape
```

# **Examples**

```
>>> df = pd.DataFrame({'col1': [1, 2], 'col2': [3, 4]})
>>> df.shape
(2, 2)
```

```
>>> df = pd.DataFrame({'col1': [1, 2], 'col2': [3, 4],
... 'col3': [5, 6]})
>>> df.shape
(2, 3)
```

## **shift** (*periods*=1, *freq*=None, *axis*=0)

Shift index by desired number of periods with an optional time freq

## **Parameters**

periods [int] Number of periods to move, can be positive or negative

**freq** [DateOffset, timedelta, or time rule string, optional] Increment to use from the tseries module or time rule (e.g. 'EOM'). See Notes.

```
axis [{0 or 'index', 1 or 'columns'}]
```

### Returns

shifted [DataFrame]

### **Notes**

If freq is specified then the index values are shifted but the data is not realigned. That is, use freq if you would like to extend the index when shifting and preserve the original data.

## size

Return an int representing the number of elements in this object.

Return the number of rows if Series. Otherwise return the number of rows times number of columns if DataFrame.

## See also:

ndarray.size

# **Examples**

```
>>> s = pd.Series({'a': 1, 'b': 2, 'c': 3})
>>> s.size
3
```

```
>>> df = pd.DataFrame({'col1': [1, 2], 'col2': [3, 4]})
>>> df.size
4
```

**skew** (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

Return unbiased skew over requested axis Normalized by N-1

# **Parameters**

```
axis [{index (0), columns (1)}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

## Returns

**skew** [Series or DataFrame (if level specified)]

## slice shift (periods=1, axis=0)

Equivalent to *shift* without copying data. The shifted data will not include the dropped periods and the shifted axis will be smaller than the original.

## **Parameters**

**periods** [int] Number of periods to move, can be positive or negative

### Returns

**shifted** [same type as caller]

## **Notes**

While the *slice\_shift* is faster than *shift*, you may pay for it later during alignment.

## **Parameters**

axis [index, columns to direct sorting]

**level** [int or level name or list of ints or list of level names] if not None, sort on values in specified index level(s)

ascending [boolean, default True] Sort ascending vs. descending

inplace [bool, default False] if True, perform operation in-place

**kind** [{'quicksort', 'mergesort', 'heapsort'}, default 'quicksort'] Choice of sorting algorithm. See also ndarray.np.sort for more information. *mergesort* is the only stable algorithm. For DataFrames, this option is only applied when sorting on a single column or label.

**na\_position** [{'first', 'last'}, default 'last'] *first* puts NaNs at the beginning, *last* puts NaNs at the end. Not implemented for MultiIndex.

**sort\_remaining** [bool, default True] if true and sorting by level and index is multilevel, sort by other levels too (in order) after sorting by specified level

## Returns

```
sorted_obj [DataFrame]
```

**sort\_values** (*by*, *axis*=0, *ascending*=*True*, *inplace*=*False*, *kind*='*quicksort*', *na\_position*='*last*') Sort by the values along either axis

## **Parameters**

by [str or list of str] Name or list of names to sort by.

- if axis is 0 or 'index' then by may contain index levels and/or column labels
- if axis is 1 or 'columns' then by may contain column levels and/or index labels

Changed in version 0.23.0: Allow specifying index or column level names.

```
axis [{0 or 'index', 1 or 'columns'}, default 0] Axis to be sorted
```

**ascending** [bool or list of bool, default True] Sort ascending vs. descending. Specify list for multiple sort orders. If this is a list of bools, must match the length of the by.

inplace [bool, default False] if True, perform operation in-place

**kind** [{'quicksort', 'mergesort', 'heapsort'}, default 'quicksort'] Choice of sorting algorithm. See also ndarray.np.sort for more information. *mergesort* is the only stable algorithm. For DataFrames, this option is only applied when sorting on a single column or label.

**na\_position** [{'first', 'last'}, default 'last'] *first* puts NaNs at the beginning, *last* puts NaNs at the end

## Returns

sorted\_obj [DataFrame]

## **Examples**

```
>>> df = pd.DataFrame({
         'col1': ['A', 'A', 'B', np.nan, 'D', 'C'], 'col2': [2, 1, 9, 8, 7, 4],
. . .
          'col3': [0, 1, 9, 4, 2, 3],
. . .
. . . })
>>> df
    col1 col2 col3
0
    Α
           2
                 0
           1
1
    Α
                 1
2
    В
           9
3
    NaN 8
                 4
4
           7
    D
                 2.
5
    С
```

## Sort by col1

```
>>> df.sort_values(by=['col1'])
    col1 col2 col3
0
    Α
         2
               \cap
1
    Α
         1
               1
2
    В
         9
               9
5
    С
          4
               3
4
          7
               2
    D
3
    NaN 8
```

## Sort by multiple columns

```
>>> df.sort_values(by=['col1', 'col2'])
    col1 col2 col3
         1
1
    Α
              1
         2
0
    Α
              0
2
    В
         9
               9
5
         4
               3
    С
4
    D
         7
               2
3
    NaN 8
               4
```

## Sort Descending

```
>>> df.sort_values(by='col1', ascending=False)
    col1 col2 col3
          7
4
    D
5
    С
          4
               3
          9
2
    В
               9
0
    Α
          2
               0
1
          1
    Α
               1
3
         8
    NaN
               4
```

## Putting NAs first

```
>>> df.sort_values(by='col1', ascending=False, na_position='first')
    col1 col2 col3
    NaN 8
               4
4
          7
    С
          4
               3
2
    В
          9
               9
          2
0
               \cap
    Α
1
    Α
          1
               1
```

```
sortlevel (level=0, axis=0, ascending=True, inplace=False, sort_remaining=True)
```

Sort multilevel index by chosen axis and primary level. Data will be lexicographically sorted by the chosen level followed by the other levels (in order).

Deprecated since version 0.20.0: Use DataFrame.sort\_index()

## **Parameters**

```
level [int]
axis [{0 or 'index', 1 or 'columns'}, default 0]
ascending [boolean, default True]
inplace [boolean, default False] Sort the DataFrame without creating a new instance
sort_remaining [boolean, default True] Sort by the other levels too.
```

# **Returns**

```
sorted [DataFrame]
```

## See also:

```
DataFrame.sort_index
```

# squeeze(axis=None)

Squeeze length 1 dimensions.

## **Parameters**

axis [None, integer or string axis name, optional] The axis to squeeze if 1-sized.

New in version 0.20.0.

## Returns

## scalar if 1-sized, else original object

```
stack (level=-1, dropna=True)
```

Stack the prescribed level(s) from columns to index.

Return a reshaped DataFrame or Series having a multi-level index with one or more new inner-most levels compared to the current DataFrame. The new inner-most levels are created by pivoting the columns of the current dataframe:

- if the columns have a single level, the output is a Series;
- if the columns have multiple levels, the new index level(s) is (are) taken from the prescribed level(s) and the output is a DataFrame.

The new index levels are sorted.

## **Parameters**

**level** [int, str, list, default -1] Level(s) to stack from the column axis onto the index axis, defined as one index or label, or a list of indices or labels.

**dropna** [bool, default True] Whether to drop rows in the resulting Frame/Series with missing values. Stacking a column level onto the index axis can create combinations of index and column values that are missing from the original dataframe. See Examples section.

### Returns

**DataFrame or Series** Stacked dataframe or series.

## See also:

DataFrame.unstack Unstack prescribed level(s) from index axis onto column axis.

**DataFrame.pivot** Reshape dataframe from long format to wide format.

**DataFrame.pivot\_table** Create a spreadsheet-style pivot table as a DataFrame.

## **Notes**

The function is named by analogy with a collection of books being re-organised from being side by side on a horizontal position (the columns of the dataframe) to being stacked vertically on top of each other (in the index of the dataframe).

## **Examples**

## Single level columns

Stacking a dataframe with a single level column axis returns a Series:

```
>>> df_single_level_cols
    weight height
       0
             1
cat
         2
                3
doa
>>> df_single_level_cols.stack()
cat weight 0
    height
             1
dog weight
             2
              3
    height
dtype: int64
```

## Multi level columns: simple case

Stacking a dataframe with a multi-level column axis:

```
>>> df_multi_level_cols1
    weight
        kg
              pounds
             2
cat
        1
        2
                 4
>>> df_multi_level_cols1.stack()
          weight
cat kg
   pounds
                2
                2.
dog kg
   pounds
                4
```

### Missing values

It is common to have missing values when stacking a dataframe with multi-level columns, as the stacked dataframe typically has more values than the original dataframe. Missing values are filled with NaNs:

```
>>> df_multi_level_cols2
   weight height
      kg
          2.0
cat
     1.0
dog
     3.0 4.0
>>> df_multi_level_cols2.stack()
     height weight
cat kg
      NaN 1.0
        2.0
               NaN
   m
        NaN
               3.0
dog kg
        4.0
               NaN
   m
```

# Prescribing the level(s) to be stacked

The first parameter controls which level or levels are stacked:

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```
dog height m 4.0 weight kg 3.0 dtype: float64
```

# **Dropping missing values**

Note that rows where all values are missing are dropped by default but this behaviour can be controlled via the dropna keyword parameter:

```
>>> df_multi_level_cols3
   weight height
       kg m
             1.0
      NaN
cat
      2.0 3.0
doa
>>> df_multi_level_cols3.stack(dropna=False)
       height weight
cat kg
         NaN
                  NaN
          1.0
                  NaN
   m
                  2.0
dog kg
         NaN
          3.0
                  NaN
>>> df_multi_level_cols3.stack(dropna=True)
       height weight
         1.0
                  NaN
cat m
          NaN
                  2.0
dog kg
          3.0
                  NaN
   m
```

**std** (*axis=None*, *skipna=None*, *level=None*, *ddof=1*, *numeric\_only=None*, \*\*kwargs)
Return sample standard deviation over requested axis.

Normalized by N-1 by default. This can be changed using the ddof argument

## **Parameters**

```
axis [\{index (0), columns (1)\}]
```

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**ddof** [int, default 1] Delta Degrees of Freedom. The divisor used in calculations is N - ddof, where N represents the number of elements.

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

**std** [Series or DataFrame (if level specified)]

## style

Property returning a Styler object containing methods for building a styled HTML representation fo the DataFrame.

## See also:

```
pandas.io.formats.style.Styler
```

sub (other, axis='columns', level=None, fill\_value=None)

Subtraction of dataframe and other, element-wise (binary operator *sub*).

Equivalent to dataframe - other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

```
other [Series, DataFrame, or constant]
```

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

### Returns

result [DataFrame]

## See also:

DataFrame.rsub

## **Notes**

Mismatched indices will be unioned together

# **Examples**

```
>>> a = pd.DataFrame([2, 1, 1, np.nan], index=['a', 'b', 'c', 'd'],
                     columns=['one'])
. . .
>>> a
  one
a 2.0
b 1.0
c 1.0
d NaN
>>> b = pd.DataFrame(dict(one=[1, np.nan, 1, np.nan],
                          two=[3, 2, np.nan, 2]),
                     index=['a', 'b', 'd', 'e'])
>>> b
  one
       two
  1.0
       3.0
       2.0
  NaN
  1.0
       NaN
d
  NaN 2.0
е
>>> a.sub(b, fill_value=0)
  one two
  1.0
       -3.0
  1.0
       -2.0
```

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```
c 1.0 NaN
d -1.0 NaN
e NaN -2.0
```

subtract (other, axis='columns', level=None, fill\_value=None)

Subtraction of dataframe and other, element-wise (binary operator *sub*).

Equivalent to dataframe - other, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

other [Series, DataFrame, or constant]

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

## Returns

result [DataFrame]

#### See also:

DataFrame.rsub

## **Notes**

Mismatched indices will be unioned together

# **Examples**

```
>>> a = pd.DataFrame([2, 1, 1, np.nan], index=['a', 'b', 'c', 'd'],
                     columns=['one'])
. . .
>>> a
   one
  2.0
b 1.0
c 1.0
d NaN
>>> b = pd.DataFrame(dict(one=[1, np.nan, 1, np.nan],
                          two=[3, 2, np.nan, 2]),
                     index=['a', 'b', 'd', 'e'])
>>> b
   one
        two
  1.0
        3.0
  NaN
        2.0
d
  1.0
        NaN
  NaN
       2.0
>>> a.sub(b, fill_value=0)
```

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```
one two
a 1.0 -3.0
b 1.0 -2.0
c 1.0 NaN
d -1.0 NaN
e NaN -2.0
```

**sum** (axis=None, skipna=None, level=None, numeric\_only=None, min\_count=0, \*\*kwargs)
Return the sum of the values for the requested axis

### **Parameters**

```
axis [{index (0), columns (1)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

min\_count [int, default 0] The required number of valid values to perform the operation. If fewer than min\_count non-NA values are present the result will be NA.

New in version 0.22.0: Added with the default being 0. This means the sum of an all-NA or empty Series is 0, and the product of an all-NA or empty Series is 1.

## **Returns**

**sum** [Series or DataFrame (if level specified)]

# **Examples**

By default, the sum of an empty or all-NA Series is 0.

```
>>> pd.Series([]).sum() # min_count=0 is the default
0.0
```

This can be controlled with the min\_count parameter. For example, if you'd like the sum of an empty series to be NaN, pass min\_count=1.

```
>>> pd.Series([]).sum(min_count=1)
nan
```

Thanks to the skipna parameter, min\_count handles all-NA and empty series identically.

```
>>> pd.Series([np.nan]).sum()
0.0
```

```
>>> pd.Series([np.nan]).sum(min_count=1)
nan
```

swapaxes (axis1, axis2, copy=True)

Interchange axes and swap values axes appropriately

## Returns

y [same as input]

```
swaplevel (i=-2, j=-1, axis=0)
```

Swap levels i and j in a MultiIndex on a particular axis

## **Parameters**

i, j [int, string (can be mixed)] Level of index to be swapped. Can pass level name as string.

### Returns

swapped [type of caller (new object)]

.. versionchanged:: 0.18.1 The indexes i and j are now optional, and default to the two innermost levels of the index.

### tail(n=5)

Return the last n rows.

This function returns last n rows from the object based on position. It is useful for quickly verifying data, for example, after sorting or appending rows.

### **Parameters**

**n** [int, default 5] Number of rows to select.

## Returns

**type of caller** The last n rows of the caller object.

See also:

pandas.DataFrame.head The first n rows of the caller object.

## **Examples**

```
>>> df = pd.DataFrame({'animal':['alligator', 'bee', 'falcon', 'lion',
                        'monkey', 'parrot', 'shark', 'whale', 'zebra']})
. . .
>>> df
      animal
0
  alligator
1
         bee
2
      falcon
3
        lion
4
      monkey
5
      parrot
6
       shark
7
       whale
8
       zebra
```

# Viewing the last 5 lines

```
>>> df.tail()
   animal
4 monkey
5 parrot
6 shark
7 whale
8 zebra
```

Viewing the last *n* lines (three in this case)

```
>>> df.tail(3)
animal
6 shark
7 whale
8 zebra
```

**take** (indices, axis=0, convert=None, is\_copy=True, \*\*kwargs)

Return the elements in the given *positional* indices along an axis.

This means that we are not indexing according to actual values in the index attribute of the object. We are indexing according to the actual position of the element in the object.

### **Parameters**

indices [array-like] An array of ints indicating which positions to take.

**axis** [{0 or 'index', 1 or 'columns', None}, default 0] The axis on which to select elements. 0 means that we are selecting rows, 1 means that we are selecting columns.

**convert** [bool, default True] Whether to convert negative indices into positive ones. For example, -1 would map to the len(axis) - 1. The conversions are similar to the behavior of indexing a regular Python list.

Deprecated since version 0.21.0: In the future, negative indices will always be converted.

is\_copy [bool, default True] Whether to return a copy of the original object or not.

\*\*kwargs For compatibility with numpy.take(). Has no effect on the output.

## Returns

taken [type of caller] An array-like containing the elements taken from the object.

## See also:

**DataFrame.loc** Select a subset of a DataFrame by labels.

DataFrame.iloc Select a subset of a DataFrame by positions.

numpy.take Take elements from an array along an axis.

# **Examples**

```
>>> df = pd.DataFrame([('falcon', 'bird',
                                          389.0),
                      ('parrot', 'bird',
                                            24.0),
                               'mammal',
                      ('lion',
                                            80.5),
. . .
                      ('monkey', 'mammal', np.nan)],
. . .
                      columns=['name', 'class', 'max_speed'],
. . .
                      index=[0, 2, 3, 1])
. . .
>>> df
    name class max_speed
0 falcon bird 389.0
 parrot bird
                      24.0
                      80.5
    lion mammal
 monkey mammal
                       NaN
```

Take elements at positions 0 and 3 along the axis 0 (default).

Note how the actual indices selected (0 and 1) do not correspond to our selected indices 0 and 3. That's because we are selecting the 0th and 3rd rows, not rows whose indices equal 0 and 3.

```
>>> df.take([0, 3])
    name class max_speed
0 falcon bird 389.0
1 monkey mammal NaN
```

Take elements at indices 1 and 2 along the axis 1 (column selection).

```
>>> df.take([1, 2], axis=1)
    class max_speed
0 bird 389.0
2 bird 24.0
3 mammal 80.5
1 mammal NaN
```

We may take elements using negative integers for positive indices, starting from the end of the object, just like with Python lists.

```
>>> df.take([-1, -2])
name class max_speed
1 monkey mammal NaN
3 lion mammal 80.5
```

# to\_clipboard(excel=True, sep=None, \*\*kwargs)

Copy object to the system clipboard.

Write a text representation of object to the system clipboard. This can be pasted into Excel, for example.

## **Parameters**

excel [bool, default True]

- True, use the provided separator, writing in a csv format for allowing easy pasting into excel.
- False, write a string representation of the object to the clipboard.

```
sep [str, default '\t'] Field delimiter.
```

\*\*kwargs These parameters will be passed to DataFrame.to\_csv.

### See also:

**DataFrame.to\_csv** Write a DataFrame to a comma-separated values (csv) file.

read\_clipboard Read text from clipboard and pass to read\_table.

## **Notes**

Requirements for your platform.

- Linux : xclip, or xsel (with gtk or PyQt4 modules)
- Windows: none
- OS X: none

# **Examples**

Copy the contents of a DataFrame to the clipboard.

```
>>> df = pd.DataFrame([[1, 2, 3], [4, 5, 6]], columns=['A', 'B', 'C'])
>>> df.to_clipboard(sep=',')
... # Wrote the following to the system clipboard:
... # ,A,B,C
... # 0,1,2,3
... # 1,4,5,6
```

We can omit the the index by passing the keyword *index* and setting it to false.

```
>>> df.to_clipboard(sep=',', index=False)
... # Wrote the following to the system clipboard:
... # A,B,C
... # 1,2,3
... # 4,5,6
```

```
to csv(*args, **kwargs)
```

Write DataFrame to a comma-separated values (csv) file

### **Parameters**

path\_or\_buf [string or file handle, default None] File path or object, if None is provided the result is returned as a string.

sep [character, default ','] Field delimiter for the output file.

na\_rep [string, default '] Missing data representation

**float\_format** [string, default None] Format string for floating point numbers

columns [sequence, optional] Columns to write

**header** [boolean or list of string, default True] Write out the column names. If a list of strings is given it is assumed to be aliases for the column names

index [boolean, default True] Write row names (index)

index\_label [string or sequence, or False, default None] Column label for index column(s) if desired. If None is given, and header and index are True, then the index names are used. A sequence should be given if the DataFrame uses MultiIndex. If False do not print fields for index names. Use index\_label=False for easier importing in R

mode [str] Python write mode, default 'w'

**encoding** [string, optional] A string representing the encoding to use in the output file, defaults to 'ascii' on Python 2 and 'utf-8' on Python 3.

**compression** [string, optional] A string representing the compression to use in the output file. Allowed values are 'gzip', 'bz2', 'zip', 'xz'. This input is only used when the first argument is a filename.

**line\_terminator** [string, default '\n'] The newline character or character sequence to use in the output file

**quoting** [optional constant from csv module] defaults to csv.QUOTE\_MINIMAL. If you have set a *float\_format* then floats are converted to strings and thus csv.QUOTE\_NONNUMERIC will treat them as non-numeric

quotechar [string (length 1), default "''] character used to quote fields

doublequote [boolean, default True] Control quoting of quotechar inside a field

**escapechar** [string (length 1), default None] character used to escape *sep* and *quotechar* when appropriate

chunksize [int or None] rows to write at a time

**tupleize\_cols** [boolean, default False] Deprecated since version 0.21.0: This argument will be removed and will always write each row of the multi-index as a separate row in the CSV file.

Write MultiIndex columns as a list of tuples (if True) or in the new, expanded format, where each MultiIndex column is a row in the CSV (if False).

date\_format [string, default None] Format string for datetime objects

**decimal: string, default '.'** Character recognized as decimal separator. E.g. use ',' for European data

### to dense()

Return dense representation of NDFrame (as opposed to sparse)

```
to_dict (orient='dict', into=<type 'dict'>)
```

Convert the DataFrame to a dictionary.

The type of the key-value pairs can be customized with the parameters (see below).

### **Parameters**

**orient** [str {'dict', 'list', 'series', 'split', 'records', 'index'}] Determines the type of the values of the dictionary.

- 'dict' (default): dict like {column -> {index -> value}}
- 'list' : dict like {column -> [values]}
- 'series' : dict like {column -> Series(values)}
- 'split': dict like { 'index' -> [index], 'columns' -> [columns], 'data' -> [values]}
- 'records': list like [{column -> value}, ..., {column -> value}]
- 'index' : dict like {index -> {column -> value}}

Abbreviations are allowed. *s* indicates *series* and *sp* indicates *split*.

**into** [class, default dict] The collections. Mapping subclass used for all Mappings in the return value. Can be the actual class or an empty instance of the mapping type you want. If you want a collections default dict, you must pass it initialized.

New in version 0.21.0.

## Returns

```
result [collections.Mapping like {column -> {index -> value}}]
```

See also:

```
DataFrame.from_dict create a DataFrame from a dictionary
```

DataFrame.to\_json convert a DataFrame to JSON format

## **Examples**

You can specify the return orientation.

```
>>> df.to_dict('split')
{'index': ['a', 'b'], 'columns': ['col1', 'col2'],
   'data': [[1.0, 0.5], [2.0, 0.75]]}
```

```
>>> df.to_dict('records')
[{'coll': 1.0, 'col2': 0.5}, {'col1': 2.0, 'col2': 0.75}]
```

```
>>> df.to_dict('index')
{'a': {'col1': 1.0, 'col2': 0.5}, 'b': {'col1': 2.0, 'col2': 0.75}}
```

You can also specify the mapping type.

If you want a *defaultdict*, you need to initialize it:

```
>>> dd = defaultdict(list)
>>> df.to_dict('records', into=dd)
[defaultdict(<class 'list'>, {'col1': 1.0, 'col2': 0.5}),
  defaultdict(<class 'list'>, {'col1': 2.0, 'col2': 0.75})]
```

to\_excel (\*args, \*\*kwargs)

Write DataFrame to an excel sheet

### **Parameters**

```
excel_writer [string or ExcelWriter object] File path or existing ExcelWriter
sheet_name [string, default 'Sheet1'] Name of sheet which will contain DataFrame
na_rep [string, default '] Missing data representation
float_format [string, default None] Format string for floating point numbers
```

columns [sequence, optional] Columns to write

**header** [boolean or list of string, default True] Write out the column names. If a list of strings is given it is assumed to be aliases for the column names

index [boolean, default True] Write row names (index)

index\_label [string or sequence, default None] Column label for index column(s) if desired. If None is given, and *header* and *index* are True, then the index names are used. A sequence should be given if the DataFrame uses MultiIndex.

startrow: upper left cell row to dump data frame

**startcol:** upper left cell column to dump data frame

engine [string, default None] write engine to use - you can also set this via the options io.excel.xlsx.writer, io.excel.xls.writer, and io.excel. xlsm.writer.

**merge\_cells** [boolean, default True] Write MultiIndex and Hierarchical Rows as merged cells.

**encoding: string, default None** encoding of the resulting excel file. Only necessary for xlwt, other writers support unicode natively.

inf\_rep [string, default 'inf'] Representation for infinity (there is no native representation
 for infinity in Excel)

**freeze\_panes** [tuple of integer (length 2), default None] Specifies the one-based bottom-most row and rightmost column that is to be frozen

New in version 0.20.0.

#### **Notes**

If passing an existing ExcelWriter object, then the sheet will be added to the existing workbook. This can be used to save different DataFrames to one workbook:

```
>>> writer = pd.ExcelWriter('output.xlsx')
>>> df1.to_excel(writer,'Sheet1')
>>> df2.to_excel(writer,'Sheet2')
>>> writer.save()
```

For compatibility with to\_csv, to\_excel serializes lists and dicts to strings before writing.

#### to\_feather(fname)

write out the binary feather-format for DataFrames

New in version 0.20.0.

### **Parameters**

fname [str] string file path

to\_gbq (destination\_table, project\_id, chunksize=None, verbose=None, reauth=False, if\_exists='fail', private\_key=None, auth\_local\_webserver=False, table\_schema=None)
Write a DataFrame to a Google BigQuery table.

This function requires the pandas-gbq package.

Authentication to the Google BigQuery service is via OAuth 2.0.

- If private\_key is provided, the library loads the JSON service account credentials and uses those
  to authenticate.
- If no private\_key is provided, the library tries application default credentials.
- If application default credentials are not found or cannot be used with BigQuery, the library authenticates with user account credentials. In this case, you will be asked to grant permissions for product name 'pandas GBQ'.

#### **Parameters**

**destination\_table** [str] Name of table to be written, in the form 'dataset.tablename'.

project\_id [str] Google BigQuery Account project ID.

**chunksize** [int, optional] Number of rows to be inserted in each chunk from the dataframe. Set to None to load the whole dataframe at once.

**reauth** [bool, default False] Force Google BigQuery to reauthenticate the user. This is useful if multiple accounts are used.

if\_exists [str, default 'fail'] Behavior when the destination table exists. Value can be one of:

'fail' If table exists, do nothing.

'replace' If table exists, drop it, recreate it, and insert data.

'append' If table exists, insert data. Create if does not exist.

private\_key [str, optional] Service account private key in JSON format. Can be file path
or string contents. This is useful for remote server authentication (eg. Jupyter/IPython
notebook on remote host).

**auth\_local\_webserver** [bool, default False] Use the local webserver flow instead of the console flow when getting user credentials.

New in version 0.2.0 of pandas-gbq.

table\_schema [list of dicts, optional] List of BigQuery table fields to which according DataFrame columns conform to, e.g. [{'name': 'coll', 'type': 'STRING'},...]. If schema is not provided, it will be generated according to dtypes of DataFrame columns. See BigQuery API documentation on available names of a field.

New in version 0.3.1 of pandas-gbq.

**verbose** [boolean, deprecated] *Deprecated in Pandas-GBQ 0.4.0.* Use the logging module to adjust verbosity instead.

### See also:

pandas\_gbq.to\_gbq This function in the pandas-gbq library.

pandas.read\_gbq Read a DataFrame from Google BigQuery.

to\_hdf (path\_or\_buf, key, \*\*kwargs)

Write the contained data to an HDF5 file using HDFStore.

Hierarchical Data Format (HDF) is self-describing, allowing an application to interpret the structure and contents of a file with no outside information. One HDF file can hold a mix of related objects which can be accessed as a group or as individual objects.

In order to add another DataFrame or Series to an existing HDF file please use append mode and a different a key.

For more information see the user guide.

#### **Parameters**

path\_or\_buf [str or pandas.HDFStore] File path or HDFStore object.

**key** [str] Identifier for the group in the store.

**mode** [{'a', 'w', 'r+'}, default 'a'] Mode to open file:

- 'w': write, a new file is created (an existing file with the same name would be deleted).
- 'a': append, an existing file is opened for reading and writing, and if the file does not exist it is created.
- 'r+': similar to 'a', but the file must already exist.

**format** [{'fixed', 'table'}, default 'fixed'] Possible values:

- 'fixed': Fixed format. Fast writing/reading. Not-appendable, nor searchable.
- 'table': Table format. Write as a PyTables Table structure which may perform worse but allow more flexible operations like searching / selecting subsets of the data.

append [bool, default False] For Table formats, append the input data to the existing.

data\_columns [list of columns or True, optional] List of columns to create as indexed data columns for on-disk queries, or True to use all columns. By default only the axes of the object are indexed. See io.hdf5-query-data-columns. Applicable only to format='table'.

**complevel** [{0-9}, optional] Specifies a compression level for data. A value of 0 disables compression.

complib [{'zlib', 'lzo', 'bzip2', 'blosc'}, default 'zlib'] Specifies the compression library to be used. As of v0.20.2 these additional compressors for Blosc are supported (default if no compressor specified: 'blosc:blosclz'): {'blosc:blosclz', 'blosc:lz4', 'blosc:lz4hc', 'blosc:snappy', 'blosc:zlib', 'blosc:zstd'}. Specifying a compression library which is not available issues a ValueError.

fletcher32 [bool, default False] If applying compression use the fletcher32 checksum.

**dropna** [bool, default False] If true, ALL nan rows will not be written to store.

**errors** [str, default 'strict'] Specifies how encoding and decoding errors are to be handled. See the errors argument for open () for a full list of options.

### See also:

DataFrame.read\_hdf Read from HDF file.

**DataFrame.to\_parquet** Write a DataFrame to the binary parquet format.

DataFrame.to\_sql Write to a sql table.

**DataFrame.to\_feather** Write out feather-format for DataFrames.

DataFrame.to\_csv Write out to a csv file.

## **Examples**

```
>>> df = pd.DataFrame({'A': [1, 2, 3], 'B': [4, 5, 6]},
... index=['a', 'b', 'c'])
>>> df.to_hdf('data.h5', key='df', mode='w')
```

We can add another object to the same file:

```
>>> s = pd.Series([1, 2, 3, 4])
>>> s.to_hdf('data.h5', key='s')
```

Reading from HDF file:

```
>>> pd.read_hdf('data.h5', 'df')
A B
  1
а
  2
b
с 3
>>> pd.read_hdf('data.h5', 's')
0
    1
1
     2.
2
     3
     4
dtype: int64
```

### Deleting file with data:

```
>>> import os
>>> os.remove('data.h5')
```

## to\_html (\*args, \*\*kwargs)

Render a DataFrame as an HTML table.

*to\_html*-specific options:

**bold\_rows** [boolean, default True] Make the row labels bold in the output

classes [str or list or tuple, default None] CSS class(es) to apply to the resulting html table

escape [boolean, default True] Convert the characters <, >, and & to HTML-safe sequences.

max\_rows [int, optional] Maximum number of rows to show before truncating. If None, show all.

max\_cols [int, optional] Maximum number of columns to show before truncating. If None, show all.

decimal [string, default '.'] Character recognized as decimal separator, e.g. ',' in Europe

New in version 0.18.0.

**border** [int] A border=border attribute is included in the opening tag. Default pd. options.html.border.

New in version 0.19.0.

**table\_id** [str, optional] A css id is included in the opening tag if specified.

New in version 0.23.0.

### **Parameters**

**buf** [StringIO-like, optional] buffer to write to

columns [sequence, optional] the subset of columns to write; default None writes all columns

col\_space [int, optional] the minimum width of each column

header [bool, optional] whether to print column labels, default True

index [bool, optional] whether to print index (row) labels, default True

na\_rep [string, optional] string representation of NAN to use, default 'NaN'

**formatters** [list or dict of one-parameter functions, optional] formatter functions to apply to columns' elements by position or name, default None. The result of each function must be a unicode string. List must be of length equal to the number of columns.

**float\_format** [one-parameter function, optional] formatter function to apply to columns' elements if they are floats, default None. The result of this function must be a unicode string.

**sparsify** [bool, optional] Set to False for a DataFrame with a hierarchical index to print every multiindex key at each row, default True

index\_names [bool, optional] Prints the names of the indexes, default True

**line width** [int, optional] Width to wrap a line in characters, default no wrap

table\_id [str, optional] id for the element create by to\_html

New in version 0.23.0.

**justify** [str, default None] How to justify the column labels. If None uses the option from the print configuration (controlled by set\_option), 'right' out of the box. Valid values are

- left
- right
- center
- · justify
- justify-all
- start
- end
- inherit
- match-parent
- initial
- unset

### Returns

**formatted** [string (or unicode, depending on data and options)]

to\_json (path\_or\_buf=None, orient=None, date\_format=None, double\_precision=10, force\_ascii=True, date\_unit='ms', default\_handler=None, lines=False, compression=None, index=True)

Convert the object to a JSON string.

Note NaN's and None will be converted to null and datetime objects will be converted to UNIX timestamps.

### **Parameters**

**path\_or\_buf** [string or file handle, optional] File path or object. If not specified, the result is returned as a string.

orient [string] Indication of expected JSON string format.

- Series
  - default is 'index'
  - allowed values are: {'split','records','index'}
- DataFrame
  - default is 'columns'
  - allowed values are: {'split','records','index','columns','values'}
- The format of the JSON string
  - 'split': dict like { 'index' -> [index], 'columns' -> [columns], 'data' -> [values]}
  - 'records': list like [{column -> value}, ..., {column -> value}]
  - 'index' : dict like {index -> {column -> value}}
  - 'columns' : dict like {column -> {index -> value}}}
  - 'values' : just the values array
  - 'table': dict like {'schema': {schema}, 'data': {data}} describing the data, and the data component is like orient='records'.

Changed in version 0.20.0.

- date\_format [{None, 'epoch', 'iso'}] Type of date conversion. 'epoch' = epoch milliseconds, 'iso' = ISO8601. The default depends on the orient. For orient='table',
  the default is 'iso'. For all other orients, the default is 'epoch'.
- **double\_precision** [int, default 10] The number of decimal places to use when encoding floating point values.
- force\_ascii [boolean, default True] Force encoded string to be ASCII.
- **date\_unit** [string, default 'ms' (milliseconds)] The time unit to encode to, governs timestamp and ISO8601 precision. One of 's', 'ms', 'us', 'ns' for second, millisecond, microsecond, and nanosecond respectively.
- **default\_handler** [callable, default None] Handler to call if object cannot otherwise be converted to a suitable format for JSON. Should receive a single argument which is the object to convert and return a serialisable object.
- **lines** [boolean, default False] If 'orient' is 'records' write out line delimited json format. Will throw ValueError if incorrect 'orient' since others are not list like.

New in version 0.19.0.

**compression** [{None, 'gzip', 'bz2', 'zip', 'xz'}] A string representing the compression to use in the output file, only used when the first argument is a filename.

New in version 0.21.0.

index [boolean, default True] Whether to include the index values in the JSON string. Not including the index (index=False) is only supported when orient is 'split' or 'table'.

New in version 0.23.0.

#### See also:

```
pandas.read_json
```

## **Examples**

Encoding/decoding a Dataframe using 'records' formatted JSON. Note that index labels are not preserved with this encoding.

```
>>> df.to_json(orient='records')
'[{"col 1":"a","col 2":"b"},{"col 1":"c","col 2":"d"}]'
```

Encoding/decoding a Dataframe using 'index' formatted JSON:

```
>>> df.to_json(orient='index')
'{"row 1":{"col 1":"a","col 2":"b"},"row 2":{"col 1":"c","col 2":"d"}}'
```

Encoding/decoding a Dataframe using 'columns' formatted JSON:

```
>>> df.to_json(orient='columns')
'{"col 1":{"row 1":"a","row 2":"c"},"col 2":{"row 1":"b","row 2":"d"}}'
```

Encoding/decoding a Dataframe using 'values' formatted JSON:

```
>>> df.to_json(orient='values')
'[["a","b"],["c","d"]]'
```

**Encoding with Table Schema** 

to\_latex (buf=None, columns=None, col\_space=None, header=True, index=True, na\_rep='NaN', formatters=None, float\_format=None, sparsify=None, index\_names=True, bold\_rows=False, column\_format=None, longtable=None, escape=None, encoding=None, decimal='.', multicolumn=None, multicolumn\_format=None, multirow=None)

Render an object to a tabular environment table. You can splice this into a LaTeX document. Requires \usepackage{booktabs}.

Changed in version 0.20.2: Added to Series

to\_latex-specific options:

**bold\_rows** [boolean, default False] Make the row labels bold in the output

column\_format [str, default None] The columns format as specified in LaTeX table format e.g 'rcl' for 3 columns

**longtable** [boolean, default will be read from the pandas config module] Default: False. Use a longtable environment instead of tabular. Requires adding a \usepackage{longtable} to your LaTeX preamble.

**escape** [boolean, default will be read from the pandas config module] Default: True. When set to False prevents from escaping latex special characters in column names.

**encoding** [str, default None] A string representing the encoding to use in the output file, defaults to 'ascii' on Python 2 and 'utf-8' on Python 3.

decimal [string, default '.'] Character recognized as decimal separator, e.g. ',' in Europe.

New in version 0.18.0.

**multicolumn** [boolean, default True] Use multicolumn to enhance MultiIndex columns. The default will be read from the config module.

New in version 0.20.0.

**multicolumn\_format** [str, default '1'] The alignment for multicolumns, similar to *column\_format* The default will be read from the config module.

New in version 0.20.0.

multirow [boolean, default False] Use multirow to enhance MultiIndex rows. Requires adding a \usepackage{multirow} to your LaTeX preamble. Will print centered labels (instead of top-aligned) across the contained rows, separating groups via clines. The default will be read from the pandas config module.

New in version 0.20.0.

to\_mo12 (filepath\_or\_buffer=None, columns=None)

update\_properties=True,

molecule column='mol',

Write DataFrame to Mol2 file.

New in version 0.3.

## **Parameters**

filepath\_or\_buffer [string or None] File path

**update\_properties** [bool, optional (default=True)] Switch to update properties from the DataFrames to the molecules while writting.

**molecule\_column** [string or None, optional (default='mol')] Name of molecule column. If None the molecules will be skipped.

**columns** [list or None, optional (default=None)] A list of columns to write to file. If None then all available fields are written.

to\_msgpack (path\_or\_buf=None, encoding='utf-8', \*\*kwargs) msgpack (serialize) object to input file path

THIS IS AN EXPERIMENTAL LIBRARY and the storage format may not be stable until a future release.

## **Parameters**

**path** [string File path, buffer-like, or None] if None, return generated string

append [boolean whether to append to an existing msgpack] (default is False)

compress [type of compressor (zlib or blosc), default to None (no] compression)

### to\_panel()

Transform long (stacked) format (DataFrame) into wide (3D, Panel) format.

Deprecated since version 0.20.0.

Currently the index of the DataFrame must be a 2-level MultiIndex. This may be generalized later

## Returns

```
panel [Panel]
```

```
to_parquet (fname, engine='auto', compression='snappy', **kwargs)
```

Write a DataFrame to the binary parquet format.

New in version 0.21.0.

This function writes the dataframe as a parquet file. You can choose different parquet backends, and have the option of compression. See the user guide for more details.

### **Parameters**

fname [str] String file path.

engine [{'auto', 'pyarrow', 'fastparquet'}, default 'auto'] Parquet library to use. If
 'auto', then the option io.parquet.engine is used. The default io.parquet.
 engine behavior is to try 'pyarrow', falling back to 'fastparquet' if 'pyarrow' is
 unavailable.

**compression** [{'snappy', 'gzip', 'brotli', None}, default 'snappy'] Name of the compression to use. Use None for no compression.

\*\*kwargs Additional arguments passed to the parquet library. See pandas io for more details.

### See also:

```
read_parquet Read a parquet file.
```

DataFrame.to\_csv Write a csv file.

DataFrame.to\_sql Write to a sql table.

DataFrame.to\_hdf Write to hdf.

### **Notes**

This function requires either the fastparquet or pyarrow library.

## **Examples**

```
>>> df = pd.DataFrame(data={'coll': [1, 2], 'col2': [3, 4]})
>>> df.to_parquet('df.parquet.gzip', compression='gzip')
>>> pd.read_parquet('df.parquet.gzip')
    col1 col2
0     1     3
1     2     4
```

## to\_period (freq=None, axis=0, copy=True)

Convert DataFrame from DatetimeIndex to PeriodIndex with desired frequency (inferred from index if not passed)

### **Parameters**

```
freq [string, default]
axis [{0 or 'index', 1 or 'columns'}, default 0] The axis to convert (the index by default)
copy [boolean, default True] If False then underlying input data is not copied
```

#### Returns

ts [TimeSeries with PeriodIndex]

to\_pickle (path, compression='infer', protocol=2) Pickle (serialize) object to file.

### **Parameters**

**path** [str] File path where the pickled object will be stored.

compression [{'infer', 'gzip', 'bz2', 'zip', 'xz', None}, default 'infer'] A string representing the compression to use in the output file. By default, infers from the file extension in specified path.

New in version 0.20.0.

protocol [int] Int which indicates which protocol should be used by the pickler, default HIGHEST\_PROTOCOL (see [1] paragraph 12.1.2). The possible values for this parameter depend on the version of Python. For Python 2.x, possible values are 0, 1, 2. For Python $\geq 3.0$ , 3 is a valid value. For Python  $\geq 3.4$ , 4 is a valid value. A negative value for the protocol parameter is equivalent to setting its value to HIGH-EST PROTOCOL.

New in version 0.21.0.

# See also:

read\_pickle Load pickled pandas object (or any object) from file.

DataFrame.to\_hdf Write DataFrame to an HDF5 file.

DataFrame.to\_sql Write DataFrame to a SQL database.

**DataFrame.to\_parquet** Write a DataFrame to the binary parquet format.

# **Examples**

```
>>> original_df = pd.DataFrame({"foo": range(5), "bar": range(5, 10)})
>>> original_df
   foo bar
0
     0
          5
1
          6
     1
          7
2
     2
3
          8
     3
>>> original_df.to_pickle("./dummy.pkl")
```

```
>>> unpickled_df = pd.read_pickle("./dummy.pkl")
>>> unpickled_df
   foo bar
    0
          5
     1
          6
```

(continues on next page)

```
2 2 7
3 3 8
4 4 9
```

```
>>> import os
>>> os.remove("./dummy.pkl")
```

## to\_records (index=True, convert\_datetime64=None)

Convert DataFrame to a NumPy record array.

Index will be put in the 'index' field of the record array if requested.

### **Parameters**

index [boolean, default True] Include index in resulting record array, stored in 'index' field.

**convert\_datetime64** [boolean, default None] Deprecated since version 0.23.0.

Whether to convert the index to datetime.datetime if it is a DatetimeIndex.

## Returns

y [numpy.recarray]

See also:

**DataFrame.from\_records** convert structured or record ndarray to DataFrame.

**numpy.recarray** ndarray that allows field access using attributes, analogous to typed columns in a spreadsheet.

# **Examples**

The index can be excluded from the record array:

By default, timestamps are converted to *datetime*. datetime:

```
>>> df.index = pd.date_range('2018-01-01 09:00', periods=2, freq='min')
>>> df

A B
2018-01-01 09:00:00 1 0.50
2018-01-01 09:01:00 2 0.75
```

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The timestamp conversion can be disabled so NumPy's datetime64 data type is used instead:

to\_sdf (filepath\_or\_buffer=None, update\_properties=True, molecule\_column=None, columns=None)
Write DataFrame to SDF file.

New in version 0.3.

### **Parameters**

filepath\_or\_buffer [string or None] File path

**update\_properties** [bool, optional (default=True)] Switch to update properties from the DataFrames to the molecules while writting.

**molecule\_column** [string or None, optional (default='mol')] Name of molecule column. If None the molecules will be skipped.

**columns** [list or None, optional (default=None)] A list of columns to write to file. If None then all available fields are written.

```
to_sparse(fill_value=None, kind='block')
```

Convert to SparseDataFrame

### **Parameters**

```
fill_value [float, default NaN]
kind [{'block', 'integer'}]
```

#### Returns

v [SparseDataFrame]

to\_sql (name, con, schema=None, if\_exists='fail', index=True, index\_label=None, chunksize=None, dtype=None)

Write records stored in a DataFrame to a SQL database.

Databases supported by SQLAlchemy [1] are supported. Tables can be newly created, appended to, or overwritten.

### **Parameters**

name [string] Name of SQL table.

**con** [sqlalchemy.engine.Engine or sqlite3.Connection] Using SQLAlchemy makes it possible to use any DB supported by that library. Legacy support is provided for sqlite3.Connection objects.

**schema** [string, optional] Specify the schema (if database flavor supports this). If None, use default schema.

if\_exists [{'fail', 'replace', 'append'}, default 'fail'] How to behave if the table already exists.

- fail: Raise a ValueError.
- replace: Drop the table before inserting new values.
- append: Insert new values to the existing table.

**index** [boolean, default True] Write DataFrame index as a column. Uses *index\_label* as the column name in the table.

**index\_label** [string or sequence, default None] Column label for index column(s). If None is given (default) and *index* is True, then the index names are used. A sequence should be given if the DataFrame uses MultiIndex.

**chunksize** [int, optional] Rows will be written in batches of this size at a time. By default, all rows will be written at once.

**dtype** [dict, optional] Specifying the datatype for columns. The keys should be the column names and the values should be the SQLAlchemy types or strings for the sqlite3 legacy mode.

### **Raises**

**ValueError** When the table already exists and *if\_exists* is 'fail' (the default).

#### See also:

pandas.read\_sql read a DataFrame from a table

#### References

[1], [2]

### **Examples**

Create an in-memory SQLite database.

```
>>> from sqlalchemy import create_engine
>>> engine = create_engine('sqlite://', echo=False)
```

Create a table from scratch with 3 rows.

```
>>> df.to_sql('users', con=engine)
>>> engine.execute("SELECT * FROM users").fetchall()
[(0, 'User 1'), (1, 'User 2'), (2, 'User 3')]
```

Overwrite the table with just df1.

```
>>> df1.to_sql('users', con=engine, if_exists='replace',
... index_label='id')
>>> engine.execute("SELECT * FROM users").fetchall()
[(0, 'User 4'), (1, 'User 5')]
```

Specify the dtype (especially useful for integers with missing values). Notice that while pandas is forced to store the data as floating point, the database supports nullable integers. When fetching the data with Python, we get back integer scalars.

```
>>> from sqlalchemy.types import Integer
>>> df.to_sql('integers', con=engine, index=False,
... dtype={"A": Integer()})
```

```
>>> engine.execute("SELECT * FROM integers").fetchall()
[(1,), (None,), (2,)]
```

to\_stata (fname, convert\_dates=None, write\_index=True, encoding='latin-1', byteorder=None, time\_stamp=None, data\_label=None, variable\_labels=None, version=114, convert\_strl=None)

Export Stata binary dta files.

## **Parameters**

**fname** [path (string), buffer or path object] string, path object (pathlib.Path or py.\_path.local.LocalPath) or object implementing a binary write() functions. If using a buffer then the buffer will not be automatically closed after the file data has been written.

convert\_dates [dict] Dictionary mapping columns containing datetime types to stata internal format to use when writing the dates. Options are 'tc', 'td', 'tm', 'tw', 'th', 'tq', 'ty'. Column can be either an integer or a name. Datetime columns that do not have a conversion type specified will be converted to 'tc'. Raises NotImplementedError if a datetime column has timezone information.

write index [bool] Write the index to Stata dataset.

encoding [str] Default is latin-1. Unicode is not supported.

**byteorder** [str] Can be ">", "<", "little", or "big". default is sys.byteorder.

**time stamp** [datetime] A datetime to use as file creation date. Default is the current time.

data label [str] A label for the data set. Must be 80 characters or smaller.

variable\_labels [dict] Dictionary containing columns as keys and variable labels as values. Each label must be 80 characters or smaller.

New in version 0.19.0.

**version** [{114, 117}] Version to use in the output dta file. Version 114 can be used read by Stata 10 and later. Version 117 can be read by Stata 13 or later. Version 114 limits

string variables to 244 characters or fewer while 117 allows strings with lengths up to 2.000,000 characters.

New in version 0.23.0.

convert\_strl [list, optional] List of column names to convert to string columns to Stata StrL format. Only available if version is 117. Storing strings in the StrL format can produce smaller dta files if strings have more than 8 characters and values are repeated.

New in version 0.23.0.

### **Raises**

## NotImplementedError

- If datetimes contain timezone information
- Column dtype is not representable in Stata

### ValueError

- Columns listed in convert\_dates are neither datetime64[ns] or datetime.datetime
- Column listed in convert\_dates is not in DataFrame
- Categorical label contains more than 32,000 characters

New in version 0.19.0.

#### See also:

```
pandas.read_stata Import Stata data files
pandas.io.stata.StataWriter low-level writer for Stata data files
pandas.io.stata.StataWriter117 low-level writer for version 117 files
```

## **Examples**

```
>>> data.to_stata('./data_file.dta')
```

## Or with dates

```
>>> data.to_stata('./date_data_file.dta', {2 : 'tw'})
```

Alternatively you can create an instance of the StataWriter class

```
>>> writer = StataWriter('./data_file.dta', data)
>>> writer.write_file()
```

# With dates:

```
>>> writer = StataWriter('./date_data_file.dta', data, {2 : 'tw'})
>>> writer.write_file()
```

to\_string (buf=None, columns=None, col\_space=None, header=True, index=True, na\_rep='NaN', formatters=None, float\_format=None, sparsify=None, index\_names=True, justify=None, line\_width=None, max\_rows=None, max\_cols=None, show\_dimensions=False)

Render a DataFrame to a console-friendly tabular output.

### **Parameters**

**buf** [StringIO-like, optional] buffer to write to

**columns** [sequence, optional] the subset of columns to write; default None writes all columns

col\_space [int, optional] the minimum width of each column

**header** [bool, optional] Write out the column names. If a list of strings is given, it is assumed to be aliases for the column names

index [bool, optional] whether to print index (row) labels, default True

**na\_rep** [string, optional] string representation of NAN to use, default 'NaN'

**formatters** [list or dict of one-parameter functions, optional] formatter functions to apply to columns' elements by position or name, default None. The result of each function must be a unicode string. List must be of length equal to the number of columns.

**float\_format** [one-parameter function, optional] formatter function to apply to columns' elements if they are floats, default None. The result of this function must be a unicode string.

**sparsify** [bool, optional] Set to False for a DataFrame with a hierarchical index to print every multiindex key at each row, default True

index\_names [bool, optional] Prints the names of the indexes, default True

line\_width [int, optional] Width to wrap a line in characters, default no wrap

table\_id [str, optional] id for the element create by to\_html

New in version 0.23.0.

**justify** [str, default None] How to justify the column labels. If None uses the option from the print configuration (controlled by set\_option), 'right' out of the box. Valid values are

- left
- right
- center
- · justify
- justify-all
- start
- end
- inherit
- match-parent
- initial
- unset

### **Returns**

**formatted** [string (or unicode, depending on data and options)]

to\_timestamp (freq=None, how='start', axis=0, copy=True)

Cast to DatetimeIndex of timestamps, at beginning of period

#### **Parameters**

```
freq [string, default frequency of PeriodIndex] Desired frequency
```

**how** [{'s', 'e', 'start', 'end'}] Convention for converting period to timestamp; start of period vs. end

axis [{0 or 'index', 1 or 'columns'}, default 0] The axis to convert (the index by default)

**copy** [boolean, default True] If false then underlying input data is not copied

## Returns

**df** [DataFrame with DatetimeIndex]

## to\_xarray()

Return an xarray object from the pandas object.

### Returns

- a DataArray for a Series
- a Dataset for a DataFrame
- a DataArray for higher dims

### **Notes**

See the xarray docs

## **Examples**

```
>>> df.to_xarray()
<xarray.Dataset>
Dimensions: (index: 3)
Coordinates:
  * index (index) int64 0 1 2
Data variables:
  A (index) int64 1 1 2
  B (index) object 'foo' 'bar' 'foo'
  C (index) float64 4.0 5.0 6.0
```

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```
bar 1 5.0
foo 2 6.0
```

```
>>> p.to_xarray()
<xarray.DataArray (items: 4, major_axis: 3, minor_axis: 2)>
array([[[ 0, 1],
        [ 2,
             3],
        [ 4,
             5]],
       [[6,
             7],
        [8, 9],
       [10, 11]],
       [[12, 13],
       [14, 15],
       [16, 17]],
       [[18, 19],
        [20, 21],
        [22, 23]])
Coordinates:
                (items) object 'A' 'B' 'C' 'D'
  \star items
  * major_axis (major_axis) datetime64[ns] 2013-01-01 2013-01-02 2013-01-03...

→ # noqa

  * minor_axis (minor_axis) object 'first' 'second'
```

## transform(func, \*args, \*\*kwargs)

Call function producing a like-indexed NDFrame and return a NDFrame with the transformed values New in version 0.20.0.

## **Parameters**

func [callable, string, dictionary, or list of string/callables] To apply to column

Accepted Combinations are:

- string function name
- function
- list of functions

• dict of column names -> functions (or list of functions)

### Returns

transformed [NDFrame]

### See also:

pandas.NDFrame.aggregate, pandas.NDFrame.apply

## **Examples**

```
>>> df = pd.DataFrame(np.random.randn(10, 3), columns=['A', 'B', 'C'], index=pd.date_range('1/1/2000', periods=10))
df.iloc[3:7] = np.nan
```

```
>>> df.transform(lambda x: (x - x.mean()) / x.std())
                   A
                             В
2000-01-01 0.579457 1.236184 0.123424
2000-01-02 0.370357 -0.605875 -1.231325
2000-01-03 1.455756 -0.277446 0.288967
2000-01-04 NaN
2000-01-05 NaN
2000-01-06 NaN
2000-01-07 NaN
                       NaN
                            NaN
                            NaN
                                      NaN
                           NaN
                                      NaN
2000-01-08 -0.498658 1.274522 1.642524
2000-01-09 -0.540524 -1.012676 -0.828968
2000-01-10 -1.366388 -0.614710 0.005378
```

## transpose(\*args, \*\*kwargs)

Transpose index and columns.

Reflect the DataFrame over its main diagonal by writing rows as columns and vice-versa. The property T is an accessor to the method transpose().

## **Parameters**

**copy** [bool, default False] If True, the underlying data is copied. Otherwise (default), no copy is made if possible.

\*args, \*\*kwargs Additional keywords have no effect but might be accepted for compatibility with numpy.

## Returns

DataFrame The transposed DataFrame.

#### See also:

numpy.transpose Permute the dimensions of a given array.

### **Notes**

Transposing a DataFrame with mixed dtypes will result in a homogeneous DataFrame with the *object* dtype. In such a case, a copy of the data is always made.

## **Examples**

## Square DataFrame with homogeneous dtype

```
>>> d1 = {'col1': [1, 2], 'col2': [3, 4]}
>>> df1 = pd.DataFrame(data=d1)
>>> df1
col1 col2
0 1 3
1 2 4
```

When the dtype is homogeneous in the original DataFrame, we get a transposed DataFrame with the same dtype:

```
>>> df1.dtypes
col1   int64
col2   int64
dtype: object
>>> df1_transposed.dtypes
0   int64
1   int64
dtype: object
```

## Non-square DataFrame with mixed dtypes

```
>>> d2 = {'name': ['Alice', 'Bob'],
... 'score': [9.5, 8],
... 'employed': [False, True],
... 'kids': [0, 0]}
>>> df2 = pd.DataFrame(data=d2)
>>> df2
    name score employed kids
0 Alice 9.5 False 0
1 Bob 8.0 True 0
```

When the DataFrame has mixed dtypes, we get a transposed DataFrame with the *object* dtype:

```
>>> df2.dtypes
name object
score float64
employed bool
kids int64
```

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```
dtype: object
>>> df2_transposed.dtypes
0    object
1    object
dtype: object
```

truediv (other, axis='columns', level=None, fill\_value=None)

Floating division of dataframe and other, element-wise (binary operator truediv).

Equivalent to dataframe / other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

```
other [Series, DataFrame, or constant]
```

axis [{0, 1, 'index', 'columns'}] For Series input, axis to match Series index on

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

**fill\_value** [None or float value, default None] Fill existing missing (NaN) values, and any new element needed for successful DataFrame alignment, with this value before computation. If data in both corresponding DataFrame locations is missing the result will be missing

### Returns

result [DataFrame]

#### See also:

DataFrame.rtruediv

#### **Notes**

Mismatched indices will be unioned together

## **Examples**

None

truncate (before=None, after=None, axis=None, copy=True)

Truncate a Series or DataFrame before and after some index value.

This is a useful shorthand for boolean indexing based on index values above or below certain thresholds.

### **Parameters**

**before** [date, string, int] Truncate all rows before this index value.

after [date, string, int] Truncate all rows after this index value.

**axis** [{0 or 'index', 1 or 'columns'}, optional] Axis to truncate. Truncates the index (rows) by default.

**copy** [boolean, default is True,] Return a copy of the truncated section.

## Returns

type of caller The truncated Series or DataFrame.

### See also:

DataFrame.loc Select a subset of a DataFrame by label.

**DataFrame.iloc** Select a subset of a DataFrame by position.

### **Notes**

If the index being truncated contains only datetime values, *before* and *after* may be specified as strings instead of Timestamps.

## **Examples**

```
>>> df = pd.DataFrame({'A': ['a', 'b', 'c', 'd', 'e'],
                      'B': ['f', 'g', 'h', 'i', 'j'],
                      'C': ['k', 'l', 'm', 'n', 'o']},
. . .
                      index=[1, 2, 3, 4, 5])
. . .
>>> df
  A B C
  a f k
2 b g
        1
3
  c h
        m
4
  d
     i
5 е ј о
```

```
>>> df.truncate(before=2, after=4)

A B C
2 b g l
3 c h m
4 d i n
```

The columns of a DataFrame can be truncated.

```
>>> df.truncate(before="A", after="B", axis="columns")
    A B
1 a f
2 b g
3 c h
4 d i
5 e j
```

For Series, only rows can be truncated.

The index values in truncate can be datetimes or string dates.

```
>>> dates = pd.date_range('2016-01-01', '2016-02-01', freq='s')
>>> df = pd.DataFrame(index=dates, data={'A': 1})
>>> df.tail()
```

(continues on next page)

```
A

2016-01-31 23:59:56 1

2016-01-31 23:59:57 1

2016-01-31 23:59:58 1

2016-01-31 23:59:59 1

2016-02-01 00:00:00 1
```

```
>>> df.truncate(before=pd.Timestamp('2016-01-05'),
... after=pd.Timestamp('2016-01-10')).tail()

A
2016-01-09 23:59:56 1
2016-01-09 23:59:57 1
2016-01-09 23:59:58 1
2016-01-09 23:59:59 1
2016-01-10 00:00:00 1
```

Because the index is a DatetimeIndex containing only dates, we can specify *before* and *after* as strings. They will be coerced to Timestamps before truncation.

```
>>> df.truncate('2016-01-05', '2016-01-10').tail()

A

2016-01-09 23:59:56 1

2016-01-09 23:59:57 1

2016-01-09 23:59:58 1

2016-01-09 23:59:59 1

2016-01-10 00:00:00 1
```

Note that truncate assumes a 0 value for any unspecified time component (midnight). This differs from partial string slicing, which returns any partially matching dates.

```
>>> df.loc['2016-01-05':'2016-01-10', :].tail()

A
2016-01-10 23:59:55 1
2016-01-10 23:59:56 1
2016-01-10 23:59:57 1
2016-01-10 23:59:58 1
2016-01-10 23:59:59 1
```

#### tshift (periods=1, freq=None, axis=0)

Shift the time index, using the index's frequency if available.

### **Parameters**

**periods** [int] Number of periods to move, can be positive or negative

**freq** [DateOffset, timedelta, or time rule string, default None] Increment to use from the tseries module or time rule (e.g. 'EOM')

axis [int or basestring] Corresponds to the axis that contains the Index

## Returns

shifted [NDFrame]

### **Notes**

If freq is not specified then tries to use the freq or inferred\_freq attributes of the index. If neither of those attributes exist, a ValueError is thrown

tz\_convert (tz, axis=0, level=None, copy=True)

Convert tz-aware axis to target time zone.

### **Parameters**

tz [string or pytz.timezone object]

axis [the axis to convert]

**level** [int, str, default None] If axis ia a MultiIndex, convert a specific level. Otherwise must be None

copy [boolean, default True] Also make a copy of the underlying data

#### Raises

**TypeError** If the axis is tz-naive.

tz\_localize(tz, axis=0, level=None, copy=True, ambiguous='raise')

Localize tz-naive TimeSeries to target time zone.

#### **Parameters**

tz [string or pytz.timezone object]

axis [the axis to localize]

**level** [int, str, default None] If axis ia a MultiIndex, localize a specific level. Otherwise must be None

copy [boolean, default True] Also make a copy of the underlying data

ambiguous ['infer', bool-ndarray, 'NaT', default 'raise']

- 'infer' will attempt to infer fall dst-transition hours based on order
- bool-ndarray where True signifies a DST time, False designates a non-DST time (note that this flag is only applicable for ambiguous times)
- 'NaT' will return NaT where there are ambiguous times
- 'raise' will raise an AmbiguousTimeError if there are ambiguous times

## Raises

**TypeError** If the TimeSeries is tz-aware and tz is not None.

```
unstack (level=-1, fill_value=None)
```

Pivot a level of the (necessarily hierarchical) index labels, returning a DataFrame having a new level of column labels whose inner-most level consists of the pivoted index labels. If the index is not a MultiIndex, the output will be a Series (the analogue of stack when the columns are not a MultiIndex). The level involved will automatically get sorted.

## **Parameters**

**level** [int, string, or list of these, default -1 (last level)] Level(s) of index to unstack, can pass level name

fill\_value [replace NaN with this value if the unstack produces] missing values

New in version 0.18.0.

#### Returns

unstacked [DataFrame or Series]

See also:

**DataFrame.pivot** Pivot a table based on column values.

DataFrame.stack Pivot a level of the column labels (inverse operation from unstack).

## **Examples**

```
>>> index = pd.MultiIndex.from_tuples([('one', 'a'), ('one', 'b'), ('two', 'a'), ('two', 'b')])
>>> s = pd.Series(np.arange(1.0, 5.0), index=index)
>>> s
one a 1.0
b 2.0
two a 3.0
b 4.0
dtype: float64
```

```
>>> s.unstack(level=-1)
    a    b
    one 1.0 2.0
    two 3.0 4.0
```

```
>>> s.unstack(level=0)
    one two
a 1.0 3.0
b 2.0 4.0
```

**update** (*other*, *join='left'*, *overwrite=True*, *filter\_func=None*, *raise\_conflict=False*) Modify in place using non-NA values from another DataFrame.

Aligns on indices. There is no return value.

## **Parameters**

**other** [DataFrame, or object coercible into a DataFrame] Should have at least one matching index/column label with the original DataFrame. If a Series is passed, its name attribute must be set, and that will be used as the column name to align with the original DataFrame.

**join** [{'left'}, default 'left'] Only left join is implemented, keeping the index and columns of the original object.

**overwrite** [bool, default True] How to handle non-NA values for overlapping keys:

• True: overwrite original DataFrame's values with values from other.

• False: only update values that are NA in the original DataFrame.

**filter\_func** [callable(1d-array) -> boolean 1d-array, optional] Can choose to replace values other than NA. Return True for values that should be updated.

**raise\_conflict** [bool, default False] If True, will raise a ValueError if the DataFrame and *other* both contain non-NA data in the same place.

#### Raises

ValueError When raise\_conflict is True and there's overlapping non-NA data.

See also:

dict.update Similar method for dictionaries.

**DataFrame.merge** For column(s)-on-columns(s) operations.

## **Examples**

```
>>> df = pd.DataFrame({'A': [1, 2, 3],
... 'B': [400, 500, 600]})
>>> new_df = pd.DataFrame({'B': [4, 5, 6],
... 'C': [7, 8, 9]})
>>> df.update(new_df)
>>> df
A B
0 1 4
1 2 5
2 3 6
```

The DataFrame's length does not increase as a result of the update, only values at matching index/column labels are updated.

For Series, it's name attribute must be set.

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If other contains NaNs the corresponding values are not updated in the original dataframe.

```
>>> df = pd.DataFrame({'A': [1, 2, 3],
                       'B': [400, 500, 600]})
>>> new_df = pd.DataFrame({'B': [4, np.nan, 6]})
>>> df.update(new_df)
>>> df
  Α
          R
0
        4.0
  1
  2
     500.0
1
2
  3
        6.0
```

### values

Return a Numpy representation of the DataFrame.

Only the values in the DataFrame will be returned, the axes labels will be removed.

#### Returns

**numpy.ndarray** The values of the DataFrame.

See also:

```
pandas.DataFrame.index Retrievie the index labels
pandas.DataFrame.columns Retrieving the column names
```

## **Notes**

The dtype will be a lower-common-denominator dtype (implicit upcasting); that is to say if the dtypes (even of numeric types) are mixed, the one that accommodates all will be chosen. Use this with care if you are not dealing with the blocks.

e.g. If the dtypes are float16 and float32, dtype will be upcast to float32. If dtypes are int32 and uint8, dtype will be upcast to int32. By numpy.find\_common\_type() convention, mixing int64 and uint64 will result in a float64 dtype.

## **Examples**

A DataFrame where all columns are the same type (e.g., int64) results in an array of the same type.

```
>>> df = pd.DataFrame({'age': [ 3, 29],
                        'height': [94, 170],
. . .
                        'weight': [31, 115]})
. . .
>>> df
  age
       height weight
            94
0
    3
                    31
           170
                   115
1
    29
```

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A DataFrame with mixed type columns(e.g., str/object, int64, float32) results in an ndarray of the broadest type that accommodates these mixed types (e.g., object).

```
>>> df2 = pd.DataFrame([('parrot',
                                      24.0, 'second'),
                         ('lion',
                                     80.5, 1),
                         ('monkey', np.nan, None)],
. . .
                      columns=('name', 'max_speed', 'rank'))
. . .
>>> df2.dtypes
             object
name
max_speed float64
rank
            object
dtype: object
>>> df2.values
array([['parrot', 24.0, 'second'],
       ['lion', 80.5, 1],
       ['monkey', nan, None]], dtype=object)
```

var (axis=None, skipna=None, level=None, ddof=1, numeric\_only=None, \*\*kwargs)
Return unbiased variance over requested axis.

Normalized by N-1 by default. This can be changed using the ddof argument

## **Parameters**

```
axis [{index (0), columns (1)}]
```

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a Series

**ddof** [int, default 1] Delta Degrees of Freedom. The divisor used in calculations is N - ddof, where N represents the number of elements.

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

var [Series or DataFrame (if level specified)]

where (cond, other=nan, inplace=False, axis=None, level=None, errors='raise', try\_cast=False, raise\_on\_error=None)

Return an object of same shape as self and whose corresponding entries are from self where *cond* is True and otherwise are from *other*.

### **Parameters**

**cond** [boolean NDFrame, array-like, or callable] Where *cond* is True, keep the original value. Where False, replace with corresponding value from *other*. If *cond* is callable,

it is computed on the NDFrame and should return boolean NDFrame or array. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as cond.

**other** [scalar, NDFrame, or callable] Entries where *cond* is False are replaced with corresponding value from *other*. If other is callable, it is computed on the NDFrame and should return scalar or NDFrame. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as other.

inplace [boolean, default False] Whether to perform the operation in place on the data

axis [alignment axis if needed, default None]

**level** [alignment level if needed, default None]

```
errors [str, {'raise', 'ignore'}, default 'raise']
```

- raise: allow exceptions to be raised
- ignore: suppress exceptions. On error return original object

Note that currently this parameter won't affect the results and will always coerce to a suitable dtype.

try\_cast [boolean, default False] try to cast the result back to the input type (if possible),

raise\_on\_error [boolean, default True] Whether to raise on invalid data types (e.g. trying to where on strings)

Deprecated since version 0.21.0.

## Returns

**wh** [same type as caller]

### See also:

```
DataFrame.mask()
```

#### **Notes**

The where method is an application of the if-then idiom. For each element in the calling DataFrame, if cond is True the element is used; otherwise the corresponding element from the DataFrame other is used.

The signature for DataFrame.where () differs from numpy.where (). Roughly df1.where (m, df2) is equivalent to np.where (m, df1, df2).

For further details and examples see the where documentation in indexing.

## **Examples**

```
>>> s = pd.Series(range(5))
>>> s.where(s > 0)
0    NaN
1    1.0
2    2.0
```

(continues on next page)

```
3 3.0
4 4.0
```

```
>>> s.mask(s > 0)
0     0.0
1     NaN
2     NaN
3     NaN
4     NaN
```

```
>>> s.where(s > 1, 10)
0    10.0
1    10.0
2    2.0
3    3.0
4    4.0
```

```
>>> df = pd.DataFrame(np.arange(10).reshape(-1, 2), columns=['A', 'B'])
>>> m = df % 3 == 0
>>> df.where(m, -df)
   A B
0 0 -1
1 -2 3
2 - 4 - 5
3 \quad 6 \quad -7
4 - 8 9
\rightarrow \rightarrow df.where(m, -df) == np.where(m, df, -df)
      Α
            В
0 True True
  True True
  True True
3
 True True
4 True True
>>> df.where(m, -df) == df.mask(\simm, -df)
      Α
            В
  True True
   True True
         True
   True
3
   True True
4
  True True
```

## **xs** (key, axis=0, level=None, drop\_level=True)

Returns a cross-section (row(s) or column(s)) from the Series/DataFrame. Defaults to cross-section on the rows (axis=0).

## **Parameters**

key [object] Some label contained in the index, or partially in a MultiIndex

axis [int, default 0] Axis to retrieve cross-section on

**level** [object, defaults to first n levels (n=1 or len(key))] In case of a key partially contained in a MultiIndex, indicate which levels are used. Levels can be referred by label or position.

**drop\_level** [boolean, default True] If False, returns object with same levels as self.

## Returns

xs [Series or DataFrame]

## **Notes**

xs is only for getting, not setting values.

MultiIndex Slicers is a generic way to get/set values on any level or levels. It is a superset of xs functionality, see MultiIndex Slicers

## **Examples**

```
>>> df
  A B C
  4 5
       2
b 4 0 9
c 9 7
>>> df.xs('a')
A
   4
В
    5
С
    2
Name: a
>>> df.xs('C', axis=1)
а
    9
b
С
    3
Name: C
```

```
>>> df
                   A B C D
first second third
bar one 1 4 1 8 9 two 1 7 5 5 0 baz one 1 6 6 8 0 three 2 5 3 5 3
>>> df.xs(('baz', 'three'))
     A B C D
third
2 5 3 5 3
>>> df.xs('one', level=1)
            A B C D
first third
bar 1
            4 1 8 9
    1
            6 6 8
                     0
>>> df.xs(('baz', 2), level=[0, 'third'])
      A B C D
second
three 5 3 5 3
```

Modified *pandas.Panel* to adopt higher dimension data than *ChemDataFrame*. Main purpose is to store molecular fingerprints in one column and keep 2D numpy array underneath.

New in version 0.3.

### **Attributes**

at Access a single value for a row/column label pair.

axes Return index label(s) of the internal NDFrame

blocks Internal property, property synonym for as\_blocks()

dtypes Return the dtypes in the DataFrame.

empty Indicator whether DataFrame is empty.

ftypes Return the ftypes (indication of sparse/dense and dtype) in DataFrame.

iat Access a single value for a row/column pair by integer position.

iloc Purely integer-location based indexing for selection by position.

### is\_copy

items items

**ix** A primarily label-location based indexer, with integer position fallback.

**loc** Access a group of rows and columns by label(s) or a boolean array.

major\_axis major\_axis

minor\_axis minor\_axis

**ndim** Return an int representing the number of axes / array dimensions.

**shape** Return a tuple of axis dimensions

size Return an int representing the number of elements in this object.

values Return a Numpy representation of the DataFrame.

## **Methods**

abs()	Return a Series/DataFrame with absolute numeric
	value of each element.
add(other[, axis])	Addition of series and other, element-wise (binary
	operator <i>add</i> ).
add_prefix(prefix)	Prefix labels with string <i>prefix</i> .
add_suffix(suffix)	Suffix labels with string <i>suffix</i> .
all([axis, bool_only, skipna, level])	Return whether all elements are True, potentially
	over an axis.
any([axis, bool_only, skipna, level])	Return whether any element is True over requested
	axis.
apply(func[, axis])	Applies function along axis (or axes) of the Panel
as_blocks([copy])	Convert the frame to a dict of dtype -> Constructor
	Types that each has a homogeneous dtype.
<pre>asfreq(freq[, method, how, normalize,])</pre>	Convert TimeSeries to specified frequency.
asof(where[, subset])	The last row without any NaN is taken (or the last
	row without NaN considering only the subset of
	columns in the case of a DataFrame)
astype(**kwargs)	Cast a pandas object to a specified dtype dtype.
at_time(time[, asof])	Select values at particular time of day (e.g.
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between_time(start_time, end_time[,])	Select values between particular times of the day (e.g., 9:00-9:30 AM).
bfill([axis, inplace, limit, downcast])	Synonym for DataFrame.
DITII([uxis, inplace, inint, downeast])	fillna (method='bfill')
bool()	Return the bool of a single element PandasObject.
clip([lower, upper, axis, inplace])	Trim values at input threshold(s).
clip_lower(threshold[, axis, inplace])	Return copy of the input with values below a thresh-
•	old truncated.
<pre>clip_upper(threshold[, axis, inplace])</pre>	Return copy of input with values above given value(s) truncated.
compound([axis, skipna, level])	Return the compound percentage of the values for the requested axis
conform(frame[, axis])	Conform input DataFrame to align with chosen axis
(	pair.
consolidate([inplace])	Compute NDFrame with "consolidated" internals
(L T)	(data of each dtype grouped together in a single ndar-
	ray).
<pre>convert_objects([convert_dates,])</pre>	Attempt to infer better dtype for object columns.
copy([deep])	Make a copy of this object's indices and data.
count([axis])	Return number of observations over requested axis.
cummax([axis, skipna])	Return cumulative maximum over a DataFrame or
(Lances, 2	Series axis.
cummin([axis, skipna])	Return cumulative minimum over a DataFrame or
(for a) a L (d)	Series axis.
cumprod([axis, skipna])	Return cumulative product over a DataFrame or Se-
1 1	ries axis.
cumsum([axis, skipna])	Return cumulative sum over a DataFrame or Series
	axis.
describe([percentiles, include, exclude])	Generates descriptive statistics that summarize the
-	central tendency, dispersion and shape of a dataset's
	distribution, excluding NaN values.
div(other[, axis])	Floating division of series and other, element-wise
	(binary operator truediv).
divide(other[, axis])	Floating division of series and other, element-wise
	(binary operator <i>truediv</i> ).
dropna([axis, how, inplace])	Drop 2D from panel, holding passed axis constant
eq(other[, axis])	Wrapper for comparison method eq
equals(other)	Determines if two NDFrame objects contain the
	same elements.
ffill([axis, inplace, limit, downcast])	Synonym for DataFrame.
	fillna(method='ffill')
fillna([value, method, axis, inplace,])	Fill NA/NaN values using the specified method
filter([items, like, regex, axis])	Subset rows or columns of dataframe according to
	labels in the specified index.
first(offset)	Convenience method for subsetting initial periods of
	time series data based on a date offset.
first_valid_index()	Return index for first non-NA/null value.
floordiv(other[, axis])	Integer division of series and other, element-wise (bi-
	nary operator floordiv).
<pre>fromDict(data[, intersect, orient, dtype])</pre>	Construct Panel from dict of DataFrame objects
<pre>from_dict(data[, intersect, orient, dtype])</pre>	Construct Panel from dict of DataFrame objects
	Continued on next page

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ge(other], axis])  get(key], default])  get(key], default])  get_dtype_counts()  get_dtype_counts()  get_titepe_counts()  get_value(*args, **kwargs)  groupby(function[, axis])  groupby(function[, axis])  groupby(function[, axis])  groupby(function[, axis])  groupdata on given axis, returning GroupBy object  g(other[, axis])  interpolare([method, axis, limit, inplace,))  isna()  isna()  isna()  isna()  isna()  isna()  isna()  iteritems()  foin(other[, how, Isuflix, rsuflix])  keys()  kurt([axis, skipna, level, numeric_only])  kurtosis([axis, skipna, level, numeric_only])  fisher's definition of kurtosis (kurtosis of normal == 0.0).  last(offset)  convenience method for subsetting final periods of time series data based on a date offset.  last_valid_index()  return index for last non-NAvanlul value.  le(other], axis])  mac([axis, skipna, level, numeric_only])  mac([axis, skipna, level])  mac([axis, skipna, level])  mac([axis, skipna, level], numeric_only])  mac([axis, skipna, level])  mac([axis, skipna, level])  mac([axis, skipna, level], numeric_only])  mac([axis, skipna, level])  mac([axis, skipna, level])  mac([axis, skipna, level], numeric_only])  mac([axis, skipna,		a from previous page
umn, Panel slice, etc.).   get_dtype_counts()   Return counts of unique dtypes in this object.   get_value(*args, **kwargs)   Quickly retrieve single value at (item, major, minor) location   get_values()   Return an indurray after converting sparse values to dense.   groupby(function[, axis])   Group data on given axis, returning GroupBy object gt(other[, axis])   Wrapper for comparison method gt     interpolate([method, axis, limit, inplace,])   Interpolate values according to different methods.     isna()   Detect missing values.     isna()   Detect missin	ge(other[, axis])	Wrapper for comparison method ge
get_dtype_counts()         Return counts of unique dtypes in this object.           get_value(*args, **kwargs)         Quickly retrieve single value at (item, major, minor) location           get_values()         Return an ndarray after converting sparse values to dense.           groupby(function[, axis])         Group data on given axis, returning GroupBy object volumes.           groupby(function[, axis])         Wrapper for comparison method gt           interpolate([method, axis, limit, inplace,])         Interpolate values according to different methods.           isnal()         Detect missing values.           iteritems()         Join items with other Panel either on major and minor axes column           keys()         Get the 'info axis' (see Indexing for more)           kurtosis([axis, skipna, level, numeric_only])         Return unbiased kurtosis (ver requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0).           kurtosis([axis, skipna, level, numeric_only])         Return index for last non-NA/null value.           i=o(other[, axis])         Wrapper for comparison	get(key[, default])	
get_ftype_counts()         Return counts of unique flypes in this object.           get_value(*args, **kwargs)         Quickly retrieve single value at (item, major, minor) location           get_values()         Return an ndarray after converting sparse values to dense.           groupby(function(, axis))         Group data on given axis, returning GroupBy object of the return of the values of the dense.           groupby(function(, axis))         Wrapper for comparison method gt the values of the values for the return of the values in the object.           groupby(function(, axis))         Attempt to infer for the return of the values of the value of values of the values for the requested axis (axis)           group for value of values of valu		<u> </u>
get_values()         Quickly retrieve single value at (item, major, minor) location           get_values()         Return an ndarray after converting sparse values to dense.           groupby(function[, axis])         Group data on given axis, returning GroupBy object of the comparison method gt.           dinter_objects()         Attempt to infer better dtypes for object columns.           interpolate([method, axis, limit, inplace,])         Interpolate values according to different methods.           isna()         Detect missing values.           isna()         Detect missing values.           isna()         Detect missing values.           isna()         Join items with other Panel either on major and minor axes column           keys()         Get the 'info axis' (see Indexing for more)           kurt([axis, skipna, level, numeric_only])         Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0).           kurtosis([axis, skipna, level, numeric_only])         Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0).           last(offset)         Convenience method for subsetting final periods of time series data based on a date offset.           last(valid, axis)         Wrapper for comparison method lt           let(other[, axis])         Wrapper for comparison method lt           max([axis, skipna, level], numeric_only])         Return the me		
location   Return an Indarray after converting sparse values to dense.   groupby(function[, axis])   Group data on given axis, returning GroupBy object gt(other[, axis])   Wrapper for comparison method gt   Attempt to infer better dtypes for object columns.   Inter_objects()   Attempt to infer better dtypes for object columns.   Interpolate (Imethod, axis, limit, inplace,  )   Interpolate values according to different methods.   isna()   Detect missing values.   Detect missing values.		
Return an ndarray after converting sparse values to dense.   Group data on given axis, returning GroupBy object	get_value(*args, **kwargs)	Quickly retrieve single value at (item, major, minor)
groupby(function[, axis])         Group data on given axis, returning GroupBy object           group by (tother[, axis])         Wrapper for comparison method gt           inter_objects()         Attempt to infer better dtypes for object columns.           interpolate([method, axis, limit, inplace,])         Interpolate values according to different methods.           isna()         Detect missing values.           iterate over (label, values) on info axis         Detect missing values.           join(other[, how, lsuffix, rsuffix])         Join items with other Panel either on major and minor axes column           keys()         Get the 'info axis' (see Indexing for more)           kurt(axis, skipna, level, numeric_only])         Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0).           kurtosis([axis, skipna, level, numeric_only])         Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0).           last_valid_index()         Convenience method for subsetting final periods of time series data based on a date offset.           last_valid_index()         Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0).           leother[, axis])         Wrapper for comparison method le           leother[, axis]         Wrapper for comparison method lt           max([axis, skipna, level])         Return the mean absolute deviatio		location
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Continued on next page	ne(other[, axis])	Wrapper for comparison method ne
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Table 20 – continued from previous page

Table 20 – continue	d from previous page
notna()	Detect existing (non-missing) values.
notnull()	Detect existing (non-missing) values.
<pre>pct_change([periods, fill_method, limit, freq])</pre>	Percentage change between the current and a prior element.
pipe(func, *args, **kwargs)	Apply func(self, *args, **kwargs)
pop(item)	Return item and drop from frame.
pow(other[, axis])	Exponential power of series and other, element-wise
	(binary operator <i>pow</i> ).
prod([axis, skipna, level, numeric_only,])	Return the product of the values for the requested axis
<pre>product([axis, skipna, level, numeric_only,])</pre>	Return the product of the values for the requested axis
radd(other[, axis])	Addition of series and other, element-wise (binary operator <i>radd</i> ).
rank([axis, method, numeric_only,])	Compute numerical data ranks (1 through n) along
(Co. 1)	axis.
rdiv(other[, axis])	Floating division of series and other, element-wise
	(binary operator <i>rtruediv</i> ).
reindex(*args, **kwargs)	Conform Panel to new index with optional filling
	logic, placing NA/NaN in locations having no value in the previous index.
reindex_axis(labels[, axis, method, level,])	Conform input object to new index with optional fill-
	ing logic, placing NA/NaN in locations having no
	value in the previous index.
<pre>reindex_like(other[, method, copy, limit,])</pre>	Return an object with matching indices to myself.
rename([items, major_axis, minor_axis])	Alter axes input function or functions.
rename_axis(mapper[, axis, copy, inplace])	Alter the name of the index or columns.
replace([to_replace, value, inplace, limit,])	Replace values given in to_replace with value.
resample(rule[, how, axis, fill_method,])	Convenience method for frequency conversion and
1 ( 1) / / = / 1)	resampling of time series.
rfloordiv(other[, axis])	Integer division of series and other, element-wise (bi-
7	nary operator <i>rfloordiv</i> ).
rmod(other[, axis])	Modulo of series and other, element-wise (binary op-
	erator <i>rmod</i> ).
rmul(other[, axis])	Multiplication of series and other, element-wise (bi-
	nary operator rmul).
round([decimals])	Round each value in Panel to a specified number of
	decimal places.
rpow(other[, axis])	Exponential power of series and other, element-wise
	(binary operator <i>rpow</i> ).
rsub(other[, axis])	Subtraction of series and other, element-wise (binary
	operator rsub).
rtruediv(other[, axis])	Floating division of series and other, element-wise
	(binary operator rtruediv).
sample([n, frac, replace, weights,])	Return a random sample of items from an axis of
	object.
select(crit[, axis])	Return data corresponding to axis labels matching
	criteria
sem([axis, skipna, level, ddof, numeric_only])	Return unbiased standard error of the mean over re-
	quested axis.
set_axis(labels[, axis, inplace])	Assign desired index to given axis.
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	d from previous page
set_value(*args, **kwargs)	Quickly set single value at (item, major, minor) location
shift([periods, freq, axis])	Shift index by desired number of periods with an op-
SHIII e([periods, freq, axis])	tional time freq.
skew([axis, skipna, level, numeric_only])	Return unbiased skew over requested axis Normal-
SAEW ([axis, skipila, level, numeric_omy])	ized by N-1
slice_shift([periods, axis])	Equivalent to <i>shift</i> without copying data.
sort_index([axis, level, ascending,])	Sort object by labels (along an axis)
sort_values([by, axis, ascending, inplace,])	NOT IMPLEMENTED: do not call this method, as
	sorting values is not supported for Panel objects and
	will raise an error.
squeeze([axis])	Squeeze length 1 dimensions.
std([axis, skipna, level, ddof, numeric_only])	Return sample standard deviation over requested
(t / 1 / / / / = 33)	axis.
sub(other[, axis])	Subtraction of series and other, element-wise (binary
( 1)	operator <i>sub</i> ).
subtract(other[, axis])	Subtraction of series and other, element-wise (binary
(- · · · · · · · · · · · · · · · ·	operator <i>sub</i> ).
sum([axis, skipna, level, numeric_only,])	Return the sum of the values for the requested axis
swapaxes(axis1, axis2[, copy])	Interchange axes and swap values axes appropriately
swaplevel([i, j, axis])	Swap levels i and j in a MultiIndex on a particular
	axis
take(indices[, axis, convert, is_copy])	Return the elements in the given <i>positional</i> indices
, , , , , = 1,,,,	along an axis.
to_clipboard([excel, sep])	Copy object to the system clipboard.
to_dense()	Return dense representation of NDFrame (as op-
_ · ·	posed to sparse)
to_excel(path[, na_rep, engine])	Write each DataFrame in Panel to a separate excel
	sheet
to_frame([filter_observations])	Transform wide format into long (stacked) format as
	DataFrame whose columns are the Panel's items and
	whose index is a MultiIndex formed of the Panel's
	major and minor axes.
to_hdf(path_or_buf, key, **kwargs)	Write the contained data to an HDF5 file using HDF-
	Store.
to_json([path_or_buf, orient, date_format,])	Convert the object to a JSON string.
<pre>to_latex([buf, columns, col_space, header,])</pre>	Render an object to a tabular environment table.
to_msgpack([path_or_buf, encoding])	msgpack (serialize) object to input file path
to_pickle(path[, compression, protocol])	Pickle (serialize) object to file.
to_sparse(*args, **kwargs)	NOT IMPLEMENTED: do not call this method, as
	sparsifying is not supported for Panel objects and
	will raise an error.
to_sql(name, con[, schema, if_exists,])	Write records stored in a DataFrame to a SQL
	database.
to_xarray()	Return an xarray object from the pandas object.
transpose(*args, **kwargs)	Permute the dimensions of the Panel
truediv(other[, axis])	Floating division of series and other, element-wise
	(binary operator <i>truediv</i> ).
truncate([before, after, axis, copy])	Truncate a Series or DataFrame before and after
	some index value.
tz_convert(tz[, axis, level, copy])	Convert tz-aware axis to target time zone.
	Continued on next page

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tz_localize(tz[, axis, level, copy, ambiguous])	Localize tz-naive TimeSeries to target time zone.
update(other[, join, overwrite,])	Modify Panel in place using non-NA values from
	passed Panel, or object coercible to Panel.
var([axis, skipna, level, ddof, numeric_only])	Return unbiased variance over requested axis.
where(cond[, other, inplace, axis, level,])	Return an object of same shape as self and whose
	corresponding entries are from self where cond is
	True and otherwise are from <i>other</i> .
xs(key[, axis])	Return slice of panel along selected axis

## abs()

Return a Series/DataFrame with absolute numeric value of each element.

This function only applies to elements that are all numeric.

## **Returns**

abs Series/DataFrame containing the absolute value of each element.

#### See also:

numpy.absolute calculate the absolute value element-wise.

## **Notes**

For complex inputs, 1.2 + 1j, the absolute value is  $\sqrt{a^2+b^2}$ .

## **Examples**

Absolute numeric values in a Series.

```
>>> s = pd.Series([-1.10, 2, -3.33, 4])
>>> s.abs()
0     1.10
1     2.00
2     3.33
3     4.00
dtype: float64
```

Absolute numeric values in a Series with complex numbers.

```
>>> s = pd.Series([1.2 + 1j])
>>> s.abs()
```

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```
0 1.56205
dtype: float64
```

Absolute numeric values in a Series with a Timedelta element.

```
>>> s = pd.Series([pd.Timedelta('1 days')])
>>> s.abs()
0  1 days
dtype: timedelta64[ns]
```

Select rows with data closest to certain value using argsort (from StackOverflow).

```
>>> df = pd.DataFrame({
        'a': [4, 5, 6, 7],
        'b': [10, 20, 30, 40],
        'c': [100, 50, -30, -50]
. . .
...})
>>> df
          b
               C
         10 100
     4
0
     5
         20
             50
1
2
     6
         30
            -30
3
     7
         40
             -50
>>> df.loc[(df.c - 43).abs().argsort()]
         b
     а
               C
         20
              50
1
     5
0
         10 100
     4
2
         30
             -30
3
         40
             -50
```

## add(other, axis=0)

Addition of series and other, element-wise (binary operator add). Equivalent to panel + other.

### **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

### **Returns**

**Panel** 

# See also:

Panel.radd

## add\_prefix (prefix)

Prefix labels with string prefix.

For Series, the row labels are prefixed. For DataFrame, the column labels are prefixed.

### **Parameters**

**prefix** [str] The string to add before each label.

## Returns

Series or DataFrame New Series or DataFrame with updated labels.

See also:

Series.add\_suffix Suffix row labels with string suffix.

DataFrame.add\_suffix Suffix column labels with string suffix.

## **Examples**

```
>>> df = pd.DataFrame({'A': [1, 2, 3, 4], 'B': [3, 4, 5, 6]})
>>> df

A B
0 1 3
1 2 4
2 3 5
3 4 6
```

## add\_suffix (suffix)

Suffix labels with string *suffix*.

For Series, the row labels are suffixed. For DataFrame, the column labels are suffixed.

#### **Parameters**

suffix [str] The string to add after each label.

# Returns

Series or DataFrame New Series or DataFrame with updated labels.

## See also:

**Series.add\_prefix** Prefix row labels with string *prefix*.

 ${\tt DataFrame.add\_prefix}\ \ Prefix\ column\ labels\ with\ string\ \textit{prefix}.$ 

## **Examples**

```
>>> df = pd.DataFrame({'A': [1, 2, 3, 4], 'B': [3, 4, 5, 6]})
>>> df

A B
0 1 3
1 2 4
2 3 5
3 4 6
```

```
agg (func, *args, **kwargs)
```

aggregate (func, \*args, \*\*kwargs)

align (other, \*\*kwargs)

Align two objects on their axes with the specified join method for each axis Index

### **Parameters**

```
other [DataFrame or Series]
```

```
join [{'outer', 'inner', 'left', 'right'}, default 'outer']
```

**axis** [allowed axis of the other object, default None] Align on index (0), columns (1), or both (None)

**level** [int or level name, default None] Broadcast across a level, matching Index values on the passed MultiIndex level

**copy** [boolean, default True] Always returns new objects. If copy=False and no reindexing is required then original objects are returned.

**fill\_value** [scalar, default np.NaN] Value to use for missing values. Defaults to NaN, but can be any "compatible" value

method [str, default None]

**limit** [int, default None]

fill\_axis [int or labels for object, default 0] Filling axis, method and limit

**broadcast\_axis** [int or labels for object, default None] Broadcast values along this axis, if aligning two objects of different dimensions

### Returns

(left, right) [(NDFrame, type of other)] Aligned objects

**all** (axis=0, bool\_only=None, skipna=True, level=None, \*\*kwargs) Return whether all elements are True, potentially over an axis.

Returns True if all elements within a series or along a Dataframe axis are non-zero, not-empty or not-False.

## **Parameters**

axis [{0 or 'index', 1 or 'columns', None}, default 0] Indicate which axis or axes should be reduced.

- 0 / 'index' : reduce the index, return a Series whose index is the original column labels.
- 1 / 'columns' : reduce the columns, return a Series whose index is the original index.
- None: reduce all axes, return a scalar.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame.

**bool\_only** [boolean, default None] Include only boolean columns. If None, will attempt to use everything, then use only boolean data. Not implemented for Series.

\*\*kwargs [any, default None] Additional keywords have no effect but might be accepted for compatibility with NumPy.

# Returns

**all** [DataFrame or Panel (if level specified)]

## See also:

```
pandas.Series.all Return True if all elements are True
pandas.DataFrame.any Return True if one (or more) elements are True
```

## **Examples**

## Series

```
>>> pd.Series([True, True]).all()
True
>>> pd.Series([True, False]).all()
False
```

## **DataFrames**

Create a dataframe from a dictionary.

```
>>> df = pd.DataFrame({'col1': [True, True], 'col2': [True, False]})
>>> df
    col1    col2
0    True    True
1    True    False
```

Default behaviour checks if column-wise values all return True.

```
>>> df.all()
col1 True
col2 False
dtype: bool
```

Specify axis='columns' to check if row-wise values all return True.

```
>>> df.all(axis='columns')
0 True
1 False
dtype: bool
```

Or axis=None for whether every value is True.

```
>>> df.all(axis=None)
False
```

any (axis=0, bool\_only=None, skipna=True, level=None, \*\*kwargs)

Return whether any element is True over requested axis.

Unlike DataFrame.all(), this performs an *or* operation. If any of the values along the specified axis is True, this will return True.

## **Parameters**

axis [{0 or 'index', 1 or 'columns', None}, default 0] Indicate which axis or axes should be reduced.

- 0 / 'index': reduce the index, return a Series whose index is the original column labels.
- 1 / 'columns' : reduce the columns, return a Series whose index is the original index.
- None: reduce all axes, return a scalar.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame.

**bool\_only** [boolean, default None] Include only boolean columns. If None, will attempt to use everything, then use only boolean data. Not implemented for Series.

\*\*kwargs [any, default None] Additional keywords have no effect but might be accepted for compatibility with NumPy.

## Returns

any [DataFrame or Panel (if level specified)]

See also:

pandas.DataFrame.all Return whether all elements are True.

# **Examples**

## **Series**

For Series input, the output is a scalar indicating whether any element is True.

```
>>> pd.Series([True, False]).any()
True
```

#### **DataFrame**

Whether each column contains at least one True element (the default).

```
>>> df = pd.DataFrame({"A": [1, 2], "B": [0, 2], "C": [0, 0]})
>>> df

A B C
0 1 0 0
1 2 2 0
```

```
>>> df.any()
A True
B True
C False
dtype: bool
```

Aggregating over the columns.

```
>>> df = pd.DataFrame({"A": [True, False], "B": [1, 2]})
>>> df

A B

0 True 1

1 False 2
```

```
>>> df.any(axis='columns')

0 True

1 True

dtype: bool
```

```
>>> df = pd.DataFrame({"A": [True, False], "B": [1, 0]})
>>> df

A B

0 True 1

1 False 0
```

```
>>> df.any(axis='columns')
0 True
1 False
dtype: bool
```

Aggregating over the entire DataFrame with axis=None.

```
>>> df.any(axis=None)
True
```

any for an empty DataFrame is an empty Series.

```
>>> pd.DataFrame([]).any()
Series([], dtype: bool)
```

```
apply (func, axis='major', **kwargs)
```

Applies function along axis (or axes) of the Panel

### **Parameters**

```
func [function] Function to apply to each combination of 'other' axes e.g. if axis = 'items', the combination of major_axis/minor_axis will each be passed as a Series; if axis = ('items', 'major'), DataFrames of items & major axis will be passed
```

```
axis [{'items', 'minor', 'major'}, or {0, 1, 2}, or a tuple with two] axes
```

## Additional keyword arguments will be passed as keywords to the function

### Returns

result [Panel, DataFrame, or Series]

## **Examples**

Returns a Panel with the square root of each element

```
>>> p = pd.Panel(np.random.rand(4,3,2))
>>> p.apply(np.sqrt)
```

Equivalent to p.sum(1), returning a DataFrame

```
>>> p.apply(lambda x: x.sum(), axis=1)
```

Equivalent to previous:

```
>>> p.apply(lambda x: x.sum(), axis='major')
```

Return the shapes of each DataFrame over axis 2 (i.e the shapes of items x major), as a Series

```
>>> p.apply(lambda x: x.shape, axis=(0,1))
```

## as\_blocks(copy=True)

Convert the frame to a dict of dtype -> Constructor Types that each has a homogeneous dtype.

Deprecated since version 0.21.0.

**NOTE:** the dtypes of the blocks WILL BE PRESERVED HERE (unlike in as\_matrix)

## **Parameters**

```
copy [boolean, default True]
```

### **Returns**

**values** [a dict of dtype -> Constructor Types]

## as\_matrix()

Convert the frame to its Numpy-array representation.

Deprecated since version 0.23.0: Use DataFrame.values() instead.

### **Parameters**

**columns: list, optional, default:None** If None, return all columns, otherwise, returns specified columns.

#### Returns

values [ndarray] If the caller is heterogeneous and contains booleans or objects, the result will be of dtype=object. See Notes.

#### See also:

pandas.DataFrame.values

#### **Notes**

Return is NOT a Numpy-matrix, rather, a Numpy-array.

The dtype will be a lower-common-denominator dtype (implicit upcasting); that is to say if the dtypes (even of numeric types) are mixed, the one that accommodates all will be chosen. Use this with care if you are not dealing with the blocks.

e.g. If the dtypes are float16 and float32, dtype will be upcast to float32. If dtypes are int32 and uint8, dtype will be upcase to int32. By numpy.find\_common\_type convention, mixing int64 and uint64 will result in a flot64 dtype.

This method is provided for backwards compatibility. Generally, it is recommended to use '.values'.

**asfreq** (freq, method=None, how=None, normalize=False, fill\_value=None) Convert TimeSeries to specified frequency.

Optionally provide filling method to pad/backfill missing values.

Returns the original data conformed to a new index with the specified frequency. resample is more appropriate if an operation, such as summarization, is necessary to represent the data at the new frequency.

### **Parameters**

**freq** [DateOffset object, or string]

**method** [{'backfill'/'bfill', 'pad'/'ffill'}, default None] Method to use for filling holes in reindexed Series (note this does not fill NaNs that already were present):

- 'pad' / 'ffill': propagate last valid observation forward to next valid
- 'backfill' / 'bfill': use NEXT valid observation to fill

**how** [{'start', 'end'}, default end] For PeriodIndex only, see PeriodIndex.asfreq

**normalize** [bool, default False] Whether to reset output index to midnight

**fill\_value: scalar, optional** Value to use for missing values, applied during upsampling (note this does not fill NaNs that already were present).

New in version 0.20.0.

### **Returns**

**converted** [type of caller]

## See also:

reindex

### **Notes**

To learn more about the frequency strings, please see this link.

## **Examples**

Start by creating a series with 4 one minute timestamps.

Upsample the series into 30 second bins.

Upsample again, providing a fill value.

```
>>> df.asfreq(freq='30S', fill_value=9.0)
                       S
2000-01-01 00:00:00
                       0.0
2000-01-01 00:00:30
                       9.0
2000-01-01 00:01:00
                       NaN
2000-01-01 00:01:30
                       9.0
2000-01-01 00:02:00
                       2.0
2000-01-01 00:02:30
                       9.0
2000-01-01 00:03:00
                       3.0
```

Upsample again, providing a method.

```
>>> df.asfreq(freq='30S', method='bfill')
                       S
2000-01-01 00:00:00
                       0.0
2000-01-01 00:00:30
                       NaN
2000-01-01 00:01:00
                       NaN
2000-01-01 00:01:30
                       2.0
2000-01-01 00:02:00
                       2.0
2000-01-01 00:02:30
                       3.0
2000-01-01 00:03:00
                       3 0
```

asof (where, subset=None)

The last row without any NaN is taken (or the last row without NaN considering only the subset of columns in the case of a DataFrame)

New in version 0.19.0: For DataFrame

If there is no good value, NaN is returned for a Series a Series of NaN values for a DataFrame

## **Parameters**

where [date or array of dates]

**subset** [string or list of strings, default None] if not None use these columns for NaN propagation

### Returns

### where is scalar

- value or NaN if input is Series
- Series if input is DataFrame

where is Index: same shape object as input

### See also:

merge asof

### **Notes**

Dates are assumed to be sorted Raises if this is not the case

```
astype (**kwargs)
```

Cast a pandas object to a specified dtype dtype.

### **Parameters**

**dtype** [data type, or dict of column name -> data type] Use a numpy.dtype or Python type to cast entire pandas object to the same type. Alternatively, use {col: dtype, ...}, where col is a column label and dtype is a numpy.dtype or Python type to cast one or more of the DataFrame's columns to column-specific types.

**copy** [bool, default True.] Return a copy when copy=True (be very careful setting copy=False as changes to values then may propagate to other pandas objects).

**errors** [{'raise', 'ignore'}, default 'raise'.] Control raising of exceptions on invalid data for provided dtype.

- raise: allow exceptions to be raised
- ignore: suppress exceptions. On error return original object

New in version 0.20.0.

raise\_on\_error [raise on invalid input] Deprecated since version 0.20.0: Use errors
instead

**kwargs** [keyword arguments to pass on to the constructor]

### **Returns**

**casted** [type of caller]

## See also:

```
pandas.to_datetime Convert argument to datetime.
```

pandas.to\_timedelta Convert argument to timedelta.

pandas.to\_numeric Convert argument to a numeric type.

numpy.ndarray.astype Cast a numpy array to a specified type.

## **Examples**

### Convert to categorical type:

```
>>> ser.astype('category')
0    1
1    2
dtype: category
Categories (2, int64): [1, 2]
```

Convert to ordered categorical type with custom ordering:

Note that using copy=False and changing data on a new pandas object may propagate changes:

```
>>> s1 = pd.Series([1,2])
>>> s2 = s1.astype('int64', copy=False)
>>> s2[0] = 10
>>> s1  # note that s1[0] has changed too
0  10
1  2
dtype: int64
```

at

Access a single value for a row/column label pair.

Similar to loc, in that both provide label-based lookups. Use at if you only need to get or set a single value in a DataFrame or Series.

Raises

**KeyError** When label does not exist in DataFrame

See also:

DataFrame.iat Access a single value for a row/column pair by integer position

**DataFrame.loc** Access a group of rows and columns by label(s)

Series.at Access a single value using a label

## **Examples**

Get value at specified row/column pair

```
>>> df.at[4, 'B']
2
```

Set value at specified row/column pair

```
>>> df.at[4, 'B'] = 10
>>> df.at[4, 'B']
10
```

Get value within a Series

```
>>> df.loc[5].at['B']
4
```

at\_time (time, asof=False)

Select values at particular time of day (e.g. 9:30AM).

## **Parameters**

**time** [datetime.time or string]

Returns

values\_at\_time [type of caller]

**Raises** 

 $\textbf{TypeError} \ \ \textbf{If the index is not a} \ \textbf{DatetimeIndex}$ 

See also:

between\_time Select values between particular times of the day

first Select initial periods of time series based on a date offset

last Select final periods of time series based on a date offset

DatetimeIndex.indexer\_at\_time Get just the index locations for values at particular time of
 the day

## **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='12H')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
```

(continues on next page)

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```
2018-04-09 00:00:00 1
2018-04-09 12:00:00 2
2018-04-10 00:00:00 3
2018-04-10 12:00:00 4
```

```
>>> ts.at_time('12:00')

A
2018-04-09 12:00:00 2
2018-04-10 12:00:00 4
```

#### axes

Return index label(s) of the internal NDFrame

 $\textbf{between\_time} \ (\textit{start\_time}, \textit{end\_time}, \textit{include\_start=True}, \textit{include\_end=True})$ 

Select values between particular times of the day (e.g., 9:00-9:30 AM).

By setting start\_time to be later than end\_time, you can get the times that are *not* between the two times.

### **Parameters**

```
start_time [datetime.time or string]
end_time [datetime.time or string]
include_start [boolean, default True]
include_end [boolean, default True]
```

### Returns

values\_between\_time [type of caller]

### **Raises**

TypeError If the index is not a DatetimeIndex

## See also:

at\_time Select values at a particular time of the day

first Select initial periods of time series based on a date offset

last Select final periods of time series based on a date offset

DatetimeIndex.indexer\_between\_time Get just the index locations for values between particular times of the day

## **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='1D20min')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 00:00:00 1
2018-04-10 00:20:00 2
2018-04-11 00:40:00 3
2018-04-12 01:00:00 4
```

```
>>> ts.between_time('0:15', '0:45')

A
2018-04-10 00:20:00 2
2018-04-11 00:40:00 3
```

You get the times that are *not* between two times by setting start\_time later than end\_time:

```
>>> ts.between_time('0:45', '0:15')

A
2018-04-09 00:00:00 1
2018-04-12 01:00:00 4
```

## **bfill** (axis=None, inplace=False, limit=None, downcast=None)

Synonym for DataFrame.fillna (method='bfill')

### blocks

Internal property, property synonym for as\_blocks()

Deprecated since version 0.21.0.

### bool()

Return the bool of a single element PandasObject.

This must be a boolean scalar value, either True or False. Raise a ValueError if the PandasObject does not have exactly 1 element, or that element is not boolean

```
clip (lower=None, upper=None, axis=None, inplace=False, *args, **kwargs)
Trim values at input threshold(s).
```

Assigns values outside boundary to boundary values. Thresholds can be singular values or array like, and in the latter case the clipping is performed element-wise in the specified axis.

### **Parameters**

**lower** [float or array\_like, default None] Minimum threshold value. All values below this threshold will be set to it.

**upper** [float or array\_like, default None] Maximum threshold value. All values above this threshold will be set to it.

**axis** [int or string axis name, optional] Align object with lower and upper along the given axis.

**inplace** [boolean, default False] Whether to perform the operation in place on the data.

New in version 0.21.0.

\*args, \*\*kwargs Additional keywords have no effect but might be accepted for compatibility with numpy.

### Returns

**Series or DataFrame** Same type as calling object with the values outside the clip boundaries replaced

## See also:

```
clip_lower Clip values below specified threshold(s).clip_upper Clip values above specified threshold(s).
```

## **Examples**

```
>>> data = {'col_0': [9, -3, 0, -1, 5], 'col_1': [-2, -7, 6, 8, -5]}
>>> df = pd.DataFrame(data)
>>> df
   col_0 col_1
       9
      -3
             -7
1
2
       0
              6
3
      -1
              8
4
       5
             -5
```

Clips per column using lower and upper thresholds:

```
>>> df.clip(-4, 6)
   col_0 col_1
0
       6
              -2
1
      -3
              -4
2
       0
               6
3
      -1
               6
4
       5
              -4
```

Clips using specific lower and upper thresholds per column element:

```
>>> df.clip(t, t + 4, axis=0)
   col_0 col_1
0
       6
               2
1
      -3
              -4
2
       0
               3
3
       6
               8
4
       5
               3
```

clip\_lower (threshold, axis=None, inplace=False)

Return copy of the input with values below a threshold truncated.

### **Parameters**

**threshold** [numeric or array-like] Minimum value allowed. All values below threshold will be set to this value.

- float : every value is compared to *threshold*.
- array-like: The shape of *threshold* should match the object it's compared to. When *self* is a Series, *threshold* should be the length. When *self* is a DataFrame, *threshold* should 2-D and the same shape as *self* for axis=None, or 1-D and the same length as the axis being compared.

**axis** [{0 or 'index', 1 or 'columns'}, default 0] Align *self* with *threshold* along the given axis.

**inplace** [boolean, default False] Whether to perform the operation in place on the data. New in version 0.21.0.

### Returns

**clipped** [same type as input]

See also:

Series.clip Return copy of input with values below and above thresholds truncated.

Series.clip\_upper Return copy of input with values above threshold truncated.

## **Examples**

Series single threshold clipping:

```
>>> s = pd.Series([5, 6, 7, 8, 9])
>>> s.clip_lower(8)
0     8
1     8
2     8
3     8
4     9
dtype: int64
```

Series clipping element-wise using an array of thresholds. *threshold* should be the same length as the Series.

```
>>> elemwise_thresholds = [4, 8, 7, 2, 5]
>>> s.clip_lower(elemwise_thresholds)
0    5
1    8
2    7
3    8
4    9
dtype: int64
```

DataFrames can be compared to a scalar.

```
>>> df = pd.DataFrame({"A": [1, 3, 5], "B": [2, 4, 6]})
>>> df

A B
0 1 2
1 3 4
2 5 6
```

```
>>> df.clip_lower(3)

A B

0 3 3

1 3 4

2 5 6
```

Or to an array of values. By default, *threshold* should be the same shape as the DataFrame.

```
>>> df.clip_lower(np.array([[3, 4], [2, 2], [6, 2]]))

A B

0 3 4

1 3 4

2 6 6
```

Control how *threshold* is broadcast with *axis*. In this case *threshold* should be the same length as the axis specified by *axis*.

```
>>> df.clip_lower(np.array([3, 3, 5]), axis='index')

A B
0 3 3
1 3 4
2 5 6
```

```
>>> df.clip_lower(np.array([4, 5]), axis='columns')

A B
0 4 5
1 4 5
2 5 6
```

# clip\_upper (threshold, axis=None, inplace=False)

Return copy of input with values above given value(s) truncated.

### **Parameters**

```
threshold [float or array_like]
```

axis [int or string axis name, optional] Align object with threshold along the given axis.

inplace [boolean, default False] Whether to perform the operation in place on the data

New in version 0.21.0.

## Returns

**clipped** [same type as input]

## See also:

clip

compound(axis=None, skipna=None, level=None)

Return the compound percentage of the values for the requested axis

## **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

**compounded** [DataFrame or Panel (if level specified)]

```
conform (frame, axis='items')
```

Conform input DataFrame to align with chosen axis pair.

### **Parameters**

frame [DataFrame]

axis [{'items', 'major', 'minor'}] Axis the input corresponds to. E.g., if axis='major', then the frame's columns would be items, and the index would be values of the minor axis

#### Returns

#### **DataFrame**

### consolidate(inplace=False)

Compute NDFrame with "consolidated" internals (data of each dtype grouped together in a single ndarray).

Deprecated since version 0.20.0: Consolidate will be an internal implementation only.

Attempt to infer better dtype for object columns.

Deprecated since version 0.21.0.

### **Parameters**

**convert\_dates** [boolean, default True] If True, convert to date where possible. If 'coerce', force conversion, with unconvertible values becoming NaT.

**convert\_numeric** [boolean, default False] If True, attempt to coerce to numbers (including strings), with unconvertible values becoming NaN.

**convert\_timedeltas** [boolean, default True] If True, convert to timedelta where possible. If 'coerce', force conversion, with unconvertible values becoming NaT.

**copy** [boolean, default True] If True, return a copy even if no copy is necessary (e.g. no conversion was done). Note: This is meant for internal use, and should not be confused with inplace.

### Returns

converted [same as input object]

See also:

pandas.to\_datetime Convert argument to datetime.

pandas.to\_timedelta Convert argument to timedelta.

pandas.to\_numeric Return a fixed frequency timedelta index, with day as the default.

### copy (deep=True)

Make a copy of this object's indices and data.

When deep=True (default), a new object will be created with a copy of the calling object's data and indices. Modifications to the data or indices of the copy will not be reflected in the original object (see notes below).

When deep=False, a new object will be created without copying the calling object's data or index (only references to the data and index are copied). Any changes to the data of the original will be reflected in the shallow copy (and vice versa).

## **Parameters**

**deep** [bool, default True] Make a deep copy, including a copy of the data and the indices. With deep=False neither the indices nor the data are copied.

#### Returns

**copy** [Series, DataFrame or Panel] Object type matches caller.

### **Notes**

When deep=True, data is copied but actual Python objects will not be copied recursively, only the reference to the object. This is in contrast to *copy.deepcopy* in the Standard Library, which recursively copies object data (see examples below).

While Index objects are copied when deep=True, the underlying numpy array is not copied for performance reasons. Since Index is immutable, the underlying data can be safely shared and a copy is not needed.

## **Examples**

```
>>> s = pd.Series([1, 2], index=["a", "b"])
>>> s
a    1
b    2
dtype: int64
```

## Shallow copy versus default (deep) copy:

```
>>> s = pd.Series([1, 2], index=["a", "b"])
>>> deep = s.copy()
>>> shallow = s.copy(deep=False)
```

Shallow copy shares data and index with original.

```
>>> s is shallow
False
>>> s.values is shallow.values and s.index is shallow.index
True
```

Deep copy has own copy of data and index.

```
>>> s is deep
False
>>> s.values is deep.values or s.index is deep.index
False
```

Updates to the data shared by shallow copy and original is reflected in both; deep copy remains unchanged.

Note that when copying an object containing Python objects, a deep copy will copy the data, but will not do so recursively. Updating a nested data object will be reflected in the deep copy.

```
>>> s = pd.Series([[1, 2], [3, 4]])
>>> deep = s.copy()
>>> s[0][0] = 10
>>> s
0    [10, 2]
1    [3, 4]
dtype: object
>>> deep
0    [10, 2]
1    [3, 4]
dtype: object
```

count (axis='major')

Return number of observations over requested axis.

### **Parameters**

```
axis [{'items', 'major', 'minor'} or {0, 1, 2}]
```

### Returns

count [DataFrame]

cummax (axis=None, skipna=True, \*args, \*\*kwargs)

Return cumulative maximum over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative maximum.

## **Parameters**

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

## **Returns**

cummax [DataFrame or Panel]

See also:

pandas.core.window.Expanding.max Similar functionality but ignores NaN values.

Panel.max Return the maximum over Panel axis.

Panel.cummax Return cumulative maximum over Panel axis.

Panel.cummin Return cumulative minimum over Panel axis.

Panel.cumsum Return cumulative sum over Panel axis.

Panel.cumprod Return cumulative product over Panel axis.

## **Examples**

### Series

By default, NA values are ignored.

To include NA values in the operation, use skipna=False

```
>>> s.cummax(skipna=False)

0 2.0

1 NaN

2 NaN

3 NaN

4 NaN

dtype: float64
```

### **DataFrame**

By default, iterates over rows and finds the maximum in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cummax()

A B

0 2.0 1.0

1 3.0 NaN

2 3.0 1.0
```

To iterate over columns and find the maximum in each row, use axis=1

```
>>> df.cummax(axis=1)

A B
0 2.0 2.0
1 3.0 NaN
2 1.0 1.0
```

```
cummin (axis=None, skipna=True, *args, **kwargs)
```

Return cumulative minimum over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative minimum.

#### **Parameters**

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

## Returns

**cummin** [DataFrame or Panel]

## See also:

pandas.core.window.Expanding.min Similar functionality but ignores NaN values.

Panel.min Return the minimum over Panel axis.

Panel.cummax Return cumulative maximum over Panel axis.

Panel.cummin Return cumulative minimum over Panel axis.

Panel.cumsum Return cumulative sum over Panel axis.

Panel.cumprod Return cumulative product over Panel axis.

# **Examples**

## Series

By default, NA values are ignored.

```
>>> s.cummin()
0 2.0
1 NaN
2 2.0
3 -1.0
4 -1.0
dtype: float64
```

To include NA values in the operation, use skipna=False

```
>>> s.cummin(skipna=False)
0 2.0
1 NaN
2 NaN
3 NaN
4 NaN
dtype: float64
```

## **DataFrame**

By default, iterates over rows and finds the minimum in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cummin()

A B

0 2.0 1.0

1 2.0 NaN

2 1.0 0.0
```

To iterate over columns and find the minimum in each row, use axis=1

```
>>> df.cummin(axis=1)

A B
0 2.0 1.0
1 3.0 NaN
2 1.0 0.0
```

cumprod (axis=None, skipna=True, \*args, \*\*kwargs)

Return cumulative product over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative product.

### **Parameters**

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

## **Returns**

**cumprod** [DataFrame or Panel]

#### See also:

pandas.core.window.Expanding.prod Similar functionality but ignores NaN values.

Panel.prod Return the product over Panel axis.

Panel.cummax Return cumulative maximum over Panel axis.

Panel.cummin Return cumulative minimum over Panel axis.

Panel.cumsum Return cumulative sum over Panel axis.

Panel.cumprod Return cumulative product over Panel axis.

## **Examples**

## **Series**

## By default, NA values are ignored.

```
>>> s.cumprod()
0 2.0
1 NaN
2 10.0
3 -10.0
4 -0.0
dtype: float64
```

To include NA values in the operation, use skipna=False

```
>>> s.cumprod(skipna=False)
0 2.0
1 NaN
2 NaN
3 NaN
4 NaN
dtype: float64
```

## **DataFrame**

```
>>> df = pd.DataFrame([[2.0, 1.0],
... [3.0, np.nan],
... [1.0, 0.0]],
... columns=list('AB'))
>>> df
A B
0 2.0 1.0
1 3.0 NaN
2 1.0 0.0
```

By default, iterates over rows and finds the product in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cumprod()

A B

0 2.0 1.0

1 6.0 NaN

2 6.0 0.0
```

To iterate over columns and find the product in each row, use axis=1

```
>>> df.cumprod(axis=1)

A B
0 2.0 2.0
1 3.0 NaN
2 1.0 0.0
```

cumsum (axis=None, skipna=True, \*args, \*\*kwargs)

Return cumulative sum over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative sum.

### **Parameters**

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

## Returns

cumsum [DataFrame or Panel]

See also:

pandas.core.window.Expanding.sum Similar functionality but ignores NaN values.

Panel.sum Return the sum over Panel axis.

Panel.cummax Return cumulative maximum over Panel axis.

Panel.cummin Return cumulative minimum over Panel axis.

Panel.cumsum Return cumulative sum over Panel axis.

Panel.cumprod Return cumulative product over Panel axis.

## **Examples**

### Series

By default, NA values are ignored.

```
>>> s.cumsum()
0 2.0
1 NaN
2 7.0
3 6.0
4 6.0
dtype: float64
```

To include NA values in the operation, use skipna=False

```
>>> s.cumsum(skipna=False)
0 2.0
1 NaN
2 NaN
3 NaN
4 NaN
dtype: float64
```

### **DataFrame**

```
>>> df = pd.DataFrame([[2.0, 1.0],
... [3.0, np.nan],
... [1.0, 0.0]],
... columns=list('AB'))
>>> df
A B
0 2.0 1.0
1 3.0 NaN
2 1.0 0.0
```

By default, iterates over rows and finds the sum in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cumsum()

A B

0 2.0 1.0

1 5.0 NaN

2 6.0 1.0
```

To iterate over columns and find the sum in each row, use axis=1

```
>>> df.cumsum(axis=1)

A B
0 2.0 3.0
1 3.0 NaN
2 1.0 1.0
```

**describe** (percentiles=None, include=None, exclude=None)

Generates descriptive statistics that summarize the central tendency, dispersion and shape of a dataset's distribution, excluding NaN values.

Analyzes both numeric and object series, as well as DataFrame column sets of mixed data types. The output will vary depending on what is provided. Refer to the notes below for more detail.

#### **Parameters**

**percentiles** [list-like of numbers, optional] The percentiles to include in the output. All should fall between 0 and 1. The default is [.25, .5, .75], which returns the 25th, 50th, and 75th percentiles.

**include** ['all', list-like of dtypes or None (default), optional] A white list of data types to include in the result. Ignored for Series. Here are the options:

- 'all' : All columns of the input will be included in the output.
- A list-like of dtypes: Limits the results to the provided data types. To limit the result to numeric types submit numpy.number. To limit it instead to object columns submit the numpy.object data type. Strings can also be used in the style of select\_dtypes (e.g. df.describe(include=['0'])). To select pandas categorical columns, use 'category'
- None (default): The result will include all numeric columns.

**exclude** [list-like of dtypes or None (default), optional,] A black list of data types to omit from the result. Ignored for Series. Here are the options:

- A list-like of dtypes: Excludes the provided data types from the result. To exclude numeric types submit numpy.number. To exclude object columns submit the data type numpy.object. Strings can also be used in the style of select\_dtypes (e.g. df.describe(include=['O'])). To exclude pandas categorical columns, use 'category'
- None (default): The result will exclude nothing.

## Returns

## summary: Series/DataFrame of summary statistics

### See also:

DataFrame.count, DataFrame.max, DataFrame.min, DataFrame.mean, DataFrame.std, DataFrame.select\_dtypes

#### **Notes**

For numeric data, the result's index will include count, mean, std, min, max as well as lower, 50 and upper percentiles. By default the lower percentile is 25 and the upper percentile is 75. The 50 percentile is the same as the median.

For object data (e.g. strings or timestamps), the result's index will include count, unique, top, and freq. The top is the most common value. The freq is the most common value's frequency. Timestamps also include the first and last items.

If multiple object values have the highest count, then the count and top results will be arbitrarily chosen from among those with the highest count.

For mixed data types provided via a DataFrame, the default is to return only an analysis of numeric columns. If the dataframe consists only of object and categorical data without any numeric columns, the default is to return an analysis of both the object and categorical columns. If include='all' is provided as an option, the result will include a union of attributes of each type.

The *include* and *exclude* parameters can be used to limit which columns in a DataFrame are analyzed for the output. The parameters are ignored when analyzing a Series.

## **Examples**

Describing a numeric Series.

```
>>> s = pd.Series([1, 2, 3])
>>> s.describe()
count 3.0
mean
         2.0
std
        1.0
min
         1.0
25%
        1.5
         2.0
50%
75%
         2.5
         3.0
max
```

Describing a categorical Series.

```
>>> s = pd.Series(['a', 'a', 'b', 'c'])
>>> s.describe()
count     4
unique     3
top          a
freq     2
dtype: object
```

Describing a timestamp Series.

```
>>> s = pd.Series([
... np.datetime64("2000-01-01"),
... np.datetime64("2010-01-01"),
     np.datetime64("2010-01-01")
...])
>>> s.describe()
                              3
count
                              2
unique
         2010-01-01 00:00:00
top
freq
       2000-01-01 00:00:00
2010-01-01 00:00:00
first
last
dtype: object
```

Describing a DataFrame. By default only numeric fields are returned.

```
>>> df = pd.DataFrame({ 'object': ['a', 'b', 'c'],
                      'numeric': [1, 2, 3],
                      'categorical': pd.Categorical(['d','e','f'])
                    })
>>> df.describe()
 numeric
count 3.0
mean
        2.0
std
        1.0
min
        1.0
25%
        1.5
50%
         2.0
75%
         2.5
          3.0
max
```

Describing all columns of a DataFrame regardless of data type.

```
>>> df.describe(include='all')
     categorical numeric object
            3
                3.0
                      3
count
            3
                 NaN
                         3
unique
           f NaN
top
                        С
           1
                 NaN
                         1
freq
          NaN
                  2.0
                      NaN
mean
std
           NaN
                  1.0
                       NaN
min
           NaN
                  1.0
                       NaN
                      NaN
25%
           NaN
                  1.5
                      NaN
50%
           NaN
                  2.0
75%
                 2.5 NaN
           NaN
           NaN
                 3.0
max
                      NaN
```

Describing a column from a DataFrame by accessing it as an attribute.

```
>>> df.numeric.describe()
count 3.0
        2.0
mean
std
       1.0
        1.0
min
       1.5
25%
       2.0
50%
       2.5
75%
       3.0
Name: numeric, dtype: float64
```

Including only numeric columns in a DataFrame description.

```
>>> df.describe(include=[np.number])
     numeric
count 3.0
mean
         2.0
std
         1.0
         1.0
min
         1.5
25%
50%
         2.0
75%
         2.5
          3.0
max
```

Including only string columns in a DataFrame description.

Including only categorical columns from a DataFrame description.

Excluding numeric columns from a DataFrame description.

Excluding object columns from a DataFrame description.

```
>>> df.describe(exclude=[np.object])
      categorical numeric
             3
                    3.0
count
             3
                   NaN
unique
                   NaN
             f
top
             1
                   NaN
freq
          NaN
mean
                    2.0
std
            NaN
                     1.0
min
            NaN
                     1.0
25%
            NaN
                     1.5
50%
                     2.0
            NaN
75%
            NaN
                     2.5
                     3.0
max
             NaN
```

## div (other, axis=0)

Floating division of series and other, element-wise (binary operator  $\it truediv$ ). Equivalent to panel / other.

# **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

## Returns

**Panel** 

# See also:

Panel.rtruediv

#### divide(other, axis=0)

Floating division of series and other, element-wise (binary operator *truediv*). Equivalent to panel / other.

### **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

#### Returns

**Panel** 

## See also:

```
Panel.rtruediv
```

drop (labels=None, axis=0, index=None, columns=None, level=None, inplace=False, errors='raise')

```
dropna (axis=0, how='any', inplace=False)
```

Drop 2D from panel, holding passed axis constant

#### **Parameters**

axis [int, default 0] Axis to hold constant. E.g. axis=1 will drop major\_axis entries having a certain amount of NA data

**how** [{'all', 'any'}, default 'any'] 'any': one or more values are NA in the DataFrame along the axis. For 'all' they all must be.

**inplace** [bool, default False] If True, do operation inplace and return None.

### **Returns**

```
dropped [Panel]
```

## dtypes

Return the dtypes in the DataFrame.

This returns a Series with the data type of each column. The result's index is the original DataFrame's columns. Columns with mixed types are stored with the object dtype. See the User Guide for more.

### Returns

**pandas.Series** The data type of each column.

See also:

pandas.DataFrame.ftypes dtype and sparsity information.

# **Examples**

### empty

Indicator whether DataFrame is empty.

True if DataFrame is entirely empty (no items), meaning any of the axes are of length 0.

#### Returns

**bool** If DataFrame is empty, return True, if not return False.

#### See also:

```
pandas.Series.dropna, pandas.DataFrame.dropna
```

### **Notes**

If DataFrame contains only NaNs, it is still not considered empty. See the example below.

## **Examples**

An example of an actual empty DataFrame. Notice the index is empty:

```
>>> df_empty = pd.DataFrame({'A' : []})
>>> df_empty
Empty DataFrame
Columns: [A]
Index: []
>>> df_empty.empty
True
```

If we only have NaNs in our DataFrame, it is not considered empty! We will need to drop the NaNs to make the DataFrame empty:

```
>>> df = pd.DataFrame({'A' : [np.nan]})
>>> df
    A
0 NaN
>>> df.empty
False
>>> df.dropna().empty
True
```

eq(other, axis=None)

Wrapper for comparison method eq

```
equals (other)
```

Determines if two NDFrame objects contain the same elements. NaNs in the same location are considered equal.

```
ffill (axis=None, inplace=False, limit=None, downcast=None)
Synonym for DataFrame.fillna (method='ffill')
```

```
fillna (value=None, method=None, axis=None, inplace=False, limit=None, downcast=None, **kwargs)
```

Fill NA/NaÑ values using the specified method

#### **Parameters**

value [scalar, dict, Series, or DataFrame] Value to use to fill holes (e.g. 0), alternately a dict/Series/DataFrame of values specifying which value to use for each index (for a Series) or column (for a DataFrame). (values not in the dict/Series/DataFrame will not be filled). This value cannot be a list.

**method** [{'backfill', 'bfill', 'pad', 'ffill', None}, default None] Method to use for filling holes in reindexed Series pad / ffill: propagate last valid observation forward to next valid backfill / bfill: use NEXT valid observation to fill gap

```
axis [{0, 1, 2, 'items', 'major_axis', 'minor_axis'}]
```

**inplace** [boolean, default False] If True, fill in place. Note: this will modify any other views on this object, (e.g. a no-copy slice for a column in a DataFrame).

**limit** [int, default None] If method is specified, this is the maximum number of consecutive NaN values to forward/backward fill. In other words, if there is a gap with more than this number of consecutive NaNs, it will only be partially filled. If method is not specified, this is the maximum number of entries along the entire axis where NaNs will be filled. Must be greater than 0 if not None.

**downcast** [dict, default is None] a dict of item->dtype of what to downcast if possible, or the string 'infer' which will try to downcast to an appropriate equal type (e.g. float64 to int64 if possible)

## **Returns**

filled [Panel]

#### See also:

interpolate Fill NaN values using interpolation.

```
reindex, asfreq
```

## **Examples**

```
>>> df = pd.DataFrame([[np.nan, 2, np.nan, 0],
                        [3, 4, np.nan, 1],
                        [np.nan, np.nan, np.nan, 5],
. . .
                        [np.nan, 3, np.nan, 4]],
. . .
                        columns=list('ABCD'))
. . .
>>> df
            C D
    Α
          В
        2.0 NaN
  NaN
                 0
1
  3.0
       4.0 NaN
                 1
2
  NaN
       NaN NaN
  NaN
       3.0 NaN
```

Replace all NaN elements with 0s.

```
>>> df.fillna(0)

A B C D

0 0.0 2.0 0.0 0

1 3.0 4.0 0.0 1

2 0.0 0.0 0.0 5

3 0.0 3.0 0.0 4
```

We can also propagate non-null values forward or backward.

```
>>> df.fillna(method='ffill')

A B C D

0 NaN 2.0 NaN 0

1 3.0 4.0 NaN 1

2 3.0 4.0 NaN 5

3 3.0 3.0 NaN 4
```

Replace all NaN elements in column 'A', 'B', 'C', and 'D', with 0, 1, 2, and 3 respectively.

```
>>> values = {'A': 0, 'B': 1, 'C': 2, 'D': 3}
>>> df.fillna(value=values)

A B C D
0 0.0 2.0 2.0 0
1 3.0 4.0 2.0 1
2 0.0 1.0 2.0 5
3 0.0 3.0 2.0 4
```

## Only replace the first NaN element.

```
>>> df.fillna(value=values, limit=1)

A B C D

0 0.0 2.0 2.0 0

1 3.0 4.0 NaN 1

2 NaN 1.0 NaN 5

3 NaN 3.0 NaN 4
```

## filter(items=None, like=None, regex=None, axis=None)

Subset rows or columns of dataframe according to labels in the specified index.

Note that this routine does not filter a dataframe on its contents. The filter is applied to the labels of the index.

## **Parameters**

```
items [list-like] List of info axis to restrict to (must not all be present)
```

**like** [string] Keep info axis where "arg in col == True"

regex [string (regular expression)] Keep info axis with re.search(regex, col) == True

axis [int or string axis name] The axis to filter on. By default this is the info axis, 'index' for Series, 'columns' for DataFrame

## Returns

## same type as input object

### See also:

```
pandas.DataFrame.loc
```

### **Notes**

The items, like, and regex parameters are enforced to be mutually exclusive.

 ${\tt axis}$  defaults to the info axis that is used when indexing with  $[\,]\,.$ 

# **Examples**

```
>>> df
one two three
mouse 1 2 3
rabbit 4 5 6
```

```
>>> # select columns by name
>>> df.filter(items=['one', 'three'])
one three
mouse 1 3
rabbit 4 6
```

```
>>> # select columns by regular expression
>>> df.filter(regex='e$', axis=1)
one three
mouse 1 3
rabbit 4 6
```

```
>>> # select rows containing 'bbi'
>>> df.filter(like='bbi', axis=0)
one two three
rabbit 4 5 6
```

### first (offset)

Convenience method for subsetting initial periods of time series data based on a date offset.

## **Parameters**

offset [string, DateOffset, dateutil.relativedelta]

# Returns

subset [type of caller]

## **Raises**

TypeError If the index is not a DatetimeIndex

### See also:

last Select final periods of time series based on a date offset

at\_time Select values at a particular time of the day

between\_time Select values between particular times of the day

# **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='2D')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 1
2018-04-11 2
2018-04-13 3
2018-04-15 4
```

Get the rows for the first 3 days:

```
>>> ts.first('3D')

A
2018-04-09 1
2018-04-11 2
```

Notice the data for 3 first calender days were returned, not the first 3 days observed in the dataset, and therefore data for 2018-04-13 was not returned.

```
first_valid_index()
```

Return index for first non-NA/null value.

#### Returns

**scalar** [type of index]

### **Notes**

If all elements are non-NA/null, returns None. Also returns None for empty NDFrame.

## floordiv (other, axis=0)

Integer division of series and other, element-wise (binary operator *floordiv*). Equivalent to panel //other.

#### **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

## Returns

**Panel** 

### See also:

```
Panel.rfloordiv
```

classmethod fromDict (data, intersect=False, orient='items', dtype=None)

Construct Panel from dict of DataFrame objects

### **Parameters**

```
data [dict] {field : DataFrame}
```

intersect [boolean] Intersect indexes of input DataFrames

orient [{'items', 'minor'}, default 'items'] The "orientation" of the data. If the keys of the passed dict should be the items of the result panel, pass 'items' (default). Otherwise if the columns of the values of the passed DataFrame objects should be the items (which in the case of mixed-dtype data you should do), instead pass 'minor'

dtype [dtype, default None] Data type to force, otherwise infer

## **Returns**

Panel

classmethod from\_dict (data, intersect=False, orient='items', dtype=None)
Construct Panel from dict of DataFrame objects

**Parameters** 

```
data [dict] {field : DataFrame}
```

intersect [boolean] Intersect indexes of input DataFrames

**orient** [{'items', 'minor'}, default 'items'] The "orientation" of the data. If the keys of the passed dict should be the items of the result panel, pass 'items' (default). Otherwise if the columns of the values of the passed DataFrame objects should be the items (which in the case of mixed-dtype data you should do), instead pass 'minor'

**dtype** [dtype, default None] Data type to force, otherwise infer

### **Returns**

#### **Panel**

## ftypes

Return the ftypes (indication of sparse/dense and dtype) in DataFrame.

This returns a Series with the data type of each column. The result's index is the original DataFrame's columns. Columns with mixed types are stored with the object dtype. See the User Guide for more.

## Returns

pandas. Series The data type and indication of sparse/dense of each column.

#### See also:

```
pandas.DataFrame.dtypes Series with just dtype information.
```

pandas.SparseDataFrame Container for sparse tabular data.

### **Notes**

Sparse data should have the same dtypes as its dense representation.

# **Examples**

```
>>> import numpy as np
>>> arr = np.random.RandomState(0).randn(100, 4)
>>> arr[arr < .8] = np.nan
>>> pd.DataFrame(arr).ftypes
0    float64:dense
1    float64:dense
2    float64:dense
3    float64:dense
dtype: object
```

```
>>> pd.SparseDataFrame(arr).ftypes
0    float64:sparse
1    float64:sparse
2    float64:sparse
3    float64:sparse
dtype: object
```

## **ge** (*other*, *axis=None*)

Wrapper for comparison method ge

## get (key, default=None)

Get item from object for given key (DataFrame column, Panel slice, etc.). Returns default value if not found.

## **Parameters**

key [object]

#### Returns

value [type of items contained in object]

## get\_dtype\_counts()

Return counts of unique dtypes in this object.

### **Returns**

**dtype** [Series] Series with the count of columns with each dtype.

### See also:

dtypes Return the dtypes in this object.

# **Examples**

```
>>> a = [['a', 1, 1.0], ['b', 2, 2.0], ['c', 3, 3.0]]
>>> df = pd.DataFrame(a, columns=['str', 'int', 'float'])
>>> df
    str int float
0 a 1 1.0
1 b 2 2.0
2 c 3 3.0
```

```
>>> df.get_dtype_counts()
float64    1
int64    1
object    1
dtype: int64
```

## get\_ftype\_counts()

Return counts of unique ftypes in this object.

Deprecated since version 0.23.0.

This is useful for SparseDataFrame or for DataFrames containing sparse arrays.

### Returns

**dtype** [Series] Series with the count of columns with each type and sparsity (dense/sparse)

# See also:

ftypes Return ftypes (indication of sparse/dense and dtype) in this object.

# **Examples**

```
>>> a = [['a', 1, 1.0], ['b', 2, 2.0], ['c', 3, 3.0]]
>>> df = pd.DataFrame(a, columns=['str', 'int', 'float'])
>>> df
str int float
0 a 1 1.0
1 b 2 2.0
2 c 3 3.0
```

```
>>> df.get_ftype_counts()
float64:dense 1
int64:dense 1
object:dense 1
dtype: int64
```

### get\_value (\*args, \*\*kwargs)

Quickly retrieve single value at (item, major, minor) location

Deprecated since version 0.21.0.

Please use .at[] or .iat[] accessors.

### **Parameters**

```
item [item label (panel item)]major [major axis label (panel item row)]minor [minor axis label (panel item column)]
```

takeable [interpret the passed labels as indexers, default False]

## **Returns**

value [scalar value]

### get\_values()

Return an ndarray after converting sparse values to dense.

This is the same as .values for non-sparse data. For sparse data contained in a *pandas.SparseArray*, the data are first converted to a dense representation.

### **Returns**

numpy.ndarray Numpy representation of DataFrame

# See also:

values Numpy representation of DataFrame.

pandas.SparseArray Container for sparse data.

# **Examples**

```
>>> df = pd.DataFrame({'a': [1, 2], 'b': [True, False],
... 'c': [1.0, 2.0]})
>>> df
a b c
0 1 True 1.0
1 2 False 2.0
```

```
>>> df.get_values()
array([[1, True, 1.0], [2, False, 2.0]], dtype=object)
```

```
>>> df = pd.DataFrame({"a": pd.SparseArray([1, None, None]),
... "c": [1.0, 2.0, 3.0]})
>>> df
    a    c
0 1.0 1.0
1 NaN 2.0
2 NaN 3.0
```

## groupby (function, axis='major')

Group data on given axis, returning GroupBy object

#### **Parameters**

```
function [callable] Mapping function for chosen access axis [{'major', 'minor', 'items'}, default 'major']
```

### Returns

grouped [PanelGroupBy]

## gt (other, axis=None)

Wrapper for comparison method gt

## head(n=5)

Return the first *n* rows.

This function returns the first n rows for the object based on position. It is useful for quickly testing if your object has the right type of data in it.

### **Parameters**

**n** [int, default 5] Number of rows to select.

# Returns

**obj\_head** [type of caller] The first *n* rows of the caller object.

See also:

pandas.DataFrame.tail Returns the last n rows.

# **Examples**

(continues on next page)

```
3 lion
4 monkey
5 parrot
6 shark
7 whale
8 zebra
```

# Viewing the first 5 lines

```
>>> df.head()
    animal
0 alligator
1 bee
2 falcon
3 lion
4 monkey
```

### Viewing the first *n* lines (three in this case)

```
>>> df.head(3)
    animal
0 alligator
1 bee
2 falcon
```

#### iat

Access a single value for a row/column pair by integer position.

Similar to iloc, in that both provide integer-based lookups. Use iat if you only need to get or set a single value in a DataFrame or Series.

## Raises

**IndexError** When integer position is out of bounds

See also:

DataFrame.at Access a single value for a row/column label pair

**DataFrame.loc** Access a group of rows and columns by label(s)

**DataFrame.iloc** Access a group of rows and columns by integer position(s)

## **Examples**

```
>>> df = pd.DataFrame([[0, 2, 3], [0, 4, 1], [10, 20, 30]],
... columns=['A', 'B', 'C'])
>>> df
    A     B     C
0    0    2    3
1    0    4    1
2    10    20    30
```

Get value at specified row/column pair

```
>>> df.iat[1, 2]
1
```

Set value at specified row/column pair

```
>>> df.iat[1, 2] = 10
>>> df.iat[1, 2]
10
```

Get value within a series

```
>>> df.loc[0].iat[1]
2
```

#### iloc

Purely integer-location based indexing for selection by position.

.iloc[] is primarily integer position based (from 0 to length-1 of the axis), but may also be used with a boolean array.

Allowed inputs are:

- An integer, e.g. 5.
- A list or array of integers, e.g. [4, 3, 0].
- A slice object with ints, e.g. 1:7.
- · A boolean array.
- A callable function with one argument (the calling Series, DataFrame or Panel) and that returns valid output for indexing (one of the above)

.iloc will raise IndexError if a requested indexer is out-of-bounds, except *slice* indexers which allow out-of-bounds indexing (this conforms with python/numpy *slice* semantics).

See more at Selection by Position

# infer\_objects()

Attempt to infer better dtypes for object columns.

Attempts soft conversion of object-dtyped columns, leaving non-object and unconvertible columns unchanged. The inference rules are the same as during normal Series/DataFrame construction.

New in version 0.21.0.

### Returns

**converted** [same type as input object]

See also:

```
pandas.to_datetime Convert argument to datetime.
```

pandas.to\_timedelta Convert argument to timedelta.

pandas.to numeric Convert argument to numeric typeR

# **Examples**

```
>>> df = pd.DataFrame({"A": ["a", 1, 2, 3]})
>>> df = df.iloc[1:]
>>> df
A
```

(continues on next page)

```
1 1
2 2
3 3
```

```
>>> df.dtypes
A object
dtype: object
```

```
>>> df.infer_objects().dtypes
A int64
dtype: object
```

Please note that only method='linear' is supported for DataFrames/Series with a MultiIndex.

### **Parameters**

- 'linear': ignore the index and treat the values as equally spaced. This is the only method supported on MultiIndexes. default
- 'time': interpolation works on daily and higher resolution data to interpolate given length of interval
- 'index', 'values': use the actual numerical values of the index
- 'nearest', 'zero', 'slinear', 'quadratic', 'cubic', 'barycentric', 'polynomial' is passed to scipy.interpolate.interpld. Both 'polynomial' and 'spline' require that you also specify an *order* (int), e.g. df.interpolate(method='polynomial', order=4). These use the actual numerical values of the index.
- 'krogh', 'piecewise\_polynomial', 'spline', 'pchip' and 'akima' are all wrappers around the scipy interpolation methods of similar names. These use the actual numerical values of the index. For more information on their behavior, see the scipy documentation and tutorial documentation
- 'from\_derivatives' refers to BPoly.from\_derivatives which replaces 'piece-wise\_polynomial' interpolation method in scipy 0.18

New in version 0.18.1: Added support for the 'akima' method Added interpolate method 'from\_derivatives' which replaces 'piecewise\_polynomial' in scipy 0.18; backwards-compatible with scipy < 0.18

```
axis [{0, 1}, default 0]
```

- 0: fill column-by-column
- 1: fill row-by-row

**limit** [int, default None.] Maximum number of consecutive NaNs to fill. Must be greater than 0.

limit\_direction [{'forward', 'backward', 'both'}, default 'forward']

limit\_area [{'inside', 'outside'}, default None]

- None: (default) no fill restriction
- 'inside' Only fill NaNs surrounded by valid values (interpolate).
- 'outside' Only fill NaNs outside valid values (extrapolate).

If limit is specified, consecutive NaNs will be filled in this direction.

New in version 0.21.0.

inplace [bool, default False] Update the NDFrame in place if possible.

downcast [optional, 'infer' or None, defaults to None] Downcast dtypes if possible.

**kwargs** [keyword arguments to pass on to the interpolating function.]

### Returns

Series or DataFrame of same shape interpolated at the NaNs

#### See also:

```
reindex, replace, fillna
```

# **Examples**

## Filling in NaNs

### is\_copy

# isna()

Detect missing values.

Return a boolean same-sized object indicating if the values are NA. NA values, such as None or numpy. NaN, gets mapped to True values. Everything else gets mapped to False values. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True).

## Returns

**NDFrame** Mask of bool values for each element in NDFrame that indicates whether an element is not an NA value.

### See also:

```
NDFrame.isnull alias of isna
```

NDFrame.notna boolean inverse of isna

NDFrame.dropna omit axes labels with missing values

isna top-level isna

## **Examples**

Show which entries in a DataFrame are NA.

```
>>> df = pd.DataFrame({'age': [5, 6, np.NaN],
                      'born': [pd.NaT, pd.Timestamp('1939-05-27'),
                               pd.Timestamp('1940-04-25')],
. . .
                      'name': ['Alfred', 'Batman', ''],
. . .
                      'toy': [None, 'Batmobile', 'Joker'] })
. . .
>>> df
           born name
  age
                                tov
            NaT Alfred
0 5.0
                              None
 6.0 1939-05-27 Batman Batmobile
 NaN 1940-04-25
                              Joker
```

```
>>> df.isna()
   age born name toy
0 False True False True
1 False False False False
2 True False False False
```

Show which entries in a Series are NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.isna()
0 False
1 False
2 True
dtype: bool
```

# isnull()

Detect missing values.

Return a boolean same-sized object indicating if the values are NA. NA values, such as None or numpy. NaN, gets mapped to True values. Everything else gets mapped to False values. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True).

# Returns

**NDFrame** Mask of bool values for each element in NDFrame that indicates whether an element is not an NA value.

See also:

```
NDFrame.isnull alias of isna
```

NDFrame.notna boolean inverse of isna

NDFrame.dropna omit axes labels with missing values

isna top-level isna

## **Examples**

Show which entries in a DataFrame are NA.

```
>>> df = pd.DataFrame({'age': [5, 6, np.NaN],
                      'born': [pd.NaT, pd.Timestamp('1939-05-27'),
                               pd.Timestamp('1940-04-25')],
. . .
                       'name': ['Alfred', 'Batman', ''],
. . .
                       'toy': [None, 'Batmobile', 'Joker']})
. . .
>>> df
  age
           born name
                                tov
  5.0
            NaT Alfred
                               None
  6.0 1939-05-27 Batman Batmobile
 NaN 1940-04-25
                              Joker
```

```
>>> df.isna()
   age born name toy
0 False True False True
1 False False False False
2 True False False False
```

Show which entries in a Series are NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.isna()
0 False
1 False
2 True
dtype: bool
```

## items

# iteritems()

Iterate over (label, values) on info axis

This is index for Series, columns for DataFrame, major\_axis for Panel, and so on.

ix

A primarily label-location based indexer, with integer position fallback.

Warning: Starting in 0.20.0, the .ix indexer is deprecated, in favor of the more strict .iloc and .loc indexers.

.ix[] supports mixed integer and label based access. It is primarily label based, but will fall back to integer positional access unless the corresponding axis is of integer type.

.ix is the most general indexer and will support any of the inputs in .loc and .iloc. .ix also supports floating point label schemes. .ix is exceptionally useful when dealing with mixed positional and label based hierarchical indexes.

However, when an axis is integer based, ONLY label based access and not positional access is supported. Thus, in such cases, it's usually better to be explicit and use .iloc or .loc.

See more at Advanced Indexing.

```
join (other, how='left', lsuffix=", rsuffix=")
```

Join items with other Panel either on major and minor axes column

### **Parameters**

**other** [Panel or list of Panels] Index should be similar to one of the columns in this one

how [{'left', 'right', 'outer', 'inner'}] How to handle indexes of the two objects. Default: 'left' for joining on index, None otherwise \* left: use calling frame's index \* right: use input frame's index \* outer: form union of indexes \* inner: use intersection of indexes

**Isuffix** [string] Suffix to use from left frame's overlapping columns

**rsuffix** [string] Suffix to use from right frame's overlapping columns

#### Returns

joined [Panel]

## keys()

Get the 'info axis' (see Indexing for more)

This is index for Series, columns for DataFrame and major\_axis for Panel.

kurt (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0). Normalized by N-1

#### **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

**kurt** [DataFrame or Panel (if level specified)]

kurtosis (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0). Normalized by N-1

## **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

## **Returns**

**kurt** [DataFrame or Panel (if level specified)]

### last (offset)

Convenience method for subsetting final periods of time series data based on a date offset.

### **Parameters**

offset [string, DateOffset, dateutil.relativedelta]

### Returns

subset [type of caller]

## **Raises**

TypeError If the index is not a DatetimeIndex

### See also:

first Select initial periods of time series based on a date offset

at\_time Select values at a particular time of the day

between\_time Select values between particular times of the day

# **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='2D')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 1
2018-04-11 2
2018-04-13 3
2018-04-15 4
```

## Get the rows for the last 3 days:

```
>>> ts.last('3D')

A
2018-04-13 3
2018-04-15 4
```

Notice the data for 3 last calender days were returned, not the last 3 observed days in the dataset, and therefore data for 2018-04-11 was not returned.

# last\_valid\_index()

Return index for last non-NA/null value.

### Returns

**scalar** [type of index]

### **Notes**

If all elements are non-NA/null, returns None. Also returns None for empty NDFrame.

## le (other, axis=None)

Wrapper for comparison method le

#### loc

Access a group of rows and columns by label(s) or a boolean array.

.loc[] is primarily label based, but may also be used with a boolean array.

Allowed inputs are:

- A single label, e.g. 5 or 'a', (note that 5 is interpreted as a *label* of the index, and **never** as an integer position along the index).
- A list or array of labels, e.g. ['a', 'b', 'c'].
- A slice object with labels, e.g. 'a': 'f'.

Warning: Note that contrary to usual python slices, both the start and the stop are included

- A boolean array of the same length as the axis being sliced, e.g. [True, False, True].
- A callable function with one argument (the calling Series, DataFrame or Panel) and that returns valid output for indexing (one of the above)

See more at Selection by Label

### Raises

**KeyError:** when any items are not found

See also:

DataFrame.at Access a single value for a row/column label pair

**DataFrame.iloc** Access group of rows and columns by integer position(s)

**DataFrame.xs** Returns a cross-section (row(s) or column(s)) from the Series/DataFrame.

Series.loc Access group of values using labels

## **Examples**

## **Getting values**

Single label. Note this returns the row as a Series.

```
>>> df.loc['viper']
max_speed 4
shield 5
Name: viper, dtype: int64
```

List of labels. Note using [ [ ] ] returns a DataFrame.

```
>>> df.loc[['viper', 'sidewinder']]

max_speed shield

viper 4 5

sidewinder 7 8
```

### Single label for row and column

```
>>> df.loc['cobra', 'shield']
2
```

Slice with labels for row and single label for column. As mentioned above, note that both the start and stop of the slice are included.

```
>>> df.loc['cobra':'viper', 'max_speed']
cobra 1
viper 4
Name: max_speed, dtype: int64
```

### Boolean list with the same length as the row axis

```
>>> df.loc[[False, False, True]]

max_speed shield
sidewinder 7 8
```

#### Conditional that returns a boolean Series

# Conditional that returns a boolean Series with column labels specified

### Callable that returns a boolean Series

```
>>> df.loc[lambda df: df['shield'] == 8]

max_speed shield
sidewinder 7 8
```

# **Setting values**

Set value for all items matching the list of labels

# Set value for an entire row

```
>>> df.loc['cobra'] = 10
>>> df
```

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```
max_speed shield
cobra 10 10
viper 4 50
sidewinder 7 50
```

Set value for an entire column

Set value for rows matching callable condition

## Getting values on a DataFrame with an index that has integer labels

Another example using integers for the index

Slice with integer labels for rows. As mentioned above, note that both the start and stop of the slice are included.

# Getting values with a MultiIndex

A number of examples using a DataFrame with a MultiIndex

```
>>> tuples = [
... ('cobra', 'mark i'), ('cobra', 'mark ii'),
... ('sidewinder', 'mark i'), ('sidewinder', 'mark ii'),
... ('viper', 'mark ii'), ('viper', 'mark iii')
... ]
>>> index = pd.MultiIndex.from_tuples(tuples)
>>> values = [[12, 2], [0, 4], [10, 20],
... [1, 4], [7, 1], [16, 36]]
>>> df = pd.DataFrame(values, columns=['max_speed', 'shield'], index=index)
```

(continues on next page)

```
>>> df
                  max_speed shield
                  12
cobra
         mark i
                        0
                                 4
         mark ii
                       10
                                20
sidewinder mark i
         mark ii
                         1
                                4
         mark ii
                         7
                                 1
viper
         mark iii
                        16
```

Single label. Note this returns a DataFrame with a single index.

```
>>> df.loc['cobra']

max_speed shield

mark i 12 2

mark ii 0 4
```

Single index tuple. Note this returns a Series.

```
>>> df.loc[('cobra', 'mark ii')]
max_speed 0
shield 4
Name: (cobra, mark ii), dtype: int64
```

Single label for row and column. Similar to passing in a tuple, this returns a Series.

```
>>> df.loc['cobra', 'mark i']
max_speed 12
shield 2
Name: (cobra, mark i), dtype: int64
```

Single tuple. Note using [ [ ] ] returns a DataFrame.

```
>>> df.loc[[('cobra', 'mark ii')]]

max_speed shield
cobra mark ii 0 4
```

Single tuple for the index with a single label for the column

```
>>> df.loc[('cobra', 'mark i'), 'shield']
2
```

Slice from index tuple to single label

```
>>> df.loc[('cobra', 'mark i'):'viper']
                 max_speed shield
                                2
cobra
         mark i
                        12
         mark ii
                         0
                                  4
sidewinder mark i
                        10
                                20
                        1
         mark ii
                                 4
         mark ii
                         7
                                 1
viper
                        16
         mark iii
```

Slice from index tuple to index tuple

```
>>> df.loc[('cobra', 'mark i'):('viper', 'mark ii')]

max_speed shield
```

(continues on next page)

	cobra	mark	i	12	2
		mark	ii	0	4
	sidewinder	mark	i	10	20
		mark	ii	1	4
	viper	mark	ii	7	1
- 1					

### **lt** (*other*, *axis=None*)

Wrapper for comparison method lt

mad (axis=None, skipna=None, level=None)

Return the mean absolute deviation of the values for the requested axis

#### **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

mad [DataFrame or Panel (if level specified)]

### major\_axis

major\_xs(key)

Return slice of panel along major axis

## **Parameters**

key [object] Major axis label

### Returns

y [DataFrame] index -> minor axis, columns -> items

### **Notes**

major\_xs is only for getting, not setting values.

MultiIndex Slicers is a generic way to get/set values on any level or levels and is a superset of major\_xs functionality, see MultiIndex Slicers

mask (cond, other=nan, inplace=False, axis=None, level=None, errors='raise', try\_cast=False,
 raise\_on\_error=None)

Return an object of same shape as self and whose corresponding entries are from self where *cond* is False and otherwise are from *other*.

### **Parameters**

**cond** [boolean NDFrame, array-like, or callable] Where *cond* is False, keep the original value. Where True, replace with corresponding value from *other*. If *cond* is callable, it is computed on the NDFrame and should return boolean NDFrame or array. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as cond.

**other** [scalar, NDFrame, or callable] Entries where *cond* is True are replaced with corresponding value from *other*. If other is callable, it is computed on the NDFrame and should return scalar or NDFrame. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as other.

**inplace** [boolean, default False] Whether to perform the operation in place on the data

axis [alignment axis if needed, default None]

level [alignment level if needed, default None]

errors [str, {'raise', 'ignore'}, default 'raise']

- raise: allow exceptions to be raised
- ignore: suppress exceptions. On error return original object

Note that currently this parameter won't affect the results and will always coerce to a suitable dtype.

try\_cast [boolean, default False] try to cast the result back to the input type (if possible),

raise\_on\_error [boolean, default True] Whether to raise on invalid data types (e.g. trying to where on strings)

Deprecated since version 0.21.0.

#### Returns

wh [same type as caller]

### See also:

DataFrame.where()

# **Notes**

The mask method is an application of the if-then idiom. For each element in the calling DataFrame, if cond is False the element is used; otherwise the corresponding element from the DataFrame other is used.

The signature for DataFrame.where() differs from numpy.where(). Roughly df1.where(m, df2) is equivalent to np.where(m, df1, df2).

For further details and examples see the mask documentation in indexing.

# **Examples**

```
>>> s = pd.Series(range(5))
>>> s.where(s > 0)
0     NaN
1     1.0
2     2.0
3     3.0
4     4.0
```

```
>>> df = pd.DataFrame(np.arange(10).reshape(-1, 2), columns=['A', 'B'])
>>> m = df % 3 == 0
>>> df.where(m, -df)
  A B
  0 -1
1 -2 3
2 -4 -5
3 6 -7
4 - 8 9
>>> df.where(m, -df) == np.where(m, df, -df)
      Α
  True
0
        True
1
  True True
  True
        True
  True
        True
  True
        True
\rightarrow \rightarrow df.where(m, -df) == df.mask(~m, -df)
      Α
  True True
0
  True True
1
  True True
3
  True True
  True True
```

max (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

This method returns the maximum of the values in the object. If you want the *index* of the maximum, use idxmax. This is the equivalent of the numpy.ndarray method argmax.

## **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

# Returns

max [DataFrame or Panel (if level specified)]

**mean** (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)
Return the mean of the values for the requested axis

#### **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

#### Returns

mean [DataFrame or Panel (if level specified)]

median (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)
Return the median of the values for the requested axis

#### **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

## Returns

```
median [DataFrame or Panel (if level specified)]
```

min (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

This method returns the minimum of the values in the object. If you want the *index* of the minimum, use idxmin. This is the equivalent of the numpy.ndarray method argmin.

### **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

min [DataFrame or Panel (if level specified)]

### minor\_axis

minor xs (key)

Return slice of panel along minor axis

## **Parameters**

```
key [object] Minor axis label
```

### Returns

```
y [DataFrame] index -> major axis, columns -> items
```

### **Notes**

minor\_xs is only for getting, not setting values.

MultiIndex Slicers is a generic way to get/set values on any level or levels and is a superset of minor\_xs functionality, see MultiIndex Slicers

```
mod(other, axis=0)
```

Modulo of series and other, element-wise (binary operator *mod*). Equivalent to panel % other.

### **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

## **Returns**

**Panel** 

### See also:

```
Panel.rmod
```

```
mul (other, axis=0)
```

Multiplication of series and other, element-wise (binary operator mul). Equivalent to panel \* other.

## **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

## Returns

Panel

# See also:

```
Panel.rmul
```

```
multiply (other, axis=0)
```

Multiplication of series and other, element-wise (binary operator mul). Equivalent to panel \* other.

## **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

### **Returns**

**Panel** 

# See also:

Panel.rmul

#### ndim

Return an int representing the number of axes / array dimensions.

Return 1 if Series. Otherwise return 2 if DataFrame.

#### See also:

ndarray.ndim

# **Examples**

```
>>> s = pd.Series({'a': 1, 'b': 2, 'c': 3})
>>> s.ndim
1
```

```
>>> df = pd.DataFrame({'col1': [1, 2], 'col2': [3, 4]})
>>> df.ndim
2
```

ne (other, axis=None)

Wrapper for comparison method ne

#### notna()

Detect existing (non-missing) values.

Return a boolean same-sized object indicating if the values are not NA. Non-missing values get mapped to True. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True). NA values, such as None or numpy.NaN, get mapped to False values.

## **Returns**

**NDFrame** Mask of bool values for each element in NDFrame that indicates whether an element is not an NA value.

See also:

NDFrame.notnull alias of notna

NDFrame.isna boolean inverse of notna

NDFrame.dropna omit axes labels with missing values

notna top-level notna

## **Examples**

Show which entries in a DataFrame are not NA.

```
>>> df = pd.DataFrame({ 'age': [5, 6, np.NaN],
                        'born': [pd.NaT, pd.Timestamp('1939-05-27'),
                                 pd.Timestamp('1940-04-25')],
. . .
                        'name': ['Alfred', 'Batman', ''],
. . .
                        'toy': [None, 'Batmobile', 'Joker']})
. . .
>>> df
             born
                                   toy
  age
                      name
  5.0
              NaT Alfred
                                  None
```

(continues on next page)

```
1 6.0 1939-05-27 Batman Batmobile
2 NaN 1940-04-25 Joker
```

```
>>> df.notna()
age born name toy
0 True False True False
1 True True True True
2 False True True True
```

Show which entries in a Series are not NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.notna()
0 True
1 True
2 False
dtype: bool
```

# notnull()

Detect existing (non-missing) values.

Return a boolean same-sized object indicating if the values are not NA. Non-missing values get mapped to True. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True). NA values, such as None or numpy.NaN, get mapped to False values.

## Returns

**NDFrame** Mask of bool values for each element in NDFrame that indicates whether an element is not an NA value.

See also:

NDFrame.notnull alias of notna

NDFrame.isna boolean inverse of notna

NDFrame.dropna omit axes labels with missing values

notna top-level notna

## **Examples**

Show which entries in a DataFrame are not NA.

```
>>> df = pd.DataFrame({'age': [5, 6, np.NaN],
... 'born': [pd.NaT, pd.Timestamp('1939-05-27'),
pd.Timestamp('1940-04-25')],
... 'name': ['Alfred', 'Batman', ''],
... 'toy': [None, 'Batmobile', 'Joker']})
```

(continues on next page)

```
>>> df.notna()
age born name toy
0 True False True False
1 True True True
2 False True True
```

Show which entries in a Series are not NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.notna()
0 True
1 True
2 False
dtype: bool
```

pct\_change (periods=1, fill\_method='pad', limit=None, freq=None, \*\*kwargs)

Percentage change between the current and a prior element.

Computes the percentage change from the immediately previous row by default. This is useful in comparing the percentage of change in a time series of elements.

## **Parameters**

periods [int, default 1] Periods to shift for forming percent change.

fill\_method [str, default 'pad'] How to handle NAs before computing percent changes.

**limit** [int, default None] The number of consecutive NAs to fill before stopping.

**freq** [DateOffset, timedelta, or offset alias string, optional] Increment to use from time series API (e.g. 'M' or BDay()).

\*\*kwargs Additional keyword arguments are passed into DataFrame.shift or Series.shift.

## Returns

**chg** [Series or DataFrame] The same type as the calling object.

# See also:

**Series.diff** Compute the difference of two elements in a Series.

**DataFrame.diff** Compute the difference of two elements in a DataFrame.

Series.shift Shift the index by some number of periods.

DataFrame.shift Shift the index by some number of periods.

## **Examples**

### Series

```
>>> s = pd.Series([90, 91, 85])

>>> s

0  90

1  91

2  85

dtype: int64
```

See the percentage change in a Series where filling NAs with last valid observation forward to next valid.

```
>>> s = pd.Series([90, 91, None, 85])
>>> s
0 90.0
1 91.0
2 NaN
3 85.0
dtype: float64
```

## **DataFrame**

Percentage change in French franc, Deutsche Mark, and Italian lira from 1980-01-01 to 1980-03-01.

```
>>> df = pd.DataFrame({
... 'FR': [4.0405, 4.0963, 4.3149],
... 'GR': [1.7246, 1.7482, 1.8519],
... 'IT': [804.74, 810.01, 860.13]},
... index=['1980-01-01', '1980-02-01', '1980-03-01'])
>>> df

FR GR IT

1980-01-01 4.0405 1.7246 804.74

1980-02-01 4.0963 1.7482 810.01

1980-03-01 4.3149 1.8519 860.13
```

```
>>> df.pct_change()

FR GR IT

1980-01-01 NaN NaN NaN

1980-02-01 0.013810 0.013684 0.006549

1980-03-01 0.053365 0.059318 0.061876
```

Percentage of change in GOOG and APPL stock volume. Shows computing the percentage change between columns.

### **Parameters**

func [function] function to apply to the NDFrame. args, and kwargs are
 passed into func. Alternatively a (callable, data\_keyword) tuple where
 data\_keyword is a string indicating the keyword of callable that expects the
 NDFrame.

args [iterable, optional] positional arguments passed into func.

**kwargs** [mapping, optional] a dictionary of keyword arguments passed into func.

## Returns

object [the return type of func.]

### See also:

pandas.DataFrame.apply, pandas.DataFrame.applymap, pandas.Series.map

### **Notes**

Use .pipe when chaining together functions that expect Series, DataFrames or GroupBy objects. Instead of writing

```
>>> f(g(h(df), arg1=a), arg2=b, arg3=c)
```

# You can write

```
>>> (df.pipe(h) ... .pipe(g, arg1=a)
```

(continues on next page)

```
... .pipe(f, arg2=b, arg3=c)
... )
```

If you have a function that takes the data as (say) the second argument, pass a tuple indicating which keyword expects the data. For example, suppose f takes its data as arg2:

```
>>> (df.pipe(h)
... .pipe(g, arg1=a)
... .pipe((f, 'arg2'), arg1=a, arg3=c)
... )
```

## pop (item)

Return item and drop from frame. Raise KeyError if not found.

### **Parameters**

item [str] Column label to be popped

#### Returns

popped [Series]

## **Examples**

```
>>> df = pd.DataFrame([('falcon', 'bird', 389.0),
                       ('parrot', 'bird',
                                              24.0),
                       ('lion', 'mammal', 80.5), ('monkey', 'mammal', np.nan)],
. . .
. . .
                      columns=('name', 'class', 'max_speed'))
. . .
>>> df
    name class max_speed
          bird
  falcon
                   389.0
1
  parrot
            bird
                       24.0
2
   lion mammal
                       80.5
3 monkey mammal
                       NaN
```

```
>>> df.pop('class')
0 bird
1 bird
2 mammal
3 mammal
Name: class, dtype: object
```

## pow (other, axis=0)

Exponential power of series and other, element-wise (binary operator pow). Equivalent to panel \*\* other.

## **Parameters**

other [DataFrame or Panel]

axis [{items, major\_axis, minor\_axis}] Axis to broadcast over

#### Returns

**Panel** 

## See also:

Panel.rpow

prod (axis=None, skipna=None, level=None, numeric\_only=None, min\_count=0, \*\*kwargs)
Return the product of the values for the requested axis

### **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

min\_count [int, default 0] The required number of valid values to perform the operation. If fewer than min\_count non-NA values are present the result will be NA.

New in version 0.22.0: Added with the default being 0. This means the sum of an all-NA or empty Series is 0, and the product of an all-NA or empty Series is 1.

### Returns

prod [DataFrame or Panel (if level specified)]

### **Examples**

By default, the product of an empty or all-NA Series is 1

```
>>> pd.Series([]).prod()
1.0
```

This can be controlled with the min\_count parameter

```
>>> pd.Series([]).prod(min_count=1)
nan
```

Thanks to the skipna parameter, min\_count handles all-NA and empty series identically.

```
>>> pd.Series([np.nan]).prod()
1.0
```

```
>>> pd.Series([np.nan]).prod(min_count=1)
nan
```

**product** (axis=None, skipna=None, level=None, numeric\_only=None, min\_count=0, \*\*kwargs)

Return the product of the values for the requested axis

### **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

**min\_count** [int, default 0] The required number of valid values to perform the operation. If fewer than min\_count non-NA values are present the result will be NA.

New in version 0.22.0: Added with the default being 0. This means the sum of an all-NA or empty Series is 0, and the product of an all-NA or empty Series is 1.

#### Returns

prod [DataFrame or Panel (if level specified)]

# **Examples**

By default, the product of an empty or all-NA Series is 1

```
>>> pd.Series([]).prod()
1.0
```

This can be controlled with the min\_count parameter

```
>>> pd.Series([]).prod(min_count=1)
nan
```

Thanks to the skipna parameter, min\_count handles all-NA and empty series identically.

```
>>> pd.Series([np.nan]).prod()
1.0
```

```
>>> pd.Series([np.nan]).prod(min_count=1)
nan
```

radd (other, axis=0)

Addition of series and other, element-wise (binary operator radd). Equivalent to other + panel.

# **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

# Returns

Panel

### See also:

Panel.add

Compute numerical data ranks (1 through n) along axis. Equal values are assigned a rank that is the average of the ranks of those values

# **Parameters**

```
axis [{0 or 'index', 1 or 'columns'}, default 0] index to direct ranking
method [{'average', 'min', 'max', 'first', 'dense'}]
```

- average: average rank of group
- min: lowest rank in group
- max: highest rank in group
- first: ranks assigned in order they appear in the array
- dense: like 'min', but rank always increases by 1 between groups

numeric\_only [boolean, default None] Include only float, int, boolean data. Valid only
for DataFrame or Panel objects

```
na_option [{'keep', 'top', 'bottom'}]
```

- keep: leave NA values where they are
- top: smallest rank if ascending
- · bottom: smallest rank if descending

**ascending** [boolean, default True] False for ranks by high (1) to low (N)

pct [boolean, default False] Computes percentage rank of data

#### Returns

ranks [same type as caller]

# rdiv (other, axis=0)

Floating division of series and other, element-wise (binary operator *rtruediv*). Equivalent to other / panel.

## **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

## Returns

Panel

### See also:

Panel.truediv

# reindex(\*args, \*\*kwargs)

Conform Panel to new index with optional filling logic, placing NA/NaN in locations having no value in the previous index. A new object is produced unless the new index is equivalent to the current one and copy=False

## **Parameters**

items, major\_axis, minor\_axis [array-like, optional (should be specified using keywords)] New labels / index to conform to. Preferably an Index object to avoid duplicating data

method [{None, 'backfill'/'bfill', 'pad'/'ffill', 'nearest'}, optional] method to use for filling holes in reindexed DataFrame. Please note: this is only applicable to DataFrames/Series with a monotonically increasing/decreasing index.

• default: don't fill gaps

- pad / ffill: propagate last valid observation forward to next valid
- backfill / bfill: use next valid observation to fill gap
- nearest: use nearest valid observations to fill gap

copy [boolean, default True] Return a new object, even if the passed indexes are the same

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

**fill\_value** [scalar, default np.NaN] Value to use for missing values. Defaults to NaN, but can be any "compatible" value

**limit** [int, default None] Maximum number of consecutive elements to forward or backward fill

tolerance [optional] Maximum distance between original and new labels for inexact
matches. The values of the index at the matching locations most satisfy the equation abs (index[indexer] - target) <= tolerance.</pre>

Tolerance may be a scalar value, which applies the same tolerance to all values, or list-like, which applies variable tolerance per element. List-like includes list, tuple, array, Series, and must be the same size as the index and its dtype must exactly match the index's type.

New in version 0.21.0: (list-like tolerance)

#### Returns

reindexed [Panel]

## **Examples**

DataFrame.reindex supports two calling conventions

- (index=index\_labels, columns=column\_labels, ...)
- (labels, axis={'index', 'columns'}, ...)

We *highly* recommend using keyword arguments to clarify your intent.

Create a dataframe with some fictional data.

```
>>> index = ['Firefox', 'Chrome', 'Safari', 'IE10', 'Konqueror']
>>> df = pd.DataFrame({
         'http_status': [200,200,404,404,301],
         'response_time': [0.04, 0.02, 0.07, 0.08, 1.0]},
. . .
         index=index)
. . .
>>> df
          http_status response_time
Firefox
                   200
                                  0.04
                   200
                                  0.02
Chrome
Safari
                    404
                                  0.07
IE10
                    404
                                  0.08
Konqueror
                    301
                                  1.00
```

Create a new index and reindex the dataframe. By default values in the new index that do not have corresponding records in the dataframe are assigned NaN.

```
>>> new_index= ['Safari', 'Iceweasel', 'Comodo Dragon', 'IE10',
               'Chrome']
. . .
>>> df.reindex(new_index)
            http_status response_time
                   404.0 0.07
Safari
Iceweasel
                    NaN
                                   NaN
Comodo Dragon
                    NaN
                                   NaN
                                   0.08
IE10
                    404.0
Chrome
                    200.0
                                   0.02
```

We can fill in the missing values by passing a value to the keyword fill\_value. Because the index is not monotonically increasing or decreasing, we cannot use arguments to the keyword method to fill the NaN values.

```
>>> df.reindex(new_index, fill_value=0)
              http_status response_time
Safari
                     404
                                    0.07
Tceweasel
                       0
                                     0.00
Comodo Dragon
                                     0.00
                        0
IE10
                       404
                                     0.08
Chrome
                       200
                                     0.02
```

```
>>> df.reindex(new_index, fill_value='missing')
            http_status response_time
Safari
                     404
                                  0.07
Iceweasel
                 missing
                               missing
                 missing
Comodo Dragon
                               missing
IE10
                                  0.08
                     404
Chrome
                     200
                                  0.02
```

We can also reindex the columns.

```
>>> df.reindex(columns=['http_status', 'user_agent'])
           http_status user_agent
Firefox
                   200
                               NaN
                   200
Chrome
                                NaN
                   404
Safari
                                NaN
IE10
                   404
                                NaN
                   301
Konqueror
                                NaN
```

Or we can use "axis-style" keyword arguments

```
>>> df.reindex(['http_status', 'user_agent'], axis="columns")
          http_status user_agent
Firefox
                  200
                  200
Chrome
                              NaN
Safari
                  404
                              NaN
IE10
                  404
                              NaN
Konqueror
                  301
                              NaN
```

To further illustrate the filling functionality in reindex, we will create a dataframe with a monotonically increasing index (for example, a sequence of dates).

```
>>> date_index = pd.date_range('1/1/2010', periods=6, freq='D')
>>> df2 = pd.DataFrame({"prices": [100, 101, np.nan, 100, 89, 88]},
... index=date_index)
>>> df2
```

(continues on next page)

```
prices
2010-01-01 100
2010-01-02 101
2010-01-03 NaN
2010-01-04 100
2010-01-05 89
2010-01-06 88
```

Suppose we decide to expand the dataframe to cover a wider date range.

```
>>> date_index2 = pd.date_range('12/29/2009', periods=10, freq='D')
>>> df2.reindex(date_index2)
            prices
2009-12-29
               NaN
2009-12-30
               NaN
2009-12-31
               NaN
2010-01-01
               100
2010-01-02
               101
2010-01-03
               NaN
2010-01-04
               100
2010-01-05
                89
                88
2010-01-06
2010-01-07
               NaN
```

The index entries that did not have a value in the original data frame (for example, '2009-12-29') are by default filled with NaN. If desired, we can fill in the missing values using one of several options.

For example, to backpropagate the last valid value to fill the NaN values, pass bfill as an argument to the method keyword.

```
>>> df2.reindex(date_index2, method='bfill')
            prices
2009-12-29
               100
2009-12-30
               100
2009-12-31
               100
2010-01-01
               100
2010-01-02
               101
2010-01-03
               NaN
2010-01-04
               100
2010-01-05
                89
2010-01-06
                88
2010-01-07
```

Please note that the NaN value present in the original dataframe (at index value 2010-01-03) will not be filled by any of the value propagation schemes. This is because filling while reindexing does not look at dataframe values, but only compares the original and desired indexes. If you do want to fill in the NaN values present in the original dataframe, use the fillna() method.

See the user guide for more.

reindex\_axis (labels, axis=0, method=None, level=None, copy=True, limit=None, fill\_value=nan)

Conform input object to new index with optional filling logic, placing NA/NaN in locations having no value in the previous index. A new object is produced unless the new index is equivalent to the current one and copy=False

### **Parameters**

labels [array-like] New labels / index to conform to. Preferably an Index object to avoid

duplicating data

**axis** [{0, 1, 2, 'items', 'major\_axis', 'minor\_axis'}]

**method** [{None, 'backfill'/'bfill', 'pad'/'ffill', 'nearest'}, optional] Method to use for filling holes in reindexed DataFrame:

- default: don't fill gaps
- pad / ffill: propagate last valid observation forward to next valid
- backfill / bfill: use next valid observation to fill gap
- nearest: use nearest valid observations to fill gap

copy [boolean, default True] Return a new object, even if the passed indexes are the same

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

**limit** [int, default None] Maximum number of consecutive elements to forward or backward fill

**tolerance** [optional] Maximum distance between original and new labels for inexact matches. The values of the index at the matching locations most satisfy the equation abs (index[indexer] - target) <= tolerance.

Tolerance may be a scalar value, which applies the same tolerance to all values, or list-like, which applies variable tolerance per element. List-like includes list, tuple, array, Series, and must be the same size as the index and its dtype must exactly match the index's type.

New in version 0.21.0: (list-like tolerance)

### Returns

reindexed [Panel]

### See also:

reindex, reindex\_like

## **Examples**

```
>>> df.reindex_axis(['A', 'B', 'C'], axis=1)
```

 $\begin{tabular}{ll} \textbf{reindex\_like} (other, method=None, copy=True, limit=None, tolerance=None) \\ \end{tabular}$ 

Return an object with matching indices to myself.

# **Parameters**

other [Object]

method [string or None]

**copy** [boolean, default True]

**limit** [int, default None] Maximum number of consecutive labels to fill for inexact matches.

**tolerance** [optional] Maximum distance between labels of the other object and this object for inexact matches. Can be list-like.

New in version 0.21.0: (list-like tolerance)

#### **Returns**

reindexed [same as input]

### **Notes**

**Like calling s.reindex(index=other.index, columns=other.columns,** method=...)

```
rename (items=None, major_axis=None, minor_axis=None, **kwargs)
```

Alter axes input function or functions. Function / dict values must be unique (1-to-1). Labels not contained in a dict / Series will be left as-is. Extra labels listed don't throw an error. Alternatively, change Series. name with a scalar value (Series only).

#### **Parameters**

**items, major\_axis, minor\_axis** [scalar, list-like, dict-like or function, optional] Scalar or list-like will alter the Series.name attribute, and raise on DataFrame or Panel. dict-like or functions are transformations to apply to that axis' values

copy [boolean, default True] Also copy underlying data

**inplace** [boolean, default False] Whether to return a new Panel. If True then value of copy is ignored.

**level** [int or level name, default None] In case of a MultiIndex, only rename labels in the specified level.

#### Returns

```
renamed [Panel (new object)]
```

## See also:

```
pandas.NDFrame.rename_axis
```

# **Examples**

```
>>> s = pd.Series([1, 2, 3])
>>> s
  1
1
     2
2.
    3
dtype: int64
>>> s.rename("my_name") # scalar, changes Series.name
    1
1
     2
Name: my_name, dtype: int64
>>> s.rename(lambda x: x ** 2) # function, changes labels
0
    1
     2
1
4
     3
dtype: int64
>>> s.rename({1: 3, 2: 5}) # mapping, changes labels
0
    1
3
     2
5
     3
dtype: int64
```

Since DataFrame doesn't have a .name attribute, only mapping-type arguments are allowed.

```
>>> df = pd.DataFrame({"A": [1, 2, 3], "B": [4, 5, 6]})
>>> df.rename(2)
Traceback (most recent call last):
...
TypeError: 'int' object is not callable
```

DataFrame.rename supports two calling conventions

- (index=index\_mapper, columns=columns\_mapper, ...)
- (mapper, axis={'index', 'columns'}, ...)

We *highly* recommend using keyword arguments to clarify your intent.

```
>>> df.rename(index=str, columns={"A": "a", "B": "c"})
    a    c
0    1    4
1    2    5
2    3    6
```

```
>>> df.rename(index=str, columns={"A": "a", "C": "c"})

a B
0 1 4
1 2 5
2 3 6
```

Using axis-style parameters

```
>>> df.rename(str.lower, axis='columns')
    a    b
0    1    4
1    2    5
2    3    6
```

```
>>> df.rename({1: 2, 2: 4}, axis='index')

A B

0 1 4

2 2 5

4 3 6
```

See the user guide for more.

rename\_axis (mapper, axis=0, copy=True, inplace=False)

Alter the name of the index or columns.

#### **Parameters**

**mapper** [scalar, list-like, optional] Value to set as the axis name attribute.

axis [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis.

copy [boolean, default True] Also copy underlying data.

**inplace** [boolean, default False] Modifies the object directly, instead of creating a new Series or DataFrame.

## Returns

**renamed** [Series, DataFrame, or None] The same type as the caller or None if *inplace* is True.

### See also:

```
pandas.Series.rename Alter Series index labels or name
pandas.DataFrame.rename Alter DataFrame index labels or name
pandas.Index.rename Set new names on index
```

### **Notes**

Prior to version 0.21.0, rename\_axis could also be used to change the axis *labels* by passing a mapping or scalar. This behavior is deprecated and will be removed in a future version. Use rename instead.

# **Examples**

### Series

```
>>> s = pd.Series([1, 2, 3])

>>> s.rename_axis("foo")

foo

0    1

1    2

2    3

dtype: int64
```

### **DataFrame**

```
>>> df.rename_axis("bar", axis="columns")
bar A B
0 1 4
1 2 5
2 3 6
```

**replace** (to\_replace=None, value=None, inplace=False, limit=None, regex=False, method='pad')
Replace values given in to\_replace with value.

Values of the NDFrame are replaced with other values dynamically. This differs from updating with .loc or .iloc, which require you to specify a location to update with some value.

## **Parameters**

**to\_replace** [str, regex, list, dict, Series, int, float, or None] How to find the values that will be replaced.

- numeric, str or regex:
  - numeric: numeric values equal to to\_replace will be replaced with value
  - str: string exactly matching to\_replace will be replaced with value

- regex: regexs matching to\_replace will be replaced with value
- list of str, regex, or numeric:
  - First, if *to\_replace* and *value* are both lists, they **must** be the same length.
  - Second, if regex=True then all of the strings in both lists will be interpreted
    as regexs otherwise they will match directly. This doesn't matter much for *value*since there are only a few possible substitution regexes you can use.
  - str, regex and numeric rules apply as above.

#### • dict:

- Dicts can be used to specify different replacement values for different existing values. For example, {'a': 'b', 'y': 'z'} replaces the value 'a' with 'b' and 'y' with 'z'. To use a dict in this way the *value* parameter should be *None*.
- For a DataFrame a dict can specify that different values should be replaced in different columns. For example, { 'a': 1, 'b': 'z'} looks for the value 1 in column 'a' and the value 'z' in column 'b' and replaces these values with whatever is specified in *value*. The *value* parameter should not be None in this case. You can treat this as a special case of passing two lists except that you are specifying the column to search in.
- For a DataFrame nested dictionaries, e.g., {'a': {'b': np.nan}}, are read as follows: look in column 'a' for the value 'b' and replace it with NaN. The *value* parameter should be None to use a nested dict in this way. You can nest regular expressions as well. Note that column names (the top-level dictionary keys in a nested dictionary) **cannot** be regular expressions.

## • None:

- This means that the *regex* argument must be a string, compiled regular expression, or list, dict, ndarray or Series of such elements. If *value* is also None then this **must** be a nested dictionary or Series.

See the examples section for examples of each of these.

value [scalar, dict, list, str, regex, default None] Value to replace any values matching to\_replace with. For a DataFrame a dict of values can be used to specify which value to use for each column (columns not in the dict will not be filled). Regular expressions, strings and lists or dicts of such objects are also allowed.

**inplace** [boolean, default False] If True, in place. Note: this will modify any other views on this object (e.g. a column from a DataFrame). Returns the caller if this is True.

**limit** [int, default None] Maximum size gap to forward or backward fill.

**regex** [bool or same types as *to\_replace*, default False] Whether to interpret *to\_replace* and/or *value* as regular expressions. If this is True then *to\_replace must* be a string. Alternatively, this could be a regular expression or a list, dict, or array of regular expressions in which case *to\_replace* must be None.

**method** [{'pad', 'ffill', 'bfill', *None*}] The method to use when for replacement, when *to\_replace* is a scalar, list or tuple and *value* is None.

Changed in version 0.23.0: Added to DataFrame.

## **Returns**

NDFrame Object after replacement.

### **Raises**

### AssertionError

• If regex is not a bool and to\_replace is not None.

# **TypeError**

- If to replace is a dict and value is not a list, dict, ndarray, or Series
- If *to\_replace* is None and *regex* is not compilable into a regular expression or is a list, dict, ndarray, or Series.
- When replacing multiple bool or datetime64 objects and the arguments to *to\_replace* does not match the type of the value being replaced

#### ValueError

• If a list or an ndarray is passed to *to\_replace* and *value* but they are not the same length.

### See also:

```
NDFrame.fillna Fill NA values
```

NDFrame.where Replace values based on boolean condition

Series.str.replace Simple string replacement.

#### **Notes**

- Regex substitution is performed under the hood with re.sub. The rules for substitution for re.sub are the same.
- Regular expressions will only substitute on strings, meaning you cannot provide, for example, a regular expression matching floating point numbers and expect the columns in your frame that have a numeric dtype to be matched. However, if those floating point numbers *are* strings, then you can do this.
- This method has *a lot* of options. You are encouraged to experiment and play with this method to gain intuition about how it works.
- When dict is used as the *to\_replace* value, it is like key(s) in the dict are the to\_replace part and value(s) in the dict are the value parameter.

# **Examples**

## Scalar 'to\_replace' and 'value'

```
>>> df = pd.DataFrame({'A': [0, 1, 2, 3, 4],

... 'B': [5, 6, 7, 8, 9],

... 'C': ['a', 'b', 'c', 'd', 'e']})
>>> df.replace(0, 5)

A B C
0 5 5 a
1 1 6 b
2 2 7 c
3 3 8 d
4 4 9 e
```

# List-like 'to\_replace'

```
>>> df.replace([0, 1, 2, 3], 4)

A B C

0 4 5 a

1 4 6 b

2 4 7 c

3 4 8 d

4 4 9 e
```

```
>>> df.replace([0, 1, 2, 3], [4, 3, 2, 1])

A B C

0 4 5 a

1 3 6 b

2 2 7 c

3 1 8 d

4 4 9 e
```

## dict-like 'to\_replace'

```
>>> df.replace({0: 10, 1: 100})

A B C

0 10 5 a

1 100 6 b

2 2 7 c

3 3 8 d

4 4 9 e
```

```
>>> df.replace({'A': 0, 'B': 5}, 100)

A B C

0 100 100 a

1 1 6 b

2 2 7 c

3 3 8 d

4 4 9 e
```

```
>>> df.replace({'A': {0: 100, 4: 400}})

A B C

0 100 5 a

1 1 6 b

2 2 7 c

3 3 8 d

4 400 9 e
```

### Regular expression 'to\_replace'

```
>>> df.replace({'A': r'^ba.$'}, {'A': 'new'}, regex=True)

A B

0 new abc

1 foo bar

2 bait xyz
```

```
>>> df.replace(regex=r'^ba.$', value='new')

A B

0 new abc

1 foo new

2 bait xyz
```

```
>>> df.replace(regex={r'^ba.$':'new', 'foo':'xyz'})
        A    B
0    new abc
1    xyz    new
2    bait    xyz
```

```
>>> df.replace(regex=[r'^ba.$', 'foo'], value='new')

A B

0 new abc

1 new new

2 bait xyz
```

Note that when replacing multiple bool or datetime64 objects, the data types in the *to\_replace* parameter must match the data type of the value being replaced:

This raises a TypeError because one of the dict keys is not of the correct type for replacement.

Compare the behavior of  $s.replace({'a': None})$  and s.replace('a', None) to understand the pecularities of the  $to\_replace$  parameter:

```
>>> s = pd.Series([10, 'a', 'a', 'b', 'a'])
```

When one uses a dict as the *to\_replace* value, it is like the value(s) in the dict are equal to the *value* parameter. s.replace({'a': None}) is equivalent to s.replace(to\_replace={'a': None}, value=None, method=None):

```
>>> s.replace({'a': None})
0 10
1 None
2 None
3 b
4 None
dtype: object
```

When value=None and to\_replace is a scalar, list or tuple, replace uses the method parameter (default 'pad') to do the replacement. So this is why the 'a' values are being replaced by 10 in rows 1 and 2 and 'b' in row 4 in this case. The command s.replace('a', None) is actually equivalent to s. replace(to\_replace='a', value=None, method='pad'):

```
>>> s.replace('a', None)
0    10
1    10
2    10
3    b
4    b
dtype: object
```

resample (rule, how=None, axis=0, fill\_method=None, closed=None, label=None, convention='start', kind=None, loffset=None, limit=None, base=0, on=None, level=None)

Convenience method for frequency conversion and resampling of time series. Object must have a datetime-like index (DatetimeIndex, PeriodIndex, or TimedeltaIndex), or pass datetime-like values to the on or level keyword.

### **Parameters**

**rule** [string] the offset string or object representing target conversion

axis [int, optional, default 0]

**closed** [{'right', 'left'}] Which side of bin interval is closed. The default is 'left' for all frequency offsets except for 'M', 'A', 'Q', 'BM', 'BA', 'BQ', and 'W' which all have a default of 'right'.

**label** [{'right', 'left'}] Which bin edge label to label bucket with. The default is 'left' for all frequency offsets except for 'M', 'A', 'Q', 'BM', 'BA', 'BQ', and 'W' which all have a default of 'right'.

**convention** [{'start', 'end', 's', 'e'}] For PeriodIndex only, controls whether to use the start or end of *rule* 

kind: {'timestamp', 'period'}, optional Pass 'timestamp' to convert the resulting index to a DateTimeIndex or 'period' to convert it to a PeriodIndex. By default the input representation is retained.

**loffset** [timedelta] Adjust the resampled time labels

**base** [int, default 0] For frequencies that evenly subdivide 1 day, the "origin" of the aggregated intervals. For example, for '5min' frequency, base could range from 0 through 4. Defaults to 0

**on** [string, optional] For a DataFrame, column to use instead of index for resampling. Column must be datetime-like.

New in version 0.19.0.

**level** [string or int, optional] For a MultiIndex, level (name or number) to use for resampling. Level must be datetime-like.

New in version 0.19.0.

#### Returns

## Resampler object

## See also:

groupby Group by mapping, function, label, or list of labels.

#### **Notes**

See the user guide for more.

To learn more about the offset strings, please see this link.

## **Examples**

Start by creating a series with 9 one minute timestamps.

```
>>> index = pd.date_range('1/1/2000', periods=9, freq='T')
>>> series = pd.Series(range(9), index=index)
>>> series
2000-01-01 00:00:00
2000-01-01 00:01:00
                       1
2000-01-01 00:02:00
                       2.
2000-01-01 00:03:00
2000-01-01 00:04:00
2000-01-01 00:05:00
2000-01-01 00:06:00
2000-01-01 00:07:00
                       7
2000-01-01 00:08:00
Freq: T, dtype: int64
```

Downsample the series into 3 minute bins and sum the values of the timestamps falling into a bin.

Downsample the series into 3 minute bins as above, but label each bin using the right edge instead of the left. Please note that the value in the bucket used as the label is not included in the bucket, which it labels. For example, in the original series the bucket 2000-01-01 00:03:00 contains the value 3, but the summed value in the resampled bucket with the label 2000-01-01 00:03:00 does not include 3 (if it did, the summed value would be 6, not 3). To include this value close the right side of the bin interval as illustrated in the example below this one.

Downsample the series into 3 minute bins as above, but close the right side of the bin interval.

Upsample the series into 30 second bins.

Upsample the series into 30 second bins and fill the NaN values using the pad method.

Upsample the series into 30 second bins and fill the NaN values using the bfill method.

Pass a custom function via apply

```
>>> def custom_resampler(array_like):
... return np.sum(array_like)+5
```

For a Series with a PeriodIndex, the keyword *convention* can be used to control whether to use the start or end of *rule*.

Resample by month using 'start' convention. Values are assigned to the first month of the period.

Resample by month using 'end' convention. Values are assigned to the last month of the period.

```
>>> s.resample('M', convention='end').asfreq()
2012-12 1.0
2013-01
         NaN
2013-02
         NaN
2013-03
         NaN
2013-04
        NaN
2013-05
        NaN
2013-06
        NaN
2013-07 NaN
2013-08
       NaN
2013-09
       NaN
2013-10
       NaN
2013-11
        NaN
2013-12
         2.0
Freq: M, dtype: float64
```

For DataFrame objects, the keyword on can be used to specify the column instead of the index for resampling.

For a DataFrame with MultiIndex, the keyword level can be used to specify on level the resampling needs to take place.

(continued from previous page)

```
a b c d
2000-01-01 00:00:00 0 6 12 18
2000-01-01 00:03:00 0 4 8 12
```

```
rfloordiv (other, axis=0)
```

Integer division of series and other, element-wise (binary operator *rfloordiv*). Equivalent to other // panel.

### **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

#### Returns

**Panel** 

## See also:

Panel.floordiv

#### rmod (other, axis=0)

Modulo of series and other, element-wise (binary operator rmod). Equivalent to other % panel.

#### **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

### **Returns**

Panel

### See also:

Panel.mod

### rmul (other, axis=0)

Multiplication of series and other, element-wise (binary operator rmul). Equivalent to other \* panel.

# **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

# Returns

**Panel** 

## See also:

Panel.mul

```
round (decimals=0, *args, **kwargs)
```

Round each value in Panel to a specified number of decimal places.

New in version 0.18.0.

## **Parameters**

**decimals** [int] Number of decimal places to round to (default: 0). If decimals is negative, it specifies the number of positions to the left of the decimal point.

# Returns

## Panel object

```
See also:
      numpy.around
rpow (other, axis=0)
     Exponential power of series and other, element-wise (binary operator rpow). Equivalent to other \star\star
          Parameters
               other [DataFrame or Panel]
               axis [{items, major_axis, minor_axis}] Axis to broadcast over
          Returns
               Panel
     See also:
     Panel.pow
rsub(other, axis=0)
     Subtraction of series and other, element-wise (binary operator rsub). Equivalent to other - panel.
          Parameters
               other [DataFrame or Panel]
               axis [{items, major_axis, minor_axis}] Axis to broadcast over
          Returns
               Panel
     See also:
     Panel.sub
rtruediv (other, axis=0)
     Floating division of series and other, element-wise (binary operator rtruediv). Equivalent to other /
     panel.
          Parameters
               other [DataFrame or Panel]
               axis [{items, major_axis, minor_axis}] Axis to broadcast over
          Returns
               Panel
     See also:
```

Panel.truediv

**sample** (n=None, frac=None, replace=False, weights=None, random\_state=None, axis=None) Return a random sample of items from an axis of object.

You can use *random\_state* for reproducibility.

## **Parameters**

**n** [int, optional] Number of items from axis to return. Cannot be used with *frac*. Default = 1 if frac = None.

**frac** [float, optional] Fraction of axis items to return. Cannot be used with n.

**replace** [boolean, optional] Sample with or without replacement. Default = False.

weights [str or ndarray-like, optional] Default 'None' results in equal probability weighting. If passed a Series, will align with target object on index. Index values in weights not found in sampled object will be ignored and index values in sampled object not in weights will be assigned weights of zero. If called on a DataFrame, will accept the name of a column when axis = 0. Unless weights are a Series, weights must be same length as axis being sampled. If weights do not sum to 1, they will be normalized to sum to 1. Missing values in the weights column will be treated as zero. inf and -inf values not allowed.

**random\_state** [int or numpy.random.RandomState, optional] Seed for the random number generator (if int), or numpy RandomState object.

**axis** [int or string, optional] Axis to sample. Accepts axis number or name. Default is stat axis for given data type (0 for Series and DataFrames, 1 for Panels).

## Returns

A new object of same type as caller.

## **Examples**

Generate an example Series and DataFrame:

```
>>> s = pd.Series(np.random.randn(50))
>>> s.head()
   -0.038497
0
    1.820773
1
2
   -0.972766
3
   -1.598270
4
   -1.095526
dtype: float64
>>> df = pd.DataFrame(np.random.randn(50, 4), columns=list('ABCD'))
>>> df.head()
                    В
                              С
         Α
 0.016443 -2.318952 -0.566372 -1.028078
1 -1.051921 0.438836 0.658280 -0.175797
2 -1.243569 -0.364626 -0.215065 0.057736
  1.768216 0.404512 -0.385604 -1.457834
  1.072446 -1.137172 0.314194 -0.046661
```

Next extract a random sample from both of these objects...

3 random elements from the Series:

```
>>> s.sample(n=3)
27   -0.994689
55   -1.049016
67   -0.224565
dtype: float64
```

And a random 10% of the DataFrame with replacement:

```
>>> df.sample(frac=0.1, replace=True)

A B C D
```

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You can use *random state* for reproducibility:

#### select (crit, axis=0)

Return data corresponding to axis labels matching criteria

Deprecated since version 0.21.0: Use df.loc[df.index.map(crit)] to select via labels

### **Parameters**

```
crit [function] To be called on each index (label). Should return True or Falseaxis [int]
```

#### Returns

**selection** [type of caller]

**sem** (axis=None, skipna=None, level=None, ddof=1, numeric\_only=None, \*\*kwargs)
Return unbiased standard error of the mean over requested axis.

Normalized by N-1 by default. This can be changed using the ddof argument

### **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

**ddof** [int, default 1] Delta Degrees of Freedom. The divisor used in calculations is N - ddof, where N represents the number of elements.

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

```
sem [DataFrame or Panel (if level specified)]
```

```
set_axis (labels, axis=0, inplace=None)
```

Assign desired index to given axis.

Indexes for column or row labels can be changed by assigning a list-like or Index.

Changed in version 0.21.0: The signature is now *labels* and *axis*, consistent with the rest of pandas API. Previously, the *axis* and *labels* arguments were respectively the first and second positional arguments.

#### **Parameters**

labels [list-like, Index] The values for the new index.

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The axis to update. The value 0 identifies the rows, and 1 identifies the columns.

**inplace** [boolean, default None] Whether to return a new %(klass)s instance.

**Warning:** inplace=None currently falls back to to True, but in a future version, will default to False. Use inplace=True explicitly rather than relying on the default.

### **Returns**

**renamed** [%(klass)s or None] An object of same type as caller if inplace=False, None otherwise.

See also:

pandas.DataFrame.rename axis Alter the name of the index or columns.

## **Examples**

### **Series**

The original object is not modified.

```
>>> s
0    1
1    2
2    3
dtype: int64
```

### **DataFrame**

```
>>> df = pd.DataFrame({"A": [1, 2, 3], "B": [4, 5, 6]})
```

Change the row labels.

```
>>> df.set_axis(['a', 'b', 'c'], axis='index', inplace=False)

A B
a 1 4
b 2 5
c 3 6
```

Change the column labels.

Now, update the labels inplace.

```
set_value(*args, **kwargs)
```

Quickly set single value at (item, major, minor) location

Deprecated since version 0.21.0.

Please use .at[] or .iat[] accessors.

### **Parameters**

```
item [item label (panel item)]
major [major axis label (panel item row)]
minor [minor axis label (panel item column)]
value [scalar]
takeable [interpret the passed labels as indexers, default False]
```

## **Returns**

**panel** [Panel] If label combo is contained, will be reference to calling Panel, otherwise a new object

## shape

Return a tuple of axis dimensions

```
shift (periods=1, freq=None, axis='major')
```

Shift index by desired number of periods with an optional time freq. The shifted data will not include the dropped periods and the shifted axis will be smaller than the original. This is different from the behavior of DataFrame.shift()

### **Parameters**

```
periods [int] Number of periods to move, can be positive or negative
freq [DateOffset, timedelta, or time rule string, optional]
axis [{'items', 'major', 'minor'} or {0, 1, 2}]
Returns
```

### **shifted** [Panel]

#### size

Return an int representing the number of elements in this object.

Return the number of rows if Series. Otherwise return the number of rows times number of columns if DataFrame.

#### See also:

ndarray.size

# **Examples**

```
>>> s = pd.Series({'a': 1, 'b': 2, 'c': 3})
>>> s.size
3
```

```
>>> df = pd.DataFrame({'col1': [1, 2], 'col2': [3, 4]})
>>> df.size
4
```

**skew** (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

Return unbiased skew over requested axis Normalized by N-1

#### **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### **Returns**

**skew** [DataFrame or Panel (if level specified)]

```
slice_shift (periods=1, axis=0)
```

Equivalent to *shift* without copying data. The shifted data will not include the dropped periods and the shifted axis will be smaller than the original.

### **Parameters**

periods [int] Number of periods to move, can be positive or negative

## Returns

**shifted** [same type as caller]

## **Notes**

While the *slice\_shift* is faster than *shift*, you may pay for it later during alignment.

#### **Parameters**

axis [axes to direct sorting]

**level** [int or level name or list of ints or list of level names] if not None, sort on values in specified index level(s)

ascending [boolean, default True] Sort ascending vs. descending

inplace [bool, default False] if True, perform operation in-place

**kind** [{'quicksort', 'mergesort', 'heapsort'}, default 'quicksort'] Choice of sorting algorithm. See also ndarray.np.sort for more information. *mergesort* is the only stable algorithm. For DataFrames, this option is only applied when sorting on a single column or label.

**na\_position** [{'first', 'last'}, default 'last'] *first* puts NaNs at the beginning, *last* puts NaNs at the end. Not implemented for MultiIndex.

**sort\_remaining** [bool, default True] if true and sorting by level and index is multilevel, sort by other levels too (in order) after sorting by specified level

#### Returns

```
sorted_obj [NDFrame]
```

NOT IMPLEMENTED: do not call this method, as sorting values is not supported for Panel objects and will raise an error.

#### squeeze (axis=None)

Squeeze length 1 dimensions.

### **Parameters**

axis [None, integer or string axis name, optional] The axis to squeeze if 1-sized.

New in version 0.20.0.

### **Returns**

## scalar if 1-sized, else original object

**std** (*axis=None*, *skipna=None*, *level=None*, *ddof=1*, *numeric\_only=None*, \*\*kwargs)
Return sample standard deviation over requested axis.

Normalized by N-1 by default. This can be changed using the ddof argument

#### **Parameters**

```
axis [{items (0), major_axis (1), minor_axis (2)}]
```

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

**ddof** [int, default 1] Delta Degrees of Freedom. The divisor used in calculations is N - ddof, where N represents the number of elements.

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

#### Returns

```
std [DataFrame or Panel (if level specified)]
sub(other, axis=0)
      Subtraction of series and other, element-wise (binary operator sub). Equivalent to panel - other.
           Parameters
               other [DataFrame or Panel]
               axis [{items, major_axis, minor_axis}] Axis to broadcast over
           Returns
               Panel
     See also:
     Panel.rsub
subtract (other, axis=0)
      Subtraction of series and other, element-wise (binary operator sub). Equivalent to panel - other.
           Parameters
               other [DataFrame or Panel]
               axis [{items, major_axis, minor_axis}] Axis to broadcast over
           Returns
               Panel
     See also:
      Panel.rsub
sum(axis=None, skipna=None, level=None, numeric_only=None, min_count=0, **kwargs)
     Return the sum of the values for the requested axis
           Parameters
               axis [{items (0), major_axis (1), minor_axis (2)}]
               skipna [boolean, default True] Exclude NA/null values when computing the result.
               level [int or level name, default None] If the axis is a MultiIndex (hierarchical), count
                   along a particular level, collapsing into a DataFrame
               numeric_only [boolean, default None] Include only float, int, boolean columns. If None,
                   will attempt to use everything, then use only numeric data. Not implemented for
                   Series.
               min_count [int, default 0] The required number of valid values to perform the operation.
                   If fewer than min count non-NA values are present the result will be NA.
                   New in version 0.22.0: Added with the default being 0. This means the sum of an
                   all-NA or empty Series is 0, and the product of an all-NA or empty Series is 1.
           Returns
               sum [DataFrame or Panel (if level specified)]
```

## **Examples**

By default, the sum of an empty or all-NA Series is 0.

```
>>> pd.Series([]).sum() # min_count=0 is the default
0.0
```

This can be controlled with the min\_count parameter. For example, if you'd like the sum of an empty series to be NaN, pass min\_count=1.

```
>>> pd.Series([]).sum(min_count=1)
nan
```

Thanks to the skipna parameter, min\_count handles all-NA and empty series identically.

```
>>> pd.Series([np.nan]).sum()
0.0
```

```
>>> pd.Series([np.nan]).sum(min_count=1)
nan
```

swapaxes (axis1, axis2, copy=True)

Interchange axes and swap values axes appropriately

### **Returns**

y [same as input]

```
swaplevel(i=-2, j=-1, axis=0)
```

Swap levels i and j in a MultiIndex on a particular axis

### **Parameters**

**i, j** [int, string (can be mixed)] Level of index to be swapped. Can pass level name as string.

## Returns

**swapped** [type of caller (new object)]

.. versionchanged:: 0.18.1 The indexes i and j are now optional, and default to the two innermost levels of the index.

# tail(n=5)

Return the last *n* rows.

This function returns last n rows from the object based on position. It is useful for quickly verifying data, for example, after sorting or appending rows.

### **Parameters**

**n** [int, default 5] Number of rows to select.

#### Returns

**type of caller** The last n rows of the caller object.

See also:

pandas.DataFrame.head The first n rows of the caller object.

# **Examples**

```
>>> df = pd.DataFrame({'animal':['alligator', 'bee', 'falcon', 'lion',
                        'monkey', 'parrot', 'shark', 'whale', 'zebra']})
>>> df
      animal
  alligator
1
         bee
2
      falcon
3
        lion
4
     monkey
5
      parrot
6
      shark
7
       whale
8
       zebra
```

## Viewing the last 5 lines

```
>>> df.tail()
   animal
4 monkey
5 parrot
6 shark
7 whale
8 zebra
```

Viewing the last *n* lines (three in this case)

```
>>> df.tail(3)
animal
6 shark
7 whale
8 zebra
```

take (indices, axis=0, convert=None, is\_copy=True, \*\*kwargs)

Return the elements in the given *positional* indices along an axis.

This means that we are not indexing according to actual values in the index attribute of the object. We are indexing according to the actual position of the element in the object.

#### **Parameters**

**indices** [array-like] An array of ints indicating which positions to take.

**axis** [{0 or 'index', 1 or 'columns', None}, default 0] The axis on which to select elements. 0 means that we are selecting rows, 1 means that we are selecting columns.

**convert** [bool, default True] Whether to convert negative indices into positive ones. For example, -1 would map to the len(axis) - 1. The conversions are similar to the behavior of indexing a regular Python list.

Deprecated since version 0.21.0: In the future, negative indices will always be converted.

is\_copy [bool, default True] Whether to return a copy of the original object or not.

\*\*kwargs For compatibility with numpy.take(). Has no effect on the output.

### Returns

taken [type of caller] An array-like containing the elements taken from the object.

### See also:

**DataFrame.loc** Select a subset of a DataFrame by labels.

**DataFrame.iloc** Select a subset of a DataFrame by positions.

numpy.take Take elements from an array along an axis.

## **Examples**

```
>>> df = pd.DataFrame([('falcon', 'bird',
                                            389.0),
                      ('parrot', 'bird',
                                             24.0).
                      ('lion',
                                 'mammal',
                                             80.5),
. . .
                      ('monkey', 'mammal', np.nan)],
. . .
                      columns=['name', 'class', 'max_speed'],
. . .
                      index=[0, 2, 3, 1])
. . .
>>> df
    name class max_speed
          bird
                  389.0
0 falcon
           bird
2
                      24.0
 parrot
3
    lion mammal
                      80.5
  monkey mammal
                       NaN
```

Take elements at positions 0 and 3 along the axis 0 (default).

Note how the actual indices selected (0 and 1) do not correspond to our selected indices 0 and 3. That's because we are selecting the 0th and 3rd rows, not rows whose indices equal 0 and 3.

```
>>> df.take([0, 3])
    name class max_speed
0 falcon bird 389.0
1 monkey mammal NaN
```

Take elements at indices 1 and 2 along the axis 1 (column selection).

We may take elements using negative integers for positive indices, starting from the end of the object, just like with Python lists.

```
>>> df.take([-1, -2])
    name class max_speed
1 monkey mammal NaN
3 lion mammal 80.5
```

## to\_clipboard(excel=True, sep=None, \*\*kwargs)

Copy object to the system clipboard.

Write a text representation of object to the system clipboard. This can be pasted into Excel, for example.

## **Parameters**

excel [bool, default True]

- True, use the provided separator, writing in a csv format for allowing easy pasting into excel.
- False, write a string representation of the object to the clipboard.

```
sep [str, default '\t'] Field delimiter.
```

\*\*kwargs These parameters will be passed to DataFrame.to\_csv.

See also:

**DataFrame.to\_csv** Write a DataFrame to a comma-separated values (csv) file.

read\_clipboard Read text from clipboard and pass to read\_table.

#### **Notes**

Requirements for your platform.

```
• Linux : xclip, or xsel (with gtk or PyQt4 modules)
```

• Windows: none

• OS X : none

## **Examples**

Copy the contents of a DataFrame to the clipboard.

```
>>> df = pd.DataFrame([[1, 2, 3], [4, 5, 6]], columns=['A', 'B', 'C'])
>>> df.to_clipboard(sep=',')
... # Wrote the following to the system clipboard:
... # ,A,B,C
... # 0,1,2,3
... # 1,4,5,6
```

We can omit the the index by passing the keyword *index* and setting it to false.

```
>>> df.to_clipboard(sep=',', index=False)
... # Wrote the following to the system clipboard:
... # A,B,C
... # 1,2,3
... # 4,5,6
```

## to\_dense()

Return dense representation of NDFrame (as opposed to sparse)

```
to_excel (path, na_rep=", engine=None, **kwargs)
```

Write each DataFrame in Panel to a separate excel sheet

### **Parameters**

```
path [string or ExcelWriter object] File path or existing ExcelWriter
```

```
na_rep [string, default ''] Missing data representation
```

```
engine [string, default None] write engine to use - you can also set this via the op-
tions io.excel.xlsx.writer, io.excel.xls.writer, and io.excel.
xlsm.writer.
```

### **Other Parameters**

**float\_format** [string, default None] Format string for floating point numbers

cols [sequence, optional] Columns to write

**header** [boolean or list of string, default True] Write out column names. If a list of string is given it is assumed to be aliases for the column names

index [boolean, default True] Write row names (index)

index\_label [string or sequence, default None] Column label for index column(s) if desired. If None is given, and *header* and *index* are True, then the index names are used. A sequence should be given if the DataFrame uses MultiIndex.

startrow [upper left cell row to dump data frame]

**startcol** [upper left cell column to dump data frame]

#### **Notes**

Keyword arguments (and na\_rep) are passed to the to\_excel method for each DataFrame written.

## to\_frame (filter\_observations=True)

Transform wide format into long (stacked) format as DataFrame whose columns are the Panel's items and whose index is a MultiIndex formed of the Panel's major and minor axes.

#### **Parameters**

**filter\_observations** [boolean, default True] Drop (major, minor) pairs without a complete set of observations across all the items

### Returns

y [DataFrame]

## to\_hdf (path\_or\_buf, key, \*\*kwargs)

Write the contained data to an HDF5 file using HDFStore.

Hierarchical Data Format (HDF) is self-describing, allowing an application to interpret the structure and contents of a file with no outside information. One HDF file can hold a mix of related objects which can be accessed as a group or as individual objects.

In order to add another DataFrame or Series to an existing HDF file please use append mode and a different a key.

For more information see the user guide.

### **Parameters**

path\_or\_buf [str or pandas.HDFStore] File path or HDFStore object.

key [str] Identifier for the group in the store.

**mode** [{'a', 'w', 'r+'}, default 'a'] Mode to open file:

- 'w': write, a new file is created (an existing file with the same name would be deleted).
- 'a': append, an existing file is opened for reading and writing, and if the file does not exist it is created.
- 'r+': similar to 'a', but the file must already exist.

**format** [{'fixed', 'table'}, default 'fixed'] Possible values:

- 'fixed': Fixed format. Fast writing/reading. Not-appendable, nor searchable.
- 'table': Table format. Write as a PyTables Table structure which may perform worse but allow more flexible operations like searching / selecting subsets of the data.

**append** [bool, default False] For Table formats, append the input data to the existing.

data\_columns [list of columns or True, optional] List of columns to create as indexed data columns for on-disk queries, or True to use all columns. By default only the axes of the object are indexed. See io.hdf5-query-data-columns. Applicable only to format='table'.

**complevel** [{0-9}, optional] Specifies a compression level for data. A value of 0 disables compression.

**complib** [{'zlib', 'lzo', 'bzip2', 'blosc'}, default 'zlib'] Specifies the compression library to be used. As of v0.20.2 these additional compressors for Blosc are supported (default if no compressor specified: 'blosc:blosclz'): {'blosc:blosclz', 'blosc:lz4', 'blosc:lz4hc', 'blosc:snappy', 'blosc:zlib', 'blosc:zstd'}. Specifying a compression library which is not available issues a ValueError.

**fletcher32** [bool, default False] If applying compression use the fletcher32 checksum.

**dropna** [bool, default False] If true, ALL nan rows will not be written to store.

**errors** [str, default 'strict'] Specifies how encoding and decoding errors are to be handled. See the errors argument for open () for a full list of options.

#### See also:

DataFrame.read\_hdf Read from HDF file.

**DataFrame.to\_parquet** Write a DataFrame to the binary parquet format.

DataFrame.to\_sql Write to a sql table.

DataFrame.to\_feather Write out feather-format for DataFrames.

DataFrame.to\_csv Write out to a csv file.

#### **Examples**

```
>>> df = pd.DataFrame({'A': [1, 2, 3], 'B': [4, 5, 6]},
... index=['a', 'b', 'c'])
>>> df.to_hdf('data.h5', key='df', mode='w')
```

We can add another object to the same file:

```
>>> s = pd.Series([1, 2, 3, 4])
>>> s.to_hdf('data.h5', key='s')
```

### Reading from HDF file:

```
>>> pd.read_hdf('data.h5', 'df')

A B
a 1 4
b 2 5
c 3 6
>>> pd.read_hdf('data.h5', 's')
```

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```
0 1
1 2
2 3
3 4
dtype: int64
```

### Deleting file with data:

```
>>> import os
>>> os.remove('data.h5')
```

to\_json (path\_or\_buf=None, orient=None, date\_format=None, double\_precision=10, force\_ascii=True, date\_unit='ms', default\_handler=None, lines=False, compression=None, index=True)

Convert the object to a JSON string.

Note NaN's and None will be converted to null and datetime objects will be converted to UNIX timestamps.

### **Parameters**

**path\_or\_buf** [string or file handle, optional] File path or object. If not specified, the result is returned as a string.

**orient** [string] Indication of expected JSON string format.

- Series
  - default is 'index'
  - allowed values are: {'split','records','index'}
- DataFrame
  - default is 'columns'
  - allowed values are: {'split','records','index','columns','values'}
- The format of the JSON string
  - 'split': dict like { 'index' -> [index], 'columns' -> [columns], 'data' -> [values]}
  - 'records': list like [{column -> value}, ..., {column -> value}]
  - 'index' : dict like {index -> {column -> value}}
  - 'columns' : dict like {column -> {index -> value}}
  - 'values': just the values array
  - 'table': dict like {'schema': {schema}, 'data': {data}} describing the data, and the data component is like orient='records'.

Changed in version 0.20.0.

date\_format [{None, 'epoch', 'iso'}] Type of date conversion. 'epoch' = epoch milliseconds, 'iso' = ISO8601. The default depends on the orient. For orient='table',
the default is 'iso'. For all other orients, the default is 'epoch'.

**double\_precision** [int, default 10] The number of decimal places to use when encoding floating point values.

force\_ascii [boolean, default True] Force encoded string to be ASCII.

- date\_unit [string, default 'ms' (milliseconds)] The time unit to encode to, governs timestamp and ISO8601 precision. One of 's', 'ms', 'us', 'ns' for second, millisecond, microsecond, and nanosecond respectively.
- **default\_handler** [callable, default None] Handler to call if object cannot otherwise be converted to a suitable format for JSON. Should receive a single argument which is the object to convert and return a serialisable object.
- **lines** [boolean, default False] If 'orient' is 'records' write out line delimited json format. Will throw ValueError if incorrect 'orient' since others are not list like.

New in version 0.19.0.

**compression** [{None, 'gzip', 'bz2', 'zip', 'xz'}] A string representing the compression to use in the output file, only used when the first argument is a filename.

New in version 0.21.0.

index [boolean, default True] Whether to include the index values in the JSON string. Not including the index (index=False) is only supported when orient is 'split' or 'table'.

New in version 0.23.0.

#### See also:

pandas.read\_json

## **Examples**

Encoding/decoding a Dataframe using 'records' formatted JSON. Note that index labels are not preserved with this encoding.

```
>>> df.to_json(orient='records')
'[{"col 1":"a","col 2":"b"},{"col 1":"c","col 2":"d"}]'
```

Encoding/decoding a Dataframe using 'index' formatted JSON:

```
>>> df.to_json(orient='index')
'{"row 1":{"col 1":"a","col 2":"b"},"row 2":{"col 1":"c","col 2":"d"}}'
```

Encoding/decoding a Dataframe using 'columns' formatted JSON:

```
>>> df.to_json(orient='columns')
'{"col 1":{"row 1":"a","row 2":"c"},"col 2":{"row 1":"b","row 2":"d"}}'
```

Encoding/decoding a Dataframe using 'values' formatted JSON:

```
>>> df.to_json(orient='values')
'[["a","b"],["c","d"]]'
```

### **Encoding with Table Schema**

to\_latex (buf=None, columns=None, col\_space=None, header=True, index=True, na\_rep='NaN', formatters=None, float\_format=None, sparsify=None, index\_names=True, bold\_rows=False, column\_format=None, longtable=None, escape=None, encoding=None, decimal='.', multicolumn=None, multicolumn\_format=None, multirow=None)

Render an object to a tabular environment table. You can splice this into a LaTeX document. Requires \usepackage{booktabs}.

Changed in version 0.20.2: Added to Series

*to\_latex*-specific options:

**bold\_rows** [boolean, default False] Make the row labels bold in the output

column\_format [str, default None] The columns format as specified in LaTeX table format e.g 'rcl' for 3 columns

**longtable** [boolean, default will be read from the pandas config module] Default: False. Use a longtable environment instead of tabular. Requires adding a \usepackage{longtable} to your LaTeX preamble.

**escape** [boolean, default will be read from the pandas config module] Default: True. When set to False prevents from escaping latex special characters in column names.

**encoding** [str, default None] A string representing the encoding to use in the output file, defaults to 'ascii' on Python 2 and 'utf-8' on Python 3.

**decimal** [string, default '.'] Character recognized as decimal separator, e.g. ',' in Europe.

New in version 0.18.0.

**multicolumn** [boolean, default True] Use multicolumn to enhance MultiIndex columns. The default will be read from the config module.

New in version 0.20.0.

**multicolumn\_format** [str, default '1'] The alignment for multicolumns, similar to *column\_format* The default will be read from the config module.

New in version 0.20.0.

multirow [boolean, default False] Use multirow to enhance MultiIndex rows. Requires adding a \usepackage{multirow} to your LaTeX preamble. Will print centered labels (instead of top-aligned) across the contained rows, separating groups via clines. The default will be read from the pandas config module.

New in version 0.20.0.

to\_msgpack (path\_or\_buf=None, encoding='utf-8', \*\*kwargs)
msgpack (serialize) object to input file path

THIS IS AN EXPERIMENTAL LIBRARY and the storage format may not be stable until a future release.

### **Parameters**

path [string File path, buffer-like, or None] if None, return generated stringappend [boolean whether to append to an existing msgpack] (default is False)compress [type of compressor (zlib or blosc), default to None (no] compression)

to\_pickle (path, compression='infer', protocol=2)
Pickle (serialize) object to file.

### **Parameters**

path [str] File path where the pickled object will be stored.

**compression** [{'infer', 'gzip', 'bz2', 'zip', 'xz', None}, default 'infer'] A string representing the compression to use in the output file. By default, infers from the file extension in specified path.

New in version 0.20.0.

**protocol** [int] Int which indicates which protocol should be used by the pickler, default HIGHEST\_PROTOCOL (see [1] paragraph 12.1.2). The possible values for this parameter depend on the version of Python. For Python 2.x, possible values are 0, 1, 2. For Python>=3.0, 3 is a valid value. For Python>= 3.4, 4 is a valid value. A negative value for the protocol parameter is equivalent to setting its value to HIGH-EST\_PROTOCOL.

New in version 0.21.0.

#### See also:

read\_pickle Load pickled pandas object (or any object) from file.

DataFrame.to\_hdf Write DataFrame to an HDF5 file.

DataFrame.to\_sql Write DataFrame to a SQL database.

**DataFrame.to\_parquet** Write a DataFrame to the binary parquet format.

# **Examples**

```
>>> original_df = pd.DataFrame({"foo": range(5), "bar": range(5, 10)})
>>> original_df
    foo bar
0     0     5
1     1     6
2     2     7
3     3     8
4     4     9
>>> original_df.to_pickle("./dummy.pkl")
```

```
>>> unpickled_df = pd.read_pickle("./dummy.pkl")
>>> unpickled_df
   foo bar
0
     0
          5
1
     1
          6
2
     2
          7
3
     3
          8
4
     4
          9
```

```
>>> import os
>>> os.remove("./dummy.pkl")
```

## to\_sparse(\*args, \*\*kwargs)

NOT IMPLEMENTED: do not call this method, as sparsifying is not supported for Panel objects and will raise an error.

Convert to SparsePanel

to\_sql (name, con, schema=None, if\_exists='fail', index=True, index\_label=None, chunksize=None, dtype=None)

Write records stored in a DataFrame to a SQL database.

Databases supported by SQLAlchemy [1] are supported. Tables can be newly created, appended to, or overwritten.

#### **Parameters**

name [string] Name of SQL table.

**con** [sqlalchemy.engine.Engine or sqlite3.Connection] Using SQLAlchemy makes it possible to use any DB supported by that library. Legacy support is provided for sqlite3.Connection objects.

**schema** [string, optional] Specify the schema (if database flavor supports this). If None, use default schema.

if\_exists [{'fail', 'replace', 'append'}, default 'fail'] How to behave if the table already exists.

- fail: Raise a ValueError.
- replace: Drop the table before inserting new values.
- append: Insert new values to the existing table.

**index** [boolean, default True] Write DataFrame index as a column. Uses *index\_label* as the column name in the table.

**index\_label** [string or sequence, default None] Column label for index column(s). If None is given (default) and *index* is True, then the index names are used. A sequence should be given if the DataFrame uses MultiIndex.

**chunksize** [int, optional] Rows will be written in batches of this size at a time. By default, all rows will be written at once.

**dtype** [dict, optional] Specifying the datatype for columns. The keys should be the column names and the values should be the SQLAlchemy types or strings for the sqlite3 legacy mode.

# Raises

**ValueError** When the table already exists and *if exists* is 'fail' (the default).

See also:

pandas.read\_sql read a DataFrame from a table

### References

[1], [2]

## **Examples**

Create an in-memory SQLite database.

```
>>> from sqlalchemy import create_engine
>>> engine = create_engine('sqlite://', echo=False)
```

Create a table from scratch with 3 rows.

```
>>> df.to_sql('users', con=engine)
>>> engine.execute("SELECT * FROM users").fetchall()
[(0, 'User 1'), (1, 'User 2'), (2, 'User 3')]
```

Overwrite the table with just df1.

```
>>> df1.to_sql('users', con=engine, if_exists='replace',
... index_label='id')
>>> engine.execute("SELECT * FROM users").fetchall()
[(0, 'User 4'), (1, 'User 5')]
```

Specify the dtype (especially useful for integers with missing values). Notice that while pandas is forced to store the data as floating point, the database supports nullable integers. When fetching the data with Python, we get back integer scalars.

```
>>> from sqlalchemy.types import Integer
>>> df.to_sql('integers', con=engine, index=False,
... dtype={"A": Integer()})
```

```
>>> engine.execute("SELECT * FROM integers").fetchall()
[(1,), (None,), (2,)]
```

#### to\_xarray()

Return an xarray object from the pandas object.

### Returns

a DataArray for a Series

- a Dataset for a DataFrame
- a DataArray for higher dims

#### **Notes**

See the xarray docs

## **Examples**

```
>>> df.to_xarray()
<xarray.Dataset>
Dimensions: (index: 3)
Coordinates:
  * index (index) int64 0 1 2
Data variables:
  A (index) int64 1 1 2
  B (index) object 'foo' 'bar' 'foo'
  C (index) float64 4.0 5.0 6.0
```

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```
>>> p.to_xarray()
<xarray.DataArray (items: 4, major_axis: 3, minor_axis: 2)>
array([[[ 0, 1],
        [ 2, 3],
        [4,5]],
       [[6, 7],
       [8, 9],
       [10, 11]],
       [[12, 13],
        [14, 15],
        [16, 17]],
       [[18, 19],
        [20, 21],
        [22, 23]])
Coordinates:
               (items) object 'A' 'B' 'C' 'D'
 * items
  * major_axis (major_axis) datetime64[ns] 2013-01-01 2013-01-02 2013-01-03_
→ # noqa
  * minor_axis (minor_axis) object 'first' 'second'
```

### transpose (\*args, \*\*kwargs)

Permute the dimensions of the Panel

#### **Parameters**

```
args [three positional arguments: each one of]
{0, 1, 2, 'items', 'major_axis', 'minor_axis'}
```

**copy** [boolean, default False] Make a copy of the underlying data. Mixed-dtype data will always result in a copy

#### **Returns**

y [same as input]

## **Examples**

```
>>> p.transpose(2, 0, 1)
>>> p.transpose(2, 0, 1, copy=True)
```

# truediv (other, axis=0)

Floating division of series and other, element-wise (binary operator *truediv*). Equivalent to panel / other.

#### **Parameters**

```
other [DataFrame or Panel]
axis [{items, major_axis, minor_axis}] Axis to broadcast over
```

### **Returns**

#### **Panel**

### See also:

```
Panel.rtruediv
```

truncate (before=None, after=None, axis=None, copy=True)

Truncate a Series or DataFrame before and after some index value.

This is a useful shorthand for boolean indexing based on index values above or below certain thresholds.

### **Parameters**

```
before [date, string, int] Truncate all rows before this index value.
```

```
after [date, string, int] Truncate all rows after this index value.
```

**axis** [{0 or 'index', 1 or 'columns'}, optional] Axis to truncate. Truncates the index (rows) by default.

**copy** [boolean, default is True,] Return a copy of the truncated section.

#### Returns

**type of caller** The truncated Series or DataFrame.

### See also:

**DataFrame.loc** Select a subset of a DataFrame by label.

**DataFrame.iloc** Select a subset of a DataFrame by position.

#### **Notes**

If the index being truncated contains only datetime values, *before* and *after* may be specified as strings instead of Timestamps.

# **Examples**

```
>>> df = pd.DataFrame({'A': ['a', 'b', 'c', 'd', 'e'],
                         'B': ['f', 'g', 'h', 'i', 'j'],
. . .
                         'C': ['k', 'l', 'm', 'n', 'o']},
. . .
                        index=[1, 2, 3, 4, 5])
. . .
>>> df
         С
  A B
     f
  b
         1
      g
3
  С
      h
         m
4
  d
      í
         n
5
  е
      j
         0
```

```
>>> df.truncate(before=2, after=4)

A B C

2 b g l

3 c h m

4 d i n
```

The columns of a DataFrame can be truncated.

```
>>> df.truncate(before="A", after="B", axis="columns")

A B

1 a f

2 b g

3 c h

4 d i

5 e j
```

For Series, only rows can be truncated.

The index values in truncate can be datetimes or string dates.

```
>>> dates = pd.date_range('2016-01-01', '2016-02-01', freq='s')
>>> df = pd.DataFrame(index=dates, data={'A': 1})
>>> df.tail()

A
2016-01-31 23:59:56 1
2016-01-31 23:59:57 1
2016-01-31 23:59:58 1
2016-01-31 23:59:59 1
2016-02-01 00:00:00 1
```

```
>>> df.truncate(before=pd.Timestamp('2016-01-05'),
... after=pd.Timestamp('2016-01-10')).tail()

A
2016-01-09 23:59:56 1
2016-01-09 23:59:57 1
2016-01-09 23:59:58 1
2016-01-09 23:59:59 1
2016-01-10 00:00:00 1
```

Because the index is a DatetimeIndex containing only dates, we can specify *before* and *after* as strings. They will be coerced to Timestamps before truncation.

```
>>> df.truncate('2016-01-05', '2016-01-10').tail()

A
2016-01-09 23:59:56 1
2016-01-09 23:59:57 1
2016-01-09 23:59:58 1
2016-01-09 23:59:59 1
2016-01-10 00:00:00 1
```

Note that truncate assumes a 0 value for any unspecified time component (midnight). This differs from partial string slicing, which returns any partially matching dates.

```
>>> df.loc['2016-01-05':'2016-01-10', :].tail()

A
2016-01-10 23:59:55 1
2016-01-10 23:59:56 1
2016-01-10 23:59:57 1
```

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```
2016-01-10 23:59:58 1
2016-01-10 23:59:59 1
```

tshift (periods=1, freq=None, axis='major')

Shift the time index, using the index's frequency if available.

#### **Parameters**

periods [int] Number of periods to move, can be positive or negative

**freq** [DateOffset, timedelta, or time rule string, default None] Increment to use from the tseries module or time rule (e.g. 'EOM')

axis [int or basestring] Corresponds to the axis that contains the Index

### **Returns**

shifted [NDFrame]

### **Notes**

If freq is not specified then tries to use the freq or inferred\_freq attributes of the index. If neither of those attributes exist, a ValueError is thrown

tz convert (tz, axis=0, level=None, copy=True)

Convert tz-aware axis to target time zone.

### **Parameters**

tz [string or pytz.timezone object]

axis [the axis to convert]

**level** [int, str, default None] If axis ia a MultiIndex, convert a specific level. Otherwise must be None

copy [boolean, default True] Also make a copy of the underlying data

#### Raises

**TypeError** If the axis is tz-naive.

tz\_localize(tz, axis=0, level=None, copy=True, ambiguous='raise')

Localize tz-naive TimeSeries to target time zone.

## **Parameters**

**tz** [string or pytz.timezone object]

axis [the axis to localize]

**level** [int, str, default None] If axis ia a MultiIndex, localize a specific level. Otherwise must be None

**copy** [boolean, default True] Also make a copy of the underlying data

ambiguous ['infer', bool-ndarray, 'NaT', default 'raise']

- 'infer' will attempt to infer fall dst-transition hours based on order
- bool-ndarray where True signifies a DST time, False designates a non-DST time (note that this flag is only applicable for ambiguous times)
- 'NaT' will return NaT where there are ambiguous times

• 'raise' will raise an AmbiguousTimeError if there are ambiguous times

#### Raises

**TypeError** If the TimeSeries is tz-aware and tz is not None.

update (other, join='left', overwrite=True, filter\_func=None, raise\_conflict=False)

Modify Panel in place using non-NA values from passed Panel, or object coercible to Panel. Aligns on items

#### **Parameters**

```
other [Panel, or object coercible to Panel]
```

**join** [How to join individual DataFrames] { 'left', 'right', 'outer', 'inner'}, default 'left'

**overwrite** [boolean, default True] If True then overwrite values for common keys in the calling panel

**filter\_func** [callable(1d-array) -> 1d-array<br/>boolean>, default None] Can choose to replace values other than NA. Return True for values that should be updated

**raise\_conflict** [bool] If True, will raise an error if a DataFrame and other both contain data in the same place.

#### values

Return a Numpy representation of the DataFrame.

Only the values in the DataFrame will be returned, the axes labels will be removed.

#### Returns

numpy.ndarray The values of the DataFrame.

See also:

```
pandas.DataFrame.index Retrievie the index labels
pandas.DataFrame.columns Retrieving the column names
```

## **Notes**

The dtype will be a lower-common-denominator dtype (implicit upcasting); that is to say if the dtypes (even of numeric types) are mixed, the one that accommodates all will be chosen. Use this with care if you are not dealing with the blocks.

e.g. If the dtypes are float16 and float32, dtype will be upcast to float32. If dtypes are int32 and uint8, dtype will be upcast to int32. By numpy.find\_common\_type() convention, mixing int64 and uint64 will result in a float64 dtype.

### **Examples**

A DataFrame where all columns are the same type (e.g., int64) results in an array of the same type.

```
>>> df = pd.DataFrame({'age': [ 3, 29], ... 'height': [94, 170], ... 'weight': [31, 115]})
>>> df
age height weight
0 3 94 31
```

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```
29
           170
                    115
1
>>> df.dtypes
          int64
age
height
          int64
weight
          int64
dtype: object
>>> df.values
array([[ 3, 94,
                   31],
       [ 29, 170, 115]], dtype=int64)
```

A DataFrame with mixed type columns(e.g., str/object, int64, float32) results in an ndarray of the broadest type that accommodates these mixed types (e.g., object).

```
>>> df2 = pd.DataFrame([('parrot',
                                       24.0, 'second'),
                         ('lion',
                                       80.5, 1),
. . .
                         ('monkey', np.nan, None)],
. . .
                       columns=('name', 'max_speed', 'rank'))
. . .
>>> df2.dtypes
              object
name
max_speed float64
rank
              object
dtype: object
>>> df2.values
array([['parrot', 24.0, 'second'],
       ['lion', 80.5, 1],
       ['monkey', nan, None]], dtype=object)
```

var (axis=None, skipna=None, level=None, ddof=1, numeric\_only=None, \*\*kwargs)
Return unbiased variance over requested axis.

Normalized by N-1 by default. This can be changed using the ddof argument

### **Parameters**

```
axis [{items (0), major axis (1), minor axis (2)}]
```

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a DataFrame

**ddof** [int, default 1] Delta Degrees of Freedom. The divisor used in calculations is N - ddof, where N represents the number of elements.

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

#### Returns

var [DataFrame or Panel (if level specified)]

**where** (cond, other=nan, inplace=False, axis=None, level=None, errors='raise', try\_cast=False, raise\_on\_error=None)

Return an object of same shape as self and whose corresponding entries are from self where *cond* is True and otherwise are from *other*.

### **Parameters**

**cond** [boolean NDFrame, array-like, or callable] Where *cond* is True, keep the original value. Where False, replace with corresponding value from *other*. If *cond* is callable, it is computed on the NDFrame and should return boolean NDFrame or array. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as cond.

**other** [scalar, NDFrame, or callable] Entries where *cond* is False are replaced with corresponding value from *other*. If other is callable, it is computed on the NDFrame and should return scalar or NDFrame. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as other.

inplace [boolean, default False] Whether to perform the operation in place on the data

axis [alignment axis if needed, default None]

level [alignment level if needed, default None]

errors [str, {'raise', 'ignore'}, default 'raise']

- raise: allow exceptions to be raised
- ignore: suppress exceptions. On error return original object

Note that currently this parameter won't affect the results and will always coerce to a suitable dtype.

try\_cast [boolean, default False] try to cast the result back to the input type (if possible),

raise\_on\_error [boolean, default True] Whether to raise on invalid data types (e.g. trying to where on strings)

Deprecated since version 0.21.0.

## Returns

**wh** [same type as caller]

## See also:

```
DataFrame.mask()
```

# **Notes**

The where method is an application of the if-then idiom. For each element in the calling DataFrame, if cond is True the element is used; otherwise the corresponding element from the DataFrame other is used.

The signature for DataFrame.where () differs from numpy.where(). Roughly df1.where(m, df2) is equivalent to np.where(m, df1, df2).

For further details and examples see the where documentation in indexing.

### **Examples**

```
>>> s = pd.Series(range(5))
>>> s.where(s > 0)
0    NaN
1    1.0
```

(continues on next page)

(continued from previous page)

```
2 2.0
3 3.0
4 4.0
```

```
>>> df = pd.DataFrame(np.arange(10).reshape(-1, 2), columns=['A', 'B'])
>>> m = df % 3 == 0
>>> df.where(m, -df)
  A B
0 \quad 0 \quad -1
1 -2 3
2 - 4 - 5
3 6 -7
4 -8 9
>>> df.where(m, -df) == np.where(m, df, -df)
     Α
0 True True
1 True True
2 True True
3 True True
4 True True
>>> df.where(m, -df) == df.mask(\sim m, -df)
     Α
        В
0 True True
  True True
  True True
3
  True True
4 True True
```

# **xs** (*key*, *axis*=1)

Return slice of panel along selected axis

# **Parameters**

```
key [object] Label
axis [{'items', 'major', 'minor}, default 1/'major']
Returns
y [ndim(self)-1]
```

### **Notes**

xs is only for getting, not setting values.

MultiIndex Slicers is a generic way to get/set values on any level or levels and is a superset of xs functionality, see MultiIndex Slicers

Bases: pandas.core.series.Series

Pandas Series modified to adapt *oddt.toolkit.Molecule* objects and apply molecular methods easily.

New in version 0.3.

### **Attributes**

```
T return the transpose, which is by definition self
asobject Return object Series which contains boxed values.
at Access a single value for a row/column label pair.
axes Return a list of the row axis labels
base return the base object if the memory of the underlying data is
blocks Internal property, property synonym for as blocks()
```

data return the data pointer of the underlying data

dtype return the dtype object of the underlying data

dtypes return the dtype object of the underlying data

#### empty

flags return the ndarray.flags for the underlying data

ftype return if the data is sparseldense

ftypes return if the data is sparseldense

hasnans return if I have any nans; enables various perf speedups

iat Access a single value for a row/column pair by integer position.

**iloc** Purely integer-location based indexing for selection by position.

## imag

**index** The index (axis labels) of the Series.

# is\_copy

is\_monotonic Return boolean if values in the object are

is\_monotonic\_decreasing Return boolean if values in the object are

is\_monotonic\_increasing Return boolean if values in the object are

is\_unique Return boolean if values in the object are unique

itemsize return the size of the dtype of the item of the underlying data

ix A primarily label-location based indexer, with integer position fallback.

**loc** Access a group of rows and columns by label(s) or a boolean array.

name

nbytes return the number of bytes in the underlying datandim return the number of dimensions of the underlying data,

# real

shape return a tuple of the shape of the underlying data
size return the number of elements in the underlying data
strides return the strides of the underlying data
values Return Series as ndarray or ndarray-like

## **Methods**

abs()	Return a Series/DataFrame with absolute numeric value of each element.	
add(other[, level, fill_value, axis])	Addition of series and other, element-wise (binary	
add(outer[, fever, fift_value, axis])	operator <i>add</i> ).	
add profiv(profiv)	Prefix labels with string <i>prefix</i> .	
<pre>add_prefix(prefix) add_suffix(suffix)</pre>		
	Suffix labels with string <i>suffix</i> .	
agg(func[, axis])	Aggregate using one or more operations over the specified axis.	
aggregate(func[, axis])	Aggregate using one or more operations over the specified axis.	
align(other[, join, axis, level, copy,])	Align two objects on their axes with the specified	
	join method for each axis Index	
all([axis, bool_only, skipna, level])	Return whether all elements are True, potentially	
	over an axis.	
any([axis, bool_only, skipna, level])	Return whether any element is True over requested	
	axis.	
append(to_append[, ignore_index,])	Concatenate two or more Series.	
<pre>apply(func[, convert_dtype, args])</pre>	Invoke function on values of Series.	
argmax(**kwargs)		
argmin(**kwargs)		
argsort([axis, kind, order])	Overrides ndarray.argsort.	
as_blocks([copy])	Convert the frame to a dict of dtype -> Constructor	
	Types that each has a homogeneous dtype.	
as_matrix([columns])	Convert the frame to its Numpy-array representation.	
asfreq(freq[, method, how, normalize,])	Convert TimeSeries to specified frequency.	
asof(where[, subset])	The last row without any NaN is taken (or the last	
	row without NaN considering only the subset of	
	columns in the case of a DataFrame)	
astype(**kwargs)	Cast a pandas object to a specified dtype dtype.	
<pre>at_time(time[, asof])</pre>	Select values at particular time of day (e.g.	
autocorr([lag])	Lag-N autocorrelation	
between(left, right[, inclusive])	Return boolean Series equivalent to left <= series <=	
	right.	
between_time(start_time, end_time[,])	Select values between particular times of the day	
	(e.g., 9:00-9:30 AM).	
bfill([axis, inplace, limit, downcast])	Synonym for DataFrame.	
•	fillna(method='bfill')	
	Continued on next page	

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	ed from previous page
bool()	Return the bool of a single element PandasObject.
calcfp(*args, **kwargs)	Helper function to map FP calculation through the
	series
cat	alias of pandas.core.arrays.
	categorical.CategoricalAccessor
clip([lower, upper, axis, inplace])	Trim values at input threshold(s).
<pre>clip_lower(threshold[, axis, inplace])</pre>	Return copy of the input with values below a thresh-
	old truncated.
<pre>clip_upper(threshold[, axis, inplace])</pre>	Return copy of input with values above given
	value(s) truncated.
combine(other, func[, fill_value])	Perform elementwise binary operation on two Series
	using given function with optional fill value when an
	index is missing from one Series or the other
combine_first(other)	Combine Series values, choosing the calling Series's
	values first.
compound([axis, skipna, level])	Return the compound percentage of the values for
	the requested axis
compress(condition, *args, **kwargs)	Return selected slices of an array along given axis as
	a Series
consolidate([inplace])	Compute NDFrame with "consolidated" internals
-	(data of each dtype grouped together in a single ndar-
	ray).
<pre>convert_objects([convert_dates,])</pre>	Attempt to infer better dtype for object columns.
copy([deep])	Make a copy of this object's indices and data.
corr(other[, method, min_periods])	Compute correlation with <i>other</i> Series, excluding
· · · · · · · · · · · · · · · · · · ·	missing values
count([level])	Return number of non-NA/null observations in the
	Series
cov(other[, min_periods])	Compute covariance with Series, excluding missing
, <u>, , , , , , , , , , , , , , , , , , </u>	values
cummax([axis, skipna])	Return cumulative maximum over a DataFrame or
	Series axis.
cummin([axis, skipna])	Return cumulative minimum over a DataFrame or
	Series axis.
cumprod([axis, skipna])	Return cumulative product over a DataFrame or Se-
- · · · · · · · · · · · · · · · · · · ·	ries axis.
cumsum([axis, skipna])	Return cumulative sum over a DataFrame or Series
V 1 2/	axis.
describe([percentiles, include, exclude])	Generates descriptive statistics that summarize the
(c)	central tendency, dispersion and shape of a dataset's
	distribution, excluding NaN values.
diff([periods])	First discrete difference of element.
div(other[, level, fill_value, axis])	Floating division of series and other, element-wise
(	(binary operator <i>truediv</i> ).
divide(other[, level, fill_value, axis])	Floating division of series and other, element-wise
(	(binary operator <i>truediv</i> ).
divmod(other[, level, fill_value, axis])	Integer division and modulo of series and other,
al mode (outer), 1000, mi_value, anisj)	element-wise (binary operator <i>divmod</i> ).
dot(other)	Matrix multiplication with DataFrame or inner-
accionici)	product with Series objects.
drop([labels, axis, index, columns, level,])	Return Series with specified index labels removed.
ατορ(μαυσιό, αλίό, πασλ, columnio, revel,])	Continued on next page
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	d from previous page
<pre>drop_duplicates([keep, inplace])</pre>	Return Series with duplicate values removed.
dropna([axis, inplace])	Return a new Series with missing values removed.
duplicated([keep])	Indicate duplicate Series values.
eq(other[, level, fill_value, axis])	Equal to of series and other, element-wise (binary
	operator $eq$ ).
equals(other)	Determines if two NDFrame objects contain the
	same elements.
ewm([com, span, halflife, alpha,])	Provides exponential weighted functions
expanding([min_periods, center, axis])	Provides expanding transformations.
factorize([sort, na_sentinel])	Encode the object as an enumerated type or categor-
	ical variable.
ffill([axis, inplace, limit, downcast])	Synonym for DataFrame.
TTTT([mins, inplace, mins, so misses])	fillna (method='ffill')
fillna([value, method, axis, inplace,])	Fill NA/NaN values using the specified method
filter([items, like, regex, axis])	Subset rows or columns of dataframe according to
rricer ([nems, nkc, regex, axis])	labels in the specified index.
first(offeet)	Convenience method for subsetting initial periods of
first(offset)	time series data based on a date offset.
	Return index for first non-NA/null value.
first_valid_index()	
floordiv(other[, level, fill_value, axis])	Integer division of series and other, element-wise (bi-
	nary operator <i>floordiv</i> ).
from_array(arr[, index, name, dtype, copy,])	Construct Series from array.
from_csv(path[, sep, parse_dates, header,])	Read CSV file.
ge(other[, level, fill_value, axis])	Greater than or equal to of series and other, element-
	wise (binary operator ge).
get(key[, default])	Get item from object for given key (DataFrame col-
	umn, Panel slice, etc.).
get_dtype_counts()	Return counts of unique dtypes in this object.
get_ftype_counts()	Return counts of unique ftypes in this object.
get_value(label[, takeable])	Quickly retrieve single value at passed index label
get_values()	same as values (but handles sparseness conversions);
	is a view
<pre>groupby([by, axis, level, as_index, sort,])</pre>	Group series using mapper (dict or key function, ap-
	ply given function to group, return result as series) or
	by a series of columns.
gt(other[, level, fill_value, axis])	Greater than of series and other, element-wise (bi-
	nary operator $gt$ ).
head([n])	Return the first <i>n</i> rows.
hist([by, ax, grid, xlabelsize, xrot,])	Draw histogram of the input series using matplotlib
idxmax([axis, skipna])	Return the row label of the maximum value.
Taxiiax([axis, skipila])	
	Return the row label of the minimum value.
idxmin([axis, skipna])	Return the row label of the minimum value.
<pre>idxmin([axis, skipna]) infer_objects()</pre>	Return the row label of the minimum value.  Attempt to infer better dtypes for object columns.
<pre>idxmin([axis, skipna]) infer_objects() interpolate([method, axis, limit, inplace,])</pre>	Return the row label of the minimum value.  Attempt to infer better dtypes for object columns.  Interpolate values according to different methods.
<pre>idxmin([axis, skipna]) infer_objects() interpolate([method, axis, limit, inplace,]) isin(values)</pre>	Return the row label of the minimum value.  Attempt to infer better dtypes for object columns.  Interpolate values according to different methods.  Check whether <i>values</i> are contained in Series.
<pre>idxmin([axis, skipna]) infer_objects() interpolate([method, axis, limit, inplace,]) isin(values) isna()</pre>	Return the row label of the minimum value.  Attempt to infer better dtypes for object columns.  Interpolate values according to different methods.  Check whether <i>values</i> are contained in Series.  Detect missing values.
<pre>idxmin([axis, skipna]) infer_objects() interpolate([method, axis, limit, inplace,]) isin(values) isna() isnull()</pre>	Return the row label of the minimum value.  Attempt to infer better dtypes for object columns.  Interpolate values according to different methods.  Check whether <i>values</i> are contained in Series.  Detect missing values.  Detect missing values.
<pre>idxmin([axis, skipna]) infer_objects() interpolate([method, axis, limit, inplace,]) isin(values) isna()</pre>	Return the row label of the minimum value.  Attempt to infer better dtypes for object columns.  Interpolate values according to different methods.  Check whether <i>values</i> are contained in Series.  Detect missing values.  Detect missing values.  return the first element of the underlying data as a
<pre>idxmin([axis, skipna]) infer_objects() interpolate([method, axis, limit, inplace,]) isin(values) isna() isnull() item()</pre>	Return the row label of the minimum value.  Attempt to infer better dtypes for object columns.  Interpolate values according to different methods.  Check whether <i>values</i> are contained in Series.  Detect missing values.  Detect missing values.  return the first element of the underlying data as a python scalar
<pre>idxmin([axis, skipna]) infer_objects() interpolate([method, axis, limit, inplace,]) isin(values) isna() isnull() item()</pre>	Return the row label of the minimum value.  Attempt to infer better dtypes for object columns.  Interpolate values according to different methods.  Check whether <i>values</i> are contained in Series.  Detect missing values.  Detect missing values.  return the first element of the underlying data as a python scalar  Lazily iterate over (index, value) tuples
<pre>idxmin([axis, skipna]) infer_objects() interpolate([method, axis, limit, inplace,]) isin(values) isna() isnull() item()  items() items()</pre>	Return the row label of the minimum value.  Attempt to infer better dtypes for object columns.  Interpolate values according to different methods.  Check whether <i>values</i> are contained in Series.  Detect missing values.  Detect missing values.  return the first element of the underlying data as a python scalar  Lazily iterate over (index, value) tuples  Lazily iterate over (index, value) tuples
<pre>idxmin([axis, skipna]) infer_objects() interpolate([method, axis, limit, inplace,]) isin(values) isna() isnull() item()</pre>	Return the row label of the minimum value.  Attempt to infer better dtypes for object columns.  Interpolate values according to different methods.  Check whether <i>values</i> are contained in Series.  Detect missing values.  Detect missing values.  return the first element of the underlying data as a python scalar  Lazily iterate over (index, value) tuples

Table 21 – continued f	from previous	page
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Table 21 – continue	d from previous page
kurt([axis, skipna, level, numeric_only])	Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0).
<pre>kurtosis([axis, skipna, level, numeric_only])</pre>	Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0).
last(offset)	Convenience method for subsetting final periods of time series data based on a date offset.
last_valid_index()	Return index for last non-NA/null value.
1e(other[, level, fill_value, axis])	Less than or equal to of series and other, elementwise (binary operator $le$ ).
1t(other[, level, fill_value, axis])	Less than of series and other, element-wise (binary operator $lt$ ).
mad([axis, skipna, level])	Return the mean absolute deviation of the values for the requested axis
map(arg[, na_action])	Map values of Series using input correspondence (a dict, Series, or function).
mask(cond[, other, inplace, axis, level,])	Return an object of same shape as self and whose corresponding entries are from self where <i>cond</i> is False and otherwise are from <i>other</i> .
<pre>max([axis, skipna, level, numeric_only])</pre>	This method returns the maximum of the values in the object.
mean([axis, skipna, level, numeric_only])	Return the mean of the values for the requested axis
median([axis, skipna, level, numeric_only])	Return the median of the values for the requested axis
<pre>memory_usage([index, deep])</pre>	Return the memory usage of the Series.
<pre>min([axis, skipna, level, numeric_only])</pre>	This method returns the minimum of the values in the object.
mod(other[, level, fill_value, axis])	Modulo of series and other, element-wise (binary operator <i>mod</i> ).
mode()	Return the mode(s) of the dataset.
<pre>mul(other[, level, fill_value, axis])</pre>	Multiplication of series and other, element-wise (binary operator <i>mul</i> ).
<pre>multiply(other[, level, fill_value, axis])</pre>	Multiplication of series and other, element-wise (binary operator <i>mul</i> ).
ne(other[, level, fill_value, axis])	Not equal to of series and other, element-wise (binary operator <i>ne</i> ).
nlargest([n, keep])	Return the largest <i>n</i> elements.
nonzero()	Return the <i>integer</i> indices of the elements that are non-zero
notna()	Detect existing (non-missing) values.
notnull()	Detect existing (non-missing) values.
nsmallest([n, keep])	Return the smallest <i>n</i> elements.
nunique([dropna])	Return number of unique elements in the object.
pct_change([periods, fill_method, limit, freq])	Percentage change between the current and a prior element.
pipe(func, *args, **kwargs)	Apply func(self, *args, **kwargs)
plot	alias of pandas.plottingcore. SeriesPlotMethods
pop(item)	Return item and drop from frame.
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	d from previous page
pow(other[, level, fill_value, axis])	Exponential power of series and other, element-wise (binary operator <i>pow</i> ).
prod([axis, skipna, level, numeric_only,])	Return the product of the values for the requested axis
product([axis, skipna, level, numeric_only,])	Return the product of the values for the requested axis
ptp([axis, skipna, level, numeric_only])	Returns the difference between the maximum value and the minimum value in the object.
put(*args, **kwargs)	Applies the <i>put</i> method to its <i>values</i> attribute if it has one.
quantile([q, interpolation])	Return value at the given quantile, a la numpy.percentile.
radd(other[, level, fill_value, axis])	Addition of series and other, element-wise (binary operator <i>radd</i> ).
rank([axis, method, numeric_only,])	Compute numerical data ranks (1 through n) along axis.
ravel([order])	Return the flattened underlying data as an ndarray
rdiv(other[, level, fill_value, axis])	Floating division of series and other, element-wise (binary operator <i>rtruediv</i> ).
reindex([index])	Conform Series to new index with optional filling logic, placing NA/NaN in locations having no value in the previous index.
reindex_axis(labels[, axis])	Conform Series to new index with optional filling logic.
reindex_like(other[, method, copy, limit,])	Return an object with matching indices to myself.
rename([index])	Alter Series index labels or name
rename_axis(mapper[, axis, copy, inplace])	Alter the name of the index or columns.
reorder_levels(order)	Rearrange index levels using input order.
repeat(**kwargs)	Repeat elements of an Series.
replace([to_replace, value, inplace, limit,])	Replace values given in to_replace with value.
resample(rule[, how, axis, fill_method,])	Convenience method for frequency conversion and resampling of time series.
reset_index([level, drop, name, inplace])	Generate a new DataFrame or Series with the index reset.
rfloordiv(other[, level, fill_value, axis])	Integer division of series and other, element-wise (binary operator <i>rfloordiv</i> ).
rmod(other[, level, fill_value, axis])	Modulo of series and other, element-wise (binary operator <i>rmod</i> ).
rmul(other[, level, fill_value, axis])	Multiplication of series and other, element-wise (binary operator <i>rmul</i> ).
rolling(window[, min_periods, center,])	Provides rolling window calculations.
round([decimals])	Round each value in a Series to the given number of decimals.
rpow(other[, level, fill_value, axis])	Exponential power of series and other, element-wise (binary operator <i>rpow</i> ).
rsub(other[, level, fill_value, axis])	Subtraction of series and other, element-wise (binary operator <i>rsub</i> ).
rtruediv(other[, level, fill_value, axis])	Floating division of series and other, element-wise (binary operator <i>rtruediv</i> ).
sample([n, frac, replace, weights,])	Return a random sample of items from an axis of object.
	Continued on next page

Table 21 – continued from previous page	
searchsorted(**kwargs)	Find indices where elements should be inserted to
	maintain order.
select(crit[, axis])	Return data corresponding to axis labels matching
	criteria
sem([axis, skipna, level, ddof, numeric_only])	Return unbiased standard error of the mean over re-
1 (E. 1) 1 P 10 11 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	quested axis.
set_axis(labels[, axis, inplace])	Assign desired index to given axis.
set_value(label, value[, takeable])	Quickly set single value at passed label.
shift([periods, freq, axis])	Shift index by desired number of periods with an op-
SHITT E([periods, freq, axis])	tional time freq
skew([axis, skipna, level, numeric_only])	Return unbiased skew over requested axis Normal-
Shew ([axis, skipila, level, numeric_omy])	ized by N-1
alian shift([nominda avial)	
slice_shift([periods, axis])	Equivalent to <i>shift</i> without copying data.
sort_index([axis, level, ascending,])	Sort Series by index labels.
sort_values([axis, ascending, inplace,])	Sort by the values.
sortlevel([level, ascending, sort_remaining])	Sort Series with MultiIndex by chosen level.
squeeze([axis])	Squeeze length 1 dimensions.
std([axis, skipna, level, ddof, numeric_only])	Return sample standard deviation over requested
	axis.
str	alias of pandas.core.strings.
	StringMethods
sub(other[, level, fill_value, axis])	Subtraction of series and other, element-wise (binary
	operator <i>sub</i> ).
subtract(other[, level, fill_value, axis])	Subtraction of series and other, element-wise (binary
	operator <i>sub</i> ).
sum([axis, skipna, level, numeric_only,])	Return the sum of the values for the requested axis
swapaxes(axis1, axis2[, copy])	Interchange axes and swap values axes appropriately
swaplevel([i, j, copy])	Swap levels i and j in a MultiIndex
tail([n])	Return the last <i>n</i> rows.
take(indices[, axis, convert, is_copy])	Return the elements in the given <i>positional</i> indices
	along an axis.
to_clipboard([excel, sep])	Copy object to the system clipboard.
to_csv([path, index, sep, na_rep,])	Write Series to a comma-separated values (csv) file
to_dense()	Return dense representation of NDFrame (as op-
	posed to sparse)
to_dict([into])	Convert Series to {label -> value} dict or dict-like
	object.
to_excel(excel_writer[, sheet_name, na_rep,	Write Series to an excel sheet
])	
to_frame([name])	Convert Series to DataFrame
to_hdf(path_or_buf, key, **kwargs)	Write the contained data to an HDF5 file using HDF-
co_nar (pani_oi_oui, key, kwaigs)	Store.
to_json([path_or_buf, orient, date_format,])	Convert the object to a JSON string.
to_latex([buf, columns, col_space, header,])	Render an object to a tabular environment table.
to_msgpack([path_or_buf, encoding])	msgpack (serialize) object to input file path
	Convert Series from DatetimeIndex to PeriodIndex
to_period([freq, copy])	
	with desired frequency (inferred from index if not
( 1.7 ( d.f	passed)
to_pickle(path[, compression, protocol])	Pickle (serialize) object to file.
to_sparse([kind, fill_value])	Convert Series to SparseSeries
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	With a serial stand in a Data France COL
to_sql(name, con[, schema, if_exists,])	Write records stored in a DataFrame to a SQL
	database.
to_string([buf, na_rep, float_format,])	Render a string representation of the Series
to_timestamp([freq, how, copy])	Cast to datetimeindex of timestamps, at beginning of
	period
to_xarray()	Return an xarray object from the pandas object.
tolist()	Return a list of the values.
transform(func, *args, **kwargs)	Call function producing a like-indexed NDFrame
	and return a NDFrame with the transformed values
transpose(*args, **kwargs)	return the transpose, which is by definition self
truediv(other[, level, fill_value, axis])	Floating division of series and other, element-wise
	(binary operator truediv).
truncate([before, after, axis, copy])	Truncate a Series or DataFrame before and after
	some index value.
tshift([periods, freq, axis])	Shift the time index, using the index's frequency if
	available.
tz_convert(tz[, axis, level, copy])	Convert tz-aware axis to target time zone.
tz_localize(tz[, axis, level, copy, ambiguous])	Localize tz-naive TimeSeries to target time zone.
unique()	Return unique values of Series object.
unstack([level, fill_value])	Unstack, a.k.a.
update(other)	Modify Series in place using non-NA values from
	passed Series.
valid([inplace])	Return Series without null values.
<pre>value_counts([normalize, sort, ascending,])</pre>	Returns object containing counts of unique values.
var([axis, skipna, level, ddof, numeric_only])	Return unbiased variance over requested axis.
view([dtype])	Create a new view of the Series.
where(cond[, other, inplace, axis, level,])	Return an object of same shape as self and whose
	corresponding entries are from self where cond is
	True and otherwise are from <i>other</i> .
xs(key[, axis, level, drop_level])	Returns a cross-section (row(s) or column(s)) from
	the Series/DataFrame.

dt	
to_mol2	
to_sdf	
to_smiles	

Т

return the transpose, which is by definition self

# abs()

Return a Series/DataFrame with absolute numeric value of each element.

This function only applies to elements that are all numeric.

## Returns

abs Series/DataFrame containing the absolute value of each element.

# See also:

numpy.absolute calculate the absolute value element-wise.

# **Notes**

For complex inputs, 1.2 + 1 j, the absolute value is  $\sqrt{a^2 + b^2}$ .

# **Examples**

Absolute numeric values in a Series.

```
>>> s = pd.Series([-1.10, 2, -3.33, 4])

>>> s.abs()

0    1.10

1    2.00

2    3.33

3    4.00

dtype: float64
```

Absolute numeric values in a Series with complex numbers.

```
>>> s = pd.Series([1.2 + 1j])
>>> s.abs()
0    1.56205
dtype: float64
```

Absolute numeric values in a Series with a Timedelta element.

```
>>> s = pd.Series([pd.Timedelta('1 days')])
>>> s.abs()
0  1 days
dtype: timedelta64[ns]
```

Select rows with data closest to certain value using argsort (from StackOverflow).

```
>>> df = pd.DataFrame({
       'a': [4, 5, 6, 7],
       'b': [10, 20, 30, 40],
       'c': [100, 50, -30, -50]
...})
>>> df
   а
       b c
      10 100
0
  4
1
   5
       20 50
2.
       30 -30
    6
3
    7
       40 -50
>>> df.loc[(df.c - 43).abs().argsort()]
       b
    а
    5
        20
            50
       10
           100
0
    4
2
    6
       30 -30
3
    7 40 -50
```

add (other, level=None, fill\_value=None, axis=0)

Addition of series and other, element-wise (binary operator add).

Equivalent to series + other, but with support to substitute a fill\_value for missing data in one of the inputs.

**Parameters** 

**other** [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### Returns

result [Series]

## See also:

Series.radd

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
b
    1.0
С
    1.0
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
    1.0
b
     NaN
     1.0
     NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
b
    1.0
С
    1.0
    1.0
    NaN
dtype: float64
```

# add\_prefix (prefix)

Prefix labels with string prefix.

For Series, the row labels are prefixed. For DataFrame, the column labels are prefixed.

# **Parameters**

**prefix** [str] The string to add before each label.

## Returns

**Series or DataFrame** New Series or DataFrame with updated labels.

## See also:

Series.add\_suffix Suffix row labels with string suffix.

DataFrame.add\_suffix Suffix column labels with string suffix.

# **Examples**

```
>>> s.add_prefix('item_')
item_0   1
item_1   2
item_2   3
item_3   4
dtype: int64
```

```
>>> df = pd.DataFrame({'A': [1, 2, 3, 4], 'B': [3, 4, 5, 6]})
>>> df
A B
0 1 3
1 2 4
2 3 5
3 4 6
```

## add\_suffix (suffix)

Suffix labels with string suffix.

For Series, the row labels are suffixed. For DataFrame, the column labels are suffixed.

#### **Parameters**

suffix [str] The string to add after each label.

# Returns

**Series or DataFrame** New Series or DataFrame with updated labels.

## See also:

Series.add\_prefix Prefix row labels with string prefix.

DataFrame.add\_prefix Prefix column labels with string prefix.

# **Examples**

```
>>> s = pd.Series([1, 2, 3, 4])
>>> s
0 1
```

(continues on next page)

(continued from previous page)

```
1 2
2 3
3 4
dtype: int64
```

agg (func, axis=0, \*args, \*\*kwargs)

Aggregate using one or more operations over the specified axis.

New in version 0.20.0.

### **Parameters**

**func** [function, string, dictionary, or list of string/functions] Function to use for aggregating the data. If a function, must either work when passed a Series or when passed to Series.apply. For a DataFrame, can pass a dict, if the keys are DataFrame column names.

Accepted combinations are:

- string function name.
- function.
- list of functions.
- dict of column names -> functions (or list of functions).

**axis** [{0 or 'index'}] Parameter needed for compatibility with DataFrame.

\*args Positional arguments to pass to func.

\*\*kwargs Keyword arguments to pass to func.

## **Returns**

aggregated [Series]

See also:

```
pandas. Series.apply, pandas. Series.transform
```

### **Notes**

agg is an alias for aggregate. Use the alias.

A passed user-defined-function will be passed a Series for evaluation.

# **Examples**

```
>>> s = Series(np.random.randn(10))
```

```
>>> s.agg('min')
-1.3018049988556679
```

```
>>> s.agg(['min', 'max'])
min -1.301805
max 1.127688
dtype: float64
```

## aggregate (func, axis=0, \*args, \*\*kwargs)

Aggregate using one or more operations over the specified axis.

New in version 0.20.0.

#### **Parameters**

**func** [function, string, dictionary, or list of string/functions] Function to use for aggregating the data. If a function, must either work when passed a Series or when passed to Series.apply. For a DataFrame, can pass a dict, if the keys are DataFrame column names.

Accepted combinations are:

- string function name.
- function.
- list of functions.
- dict of column names -> functions (or list of functions).

axis [{0 or 'index'}] Parameter needed for compatibility with DataFrame.

\*args Positional arguments to pass to func.

\*\*kwargs Keyword arguments to pass to func.

# Returns

```
aggregated [Series]
```

## See also:

```
pandas.Series.apply, pandas.Series.transform
```

### **Notes**

agg is an alias for aggregate. Use the alias.

A passed user-defined-function will be passed a Series for evaluation.

# **Examples**

```
>>> s = Series(np.random.randn(10))
```

```
>>> s.agg('min')
-1.3018049988556679
```

```
>>> s.agg(['min', 'max'])
min -1.301805
max 1.127688
dtype: float64
```

Align two objects on their axes with the specified join method for each axis Index

#### **Parameters**

```
other [DataFrame or Series]
```

```
join [{'outer', 'inner', 'left', 'right'}, default 'outer']
```

**axis** [allowed axis of the other object, default None] Align on index (0), columns (1), or both (None)

**level** [int or level name, default None] Broadcast across a level, matching Index values on the passed MultiIndex level

**copy** [boolean, default True] Always returns new objects. If copy=False and no reindexing is required then original objects are returned.

**fill\_value** [scalar, default np.NaN] Value to use for missing values. Defaults to NaN, but can be any "compatible" value

```
method [str, default None]
```

**limit** [int, default None]

fill\_axis [{0 or 'index'}, default 0] Filling axis, method and limit

**broadcast\_axis** [{0 or 'index'}, default None] Broadcast values along this axis, if aligning two objects of different dimensions

### Returns

```
(left, right) [(Series, type of other)] Aligned objects
```

all (axis=0, bool\_only=None, skipna=True, level=None, \*\*kwargs)

Return whether all elements are True, potentially over an axis.

Returns True if all elements within a series or along a Dataframe axis are non-zero, not-empty or not-False.

## **Parameters**

**axis** [{0 or 'index', 1 or 'columns', None}, default 0] Indicate which axis or axes should be reduced.

- 0 / 'index': reduce the index, return a Series whose index is the original column labels.
- 1 / 'columns' : reduce the columns, return a Series whose index is the original index.
- None: reduce all axes, return a scalar.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar.

**bool\_only** [boolean, default None] Include only boolean columns. If None, will attempt to use everything, then use only boolean data. Not implemented for Series.

\*\*kwargs [any, default None] Additional keywords have no effect but might be accepted for compatibility with NumPy.

### **Returns**

all [scalar or Series (if level specified)]

#### See also:

```
pandas.Series.all Return True if all elements are True
pandas.DataFrame.any Return True if one (or more) elements are True
```

## **Examples**

# Series

```
>>> pd.Series([True, True]).all()
True
>>> pd.Series([True, False]).all()
False
```

# **DataFrames**

Create a dataframe from a dictionary.

```
>>> df = pd.DataFrame({'col1': [True, True], 'col2': [True, False]})
>>> df
    col1    col2
0    True    True
1    True    False
```

Default behaviour checks if column-wise values all return True.

```
>>> df.all()
col1 True
col2 False
dtype: bool
```

Specify axis='columns' to check if row-wise values all return True.

```
>>> df.all(axis='columns')

0    True

1    False
dtype: bool
```

Or axis=None for whether every value is True.

```
>>> df.all(axis=None)
False
```

any (axis=0, bool\_only=None, skipna=True, level=None, \*\*kwargs)

Return whether any element is True over requested axis.

Unlike DataFrame.all(), this performs an *or* operation. If any of the values along the specified axis is True, this will return True.

## **Parameters**

axis [{0 or 'index', 1 or 'columns', None}, default 0] Indicate which axis or axes should be reduced.

- 0 / 'index': reduce the index, return a Series whose index is the original column labels.
- 1 / 'columns': reduce the columns, return a Series whose index is the original index.
- None: reduce all axes, return a scalar.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar.

**bool\_only** [boolean, default None] Include only boolean columns. If None, will attempt to use everything, then use only boolean data. Not implemented for Series.

\*\*kwargs [any, default None] Additional keywords have no effect but might be accepted for compatibility with NumPy.

### **Returns**

any [scalar or Series (if level specified)]

See also:

pandas.DataFrame.all Return whether all elements are True.

# **Examples**

# **Series**

For Series input, the output is a scalar indicating whether any element is True.

```
>>> pd.Series([True, False]).any()
True
```

## **DataFrame**

Whether each column contains at least one True element (the default).

```
>>> df = pd.DataFrame({"A": [1, 2], "B": [0, 2], "C": [0, 0]})
>>> df

A B C
0 1 0 0
1 2 2 0
```

```
>>> df.any()
A True
B True
C False
dtype: bool
```

Aggregating over the columns.

```
>>> df = pd.DataFrame({"A": [True, False], "B": [1, 2]})
>>> df

A B

True 1

False 2
```

```
>>> df.any(axis='columns')
0 True
1 True
dtype: bool
```

```
>>> df = pd.DataFrame({"A": [True, False], "B": [1, 0]})
>>> df

A B

0 True 1

1 False 0
```

```
>>> df.any(axis='columns')

0 True

1 False
dtype: bool
```

Aggregating over the entire DataFrame with axis=None.

```
>>> df.any(axis=None)
True
```

any for an empty DataFrame is an empty Series.

```
>>> pd.DataFrame([]).any()
Series([], dtype: bool)
```

append (to\_append, ignore\_index=False, verify\_integrity=False)

Concatenate two or more Series.

## **Parameters**

to\_append [Series or list/tuple of Series]

**ignore\_index** [boolean, default False] If True, do not use the index labels.

New in version 0.19.0.

**verify\_integrity** [boolean, default False] If True, raise Exception on creating index with duplicates

#### Returns

appended [Series]

See also:

pandas.concat General function to concatenate DataFrame, Series or Panel objects

# **Notes**

Iteratively appending to a Series can be more computationally intensive than a single concatenate. A better solution is to append values to a list and then concatenate the list with the original Series all at once.

# **Examples**

```
>>> s1 = pd.Series([1, 2, 3])
>>> s2 = pd.Series([4, 5, 6])
>>> s3 = pd.Series([4, 5, 6], index=[3, 4, 5])
>>> s1.append(s2)
0
    1
1
     2
2
     3
0
     4
1
     5
2
     6
dtype: int64
```

With *ignore\_index* set to True:

With verify\_integrity set to True:

```
>>> s1.append(s2, verify_integrity=True)
Traceback (most recent call last):
...
ValueError: Indexes have overlapping values: [0, 1, 2]
```

```
apply (func, convert_dtype=True, args=(), **kwds)
```

Invoke function on values of Series. Can be ufunc (a NumPy function that applies to the entire Series) or a Python function that only works on single values

### **Parameters**

```
func [function]
```

**convert\_dtype** [boolean, default True] Try to find better dtype for elementwise function results. If False, leave as dtype=object

args [tuple] Positional arguments to pass to function in addition to the value

Additional keyword arguments will be passed as keywords to the function

#### Returns

y [Series or DataFrame if func returns a Series]

See also:

Series.map For element-wise operations

**Series.agg** only perform aggregating type operations

Series.transform only perform transformating type operations

# **Examples**

Create a series with typical summer temperatures for each city.

Square the values by defining a function and passing it as an argument to apply ().

```
>>> def square(x):
... return x**2
>>> series.apply(square)
London 400
New York 441
Helsinki 144
dtype: int64
```

Square the values by passing an anonymous function as an argument to apply ().

```
>>> series.apply(lambda x: x**2)
London 400
New York 441
Helsinki 144
dtype: int64
```

Define a custom function that needs additional positional arguments and pass these additional arguments using the args keyword.

```
>>> def subtract_custom_value(x, custom_value):
... return x-custom_value
```

```
>>> series.apply(subtract_custom_value, args=(5,))
London 15
New York 16
Helsinki 7
dtype: int64
```

Define a custom function that takes keyword arguments and pass these arguments to apply.

```
>>> series.apply(add_custom_values, june=30, july=20, august=25)
London 95
New York 96
Helsinki 87
dtype: int64
```

Use a function from the Numpy library.

# argmax (\*\*kwargs)

Deprecated since version 0.21.0:will be corrected to return the positional maximum in the future. Use 'series.values.argmax' to get the position of the maximum now.

Return the row label of the maximum value.

If multiple values equal the maximum, the first row label with that value is returned.

# 'argmax' is deprecated, use 'idxmax' instead. The behavior of 'argmax' Parameters

**skipna** [boolean, default True] Exclude NA/null values. If the entire Series is NA, the result will be NA.

**axis** [int, default 0] For compatibility with DataFrame.idxmax. Redundant for application on Series.

\*args, \*\*kwargs Additional keywors have no effect but might be accepted for compatibility with NumPy.

## **Returns**

idxmax [Index of maximum of values.]

# Raises

ValueError If the Series is empty.

See also:

numpy.argmax Return indices of the maximum values along the given axis.

**DataFrame.idxmax** Return index of first occurrence of maximum over requested axis.

Series.idxmin Return index label of the first occurrence of minimum of values.

#### **Notes**

This method is the Series version of ndarray.argmax. This method returns the label of the maximum, while ndarray.argmax returns the position. To get the position, use series.values.argmax().

# **Examples**

```
>>> s.idxmax()
'C'
```

If *skipna* is False and there is an NA value in the data, the function returns nan.

```
>>> s.idxmax(skipna=False)
nan
```

```
argmin(**kwargs)
```

Deprecated since version 0.21.0:will be corrected to return the positional minimum in the future. Use 'series.values.argmin' to get the position of the minimum now.

Return the row label of the minimum value.

If multiple values equal the minimum, the first row label with that value is returned.

# 'argmin' is deprecated, use 'idxmin' instead. The behavior of 'argmin' Parameters

**skipna** [boolean, default True] Exclude NA/null values. If the entire Series is NA, the result will be NA.

**axis** [int, default 0] For compatibility with DataFrame.idxmin. Redundant for application on Series.

\*args, \*\*kwargs Additional keywors have no effect but might be accepted for compatibility with NumPy.

## Returns

**idxmin** [Index of minimum of values.]

## **Raises**

**ValueError** If the Series is empty.

See also:

numpy.argmin Return indices of the minimum values along the given axis.

DataFrame.idxmin Return index of first occurrence of minimum over requested axis.

Series.idxmax Return index label of the first occurrence of maximum of values.

#### **Notes**

This method is the Series version of ndarray.argmin. This method returns the label of the minimum, while ndarray.argmin returns the position. To get the position, use series.values.argmin().

# **Examples**

```
>>> s = pd.Series(data=[1, None, 4, 1],
... index=['A','B','C','D'])
>>> s
A 1.0
B NaN
C 4.0
D 1.0
dtype: float64
```

```
>>> s.idxmin()
'A'
```

If *skipna* is False and there is an NA value in the data, the function returns nan.

```
>>> s.idxmin(skipna=False)
nan
```

```
argsort (axis=0, kind='quicksort', order=None)
```

Overrides ndarray.argsort. Argsorts the value, omitting NA/null values, and places the result in the same locations as the non-NA values

# **Parameters**

```
    axis [int (can only be zero)]
    kind [{'mergesort', 'quicksort', 'heapsort'}, default 'quicksort'] Choice of sorting algorithm. See np.sort for more information. 'mergesort' is the only stable algorithm
    order [ignored]
```

# Returns

argsorted [Series, with -1 indicated where nan values are present]

## See also:

```
numpy.ndarray.argsort
```

## as\_blocks (copy=True)

Convert the frame to a dict of dtype -> Constructor Types that each has a homogeneous dtype.

Deprecated since version 0.21.0.

NOTE: the dtypes of the blocks WILL BE PRESERVED HERE (unlike in as\_matrix)

#### **Parameters**

**copy** [boolean, default True]

#### Returns

values [a dict of dtype -> Constructor Types]

### as\_matrix(columns=None)

Convert the frame to its Numpy-array representation.

Deprecated since version 0.23.0: Use DataFrame.values() instead.

#### **Parameters**

**columns:** list, optional, default:None If None, return all columns, otherwise, returns specified columns.

#### Returns

**values** [ndarray] If the caller is heterogeneous and contains booleans or objects, the result will be of dtype=object. See Notes.

#### See also:

pandas.DataFrame.values

### **Notes**

Return is NOT a Numpy-matrix, rather, a Numpy-array.

The dtype will be a lower-common-denominator dtype (implicit upcasting); that is to say if the dtypes (even of numeric types) are mixed, the one that accommodates all will be chosen. Use this with care if you are not dealing with the blocks.

e.g. If the dtypes are float16 and float32, dtype will be upcast to float32. If dtypes are int32 and uint8, dtype will be upcase to int32. By numpy.find\_common\_type convention, mixing int64 and uint64 will result in a flot64 dtype.

This method is provided for backwards compatibility. Generally, it is recommended to use '.values'.

asfreq(freq, method=None, how=None, normalize=False, fill\_value=None)

Convert TimeSeries to specified frequency.

Optionally provide filling method to pad/backfill missing values.

Returns the original data conformed to a new index with the specified frequency. resample is more appropriate if an operation, such as summarization, is necessary to represent the data at the new frequency.

## **Parameters**

**freq** [DateOffset object, or string]

**method** [{'backfill'/'bfill', 'pad'/'ffill'}, default None] Method to use for filling holes in reindexed Series (note this does not fill NaNs that already were present):

- 'pad' / 'ffill': propagate last valid observation forward to next valid
- 'backfill' / 'bfill': use NEXT valid observation to fill

**how** [{'start', 'end'}, default end] For PeriodIndex only, see PeriodIndex.asfreq

normalize [bool, default False] Whether to reset output index to midnight

**fill\_value: scalar, optional** Value to use for missing values, applied during upsampling (note this does not fill NaNs that already were present).

New in version 0.20.0.

### Returns

**converted** [type of caller]

#### See also:

reindex

### **Notes**

To learn more about the frequency strings, please see this link.

# **Examples**

Start by creating a series with 4 one minute timestamps.

Upsample the series into 30 second bins.

Upsample again, providing a fill value.

Upsample again, providing a method.

```
>>> df.asfreq(freq='30S', method='bfill')

S

2000-01-01 00:00:00 0.0

2000-01-01 00:00:30 NaN

2000-01-01 00:01:00 NaN

2000-01-01 00:01:30 2.0

2000-01-01 00:02:00 2.0

2000-01-01 00:02:30 3.0

2000-01-01 00:03:00 3.0
```

## asobject

Return object Series which contains boxed values.

Deprecated since version 0.23.0: Use astype (object) instead.

this is an internal non-public method

```
asof (where, subset=None)
```

The last row without any NaN is taken (or the last row without NaN considering only the subset of columns in the case of a DataFrame)

New in version 0.19.0: For DataFrame

If there is no good value, NaN is returned for a Series a Series of NaN values for a DataFrame

#### **Parameters**

```
where [date or array of dates]
```

**subset** [string or list of strings, default None] if not None use these columns for NaN propagation

## Returns

# where is scalar

- value or NaN if input is Series
- Series if input is DataFrame

where is Index: same shape object as input

## See also:

```
merge_asof
```

# Notes

Dates are assumed to be sorted Raises if this is not the case

## astype(\*\*kwargs)

Cast a pandas object to a specified dtype dtype.

## **Parameters**

**dtype** [data type, or dict of column name -> data type] Use a numpy.dtype or Python type to cast entire pandas object to the same type. Alternatively, use {col: dtype, ...}, where col is a column label and dtype is a numpy.dtype or Python type to cast one or more of the DataFrame's columns to column-specific types.

**copy** [bool, default True.] Return a copy when copy=True (be very careful setting copy=False as changes to values then may propagate to other pandas objects).

**errors** [{ 'raise', 'ignore'}, default 'raise'.] Control raising of exceptions on invalid data for provided dtype.

- raise: allow exceptions to be raised
- ignore: suppress exceptions. On error return original object

New in version 0.20.0.

raise\_on\_error [raise on invalid input] Deprecated since version 0.20.0: Use errors
instead

**kwargs** [keyword arguments to pass on to the constructor]

## **Returns**

casted [type of caller]

See also:

pandas.to\_datetime Convert argument to datetime.

pandas.to\_timedelta Convert argument to timedelta.

pandas.to\_numeric Convert argument to a numeric type.

numpy.ndarray.astype Cast a numpy array to a specified type.

# **Examples**

```
>>> ser = pd.Series([1, 2], dtype='int32')
>>> ser
0    1
1    2
dtype: int32
>>> ser.astype('int64')
0    1
1    2
dtype: int64
```

### Convert to categorical type:

```
>>> ser.astype('category')
0 1
1 2
dtype: category
Categories (2, int64): [1, 2]
```

Convert to ordered categorical type with custom ordering:

Note that using copy=False and changing data on a new pandas object may propagate changes:

```
>>> s1 = pd.Series([1,2])
>>> s2 = s1.astype('int64', copy=False)
>>> s2[0] = 10
>>> s1  # note that s1[0] has changed too
0    10
1    2
dtype: int64
```

at

Access a single value for a row/column label pair.

Similar to loc, in that both provide label-based lookups. Use at if you only need to get or set a single value in a DataFrame or Series.

#### Raises

KeyError When label does not exist in DataFrame

See also:

DataFrame.iat Access a single value for a row/column pair by integer position

DataFrame.loc Access a group of rows and columns by label(s)

Series.at Access a single value using a label

# **Examples**

Get value at specified row/column pair

```
>>> df.at[4, 'B']
2
```

Set value at specified row/column pair

```
>>> df.at[4, 'B'] = 10
>>> df.at[4, 'B']
10
```

Get value within a Series

```
>>> df.loc[5].at['B']
4
```

at\_time (time, asof=False)

Select values at particular time of day (e.g. 9:30AM).

## **Parameters**

time [datetime.time or string]

#### Returns

```
values_at_time [type of caller]
```

### **Raises**

TypeError If the index is not a DatetimeIndex

#### See also:

between\_time Select values between particular times of the day

first Select initial periods of time series based on a date offset

last Select final periods of time series based on a date offset

DatetimeIndex.indexer\_at\_time Get just the index locations for values at particular time of
the day

# **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='12H')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 00:00:00 1
2018-04-09 12:00:00 2
2018-04-10 00:00:00 3
2018-04-10 12:00:00 4
```

```
>>> ts.at_time('12:00')

A
2018-04-09 12:00:00 2
2018-04-10 12:00:00 4
```

# autocorr (lag=1)

Lag-N autocorrelation

### **Parameters**

lag [int, default 1] Number of lags to apply before performing autocorrelation.

# Returns

autocorr [float]

# axes

Return a list of the row axis labels

### base

return the base object if the memory of the underlying data is shared

between (left, right, inclusive=True)

Return boolean Series equivalent to left <= series <= right.

This function returns a boolean vector containing *True* wherever the corresponding Series element is between the boundary values *left* and *right*. NA values are treated as *False*.

## **Parameters**

**left** [scalar] Left boundary.

```
right [scalar] Right boundary.
```

inclusive [bool, default True] Include boundaries.

### Returns

Series Each element will be a boolean.

See also:

```
pandas.Series.gt Greater than of series and other
pandas.Series.lt Less than of series and other
```

#### **Notes**

This function is equivalent to (left <= ser) & (ser <= right)

# **Examples**

```
>>> s = pd.Series([2, 0, 4, 8, np.nan])
```

Boundary values are included by default:

```
>>> s.between(1, 4)

0    True

1    False

2    True

3    False

4    False

dtype: bool
```

With *inclusive* set to False boundary values are excluded:

```
>>> s.between(1, 4, inclusive=False)

0    True

1    False

2    False

3    False

4    False
dtype: bool
```

left and right can be any scalar value:

```
>>> s = pd.Series(['Alice', 'Bob', 'Carol', 'Eve'])
>>> s.between('Anna', 'Daniel')
0    False
1    True
2    True
3    False
dtype: bool
```

 $\textbf{between\_time} \ (\textit{start\_time}, \textit{end\_time}, \textit{include\_start=True}, \textit{include\_end=True})$ 

Select values between particular times of the day (e.g., 9:00-9:30 AM).

By setting start\_time to be later than end\_time, you can get the times that are *not* between the two times.

### **Parameters**

```
start_time [datetime.time or string]
end_time [datetime.time or string]
include_start [boolean, default True]
```

include\_end [boolean, default True]

#### Returns

values\_between\_time [type of caller]

### **Raises**

TypeError If the index is not a DatetimeIndex

#### See also:

at\_time Select values at a particular time of the day

first Select initial periods of time series based on a date offset

last Select final periods of time series based on a date offset

DatetimeIndex.indexer\_between\_time Get just the index locations for values between particular times of the day

# **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='1D20min')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 00:00:00 1
2018-04-10 00:20:00 2
2018-04-11 00:40:00 3
2018-04-12 01:00:00 4
```

```
>>> ts.between_time('0:15', '0:45')

A
2018-04-10 00:20:00 2
2018-04-11 00:40:00 3
```

You get the times that are *not* between two times by setting start\_time later than end\_time:

```
>>> ts.between_time('0:45', '0:15')

A
2018-04-09 00:00:00 1
2018-04-12 01:00:00 4
```

**bfill** (axis=None, inplace=False, limit=None, downcast=None)

Synonym for DataFrame.fillna (method='bfill')

### blocks

Internal property, property synonym for as\_blocks()

Deprecated since version 0.21.0.

#### bool()

Return the bool of a single element PandasObject.

This must be a boolean scalar value, either True or False. Raise a ValueError if the PandasObject does not have exactly 1 element, or that element is not boolean

```
calcfp (*args, **kwargs)
```

Helper function to map FP calculation through the series

#### cat

```
alias of pandas.core.arrays.categorical.CategoricalAccessor
```

```
clip (lower=None, upper=None, axis=None, inplace=False, *args, **kwargs)
```

Trim values at input threshold(s).

Assigns values outside boundary to boundary values. Thresholds can be singular values or array like, and in the latter case the clipping is performed element-wise in the specified axis.

### **Parameters**

**lower** [float or array\_like, default None] Minimum threshold value. All values below this threshold will be set to it.

**upper** [float or array\_like, default None] Maximum threshold value. All values above this threshold will be set to it.

axis [int or string axis name, optional] Align object with lower and upper along the given axis.

**inplace** [boolean, default False] Whether to perform the operation in place on the data.

New in version 0.21.0.

\*args, \*\*kwargs Additional keywords have no effect but might be accepted for compatibility with numpy.

## **Returns**

**Series or DataFrame** Same type as calling object with the values outside the clip boundaries replaced

## See also:

```
clip_lower Clip values below specified threshold(s).
```

*clip\_upper* Clip values above specified threshold(s).

# **Examples**

```
>>> data = {'col_0': [9, -3, 0, -1, 5], 'col_1': [-2, -7, 6, 8, -5]}
>>> df = pd.DataFrame(data)
>>> df
   col_0 col_1
0
       9
             -2
1
      -3
2
       0
              6
3
      -1
              8
4
             -5
```

Clips per column using lower and upper thresholds:

```
>>> df.clip(-4, 6)
   col_0 col_1
0
        6
               -2
       -3
               -4
1
2
        0
                6
3
       -1
                6
4
        5
               -4
```

Clips using specific lower and upper thresholds per column element:

```
>>> df.clip(t, t + 4, axis=0)
   col_0 col_1
0
       6
               2
       -3
1
               -4
2
       0
                3
3
       6
                8
4
        5
                3
```

### clip\_lower (threshold, axis=None, inplace=False)

Return copy of the input with values below a threshold truncated.

### **Parameters**

**threshold** [numeric or array-like] Minimum value allowed. All values below threshold will be set to this value.

- float : every value is compared to *threshold*.
- array-like: The shape of *threshold* should match the object it's compared to. When *self* is a Series, *threshold* should be the length. When *self* is a DataFrame, *threshold* should 2-D and the same shape as *self* for axis=None, or 1-D and the same length as the axis being compared.

**axis** [{0 or 'index', 1 or 'columns'}, default 0] Align *self* with *threshold* along the given axis.

**inplace** [boolean, default False] Whether to perform the operation in place on the data.

New in version 0.21.0.

## Returns

clipped [same type as input]

See also:

Series.clip Return copy of input with values below and above thresholds truncated.

Series.clip\_upper Return copy of input with values above threshold truncated.

# **Examples**

Series single threshold clipping:

```
>>> s = pd.Series([5, 6, 7, 8, 9])
>>> s.clip_lower(8)
0     8
1     8
2     8
3     8
4     9
dtype: int64
```

Series clipping element-wise using an array of thresholds. *threshold* should be the same length as the Series.

```
>>> elemwise_thresholds = [4, 8, 7, 2, 5]
>>> s.clip_lower(elemwise_thresholds)
0    5
1    8
2    7
3    8
4    9
dtype: int64
```

DataFrames can be compared to a scalar.

```
>>> df = pd.DataFrame({"A": [1, 3, 5], "B": [2, 4, 6]})
>>> df

A B
0 1 2
1 3 4
2 5 6
```

```
>>> df.clip_lower(3)

A B

0 3 3

1 3 4

2 5 6
```

Or to an array of values. By default, *threshold* should be the same shape as the DataFrame.

```
>>> df.clip_lower(np.array([[3, 4], [2, 2], [6, 2]]))

A B

0 3 4

1 3 4

2 6 6
```

Control how *threshold* is broadcast with *axis*. In this case *threshold* should be the same length as the axis specified by *axis*.

```
>>> df.clip_lower(np.array([3, 3, 5]), axis='index')

A B
0 3 3
1 3 4
2 5 6
```

```
>>> df.clip_lower(np.array([4, 5]), axis='columns')

A B
0 4 5
1 4 5
2 5 6
```

### clip\_upper (threshold, axis=None, inplace=False)

Return copy of input with values above given value(s) truncated.

### **Parameters**

```
threshold [float or array_like]
```

axis [int or string axis name, optional] Align object with threshold along the given axis.

inplace [boolean, default False] Whether to perform the operation in place on the data

New in version 0.21.0.

#### Returns

**clipped** [same type as input]

## See also:

clip

# combine (other, func, fill\_value=nan)

Perform elementwise binary operation on two Series using given function with optional fill value when an index is missing from one Series or the other

### **Parameters**

```
other [Series or scalar value]
```

func [function] Function that takes two scalars as inputs and return a scalar

fill\_value [scalar value]

# Returns

result [Series]

See also:

Series.combine\_first Combine Series values, choosing the calling Series's values first

# **Examples**

## combine\_first (other)

Combine Series values, choosing the calling Series's values first. Result index will be the union of the two indexes

### **Parameters**

other [Series]

Returns

combined [Series]

See also:

Series.combine Perform elementwise operation on two Series using a given function

## **Examples**

```
>>> s1 = pd.Series([1, np.nan])
>>> s2 = pd.Series([3, 4])
>>> s1.combine_first(s2)
0     1.0
1     4.0
dtype: float64
```

compound (axis=None, skipna=None, level=None)

Return the compound percentage of the values for the requested axis

#### **Parameters**

```
axis [\{index (0)\}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### **Returns**

**compounded** [scalar or Series (if level specified)]

```
compress (condition, *args, **kwargs)
```

Return selected slices of an array along given axis as a Series

### See also:

```
numpy.ndarray.compress
```

```
consolidate(inplace=False)
```

Compute NDFrame with "consolidated" internals (data of each dtype grouped together in a single ndarray).

Deprecated since version 0.20.0: Consolidate will be an internal implementation only.

Attempt to infer better dtype for object columns.

Deprecated since version 0.21.0.

#### **Parameters**

**convert\_dates** [boolean, default True] If True, convert to date where possible. If 'coerce', force conversion, with unconvertible values becoming NaT.

**convert\_numeric** [boolean, default False] If True, attempt to coerce to numbers (including strings), with unconvertible values becoming NaN.

**convert\_timedeltas** [boolean, default True] If True, convert to timedelta where possible. If 'coerce', force conversion, with unconvertible values becoming NaT.

**copy** [boolean, default True] If True, return a copy even if no copy is necessary (e.g. no conversion was done). Note: This is meant for internal use, and should not be confused with inplace.

## **Returns**

converted [same as input object]

#### See also:

```
pandas.to_datetime Convert argument to datetime.
```

pandas.to\_timedelta Convert argument to timedelta.

pandas.to\_numeric Return a fixed frequency timedelta index, with day as the default.

```
copy (deep=True)
```

Make a copy of this object's indices and data.

When deep=True (default), a new object will be created with a copy of the calling object's data and indices. Modifications to the data or indices of the copy will not be reflected in the original object (see notes below).

When deep=False, a new object will be created without copying the calling object's data or index (only references to the data and index are copied). Any changes to the data of the original will be reflected in the shallow copy (and vice versa).

# **Parameters**

**deep** [bool, default True] Make a deep copy, including a copy of the data and the indices. With deep=False neither the indices nor the data are copied.

#### Returns

**copy** [Series, DataFrame or Panel] Object type matches caller.

## **Notes**

When deep=True, data is copied but actual Python objects will not be copied recursively, only the reference to the object. This is in contrast to *copy.deepcopy* in the Standard Library, which recursively copies object data (see examples below).

While Index objects are copied when deep=True, the underlying numpy array is not copied for performance reasons. Since Index is immutable, the underlying data can be safely shared and a copy is not needed.

## **Examples**

```
>>> s = pd.Series([1, 2], index=["a", "b"])
>>> s
a    1
b    2
dtype: int64
```

### Shallow copy versus default (deep) copy:

```
>>> s = pd.Series([1, 2], index=["a", "b"])
>>> deep = s.copy()
>>> shallow = s.copy(deep=False)
```

Shallow copy shares data and index with original.

```
>>> s is shallow
False
>>> s.values is shallow.values and s.index is shallow.index
True
```

Deep copy has own copy of data and index.

```
>>> s is deep
False
>>> s.values is deep.values or s.index is deep.index
False
```

Updates to the data shared by shallow copy and original is reflected in both; deep copy remains unchanged.

```
>>> s[0] = 3
>>> shallow[1] = 4
>>> s
а
    3
dtype: int64
>>> shallow
    3
а
b
    4
dtype: int64
>>> deep
  1
а
b
    2
dtype: int64
```

Note that when copying an object containing Python objects, a deep copy will copy the data, but will not do so recursively. Updating a nested data object will be reflected in the deep copy.

```
corr (other, method='pearson', min_periods=None)
```

Compute correlation with *other* Series, excluding missing values

#### **Parameters**

other [Series]

**method** [{'pearson', 'kendall', 'spearman'}]

- pearson: standard correlation coefficient
- kendall : Kendall Tau correlation coefficient
- spearman: Spearman rank correlation

min\_periods [int, optional] Minimum number of observations needed to have a valid result

### **Returns**

correlation [float]

count (level=None)

Return number of non-NA/null observations in the Series

#### **Parameters**

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a smaller Series

#### Returns

**nobs** [int or Series (if level specified)]

cov (other, min\_periods=None)

Compute covariance with Series, excluding missing values

### **Parameters**

other [Series]

min\_periods [int, optional] Minimum number of observations needed to have a valid result

# Returns

covariance [float]

Normalized by N-1 (unbiased estimator).

cummax (axis=None, skipna=True, \*args, \*\*kwargs)

Return cumulative maximum over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative maximum.

### **Parameters**

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

#### Returns

cummax [scalar or Series]

See also:

pandas.core.window.Expanding.max Similar functionality but ignores NaN values.

**Series.max** Return the maximum over Series axis.

Series.cummax Return cumulative maximum over Series axis.

Series.cummin Return cumulative minimum over Series axis.

Series.cumsum Return cumulative sum over Series axis.

Series.cumprod Return cumulative product over Series axis.

# **Examples**

### Series

By default, NA values are ignored.

```
>>> s.cummax()
0 2.0
1 NaN
2 5.0
3 5.0
4 5.0
dtype: float64
```

To include NA values in the operation, use skipna=False

```
>>> s.cummax(skipna=False)
0 2.0
1 NaN
2 NaN
3 NaN
4 NaN
dtype: float64
```

### **DataFrame**

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```
1 3.0 NaN
2 1.0 0.0
```

By default, iterates over rows and finds the maximum in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cummax()

A B

0 2.0 1.0

1 3.0 NaN

2 3.0 1.0
```

To iterate over columns and find the maximum in each row, use axis=1

```
>>> df.cummax(axis=1)

A B
0 2.0 2.0
1 3.0 NaN
2 1.0 1.0
```

```
cummin (axis=None, skipna=True, *args, **kwargs)
```

Return cumulative minimum over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative minimum.

### **Parameters**

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

### Returns

cummin [scalar or Series]

See also:

pandas.core.window.Expanding.min Similar functionality but ignores NaN values.

Series.min Return the minimum over Series axis.

Series.cummax Return cumulative maximum over Series axis.

Series.cummin Return cumulative minimum over Series axis.

Series.cumsum Return cumulative sum over Series axis.

Series.cumprod Return cumulative product over Series axis.

### **Examples**

### **Series**

By default, NA values are ignored.

```
>>> s.cummin()
0 2.0
1 NaN
2 2.0
3 -1.0
4 -1.0
dtype: float64
```

To include NA values in the operation, use skipna=False

```
>>> s.cummin(skipna=False)
0 2.0
1 NaN
2 NaN
3 NaN
4 NaN
dtype: float64
```

## **DataFrame**

```
>>> df = pd.DataFrame([[2.0, 1.0],
... [3.0, np.nan],
... [1.0, 0.0]],
... columns=list('AB'))
>>> df
A B
0 2.0 1.0
1 3.0 NaN
2 1.0 0.0
```

By default, iterates over rows and finds the minimum in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cummin()

A B

0 2.0 1.0
1 2.0 NaN
2 1.0 0.0
```

To iterate over columns and find the minimum in each row, use axis=1

```
>>> df.cummin(axis=1)

A B
0 2.0 1.0
1 3.0 NaN
2 1.0 0.0
```

```
cumprod (axis=None, skipna=True, *args, **kwargs)
```

Return cumulative product over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative product.

### **Parameters**

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

#### Returns

```
cumprod [scalar or Series]
```

### See also:

pandas.core.window.Expanding.prod Similar functionality but ignores NaN values.

Series.prod Return the product over Series axis.

Series.cummax Return cumulative maximum over Series axis.

Series.cummin Return cumulative minimum over Series axis.

Series.cumsum Return cumulative sum over Series axis.

Series.cumprod Return cumulative product over Series axis.

## **Examples**

#### **Series**

# By default, NA values are ignored.

To include NA values in the operation, use skipna=False

```
>>> s.cumprod(skipna=False)
0 2.0
1 NaN
2 NaN
3 NaN
4 NaN
dtype: float64
```

#### **DataFrame**

By default, iterates over rows and finds the product in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cumprod()

A B

0 2.0 1.0

1 6.0 NaN

2 6.0 0.0
```

To iterate over columns and find the product in each row, use axis=1

```
>>> df.cumprod(axis=1)

A B
0 2.0 2.0
1 3.0 NaN
2 1.0 0.0
```

cumsum (axis=None, skipna=True, \*args, \*\*kwargs)

Return cumulative sum over a DataFrame or Series axis.

Returns a DataFrame or Series of the same size containing the cumulative sum.

### **Parameters**

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis. 0 is equivalent to None or 'index'.

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA.

\*args, \*\*kwargs: Additional keywords have no effect but might be accepted for compatibility with NumPy.

#### Returns

**cumsum** [scalar or Series]

See also:

pandas.core.window.Expanding.sum Similar functionality but ignores NaN values.

Series.sum Return the sum over Series axis.

Series.cummax Return cumulative maximum over Series axis.

Series.cummin Return cumulative minimum over Series axis.

Series.cumsum Return cumulative sum over Series axis.

Series.cumprod Return cumulative product over Series axis.

## **Examples**

## **Series**

By default, NA values are ignored.

```
>>> s.cumsum()
0 2.0
1 NaN
2 7.0
3 6.0
4 6.0
dtype: float64
```

To include NA values in the operation, use skipna=False

```
>>> s.cumsum(skipna=False)
0 2.0
1 NaN
2 NaN
3 NaN
4 NaN
dtype: float64
```

# **DataFrame**

```
>>> df = pd.DataFrame([[2.0, 1.0],
                        [3.0, np.nan],
. . .
                        [1.0, 0.0]],
. . .
                        columns=list('AB'))
. . .
>>> df
        В
  2.0
        1.0
  3.0
        NaN
2
  1.0
        0.0
```

By default, iterates over rows and finds the sum in each column. This is equivalent to axis=None or axis='index'.

```
>>> df.cumsum()

A B

0 2.0 1.0

1 5.0 NaN

2 6.0 1.0
```

To iterate over columns and find the sum in each row, use axis=1

```
>>> df.cumsum(axis=1)

A B
0 2.0 3.0
1 3.0 NaN
2 1.0 1.0
```

#### data

return the data pointer of the underlying data

**describe** (percentiles=None, include=None, exclude=None)

Generates descriptive statistics that summarize the central tendency, dispersion and shape of a dataset's distribution, excluding NaN values.

Analyzes both numeric and object series, as well as DataFrame column sets of mixed data types. The output will vary depending on what is provided. Refer to the notes below for more detail.

#### **Parameters**

**percentiles** [list-like of numbers, optional] The percentiles to include in the output. All should fall between 0 and 1. The default is [.25, .5, .75], which returns the 25th, 50th, and 75th percentiles.

**include** ['all', list-like of dtypes or None (default), optional] A white list of data types to include in the result. Ignored for Series. Here are the options:

- 'all' : All columns of the input will be included in the output.
- A list-like of dtypes: Limits the results to the provided data types. To limit the result to numeric types submit numpy.number. To limit it instead to object columns submit the numpy.object data type. Strings can also be used in the style of select\_dtypes (e.g. df.describe(include=['O'])). To select pandas categorical columns, use 'category'
- None (default): The result will include all numeric columns.

**exclude** [list-like of dtypes or None (default), optional,] A black list of data types to omit from the result. Ignored for Series. Here are the options:

- A list-like of dtypes: Excludes the provided data types from the result. To exclude numeric types submit numpy.number. To exclude object columns submit the data type numpy.object. Strings can also be used in the style of select\_dtypes (e.g. df.describe(include=['0'])). To exclude pandas categorical columns, use 'category'
- None (default): The result will exclude nothing.

#### Returns

### summary: Series/DataFrame of summary statistics

# See also:

DataFrame.count, DataFrame.max, DataFrame.min, DataFrame.mean, DataFrame.std, DataFrame.select\_dtypes

### **Notes**

For numeric data, the result's index will include count, mean, std, min, max as well as lower, 50 and upper percentiles. By default the lower percentile is 25 and the upper percentile is 75. The 50 percentile is the same as the median.

For object data (e.g. strings or timestamps), the result's index will include count, unique, top, and freq. The top is the most common value. The freq is the most common value's frequency. Timestamps also include the first and last items.

If multiple object values have the highest count, then the count and top results will be arbitrarily chosen from among those with the highest count.

For mixed data types provided via a DataFrame, the default is to return only an analysis of numeric columns. If the dataframe consists only of object and categorical data without any numeric columns, the default is to return an analysis of both the object and categorical columns. If include='all' is provided as an option, the result will include a union of attributes of each type.

The *include* and *exclude* parameters can be used to limit which columns in a DataFrame are analyzed for the output. The parameters are ignored when analyzing a Series.

## **Examples**

Describing a numeric Series.

```
>>> s = pd.Series([1, 2, 3])
>>> s.describe()
count
         3.0
         2.0
mean
std
         1.0
         1.0
2.5%
         1.5
50%
         2.0
75%
         2.5
         3.0
max
```

Describing a categorical Series.

```
>>> s = pd.Series(['a', 'a', 'b', 'c'])
>>> s.describe()
count    4
unique    3
top     a
freq    2
dtype: object
```

Describing a timestamp Series.

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```
freq 2
first 2000-01-01 00:00:00
last 2010-01-01 00:00:00
dtype: object
```

Describing a DataFrame. By default only numeric fields are returned.

```
>>> df = pd.DataFrame({ 'object': ['a', 'b', 'c'],
                         'numeric': [1, 2, 3],
                         'categorical': pd.Categorical(['d','e','f'])
. . .
. . .
>>> df.describe()
       numeric
         3.0
count
          2.0
mean
           1.0
std
           1.0
min
25%
           1.5
50%
           2.0
75%
           2.5
           3.0
max
```

Describing all columns of a DataFrame regardless of data type.

```
>>> df.describe(include='all')
       categorical numeric object
                     3.0
                              3
              3
count
               3
                              3
unique
                      NaN
top
              f
                      NaN
                              С
freq
               1
                      NaN
             NaN
                      2.0
                           NaN
mean
             NaN
                      1.0
                           NaN
std
                      1.0
min
             NaN
                           NaN
25%
             NaN
                      1.5
                           NaN
50%
             NaN
                      2.0
                          NaN
75%
              NaN
                      2.5
                            NaN
              NaN
                      3.0
max
                            NaN
```

Describing a column from a DataFrame by accessing it as an attribute.

```
>>> df.numeric.describe()
count
         3.0
         2.0
mean
         1.0
std
         1.0
min
25%
         1.5
50%
         2.0
75%
         2.5
         3.0
max
Name: numeric, dtype: float64
```

Including only numeric columns in a DataFrame description.

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```
mean 2.0
std 1.0
min 1.0
25% 1.5
50% 2.0
75% 2.5
max 3.0
```

Including only string columns in a DataFrame description.

Including only categorical columns from a DataFrame description.

Excluding numeric columns from a DataFrame description.

Excluding object columns from a DataFrame description.

```
>>> df.describe(exclude=[np.object])
       categorical numeric
               3
                      3.0
count
               3
unique
                      NaN
              f
top
                      NaN
              1
                      NaN
freq
                    2.0
             NaN
std
            NaN
                     1.0
            NaN
                     1.0
min
25%
             NaN
                      1.5
50%
             NaN
                      2.0
75%
                      2.5
              NaN
                      3.0
max
              NaN
```

### diff(periods=1)

First discrete difference of element.

Calculates the difference of a Series element compared with another element in the Series (default is element in previous row).

**Parameters** 

**periods** [int, default 1] Periods to shift for calculating difference, accepts negative values.

### Returns

diffed [Series]

See also:

Series.pct\_change Percent change over given number of periods.

**Series.shift** Shift index by desired number of periods with an optional time freq.

DataFrame.diff First discrete difference of object

# **Examples**

Difference with previous row

```
>>> s = pd.Series([1, 1, 2, 3, 5, 8])
>>> s.diff()
0    NaN
1    0.0
2    1.0
3    1.0
4    2.0
5    3.0
dtype: float64
```

## Difference with 3rd previous row

# Difference with following row

div (other, level=None, fill\_value=None, axis=0)

Floating division of series and other, element-wise (binary operator truediv).

Equivalent to series / other, but with support to substitute a fill\_value for missing data in one of the inputs.

# **Parameters**

other [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

#### Returns

```
result [Series]
```

### See also:

Series.rtruediv

## **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
а
    1.0
b
     1.0
     1.0
С
     NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> h
а
    1.0
b
     NaN
d
     1.0
     NaN
dtype: float64
>>> a.add(b, fill_value=0)
     2.0
     1.0
h
С
     1.0
d
     1.0
e
     NaN
dtype: float64
```

# divide (other, level=None, fill\_value=None, axis=0)

Floating division of series and other, element-wise (binary operator truediv).

Equivalent to series / other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

other [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### Returns

# result [Series]

### See also:

Series.rtruediv

## **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
    1.0
b
    1.0
С
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
    1.0
b
    NaN
d
    1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
а
    1.0
b
    1.0
С
d
     1.0
     NaN
dtype: float64
```

# ${\tt divmod} \ (other, \ level=None, \ fill\_value=None, \ axis=0)$

Integer division and modulo of series and other, element-wise (binary operator *divmod*).

Equivalent to series divmod other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

other [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

# Returns

result [Series]

### See also:

Series.None

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
b
     1.0
С
     1.0
d
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> h
    1.0
b
    NaN
d
    1.0
е
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
b
     1.0
С
     1.0
d
     1.0
    NaN
dtype: float64
```

### dot (other)

Matrix multiplication with DataFrame or inner-product with Series objects. Can also be called using *self* @ *other* in Python >= 3.5.

### **Parameters**

other [Series or DataFrame]

### Returns

```
dot_product [scalar or Series]
```

**drop** (*labels=None*, *axis=0*, *index=None*, *columns=None*, *level=None*, *inplace=False*, *errors='raise'*) Return Series with specified index labels removed.

Remove elements of a Series based on specifying the index labels. When using a multi-index, labels on different levels can be removed by specifying the level.

### **Parameters**

labels [single label or list-like] Index labels to drop.

axis [0, default 0] Redundant for application on Series.

index, columns [None] Redundant for application on Series, but index can be used instead of labels.

New in version 0.21.0.

**level** [int or level name, optional] For MultiIndex, level for which the labels will be removed.

**inplace** [bool, default False] If True, do operation inplace and return None.

**errors** [{'ignore', 'raise'}, default 'raise'] If 'ignore', suppress error and only existing labels are dropped.

## Returns

dropped [pandas.Series]

Raises

**KeyError** If none of the labels are found in the index.

See also:

**Series.reindex** Return only specified index labels of Series.

Series.dropna Return series without null values.

Series.drop\_duplicates Return Series with duplicate values removed.

DataFrame.drop Drop specified labels from rows or columns.

# **Examples**

```
>>> s = pd.Series(data=np.arange(3), index=['A','B','C'])
>>> s
A 0
B 1
C 2
dtype: int64
```

## Drop labels B en C

```
>>> s.drop(labels=['B','C'])
A 0
dtype: int64
```

# Drop 2nd level label in MultiIndex Series

```
>>> midx = pd.MultiIndex(levels=[['lama', 'cow', 'falcon'],
                                ['speed', 'weight', 'length']],
                        labels=[[0, 0, 0, 1, 1, 1, 2, 2, 2],
                                [0, 1, 2, 0, 1, 2, 0, 1, 2]])
>>> s = pd.Series([45, 200, 1.2, 30, 250, 1.5, 320, 1, 0.3],
                 index=midx)
>>> s
                 45.0
lama
       speed
       weight 200.0
                  1.2
       length
                 30.0
COW
       speed
               250.0
       weight
       length
                  1.5
               320.0
falcon speed
                1.0
       weight
       length
                   0.3
dtype: float64
```

```
>>> s.drop(labels='weight', level=1)

lama speed 45.0
    length 1.2

cow speed 30.0
    length 1.5

falcon speed 320.0
```

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```
length 0.3 dtype: float64
```

# drop\_duplicates (keep='first', inplace=False)

Return Series with duplicate values removed.

#### **Parameters**

keep [{'first', 'last', False}, default 'first']

- 'first': Drop duplicates except for the first occurrence.
- 'last': Drop duplicates except for the last occurrence.
- False: Drop all duplicates.

inplace [boolean, default False] If True, performs operation inplace and returns None.

# Returns

```
deduplicated [Series]
```

See also:

Index.drop\_duplicates equivalent method on Index

DataFrame.drop\_duplicates equivalent method on DataFrame

Series.duplicated related method on Series, indicating duplicate Series values.

## **Examples**

Generate an Series with duplicated entries.

```
>>> s = pd.Series(['lama', 'cow', 'lama', 'beetle', 'lama', 'hippo'],
                  name='animal')
>>> s
0
       lama
1
        COW
2
       lama
3
     beetle
4
       lama
5
      hippo
Name: animal, dtype: object
```

With the 'keep' parameter, the selection behaviour of duplicated values can be changed. The value 'first' keeps the first occurrence for each set of duplicated entries. The default value of keep is 'first'.

```
>>> s.drop_duplicates()
0 lama
1 cow
3 beetle
5 hippo
Name: animal, dtype: object
```

The value 'last' for parameter 'keep' keeps the last occurrence for each set of duplicated entries.

The value False for parameter 'keep' discards all sets of duplicated entries. Setting the value of 'inplace' to True performs the operation inplace and returns None.

```
>>> s.drop_duplicates(keep=False, inplace=True)
>>> s
1     cow
3     beetle
5     hippo
Name: animal, dtype: object
```

# dropna (axis=0, inplace=False, \*\*kwargs)

Return a new Series with missing values removed.

See the User Guide for more on which values are considered missing, and how to work with missing data.

#### **Parameters**

```
axis [{0 or 'index'}, default 0] There is only one axis to drop values from.
```

inplace [bool, default False] If True, do operation inplace and return None.

\*\*kwargs Not in use.

### **Returns**

Series Series with NA entries dropped from it.

## See also:

Series.isna Indicate missing values.

Series.notna Indicate existing (non-missing) values.

Series.fillna Replace missing values.

DataFrame.dropna Drop rows or columns which contain NA values.

Index.dropna Drop missing indices.

### **Examples**

```
>>> ser = pd.Series([1., 2., np.nan])
>>> ser
0    1.0
1    2.0
2    NaN
dtype: float64
```

Drop NA values from a Series.

```
>>> ser.dropna()
0 1.0
1 2.0
dtype: float64
```

Keep the Series with valid entries in the same variable.

```
>>> ser.dropna(inplace=True)
>>> ser
0 1.0
1 2.0
dtype: float64
```

Empty strings are not considered NA values. None is considered an NA value.

```
>>> ser = pd.Series([np.NaN, 2, pd.NaT, '', None, 'I stay'])
>>> ser
0
        NaN
1
          2
2
        NaT
3
4
       None
    I stay
dtype: object
>>> ser.dropna()
1
          2
3
5
    I stay
dtype: object
```

dt

alias of pandas.core.indexes.accessors.CombinedDatetimelikeProperties

# dtype

return the dtype object of the underlying data

# dtypes

return the dtype object of the underlying data

## duplicated(keep='first')

Indicate duplicate Series values.

Duplicated values are indicated as True values in the resulting Series. Either all duplicates, all except the first or all except the last occurrence of duplicates can be indicated.

### **Parameters**

```
keep [{'first', 'last', False}, default 'first']
```

- 'first': Mark duplicates as True except for the first occurrence.
- 'last' : Mark duplicates as True except for the last occurrence.
- $\bullet$  False: Mark all duplicates as  ${\tt True}.$

### Returns

pandas.core.series.Series

See also:

 ${\tt pandas.Index.duplicated} \ \ Equivalent\ method\ on\ pandas.Index$ 

pandas.DataFrame.duplicated Equivalent method on pandas.DataFrame
pandas.Series.drop\_duplicates Remove duplicate values from Series

## **Examples**

By default, for each set of duplicated values, the first occurrence is set on False and all others on True:

```
>>> animals = pd.Series(['lama', 'cow', 'lama', 'beetle', 'lama'])
>>> animals.duplicated()
0   False
1   False
2   True
3   False
4   True
dtype: bool
```

## which is equivalent to

```
>>> animals.duplicated(keep='first')

0 False
1 False
2 True
3 False
4 True
dtype: bool
```

By using 'last', the last occurrence of each set of duplicated values is set on False and all others on True:

```
>>> animals.duplicated(keep='last')

0    True

1    False

2    True

3    False

4    False
dtype: bool
```

By setting keep on False, all duplicates are True:

```
>>> animals.duplicated(keep=False)

0    True

1    False

2    True

3    False

4    True

dtype: bool
```

### empty

```
eq (other, level=None, fill_value=None, axis=0)
```

Equal to of series and other, element-wise (binary operator eq).

Equivalent to series == other, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

**other** [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

#### Returns

result [Series]

### See also:

Series.None

## **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
а
    1.0
b
     1.0
     1.0
С
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> h
а
    1.0
    NaN
d
     1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
     1.0
h
С
     1.0
d
     1.0
    NaN
dtype: float64
```

### equals (other)

Determines if two NDFrame objects contain the same elements. NaNs in the same location are considered equal.

```
ewm (com=None, span=None, halflife=None, alpha=None, min_periods=0, adjust=True, ignore_na=False, axis=0)
Provides exponential weighted functions
```

New in version 0.18.0.

# **Parameters**

**alpha** [float, optional] Specify smoothing factor  $\alpha$  directly,  $0 < \alpha \le 1$ 

New in version 0.18.0.

**min\_periods** [int, default 0] Minimum number of observations in window required to have a value (otherwise result is NA).

**adjust** [boolean, default True] Divide by decaying adjustment factor in beginning periods to account for imbalance in relative weightings (viewing EWMA as a moving average)

**ignore\_na** [boolean, default False] Ignore missing values when calculating weights; specify True to reproduce pre-0.15.0 behavior

### **Returns**

a Window sub-classed for the particular operation

#### See also:

rolling Provides rolling window calculations

**expanding** Provides expanding transformations.

#### **Notes**

Exactly one of center of mass, span, half-life, and alpha must be provided. Allowed values and relationship between the parameters are specified in the parameter descriptions above; see the link at the end of this section for a detailed explanation.

When adjust is True (default), weighted averages are calculated using weights (1-alpha)\*\*(n-1), (1-alpha)\*\*(n-2), ..., 1-alpha, 1.

```
When adjust is False, weighted averages are calculated recursively as: weighted_average[0] = arg[0]; weighted_average[i] = (1-alpha)*weighted_average[i-1] + alpha*arg[i].
```

When ignore\_na is False (default), weights are based on absolute positions. For example, the weights of x and y used in calculating the final weighted average of [x, None, y] are (1-alpha)\*\*2 and 1 (if adjust is True), and (1-alpha)\*\*2 and alpha (if adjust is False).

When ignore\_na is True (reproducing pre-0.15.0 behavior), weights are based on relative positions. For example, the weights of x and y used in calculating the final weighted average of [x, None, y] are 1-alpha and 1 (if adjust is True), and 1-alpha and alpha (if adjust is False).

More details can be found at http://pandas.pydata.org/pandas-docs/stable/computation.html# exponentially-weighted-windows

# **Examples**

```
>>> df = DataFrame({'B': [0, 1, 2, np.nan, 4]})
B
0 0.0
1 1.0
2 2.0
3 NaN
4 4.0
```

```
>>> df.ewm(com=0.5).mean()

B
0 0.000000
1 0.750000
2 1.615385
3 1.615385
4 3.670213
```

## **expanding** (*min\_periods=1*, *center=False*, *axis=0*)

Provides expanding transformations.

New in version 0.18.0.

#### **Parameters**

**min\_periods** [int, default 1] Minimum number of observations in window required to have a value (otherwise result is NA).

center [boolean, default False] Set the labels at the center of the window.

axis [int or string, default 0]

### Returns

a Window sub-classed for the particular operation

## See also:

rolling Provides rolling window calculations

ewm Provides exponential weighted functions

### **Notes**

By default, the result is set to the right edge of the window. This can be changed to the center of the window by setting center=True.

## **Examples**

```
>>> df = DataFrame({'B': [0, 1, 2, np.nan, 4]})

B
0 0.0
1 1.0
2 2.0
3 NaN
4 4.0
```

```
>>> df.expanding(2).sum()
B
0 NaN
1 1.0
2 3.0
3 3.0
4 7.0
```

# factorize (sort=False, na\_sentinel=-1)

Encode the object as an enumerated type or categorical variable.

This method is useful for obtaining a numeric representation of an array when all that matters is identifying distinct values. *factorize* is available as both a top-level function pandas.factorize(), and as a method Series.factorize() and Index.factorize().

### **Parameters**

**sort** [boolean, default False] Sort *uniques* and shuffle *labels* to maintain the relationship. **na\_sentinel** [int, default -1] Value to mark "not found".

#### Returns

**labels** [ndarray] An integer ndarray that's an indexer into *uniques*. uniques. take(labels) will have the same values as *values*.

**uniques** [ndarray, Index, or Categorical] The unique valid values. When *values* is Categorical, *uniques* is a Categorical. When *values* is some other pandas object, an *Index* is returned. Otherwise, a 1-D ndarray is returned.

**Note:** Even if there's a missing value in *values*, *uniques* will *not* contain an entry for it.

#### See also:

pandas.cut Discretize continuous-valued array.
pandas.unique Find the unique valuse in an array.

## **Examples**

These examples all show factorize as a top-level method like pd.factorize (values). The results are identical for methods like Series.factorize().

```
>>> labels, uniques = pd.factorize(['b', 'b', 'a', 'c', 'b'])
>>> labels
array([0, 0, 1, 2, 0])
>>> uniques
array(['b', 'a', 'c'], dtype=object)
```

With sort=True, the *uniques* will be sorted, and *labels* will be shuffled so that the relationship is the maintained.

```
>>> labels, uniques = pd.factorize(['b', 'b', 'a', 'c', 'b'], sort=True)
>>> labels
array([1, 1, 0, 2, 1])
>>> uniques
array(['a', 'b', 'c'], dtype=object)
```

Missing values are indicated in *labels* with  $na\_sentinel$  (-1 by default). Note that missing values are never included in *uniques*.

```
>>> labels, uniques = pd.factorize(['b', None, 'a', 'c', 'b'])
>>> labels
array([ 0, -1,  1,  2,  0])
>>> uniques
array(['b', 'a', 'c'], dtype=object)
```

Thus far, we've only factorized lists (which are internally coerced to NumPy arrays). When factorizing pandas objects, the type of *uniques* will differ. For Categoricals, a *Categorical* is returned.

```
>>> cat = pd.Categorical(['a', 'a', 'c'], categories=['a', 'b', 'c'])
>>> labels, uniques = pd.factorize(cat)
>>> labels
array([0, 0, 1])
>>> uniques
[a, c]
Categories (3, object): [a, b, c]
```

Notice that 'b' is in uniques.categories, desipite not being present in cat.values.

For all other pandas objects, an Index of the appropriate type is returned.

```
>>> cat = pd.Series(['a', 'a', 'c'])
>>> labels, uniques = pd.factorize(cat)
>>> labels
array([0, 0, 1])
>>> uniques
Index(['a', 'c'], dtype='object')
```

**ffill** (axis=None, inplace=False, limit=None, downcast=None)

Synonym for DataFrame.fillna (method='ffill')

**fillna** (value=None, method=None, axis=None, inplace=False, limit=None, downcast=None, \*\*kwargs)

Fill NA/NaN values using the specified method

#### **Parameters**

value [scalar, dict, Series, or DataFrame] Value to use to fill holes (e.g. 0), alternately a dict/Series/DataFrame of values specifying which value to use for each index (for a Series) or column (for a DataFrame). (values not in the dict/Series/DataFrame will not be filled). This value cannot be a list.

**method** [{'backfill', 'bfill', 'pad', 'ffill', None}, default None] Method to use for filling holes in reindexed Series pad / ffill: propagate last valid observation forward to next valid backfill / bfill: use NEXT valid observation to fill gap

```
axis [{0 or 'index'}]
```

**inplace** [boolean, default False] If True, fill in place. Note: this will modify any other views on this object, (e.g. a no-copy slice for a column in a DataFrame).

**limit** [int, default None] If method is specified, this is the maximum number of consecutive NaN values to forward/backward fill. In other words, if there is a gap with more than this number of consecutive NaNs, it will only be partially filled. If method is not specified, this is the maximum number of entries along the entire axis where NaNs will be filled. Must be greater than 0 if not None.

**downcast** [dict, default is None] a dict of item->dtype of what to downcast if possible, or the string 'infer' which will try to downcast to an appropriate equal type (e.g. float64 to int64 if possible)

### Returns

filled [Series]

### See also:

interpolate Fill NaN values using interpolation.

```
reindex, asfreq
```

# **Examples**

Replace all NaN elements with 0s.

```
>>> df.fillna(0)

A B C D

0 0.0 2.0 0.0 0

1 3.0 4.0 0.0 1

2 0.0 0.0 0.0 5

3 0.0 3.0 0.0 4
```

We can also propagate non-null values forward or backward.

```
>>> df.fillna(method='ffill')

A B C D

0 NaN 2.0 NaN 0

1 3.0 4.0 NaN 1

2 3.0 4.0 NaN 5

3 3.0 3.0 NaN 4
```

Replace all NaN elements in column 'A', 'B', 'C', and 'D', with 0, 1, 2, and 3 respectively.

Only replace the first NaN element.

```
>>> df.fillna(value=values, limit=1)

A B C D
0 0.0 2.0 2.0 0
1 3.0 4.0 NaN 1
2 NaN 1.0 NaN 5
3 NaN 3.0 NaN 4
```

**filter** (*items=None*, *like=None*, *regex=None*, *axis=None*)

Subset rows or columns of dataframe according to labels in the specified index.

Note that this routine does not filter a dataframe on its contents. The filter is applied to the labels of the index.

#### **Parameters**

```
items [list-like] List of info axis to restrict to (must not all be present)
```

```
like [string] Keep info axis where "arg in col == True"
```

```
regex [string (regular expression)] Keep info axis with re.search(regex, col) == True
```

**axis** [int or string axis name] The axis to filter on. By default this is the info axis, 'index' for Series, 'columns' for DataFrame

## Returns

# same type as input object

### See also:

```
pandas.DataFrame.loc
```

### **Notes**

The items, like, and regex parameters are enforced to be mutually exclusive.

axis defaults to the info axis that is used when indexing with [].

# **Examples**

```
>>> df
one two three
mouse 1 2 3
rabbit 4 5 6
```

```
>>> # select columns by name
>>> df.filter(items=['one', 'three'])
one three
mouse 1 3
rabbit 4 6
```

```
>>> # select columns by regular expression
>>> df.filter(regex='e$', axis=1)
one three
mouse 1 3
rabbit 4 6
```

```
>>> # select rows containing 'bbi'
>>> df.filter(like='bbi', axis=0)
one two three
rabbit 4 5 6
```

### first (offset)

Convenience method for subsetting initial periods of time series data based on a date offset.

### **Parameters**

offset [string, DateOffset, dateutil.relativedelta]

#### **Returns**

subset [type of caller]

### **Raises**

TypeError If the index is not a DatetimeIndex

#### See also:

last Select final periods of time series based on a date offset

at\_time Select values at a particular time of the day

between\_time Select values between particular times of the day

# **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='2D')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 1
2018-04-11 2
2018-04-13 3
2018-04-15 4
```

## Get the rows for the first 3 days:

```
>>> ts.first('3D')

A
2018-04-09 1
2018-04-11 2
```

Notice the data for 3 first calender days were returned, not the first 3 days observed in the dataset, and therefore data for 2018-04-13 was not returned.

## first\_valid\_index()

Return index for first non-NA/null value.

## Returns

scalar [type of index]

### **Notes**

If all elements are non-NA/null, returns None. Also returns None for empty NDFrame.

# flags

return the ndarray.flags for the underlying data

```
floordiv (other, level=None, fill_value=None, axis=0)
```

Integer division of series and other, element-wise (binary operator *floordiv*).

Equivalent to series // other, but with support to substitute a fill\_value for missing data in one of the inputs.

#### **Parameters**

other [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

#### Returns

result [Series]

### See also:

Series.rfloordiv

## **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
а
    1.0
b
    1.0
С
    1.0
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> h
а
    1.0
    NaN
d
    1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
    1.0
С
    1.0
d
    1.0
e
    NaN
dtype: float64
```

 $\begin{tabular}{ll} \textbf{classmethod from\_array} (arr, index=None, name=None, dtype=None, copy=False, fast-path=False) \\ \end{tabular}$ 

Construct Series from array.

Deprecated since version 0.23.0: Use pd.Series(..) constructor instead.

classmethod from\_csv (path, sep=', ',  $parse\_dates=True$ , header=None,  $index\_col=0$ , encod-ing=None,  $infer\_datetime\_format=False$ )

Read CSV file.

Deprecated since version 0.21.0: Use pandas.read\_csv() instead.

It is preferable to use the more powerful pandas.read\_csv() for most general purposes, but from\_csv makes for an easy roundtrip to and from a file (the exact counterpart of to\_csv), especially with a time Series.

This method only differs from pandas.read\_csv() in some defaults:

- index\_col is 0 instead of None (take first column as index by default)
- header is None instead of 0 (the first row is not used as the column names)

• parse\_dates is True instead of False (try parsing the index as datetime by default)

With pandas.read\_csv(), the option squeeze=True can be used to return a Series like from\_csv.

#### **Parameters**

path [string file path or file handle / StringIO]

sep [string, default ','] Field delimiter

parse\_dates [boolean, default True] Parse dates. Different default from read\_table

**header** [int, default None] Row to use as header (skip prior rows)

index\_col [int or sequence, default 0] Column to use for index. If a sequence is given, a
 MultiIndex is used. Different default from read\_table

**encoding** [string, optional] a string representing the encoding to use if the contents are non-ascii, for python versions prior to 3

**infer\_datetime\_format: boolean, default False** If True and *parse\_dates* is True for a column, try to infer the datetime format based on the first datetime string. If the format can be inferred, there often will be a large parsing speed-up.

#### Returns

y [Series]

#### See also:

pandas.read\_csv

## ftype

return if the data is sparseldense

### ftypes

return if the data is sparseldense

ge (other, level=None, fill\_value=None, axis=0)

Greater than or equal to of series and other, element-wise (binary operator ge).

Equivalent to series >= other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

other [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### **Returns**

result [Series]

# See also:

Series.None

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
b
     1.0
С
     1.0
    NaN
d
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
    1.0
b
    NaN
d
    1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
b
     1.0
С
     1.0
d
     1.0
0
     NaN
dtype: float64
```

## get (key, default=None)

Get item from object for given key (DataFrame column, Panel slice, etc.). Returns default value if not found.

### **Parameters**

key [object]

### Returns

value [type of items contained in object]

## get\_dtype\_counts()

Return counts of unique dtypes in this object.

## Returns

**dtype** [Series] Series with the count of columns with each dtype.

# See also:

dtypes Return the dtypes in this object.

# **Examples**

```
>>> a = [['a', 1, 1.0], ['b', 2, 2.0], ['c', 3, 3.0]]
>>> df = pd.DataFrame(a, columns=['str', 'int', 'float'])
>>> df
    str int float
0 a 1 1.0
1 b 2 2.0
2 c 3 3.0
```

```
>>> df.get_dtype_counts()
float64 1
int64 1
object 1
dtype: int64
```

### get\_ftype\_counts()

Return counts of unique ftypes in this object.

Deprecated since version 0.23.0.

This is useful for SparseDataFrame or for DataFrames containing sparse arrays.

#### Returns

**dtype** [Series] Series with the count of columns with each type and sparsity (dense/sparse)

### See also:

ftypes Return ftypes (indication of sparse/dense and dtype) in this object.

# **Examples**

```
>>> a = [['a', 1, 1.0], ['b', 2, 2.0], ['c', 3, 3.0]]
>>> df = pd.DataFrame(a, columns=['str', 'int', 'float'])
>>> df
    str int float
0 a 1 1.0
1 b 2 2.0
2 c 3 3.0
```

```
>>> df.get_ftype_counts()
float64:dense 1
int64:dense 1
object:dense 1
dtype: int64
```

## get\_value (label, takeable=False)

Quickly retrieve single value at passed index label

Deprecated since version 0.21.0: Please use .at[] or .iat[] accessors.

## **Parameters**

```
label [object]
```

takeable [interpret the index as indexers, default False]

### Returns

value [scalar value]

### get\_values()

same as values (but handles sparseness conversions); is a view

```
groupby (by=None, axis=0, level=None, as_index=True, squeeze=False, observed=False, **kwargs)
```

Group series using mapper (dict or key function, apply given function to group, return result as series) or by a series of columns.

### **Parameters**

by [mapping, function, label, or list of labels] Used to determine the groups for the groupby. If by is a function, it's called on each value of the object's index. If a dict or Series is passed, the Series or dict VALUES will be used to determine the groups (the Series' values are first aligned; see .align() method). If an ndarray is passed, the values are used as-is determine the groups. A label or list of labels may be passed to group by the columns in self. Notice that a tuple is interpreted a (single) key.

axis [int, default 0]

- **level** [int, level name, or sequence of such, default None] If the axis is a MultiIndex (hierarchical), group by a particular level or levels
- as\_index [boolean, default True] For aggregated output, return object with group labels as the index. Only relevant for DataFrame input. as\_index=False is effectively "SQLstyle" grouped output
- **sort** [boolean, default True] Sort group keys. Get better performance by turning this off. Note this does not influence the order of observations within each group. groupby preserves the order of rows within each group.
- **group\_keys** [boolean, default True] When calling apply, add group keys to index to identify pieces
- **squeeze** [boolean, default False] reduce the dimensionality of the return type if possible, otherwise return a consistent type
- **observed** [boolean, default False] This only applies if any of the groupers are Categoricals If True: only show observed values for categorical groupers. If False: show all values for categorical groupers.

New in version 0.23.0.

### **Returns**

# GroupBy object

## See also:

**resample** Convenience method for frequency conversion and resampling of time series.

## **Notes**

See the user guide for more.

## **Examples**

### DataFrame results

```
>>> data.groupby(func, axis=0).mean()
>>> data.groupby(['col1', 'col2'])['col3'].mean()
```

### DataFrame with hierarchical index

```
>>> data.groupby(['col1', 'col2']).mean()
```

```
gt (other, level=None, fill_value=None, axis=0)
```

Greater than of series and other, element-wise (binary operator gt).

Equivalent to series > other, but with support to substitute a fill\_value for missing data in one of the inputs.

# **Parameters**

**other** [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### **Returns**

result [Series]

#### See also:

Series.None

## **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
b
     1.0
     1.0
С
d
     NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
     1.0
а
b
     NaN
d
     1.0
     NaN
dtype: float64
>>> a.add(b, fill_value=0)
     2.0
b
     1.0
     1.0
С
d
     1.0
     NaN
dtype: float64
```

## hasnans

return if I have any nans; enables various perf speedups

#### head(n=5)

Return the first *n* rows.

This function returns the first n rows for the object based on position. It is useful for quickly testing if your object has the right type of data in it.

## **Parameters**

**n** [int, default 5] Number of rows to select.

### Returns

**obj\_head** [type of caller] The first *n* rows of the caller object.

See also:

pandas.DataFrame.tail Returns the last n rows.

# **Examples**

```
>>> df = pd.DataFrame({'animal':['alligator', 'bee', 'falcon', 'lion',
                        'monkey', 'parrot', 'shark', 'whale', 'zebra']})
>>> df
      animal
0
  alligator
1
         bee
2
     falcon
3
       lion
4
      monkey
5
      parrot
6
       shark
7
       whale
8
       zebra
```

## Viewing the first 5 lines

```
>>> df.head()
    animal
0 alligator
1 bee
2 falcon
3 lion
4 monkey
```

## Viewing the first *n* lines (three in this case)

```
>>> df.head(3)
animal
0 alligator
1 bee
2 falcon
```

hist (by=None, ax=None, grid=True, xlabelsize=None, xrot=None, ylabelsize=None, yrot=None, figsize=None, bins=10, \*\*kwds)

Draw histogram of the input series using matplotlib

### **Parameters**

```
by [object, optional] If passed, then used to form histograms for separate groups
ax [matplotlib axis object] If not passed, uses gca()
grid [boolean, default True] Whether to show axis grid lines
xlabelsize [int, default None] If specified changes the x-axis label size
xrot [float, default None] rotation of x axis labels
```

ylabelsize [int, default None] If specified changes the y-axis label size

yrot [float, default None] rotation of y axis labels

figsize [tuple, default None] figure size in inches by default

bins [integer or sequence, default 10] Number of histogram bins to be used. If an integer is given, bins + 1 bin edges are calculated and returned. If bins is a sequence, gives bin edges, including left edge of first bin and right edge of last bin. In this case, bins is returned unmodified.

bins: integer, default 10 Number of histogram bins to be used

**\*\*\*kwds\*** [keywords] To be passed to the actual plotting function

See also:

matplotlib.axes.Axes.hist Plot a histogram using matplotlib.

iat

Access a single value for a row/column pair by integer position.

Similar to iloc, in that both provide integer-based lookups. Use iat if you only need to get or set a single value in a DataFrame or Series.

### **Raises**

**IndexError** When integer position is out of bounds

See also:

DataFrame.at Access a single value for a row/column label pair

DataFrame.loc Access a group of rows and columns by label(s)

**DataFrame.iloc** Access a group of rows and columns by integer position(s)

# **Examples**

Get value at specified row/column pair

```
>>> df.iat[1, 2]
1
```

Set value at specified row/column pair

```
>>> df.iat[1, 2] = 10
>>> df.iat[1, 2]
10
```

Get value within a series

```
>>> df.loc[0].iat[1]
2
```

idxmax (axis=0, skipna=True, \*args, \*\*kwargs)

Return the row label of the maximum value.

If multiple values equal the maximum, the first row label with that value is returned.

### **Parameters**

**skipna** [boolean, default True] Exclude NA/null values. If the entire Series is NA, the result will be NA.

axis [int, default 0] For compatibility with DataFrame.idxmax. Redundant for application on Series.

\*args, \*\*kwargs Additional keywors have no effect but might be accepted for compatibility with NumPy.

### **Returns**

idxmax [Index of maximum of values.]

#### Raises

ValueError If the Series is empty.

### See also:

numpy.argmax Return indices of the maximum values along the given axis.

**DataFrame.idxmax** Return index of first occurrence of maximum over requested axis.

Series.idxmin Return index label of the first occurrence of minimum of values.

# **Notes**

This method is the Series version of ndarray.argmax. This method returns the label of the maximum, while ndarray.argmax returns the position. To get the position, use series.values.argmax().

## **Examples**

```
>>> s.idxmax()
'C'
```

If *skipna* is False and there is an NA value in the data, the function returns nan.

```
>>> s.idxmax(skipna=False)
nan
```

idxmin (axis=None, skipna=True, \*args, \*\*kwargs)

Return the row label of the minimum value.

If multiple values equal the minimum, the first row label with that value is returned.

#### **Parameters**

**skipna** [boolean, default True] Exclude NA/null values. If the entire Series is NA, the result will be NA.

axis [int, default 0] For compatibility with DataFrame.idxmin. Redundant for application on Series.

\*args, \*\*kwargs Additional keywors have no effect but might be accepted for compatibility with NumPy.

### **Returns**

idxmin [Index of minimum of values.]

#### Raises

**ValueError** If the Series is empty.

### See also:

numpy.argmin Return indices of the minimum values along the given axis.

**DataFrame.idxmin** Return index of first occurrence of minimum over requested axis.

Series.idxmax Return index label of the first occurrence of maximum of values.

## **Notes**

This method is the Series version of ndarray.argmin. This method returns the label of the minimum, while ndarray.argmin returns the position. To get the position, use series.values.argmin().

## **Examples**

```
>>> s.idxmin()
'A'
```

If skipna is False and there is an NA value in the data, the function returns nan.

```
>>> s.idxmin(skipna=False)
nan
```

#### iloc

Purely integer-location based indexing for selection by position.

.iloc[] is primarily integer position based (from 0 to length-1 of the axis), but may also be used with a boolean array.

Allowed inputs are:

- An integer, e.g. 5.
- A list or array of integers, e.g. [4, 3, 0].
- A slice object with ints, e.g. 1:7.
- A boolean array.
- A callable function with one argument (the calling Series, DataFrame or Panel) and that returns valid output for indexing (one of the above)

.iloc will raise IndexError if a requested indexer is out-of-bounds, except *slice* indexers which allow out-of-bounds indexing (this conforms with python/numpy *slice* semantics).

See more at Selection by Position

#### imag

### index

The index (axis labels) of the Series.

## infer objects()

Attempt to infer better dtypes for object columns.

Attempts soft conversion of object-dtyped columns, leaving non-object and unconvertible columns unchanged. The inference rules are the same as during normal Series/DataFrame construction.

New in version 0.21.0.

#### Returns

converted [same type as input object]

See also:

```
pandas.to_datetime Convert argument to datetime.
```

pandas.to\_timedelta Convert argument to timedelta.

pandas.to\_numeric Convert argument to numeric typeR

# **Examples**

```
>>> df = pd.DataFrame({"A": ["a", 1, 2, 3]})
>>> df = df.iloc[1:]
>>> df
    A
1    1
2    2
3    3
```

```
>>> df.dtypes
A object
dtype: object
```

```
>>> df.infer_objects().dtypes
A int64
dtype: object
```

Please note that only method='linear' is supported for DataFrames/Series with a MultiIndex.

#### **Parameters**

- 'linear': ignore the index and treat the values as equally spaced. This is the only method supported on MultiIndexes. default
- 'time': interpolation works on daily and higher resolution data to interpolate given length of interval
- 'index', 'values': use the actual numerical values of the index
- 'nearest', 'zero', 'slinear', 'quadratic', 'cubic', 'barycentric', 'polynomial' is passed to scipy.interpolate.interpld. Both 'polynomial' and 'spline' require that you also specify an *order* (int), e.g. df.interpolate(method='polynomial', order=4). These use the actual numerical values of the index.
- 'krogh', 'piecewise\_polynomial', 'spline', 'pchip' and 'akima' are all wrappers around the scipy interpolation methods of similar names. These use the actual numerical values of the index. For more information on their behavior, see the scipy documentation and tutorial documentation
- 'from\_derivatives' refers to BPoly.from\_derivatives which replaces 'piece-wise\_polynomial' interpolation method in scipy 0.18

New in version 0.18.1: Added support for the 'akima' method Added interpolate method 'from\_derivatives' which replaces 'piecewise\_polynomial' in scipy 0.18; backwards-compatible with scipy < 0.18

```
axis [{0, 1}, default 0]
```

- 0: fill column-by-column
- 1: fill row-by-row

**limit** [int, default None.] Maximum number of consecutive NaNs to fill. Must be greater than 0.

```
limit_direction [{'forward', 'backward', 'both'}, default 'forward']
```

limit\_area [{'inside', 'outside'}, default None]

- None: (default) no fill restriction
- 'inside' Only fill NaNs surrounded by valid values (interpolate).
- 'outside' Only fill NaNs outside valid values (extrapolate).

If limit is specified, consecutive NaNs will be filled in this direction.

New in version 0.21.0.

inplace [bool, default False] Update the NDFrame in place if possible.

downcast [optional, 'infer' or None, defaults to None] Downcast dtypes if possible.

**kwargs** [keyword arguments to pass on to the interpolating function.]

### **Returns**

Series or DataFrame of same shape interpolated at the NaNs

## See also:

```
reindex, replace, fillna
```

# **Examples**

## Filling in NaNs

### is\_copy

## is\_monotonic

Return boolean if values in the object are monotonic\_increasing

New in version 0.19.0.

## **Returns**

is\_monotonic [boolean]

# is\_monotonic\_decreasing

Return boolean if values in the object are monotonic\_decreasing

New in version 0.19.0.

### **Returns**

is\_monotonic\_decreasing [boolean]

# is\_monotonic\_increasing

Return boolean if values in the object are monotonic\_increasing

New in version 0.19.0.

### **Returns**

is\_monotonic [boolean]

## is\_unique

Return boolean if values in the object are unique

### Returns

is\_unique [boolean]

#### isin (values)

Check whether values are contained in Series.

Return a boolean Series showing whether each element in the Series matches an element in the passed sequence of *values* exactly.

## **Parameters**

**values** [set or list-like] The sequence of values to test. Passing in a single string will raise a TypeError. Instead, turn a single string into a list of one element.

New in version 0.18.1: Support for values as a set.

### **Returns**

isin [Series (bool dtype)]

### Raises

## **TypeError**

• If values is a string

See also:

pandas.DataFrame.isin equivalent method on DataFrame

## **Examples**

Passing a single string as s.isin('lama') will raise an error. Use a list of one element instead:

```
>>> s.isin(['lama'])
0    True
1    False
2    True
3    False
4    True
5    False
Name: animal, dtype: bool
```

# isna()

Detect missing values.

Return a boolean same-sized object indicating if the values are NA. NA values, such as None or numpy. NaN, gets mapped to True values. Everything else gets mapped to False values. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True).

Returns

**Series** Mask of bool values for each element in Series that indicates whether an element is not an NA value.

See also:

Series.isnull alias of isna

Series.notna boolean inverse of isna

Series.dropna omit axes labels with missing values

isna top-level isna

## **Examples**

Show which entries in a DataFrame are NA.

```
>>> df = pd.DataFrame({'age': [5, 6, np.NaN],
                      'born': [pd.NaT, pd.Timestamp('1939-05-27'),
                              pd.Timestamp('1940-04-25')],
                      'name': ['Alfred', 'Batman', ''],
                      'toy': [None, 'Batmobile', 'Joker']})
. . .
>>> df
  age
           born
                 name
                               toy
           NaT Alfred
  5.0
                             None
 6.0 1939-05-27 Batman Batmobile
2 NaN 1940-04-25
                             Joker
```

```
>>> df.isna()
age born name toy
0 False True False True
1 False False False
2 True False False
```

Show which entries in a Series are NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.isna()
0 False
1 False
2 True
dtype: bool
```

## isnull()

Detect missing values.

Return a boolean same-sized object indicating if the values are NA. NA values, such as None or numpy. NaN, gets mapped to True values. Everything else gets mapped to False values. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True).

Returns

**Series** Mask of bool values for each element in Series that indicates whether an element is not an NA value.

See also:

Series.isnull alias of isna

Series.notna boolean inverse of isna

Series.dropna omit axes labels with missing values

isna top-level isna

# **Examples**

Show which entries in a DataFrame are NA.

```
>>> df = pd.DataFrame({'age': [5, 6, np.NaN],
                      'born': [pd.NaT, pd.Timestamp('1939-05-27'),
. . .
                              pd.Timestamp('1940-04-25')],
                      'name': ['Alfred', 'Batman', ''],
                      'toy': [None, 'Batmobile', 'Joker']})
. . .
>>> df
  age
           born
                  name
                               toy
           NaT Alfred None
  5.0
 6.0 1939-05-27 Batman Batmobile
2 NaN 1940-04-25
                              Joker
```

```
>>> df.isna()
age born name toy
0 False True False True
1 False False False False
2 True False False
```

Show which entries in a Series are NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.isna()
0 False
1 False
2 True
dtype: bool
```

## item()

return the first element of the underlying data as a python scalar

### items()

Lazily iterate over (index, value) tuples

#### itemsize

return the size of the dtype of the item of the underlying data

#### iteritems()

Lazily iterate over (index, value) tuples

ix

A primarily label-location based indexer, with integer position fallback.

Warning: Starting in 0.20.0, the .ix indexer is deprecated, in favor of the more strict .iloc and .loc indexers.

.ix[] supports mixed integer and label based access. It is primarily label based, but will fall back to integer positional access unless the corresponding axis is of integer type.

.ix is the most general indexer and will support any of the inputs in .loc and .iloc. .ix also supports floating point label schemes. .ix is exceptionally useful when dealing with mixed positional and label based hierarchical indexes.

However, when an axis is integer based, ONLY label based access and not positional access is supported. Thus, in such cases, it's usually better to be explicit and use .iloc or .loc.

See more at Advanced Indexing.

### keys()

Alias for index

kurt (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0). Normalized by N-1

### **Parameters**

```
axis [\{index (0)\}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

## Returns

**kurt** [scalar or Series (if level specified)]

kurtosis (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

Return unbiased kurtosis over requested axis using Fisher's definition of kurtosis (kurtosis of normal == 0.0). Normalized by N-1

### **Parameters**

```
axis [\{index (0)\}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

# Returns

kurt [scalar or Series (if level specified)]

### last (offset)

Convenience method for subsetting final periods of time series data based on a date offset.

#### **Parameters**

offset [string, DateOffset, dateutil.relativedelta]

### Returns

subset [type of caller]

## **Raises**

TypeError If the index is not a DatetimeIndex

### See also:

first Select initial periods of time series based on a date offset

at\_time Select values at a particular time of the day

between\_time Select values between particular times of the day

# **Examples**

```
>>> i = pd.date_range('2018-04-09', periods=4, freq='2D')
>>> ts = pd.DataFrame({'A': [1,2,3,4]}, index=i)
>>> ts

A
2018-04-09 1
2018-04-11 2
2018-04-13 3
2018-04-15 4
```

## Get the rows for the last 3 days:

```
>>> ts.last('3D')

A
2018-04-13 3
2018-04-15 4
```

Notice the data for 3 last calender days were returned, not the last 3 observed days in the dataset, and therefore data for 2018-04-11 was not returned.

# last\_valid\_index()

Return index for last non-NA/null value.

### Returns

**scalar** [type of index]

### **Notes**

If all elements are non-NA/null, returns None. Also returns None for empty NDFrame.

### **le** (*other*, *level=None*, *fill\_value=None*, *axis=0*)

Less than or equal to of series and other, element-wise (binary operator *le*).

Equivalent to series <= other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

other [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### **Returns**

result [Series]

### See also:

Series.None

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
а
    1.0
    1.0
b
     1.0
С
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
    1.0
а
b
    NaN
d
    1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
а
     1.0
b
    1.0
С
d
     1.0
    NaN
dtype: float64
```

## loc

Access a group of rows and columns by label(s) or a boolean array.

.loc[] is primarily label based, but may also be used with a boolean array.

Allowed inputs are:

- A single label, e.g. 5 or 'a', (note that 5 is interpreted as a *label* of the index, and **never** as an integer position along the index).
- A list or array of labels, e.g. ['a', 'b', 'c'].
- A slice object with labels, e.g. 'a':'f'.

Warning: Note that contrary to usual python slices, both the start and the stop are included

- A boolean array of the same length as the axis being sliced, e.g. [True, False, True].
- A callable function with one argument (the calling Series, DataFrame or Panel) and that returns valid output for indexing (one of the above)

See more at Selection by Label

#### Raises

**KeyError:** when any items are not found

### See also:

DataFrame.at Access a single value for a row/column label pair

**DataFrame.iloc** Access group of rows and columns by integer position(s)

**DataFrame.xs** Returns a cross-section (row(s) or column(s)) from the Series/DataFrame.

Series.loc Access group of values using labels

# **Examples**

# **Getting values**

Single label. Note this returns the row as a Series.

```
>>> df.loc['viper']
max_speed 4
shield 5
Name: viper, dtype: int64
```

List of labels. Note using [ [ ] ] returns a DataFrame.

```
>>> df.loc[['viper', 'sidewinder']]

max_speed shield

viper 4 5

sidewinder 7 8
```

Single label for row and column

```
>>> df.loc['cobra', 'shield']
2
```

Slice with labels for row and single label for column. As mentioned above, note that both the start and stop of the slice are included.

```
>>> df.loc['cobra':'viper', 'max_speed']
cobra 1
viper 4
Name: max_speed, dtype: int64
```

## Boolean list with the same length as the row axis

```
>>> df.loc[[False, False, True]]

max_speed shield
sidewinder 7 8
```

### Conditional that returns a boolean Series

## Conditional that returns a boolean Series with column labels specified

#### Callable that returns a boolean Series

```
>>> df.loc[lambda df: df['shield'] == 8]

max_speed shield
sidewinder 7 8
```

# **Setting values**

Set value for all items matching the list of labels

## Set value for an entire row

## Set value for an entire column

Set value for rows matching callable condition

## Getting values on a DataFrame with an index that has integer labels

Another example using integers for the index

Slice with integer labels for rows. As mentioned above, note that both the start and stop of the slice are included.

# Getting values with a MultiIndex

A number of examples using a DataFrame with a MultiIndex

```
>>> tuples = [
       ('cobra', 'mark i'), ('cobra', 'mark ii'),
. . .
       ('sidewinder', 'mark i'), ('sidewinder', 'mark ii'),
. . .
       ('viper', 'mark ii'), ('viper', 'mark iii')
...]
>>> index = pd.MultiIndex.from_tuples(tuples)
>>> values = [[12, 2], [0, 4], [10, 20],
            [1, 4], [7, 1], [16, 36]]
>>> df = pd.DataFrame(values, columns=['max_speed', 'shield'], index=index)
>>> df
                     max_speed shield
          mark i
                            12
                                     2
cobra
          mark ii
                            0
                                     4
sidewinder mark i
                            10
                                    20
           mark ii
                            1
                                     4
                            7
           mark ii
                                     1
viper
                           16
          mark iii
```

Single label. Note this returns a DataFrame with a single index.

Single index tuple. Note this returns a Series.

```
>>> df.loc[('cobra', 'mark ii')]
max_speed 0
shield 4
Name: (cobra, mark ii), dtype: int64
```

Single label for row and column. Similar to passing in a tuple, this returns a Series.

```
>>> df.loc['cobra', 'mark i']
max_speed 12
shield 2
Name: (cobra, mark i), dtype: int64
```

Single tuple. Note using [ [ ] ] returns a DataFrame.

```
>>> df.loc[[('cobra', 'mark ii')]]

max_speed shield
cobra mark ii 0 4
```

Single tuple for the index with a single label for the column

```
>>> df.loc[('cobra', 'mark i'), 'shield']
2
```

Slice from index tuple to single label

```
>>> df.loc[('cobra', 'mark i'):'viper']
                   max_speed shield
cobra
          mark i
                          12
          mark ii
                           0
                                   4
                          10
                                  20
sidewinder mark i
                          1
          mark ii
                                   4
                           7
viper
          mark ii
                                   1
          mark iii
                          16
                                  36
```

Slice from index tuple to index tuple

```
>>> df.loc[('cobra', 'mark i'):('viper', 'mark ii')]
                  max_speed shield
                    12
cobra
          mark i
          mark ii
                          0
                                  4
sidewinder mark i
                         10
                                 20
                                  4
          mark ii
                          1
                          7
viper
          mark ii
                                  1
```

**lt** (other, level=None, fill\_value=None, axis=0)

Less than of series and other, element-wise (binary operator *lt*).

Equivalent to series < other, but with support to substitute a fill\_value for missing data in one of the inputs.

**Parameters** 

**other** [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

#### Returns

result [Series]

### See also:

Series.None

## **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
     1.0
h
     1.0
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
    1.0
а
b
    NaN
d
    1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
     2.0
h
     1.0
     1.0
С
     1.0
     NaN
dtype: float64
```

mad (axis=None, skipna=None, level=None)

Return the mean absolute deviation of the values for the requested axis

### **Parameters**

```
axis [\{index (0)\}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

mad [scalar or Series (if level specified)]

map (arg, na\_action=None)

Map values of Series using input correspondence (a dict, Series, or function).

### **Parameters**

arg [function, dict, or Series] Mapping correspondence.

**na\_action** [{None, 'ignore'}] If 'ignore', propagate NA values, without passing them to the mapping correspondence.

#### Returns

y [Series] Same index as caller.

See also:

**Series.apply** For applying more complex functions on a Series.

DataFrame.apply Apply a function row-/column-wise.

**DataFrame**. applymap Apply a function elementwise on a whole DataFrame.

### **Notes**

When *arg* is a dictionary, values in Series that are not in the dictionary (as keys) are converted to NaN. However, if the dictionary is a dict subclass that defines \_\_missing\_\_ (i.e. provides a method for default values), then this default is used rather than NaN:

```
>>> from collections import Counter
>>> counter = Counter()
>>> counter['bar'] += 1
>>> y.map(counter)
1     0
2     1
3     0
dtype: int64
```

# **Examples**

Map inputs to outputs (both of type Series):

```
>>> x = pd.Series([1,2,3], index=['one', 'two', 'three'])
>>> x
one    1
two    2
three    3
dtype: int64
```

```
>>> y = pd.Series(['foo', 'bar', 'baz'], index=[1,2,3])
>>> y
1    foo
2    bar
3    baz
```

```
>>> x.map(y)
one foo
two bar
three baz
```

If arg is a dictionary, return a new Series with values converted according to the dictionary's mapping:

```
>>> z = {1: 'A', 2: 'B', 3: 'C'}
```

```
>>> x.map(z)
one A
two B
three C
```

Use na\_action to control whether NA values are affected by the mapping function.

```
>>> s = pd.Series([1, 2, 3, np.nan])
```

```
>>> s2 = s.map('this is a string {}'.format, na_action=None)
0 this is a string 1.0
1 this is a string 2.0
2 this is a string 3.0
3 this is a string nan
dtype: object
```

```
>>> s3 = s.map('this is a string {}'.format, na_action='ignore')

0 this is a string 1.0

1 this is a string 2.0

2 this is a string 3.0

3 NaN
dtype: object
```

mask (cond, other=nan, inplace=False, axis=None, level=None, errors='raise', try\_cast=False, raise on error=None)

Return an object of same shape as self and whose corresponding entries are from self where *cond* is False and otherwise are from *other*.

#### **Parameters**

**cond** [boolean NDFrame, array-like, or callable] Where *cond* is False, keep the original value. Where True, replace with corresponding value from *other*. If *cond* is callable, it is computed on the NDFrame and should return boolean NDFrame or array. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as cond.

**other** [scalar, NDFrame, or callable] Entries where *cond* is True are replaced with corresponding value from *other*. If other is callable, it is computed on the NDFrame and should return scalar or NDFrame. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as other.

**inplace** [boolean, default False] Whether to perform the operation in place on the data

axis [alignment axis if needed, default None]

**level** [alignment level if needed, default None]

```
errors [str, {'raise', 'ignore'}, default 'raise']
```

- raise: allow exceptions to be raised
- ignore: suppress exceptions. On error return original object

Note that currently this parameter won't affect the results and will always coerce to a suitable dtype.

try\_cast [boolean, default False] try to cast the result back to the input type (if possible),

raise\_on\_error [boolean, default True] Whether to raise on invalid data types (e.g. trying to where on strings)

Deprecated since version 0.21.0.

### Returns

**wh** [same type as caller]

### See also:

```
DataFrame.where()
```

#### **Notes**

The mask method is an application of the if-then idiom. For each element in the calling DataFrame, if cond is False the element is used; otherwise the corresponding element from the DataFrame other is used.

The signature for DataFrame.where() differs from numpy.where(). Roughly dfl.where(m, df2) is equivalent to np.where(m, df1, df2).

For further details and examples see the mask documentation in indexing.

# **Examples**

```
>>> s = pd.Series(range(5))
>>> s.where(s > 0)

0     NaN
1     1.0
2     2.0
3     3.0
4     4.0
```

```
>>> df = pd.DataFrame(np.arange(10).reshape(-1, 2), columns=['A', 'B'])
>>> m = df % 3 == 0
>>> df.where(m, -df)

A B
0 0 -1
1 -2 3
2 -4 -5
3 6 -7
```

(continues on next page)

```
4 -8
>>> df.where(m, -df) == np.where(m, df, -df)
     A
\cap
  True True
  True
        True
  True
        True
  True
        True
  True
        True
>>> df.where(m, -df) == df.mask(\sim m, -df)
     Α
0
  True True
  True True
2
  True True
3
  True True
  True True
```

max (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

This method returns the maximum of the values in the object. If you want the *index* of the maximum, use idxmax. This is the equivalent of the numpy.ndarray method argmax.

#### **Parameters**

```
axis [\{index (0)\}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

#### Returns

```
max [scalar or Series (if level specified)]
```

mean (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)
Return the mean of the values for the requested axis

#### **Parameters**

```
axis [\{index (0)\}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

```
mean [scalar or Series (if level specified)]
```

**median** (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)
Return the median of the values for the requested axis

#### **Parameters**

```
axis [\{index (0)\}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### **Returns**

**median** [scalar or Series (if level specified)]

```
memory_usage (index=True, deep=False)
```

Return the memory usage of the Series.

The memory usage can optionally include the contribution of the index and of elements of *object* dtype.

### **Parameters**

index [bool, default True] Specifies whether to include the memory usage of the Series index.

**deep** [bool, default False] If True, introspect the data deeply by interrogating *object* dtypes for system-level memory consumption, and include it in the returned value.

### Returns

int Bytes of memory consumed.

#### See also:

numpy.ndarray.nbytes Total bytes consumed by the elements of the array.

**DataFrame.memory\_usage** Bytes consumed by a DataFrame.

# **Examples**

```
>>> s = pd.Series(range(3))
>>> s.memory_usage()
104
```

Not including the index gives the size of the rest of the data, which is necessarily smaller:

```
>>> s.memory_usage(index=False)
24
```

The memory footprint of *object* values is ignored by default:

```
>>> s = pd.Series(["a", "b"])
>>> s.values
array(['a', 'b'], dtype=object)
>>> s.memory_usage()
96
>>> s.memory_usage(deep=True)
212
```

min (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

This method returns the minimum of the values in the object. If you want the *index* of the minimum, use idxmin. This is the equivalent of the numpy.ndarray method argmin.

## **Parameters**

```
axis [\{index (0)\}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series

#### Returns

**min** [scalar or Series (if level specified)]

```
mod (other, level=None, fill value=None, axis=0)
```

Modulo of series and other, element-wise (binary operator *mod*).

Equivalent to series % other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

**other** [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### **Returns**

result [Series]

## See also:

Series.rmod

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
а
    1.0
b
     1.0
     1.0
С
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
     1.0
а
b
     NaN
d
     1.0
```

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```
e NaN
dtype: float64
>>> a.add(b, fill_value=0)
a 2.0
b 1.0
c 1.0
d 1.0
e NaN
dtype: float64
```

## mode()

Return the mode(s) of the dataset.

Always returns Series even if only one value is returned.

### **Returns**

```
modes [Series (sorted)]
```

**mul** (other, level=None, fill value=None, axis=0)

Multiplication of series and other, element-wise (binary operator mul).

Equivalent to series \* other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

other [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### **Returns**

result [Series]

### See also:

Series.rmul

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
     1.0
b
    1.0
С
d
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
а
    1.0
     NaN
b
d
     1.0
```

(continues on next page)

```
e NaN
dtype: float64
>>> a.add(b, fill_value=0)
a 2.0
b 1.0
c 1.0
d 1.0
e NaN
dtype: float64
```

# multiply (other, level=None, fill\_value=None, axis=0)

Multiplication of series and other, element-wise (binary operator mul).

Equivalent to series \* other, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

**other** [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

#### Returns

result [Series]

## See also:

Series.rmul

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
     1.0
b
С
     1.0
     NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
    1.0
а
b
     NaN
d
    1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
     2.0
     1.0
b
С
     1.0
d
     1.0
```

(continues on next page)

```
e NaN
dtype: float64
```

#### name

### nbytes

return the number of bytes in the underlying data

#### ndim

return the number of dimensions of the underlying data, by definition 1

```
ne (other, level=None, fill value=None, axis=0)
```

Not equal to of series and other, element-wise (binary operator *ne*).

Equivalent to series != other, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

**other** [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### Returns

result [Series]

## See also:

Series.None

## **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
     1.0
b
С
     1.0
     NaN
d
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
а
     1.0
b
     NaN
d
     1.0
     NaN
dtype: float64
>>> a.add(b, fill_value=0)
     2.0
а
     1.0
b
С
     1.0
d
     1.0
```

(continues on next page)

```
e NaN dtype: float64
```

### nlargest (n=5, keep='first')

Return the largest n elements.

#### **Parameters**

n [int] Return this many descending sorted values

**keep** [{'first', 'last'}, default 'first'] Where there are duplicate values: - first: take the first occurrence. - last: take the last occurrence.

#### Returns

top\_n [Series] The n largest values in the Series, in sorted order

### See also:

Series.nsmallest

### **Notes**

Faster than  $.sort_values(ascending=False).head(n)$  for small n relative to the size of the Series object.

## **Examples**

```
>>> import pandas as pd
>>> import numpy as np
>>> s = pd.Series(np.random.randn(10**6))
>>> s.nlargest(10) # only sorts up to the N requested
219921 4.644710
        4.608745
82124
421689 4.564644
425277
       4.447014
718691
      4.414137
43154
        4.403520
283187
       4.313922
595519
       4.273635
503969
       4.250236
       4.240952
121637
dtype: float64
```

### nonzero()

Return the *integer* indices of the elements that are non-zero

This method is equivalent to calling *numpy.nonzero* on the series data. For compatibility with NumPy, the return value is the same (a tuple with an array of indices for each dimension), but it will always be a one-item tuple because series only have one dimension.

### See also:

```
numpy.nonzero
```

# **Examples**

```
>>> s = pd.Series([0, 3, 0, 4], index=['a', 'b', 'c', 'd'])
# same return although index of s is different
>>> s.nonzero()
(array([1, 3]),)
>>> s.iloc[s.nonzero()[0]]
b     3
d     4
dtype: int64
```

## notna()

Detect existing (non-missing) values.

Return a boolean same-sized object indicating if the values are not NA. Non-missing values get mapped to True. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True). NA values, such as None or numpy.NaN, get mapped to False values.

### Returns

**Series** Mask of bool values for each element in Series that indicates whether an element is not an NA value.

See also:

Series.notnull alias of notna

Series.isna boolean inverse of notna

Series.dropna omit axes labels with missing values

notna top-level notna

# **Examples**

Show which entries in a DataFrame are not NA.

```
>>> df = pd.DataFrame({ 'age': [5, 6, np.NaN],
                      'born': [pd.NaT, pd.Timestamp('1939-05-27'),
                               pd.Timestamp('1940-04-25')],
                      'name': ['Alfred', 'Batman', ''],
. . .
                      'toy': [None, 'Batmobile', 'Joker'] })
. . .
>>> df
  age
           born name
                                t.ov
            NaT Alfred
0 5.0
                              None
1 6.0 1939-05-27 Batman Batmobile
2 NaN 1940-04-25
                              Joker
```

```
>>> df.notna()
age born name toy

0 True False True False

1 True True True True

2 False True True True
```

Show which entries in a Series are not NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.notna()
0 True
1 True
2 False
dtype: bool
```

#### notnull()

Detect existing (non-missing) values.

Return a boolean same-sized object indicating if the values are not NA. Non-missing values get mapped to True. Characters such as empty strings '' or numpy.inf are not considered NA values (unless you set pandas.options.mode.use\_inf\_as\_na = True). NA values, such as None or numpy.NaN, get mapped to False values.

### **Returns**

**Series** Mask of bool values for each element in Series that indicates whether an element is not an NA value.

See also:

Series.notnull alias of notna

Series.isna boolean inverse of notna

Series.dropna omit axes labels with missing values

notna top-level notna

## **Examples**

Show which entries in a DataFrame are not NA.

```
>>> df = pd.DataFrame({'age': [5, 6, np.NaN],
                       'born': [pd.NaT, pd.Timestamp('1939-05-27'),
. . .
                               pd.Timestamp('1940-04-25')],
. . .
                       'name': ['Alfred', 'Batman', ''],
. . .
                       'toy': [None, 'Batmobile', 'Joker']})
. . .
>>> df
  age
           born name
                               toy
 5.0
            NaT Alfred
1 6.0 1939-05-27 Batman Batmobile
2 NaN 1940-04-25
                               Joker
```

```
>>> df.notna()
age born name toy

0 True False True False

1 True True True True

2 False True True True
```

Show which entries in a Series are not NA.

```
>>> ser = pd.Series([5, 6, np.NaN])
>>> ser
0    5.0
1    6.0
2    NaN
dtype: float64
```

```
>>> ser.notna()
0 True
1 True
2 False
dtype: bool
```

# nsmallest (n=5, keep='first')

Return the smallest *n* elements.

### **Parameters**

n [int] Return this many ascending sorted values

**keep** [{'first', 'last'}, default 'first'] Where there are duplicate values: - first: take the first occurrence. - last: take the last occurrence.

### **Returns**

**bottom\_n** [Series] The n smallest values in the Series, in sorted order

## See also:

Series.nlargest

## **Notes**

Faster than .sort\_values().head(n) for small *n* relative to the size of the Series object.

## **Examples**

```
>>> import pandas as pd
>>> import numpy as np
>>> s = pd.Series(np.random.randn(10**6))
>>> s.nsmallest(10) # only sorts up to the N requested
288532 -4.954580
732345
       -4.835960
64803
        -4.812550
       -4.609998
446457
501225
        -4.483945
669476
       -4.472935
973615
       -4.401699
```

(continues on next page)

```
621279 -4.355126
773916 -4.347355
359919 -4.331927
dtype: float64
```

## nunique (dropna=True)

Return number of unique elements in the object.

Excludes NA values by default.

### **Parameters**

dropna [boolean, default True] Don't include NaN in the count.

#### Returns

```
nunique [int]
```

 $\verb|pct_change| (periods=1, fill_method='pad', limit=None, freq=None, **kwargs)|$ 

Percentage change between the current and a prior element.

Computes the percentage change from the immediately previous row by default. This is useful in comparing the percentage of change in a time series of elements.

### **Parameters**

periods [int, default 1] Periods to shift for forming percent change.

fill\_method [str, default 'pad'] How to handle NAs before computing percent changes.

**limit** [int, default None] The number of consecutive NAs to fill before stopping.

**freq** [DateOffset, timedelta, or offset alias string, optional] Increment to use from time series API (e.g. 'M' or BDay()).

\*\*kwargs Additional keyword arguments are passed into DataFrame.shift or Series.shift.

### Returns

**chg** [Series or DataFrame] The same type as the calling object.

### See also:

**Series.diff** Compute the difference of two elements in a Series.

**DataFrame**. **diff** Compute the difference of two elements in a DataFrame.

Series.shift Shift the index by some number of periods.

**DataFrame.shift** Shift the index by some number of periods.

# **Examples**

# **Series**

```
>>> s = pd.Series([90, 91, 85])
>>> s
0 90
1 91
2 85
dtype: int64
```

See the percentage change in a Series where filling NAs with last valid observation forward to next valid.

## **DataFrame**

Percentage change in French franc, Deutsche Mark, and Italian lira from 1980-01-01 to 1980-03-01.

```
>>> df = pd.DataFrame({
... 'FR': [4.0405, 4.0963, 4.3149],
... 'GR': [1.7246, 1.7482, 1.8519],
... 'IT': [804.74, 810.01, 860.13]},
... index=['1980-01-01', '1980-02-01', '1980-03-01'])
>>> df

FR GR IT

1980-01-01 4.0405 1.7246 804.74

1980-02-01 4.0963 1.7482 810.01

1980-03-01 4.3149 1.8519 860.13
```

```
>>> df.pct_change()

FR GR IT

1980-01-01 NaN NaN NaN

1980-02-01 0.013810 0.013684 0.006549

1980-03-01 0.053365 0.059318 0.061876
```

Percentage of change in GOOG and APPL stock volume. Shows computing the percentage change between columns.

```
>>> df = pd.DataFrame({
... '2016': [1769950, 30586265],
... '2015': [1500923, 40912316],
```

(continues on next page)

## **Parameters**

func [function] function to apply to the NDFrame. args, and kwargs are
 passed into func. Alternatively a (callable, data\_keyword) tuple where
 data\_keyword is a string indicating the keyword of callable that expects the
 NDFrame.

args [iterable, optional] positional arguments passed into func.

**kwargs** [mapping, optional] a dictionary of keyword arguments passed into func.

## Returns

object [the return type of func.]

## See also:

pandas.DataFrame.apply, pandas.DataFrame.applymap, pandas.Series.map

## **Notes**

Use .pipe when chaining together functions that expect Series, DataFrames or GroupBy objects. Instead of writing

```
>>> f(g(h(df), arg1=a), arg2=b, arg3=c)
```

You can write

```
>>> (df.pipe(h)
... .pipe(g, arg1=a)
... .pipe(f, arg2=b, arg3=c)
... )
```

If you have a function that takes the data as (say) the second argument, pass a tuple indicating which keyword expects the data. For example, suppose f takes its data as arg2:

```
>>> (df.pipe(h)
... .pipe(g, arg1=a)
... .pipe((f, 'arg2'), arg1=a, arg3=c)
... )
```

# plot

alias of pandas.plotting.\_core.SeriesPlotMethods

```
pop (item)
```

Return item and drop from frame. Raise KeyError if not found.

### **Parameters**

item [str] Column label to be popped

#### Returns

popped [Series]

# **Examples**

```
>>> df = pd.DataFrame([('falcon', 'bird',
                                              389.0),
                        ('parrot', 'bird',
                                               24.0),
. . .
                        ('lion', 'mammal', 80.5), ('monkey', 'mammal', np.nan)],
. . .
. . .
                       columns=('name', 'class', 'max_speed'))
>>> df
     name class max_speed
           bird
                    389.0
  falcon
  parrot
            bird
                         24.0
2
    lion mammal
                         80.5
3
  monkey mammal
                         NaN
```

```
>>> df.pop('class')
0 bird
1 bird
2 mammal
3 mammal
Name: class, dtype: object
```

```
>>> df
    name max_speed
0 falcon 389.0
1 parrot 24.0
2 lion 80.5
3 monkey NaN
```

**pow** (other, level=None, fill\_value=None, axis=0)

Exponential power of series and other, element-wise (binary operator pow).

Equivalent to series \*\* other, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

**other** [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

# Returns

result [Series]

### See also:

```
Series.rpow
```

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
b
    1.0
    1.0
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
а
    1.0
b
    NaN
d
     1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
    1.0
b
С
    1.0
d
    1.0
    NaN
dtype: float64
```

prod (axis=None, skipna=None, level=None, numeric\_only=None, min\_count=0, \*\*kwargs)
Return the product of the values for the requested axis

# **Parameters**

```
axis [\{index (0)\}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

**min\_count** [int, default 0] The required number of valid values to perform the operation. If fewer than min\_count\_non-NA values are present the result will be NA.

New in version 0.22.0: Added with the default being 0. This means the sum of an all-NA or empty Series is 0, and the product of an all-NA or empty Series is 1.

## Returns

prod [scalar or Series (if level specified)]

# **Examples**

By default, the product of an empty or all-NA Series is 1

```
>>> pd.Series([]).prod()
1.0
```

This can be controlled with the min\_count parameter

```
>>> pd.Series([]).prod(min_count=1)
nan
```

Thanks to the skipna parameter, min\_count handles all-NA and empty series identically.

```
>>> pd.Series([np.nan]).prod()
1.0
```

```
>>> pd.Series([np.nan]).prod(min_count=1)
nan
```

**product** (axis=None, skipna=None, level=None, numeric\_only=None, min\_count=0, \*\*kwargs)

Return the product of the values for the requested axis

## **Parameters**

```
axis [\{index (0)\}]
```

**skipna** [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

min\_count [int, default 0] The required number of valid values to perform the operation. If fewer than min\_count non-NA values are present the result will be NA.

New in version 0.22.0: Added with the default being 0. This means the sum of an all-NA or empty Series is 0, and the product of an all-NA or empty Series is 1.

## Returns

**prod** [scalar or Series (if level specified)]

# **Examples**

By default, the product of an empty or all-NA Series is 1

```
>>> pd.Series([]).prod()
1.0
```

This can be controlled with the min\_count parameter

```
>>> pd.Series([]).prod(min_count=1)
nan
```

Thanks to the skipna parameter, min\_count handles all-NA and empty series identically.

```
>>> pd.Series([np.nan]).prod()
1.0
```

```
>>> pd.Series([np.nan]).prod(min_count=1)
nan
```

ptp (axis=None, skipna=None, level=None, numeric\_only=None, \*\*kwargs)

**Returns the difference between the maximum value and the** minimum value in the object. This is the equivalent of the numpy.ndarray method ptp.

#### **Parameters**

```
axis [\{index (0)\}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

## Returns

```
ptp [scalar or Series (if level specified)]
```

```
put (*args, **kwargs)
```

Applies the *put* method to its *values* attribute if it has one.

### See also:

```
numpy.ndarray.put
```

```
quantile (q=0.5, interpolation='linear')
```

Return value at the given quantile, a la numpy.percentile.

## **Parameters**

**q** [float or array-like, default 0.5 (50% quantile)]  $0 \le q \le 1$ , the quantile(s) to compute **interpolation** [{'linear', 'lower', 'higher', 'midpoint', 'nearest'}] New in version 0.18.0.

This optional parameter specifies the interpolation method to use, when the desired quantile lies between two data points i and j:

- linear: i + (j i) \* fraction, where fraction is the fractional part of the index surrounded by i and j.
- lower: *i*.
- higher: *j*.
- nearest: i or j whichever is nearest.
- midpoint: (i + j) / 2.

## **Returns**

**quantile** [float or Series] if q is an array, a Series will be returned where the index is q and the values are the quantiles.

### See also:

```
pandas.core.window.Rolling.quantile
```

# **Examples**

**radd** (other, level=None, fill\_value=None, axis=0)

Addition of series and other, element-wise (binary operator radd).

Equivalent to other + series, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

**other** [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### **Returns**

result [Series]

## See also:

Series.add

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
а
    1.0
b
     1.0
     1.0
С
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
    1.0
а
    NaN
b
d
    1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
а
b
     1.0
     1.0
С
d
     1.0
```

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```
e NaN dtype: float64
```

Compute numerical data ranks (1 through n) along axis. Equal values are assigned a rank that is the average of the ranks of those values

#### **Parameters**

axis [{0 or 'index', 1 or 'columns'}, default 0] index to direct ranking
method [{'average', 'min', 'max', 'first', 'dense'}]

- average: average rank of group
- min: lowest rank in group
- max: highest rank in group
- first: ranks assigned in order they appear in the array
- dense: like 'min', but rank always increases by 1 between groups

numeric\_only [boolean, default None] Include only float, int, boolean data. Valid only
for DataFrame or Panel objects

na\_option [{'keep', 'top', 'bottom'}]

- keep: leave NA values where they are
- · top: smallest rank if ascending
- bottom: smallest rank if descending

ascending [boolean, default True] False for ranks by high (1) to low (N)

pct [boolean, default False] Computes percentage rank of data

# Returns

ranks [same type as caller]

ravel (order='C')

Return the flattened underlying data as an ndarray

### See also:

```
numpy.ndarray.ravel
```

rdiv (other, level=None, fill\_value=None, axis=0)

Floating division of series and other, element-wise (binary operator *rtruediv*).

Equivalent to other / series, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

other [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

#### Returns

result [Series]

### See also:

Series.truediv

## **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
     1.0
h
     1.0
С
     NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
    1.0
а
    NaN
b
d
    1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
h
     1.0
     1.0
С
d
     1.0
     NaN
dtype: float64
```

### real

## reindex (index=None, \*\*kwargs)

Conform Series to new index with optional filling logic, placing NA/NaN in locations having no value in the previous index. A new object is produced unless the new index is equivalent to the current one and copy=False

## **Parameters**

**index** [array-like, optional (should be specified using keywords)] New labels / index to conform to. Preferably an Index object to avoid duplicating data

**method** [{None, 'backfill'/'bfill', 'pad'/'ffill', 'nearest'}, optional] method to use for filling holes in reindexed DataFrame. Please note: this is only applicable to DataFrames/Series with a monotonically increasing/decreasing index.

- default: don't fill gaps
- pad / ffill: propagate last valid observation forward to next valid
- backfill / bfill: use next valid observation to fill gap
- nearest: use nearest valid observations to fill gap

copy [boolean, default True] Return a new object, even if the passed indexes are the same

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

**fill\_value** [scalar, default np.NaN] Value to use for missing values. Defaults to NaN, but can be any "compatible" value

**limit** [int, default None] Maximum number of consecutive elements to forward or backward fill

**tolerance** [optional] Maximum distance between original and new labels for inexact matches. The values of the index at the matching locations most satisfy the equation abs (index[indexer] - target) <= tolerance.

Tolerance may be a scalar value, which applies the same tolerance to all values, or list-like, which applies variable tolerance per element. List-like includes list, tuple, array, Series, and must be the same size as the index and its dtype must exactly match the index's type.

New in version 0.21.0: (list-like tolerance)

#### Returns

reindexed [Series]

## **Examples**

DataFrame.reindex supports two calling conventions

- (index=index\_labels, columns=column\_labels, ...)
- (labels, axis={'index', 'columns'}, ...)

We highly recommend using keyword arguments to clarify your intent.

Create a dataframe with some fictional data.

```
>>> index = ['Firefox', 'Chrome', 'Safari', 'IE10', 'Konqueror']
>>> df = pd.DataFrame({
         'http_status': [200,200,404,404,301],
         'response_time': [0.04, 0.02, 0.07, 0.08, 1.0]},
. . .
          index=index)
. . .
>>> df
          http_status response_time
                  200
                                 0.04
Firefox
                   200
                                  0.02
Chrome
                                  0.07
Safari
                   404
IE10
                   404
                                  0.08
Konqueror
                   301
                                  1.00
```

Create a new index and reindex the dataframe. By default values in the new index that do not have corresponding records in the dataframe are assigned NaN.

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IE10	404.0	0.08
Chrome	200.0	0.02

We can fill in the missing values by passing a value to the keyword fill\_value. Because the index is not monotonically increasing or decreasing, we cannot use arguments to the keyword method to fill the NaN values.

```
>>> df.reindex(new_index, fill_value=0)
             http_status response_time
                   404
Safari
                             0.07
                      0
                                  0.00
Iceweasel
                      0
Comodo Dragon
                                  0.00
IE10
                     404
                                  0.08
Chrome
                     200
                                   0.02
```

We can also reindex the columns.

```
>>> df.reindex(columns=['http_status', 'user_agent'])
          http_status user_agent
Firefox
                  200
                              NaN
                  200
                              NaN
Chrome
Safari
                  404
                              NaN
IE10
                  404
                              NaN
Konqueror
                  301
                               NaN
```

Or we can use "axis-style" keyword arguments

```
>>> df.reindex(['http_status', 'user_agent'], axis="columns")
         http_status user_agent
Firefox
                 200
                          NaN
                  200
Chrome
                             NaN
Safari
                  404
                             NaN
                  404
TE10
                             NaN
                  301
Konqueror
                             NaN
```

To further illustrate the filling functionality in reindex, we will create a dataframe with a monotonically increasing index (for example, a sequence of dates).

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```
2010-01-05 89
2010-01-06 88
```

Suppose we decide to expand the dataframe to cover a wider date range.

```
>>> date_index2 = pd.date_range('12/29/2009', periods=10, freq='D')
>>> df2.reindex(date_index2)
           prices
2009-12-29
             NaN
2009-12-30
              NaN
2009-12-31
              NaN
2010-01-01
              100
2010-01-02
              101
2010-01-03
              NaN
2010-01-04
               100
2010-01-05
               89
2010-01-06
                88
2010-01-07
               NaN
```

The index entries that did not have a value in the original data frame (for example, '2009-12-29') are by default filled with NaN. If desired, we can fill in the missing values using one of several options.

For example, to backpropagate the last valid value to fill the NaN values, pass bfill as an argument to the method keyword.

```
>>> df2.reindex(date_index2, method='bfill')
            prices
2009-12-29
               100
2009-12-30
               100
2009-12-31
               100
2010-01-01
               100
2010-01-02
               101
2010-01-03
               NaN
2010-01-04
               100
2010-01-05
                89
2010-01-06
                88
2010-01-07
               NaN
```

Please note that the NaN value present in the original dataframe (at index value 2010-01-03) will not be filled by any of the value propagation schemes. This is because filling while reindexing does not look at dataframe values, but only compares the original and desired indexes. If you do want to fill in the NaN values present in the original dataframe, use the fillna() method.

See the user guide for more.

```
reindex_axis (labels, axis=0, **kwargs)
```

Conform Series to new index with optional filling logic.

Deprecated since version 0.21.0: Use Series . reindex instead.

 $\verb"reindex_like" (other, method=None, copy=True, limit=None, tolerance=None)$ 

Return an object with matching indices to myself.

### **Parameters**

```
other [Object]method [string or None]copy [boolean, default True]
```

**limit** [int, default None] Maximum number of consecutive labels to fill for inexact matches.

**tolerance** [optional] Maximum distance between labels of the other object and this object for inexact matches. Can be list-like.

New in version 0.21.0: (list-like tolerance)

#### Returns

reindexed [same as input]

#### **Notes**

Like calling s.reindex(index=other.index, columns=other.columns, method=...)

```
rename (index=None, **kwargs)
```

Alter Series index labels or name

Function / dict values must be unique (1-to-1). Labels not contained in a dict / Series will be left as-is. Extra labels listed don't throw an error.

Alternatively, change Series.name with a scalar value.

See the user guide for more.

### **Parameters**

index [scalar, hashable sequence, dict-like or function, optional] dict-like or functions are transformations to apply to the index. Scalar or hashable sequence-like will alter the Series.name attribute.

copy [boolean, default True] Also copy underlying data

**inplace** [boolean, default False] Whether to return a new Series. If True then value of copy is ignored.

**level** [int or level name, default None] In case of a MultiIndex, only rename labels in the specified level.

# Returns

```
renamed [Series (new object)]
```

### See also:

```
pandas.Series.rename_axis
```

## **Examples**

(continues on next page)

```
Name: my_name, dtype: int64
>>> s.rename(lambda x: x ** 2)  # function, changes labels
0    1
1    2
4    3
dtype: int64
>>> s.rename({1: 3, 2: 5})  # mapping, changes labels
0    1
3    2
5    3
dtype: int64
```

rename\_axis (mapper, axis=0, copy=True, inplace=False)

Alter the name of the index or columns.

#### **Parameters**

mapper [scalar, list-like, optional] Value to set as the axis name attribute.

axis [{0 or 'index', 1 or 'columns'}, default 0] The index or the name of the axis.

copy [boolean, default True] Also copy underlying data.

**inplace** [boolean, default False] Modifies the object directly, instead of creating a new Series or DataFrame.

### Returns

**renamed** [Series, DataFrame, or None] The same type as the caller or None if *inplace* is True.

### See also:

```
pandas.Series.rename Alter Series index labels or name
pandas.DataFrame.rename Alter DataFrame index labels or name
pandas.Index.rename Set new names on index
```

## **Notes**

Prior to version 0.21.0, rename\_axis could also be used to change the axis *labels* by passing a mapping or scalar. This behavior is deprecated and will be removed in a future version. Use rename instead.

# **Examples**

## **Series**

```
>>> s = pd.Series([1, 2, 3])
>>> s.rename_axis("foo")
foo
0    1
1    2
2    3
dtype: int64
```

## **DataFrame**

```
>>> df.rename_axis("bar", axis="columns")
bar A B
0 1 4
1 2 5
2 3 6
```

### reorder levels(order)

Rearrange index levels using input order. May not drop or duplicate levels

#### **Parameters**

order [list of int representing new level order.] (reference level by number or key)axis [where to reorder levels]

### Returns

type of caller (new object)

### repeat (\*\*kwargs)

Repeat elements of an Series. Refer to *numpy.ndarray.repeat* for more information about the *repeats* argument.

### See also:

```
numpy.ndarray.repeat
```

**replace** (to\_replace=None, value=None, inplace=False, limit=None, regex=False, method='pad')
Replace values given in to\_replace with value.

Values of the Series are replaced with other values dynamically. This differs from updating with .loc or .iloc, which require you to specify a location to update with some value.

### **Parameters**

**to\_replace** [str, regex, list, dict, Series, int, float, or None] How to find the values that will be replaced.

- numeric, str or regex:
  - numeric: numeric values equal to to\_replace will be replaced with value
  - str: string exactly matching to\_replace will be replaced with value
  - regex: regexs matching to\_replace will be replaced with value
- list of str, regex, or numeric:
  - First, if *to\_replace* and *value* are both lists, they **must** be the same length.
  - Second, if regex=True then all of the strings in **both** lists will be interpreted
    as regexs otherwise they will match directly. This doesn't matter much for *value*since there are only a few possible substitution regexes you can use.
  - str, regex and numeric rules apply as above.
- dict:

- Dicts can be used to specify different replacement values for different existing values. For example, {'a': 'b', 'y': 'z'} replaces the value 'a' with 'b' and 'y' with 'z'. To use a dict in this way the *value* parameter should be *None*.
- For a DataFrame a dict can specify that different values should be replaced in different columns. For example, { 'a': 1, 'b': 'z'} looks for the value 1 in column 'a' and the value 'z' in column 'b' and replaces these values with whatever is specified in *value*. The *value* parameter should not be None in this case. You can treat this as a special case of passing two lists except that you are specifying the column to search in.
- For a DataFrame nested dictionaries, e.g., { 'a': {'b': np.nan}}, are read as follows: look in column 'a' for the value 'b' and replace it with NaN. The value parameter should be None to use a nested dict in this way. You can nest regular expressions as well. Note that column names (the top-level dictionary keys in a nested dictionary) cannot be regular expressions.

#### • None:

This means that the *regex* argument must be a string, compiled regular expression, or list, dict, ndarray or Series of such elements. If *value* is also None then this **must** be a nested dictionary or Series.

See the examples section for examples of each of these.

value [scalar, dict, list, str, regex, default None] Value to replace any values matching to\_replace with. For a DataFrame a dict of values can be used to specify which value to use for each column (columns not in the dict will not be filled). Regular expressions, strings and lists or dicts of such objects are also allowed.

**inplace** [boolean, default False] If True, in place. Note: this will modify any other views on this object (e.g. a column from a DataFrame). Returns the caller if this is True.

**limit** [int, default None] Maximum size gap to forward or backward fill.

**regex** [bool or same types as *to\_replace*, default False] Whether to interpret *to\_replace* and/or *value* as regular expressions. If this is True then *to\_replace must* be a string. Alternatively, this could be a regular expression or a list, dict, or array of regular expressions in which case *to\_replace* must be None.

**method** [{'pad', 'ffill', 'bfill', *None*}] The method to use when for replacement, when *to\_replace* is a scalar, list or tuple and *value* is None.

Changed in version 0.23.0: Added to DataFrame.

# Returns

**Series** Object after replacement.

## Raises

### AssertionError

• If regex is not a bool and to\_replace is not None.

### **TypeError**

- If to\_replace is a dict and value is not a list, dict, ndarray, or Series
- If *to\_replace* is None and *regex* is not compilable into a regular expression or is a list, dict, ndarray, or Series.

• When replacing multiple bool or datetime64 objects and the arguments to *to\_replace* does not match the type of the value being replaced

#### ValueError

• If a list or an ndarray is passed to *to\_replace* and *value* but they are not the same length.

### See also:

```
Series.fillna Fill NA values

Series.where Replace values based on boolean condition

Series.str.replace Simple string replacement.
```

## **Notes**

- Regex substitution is performed under the hood with re.sub. The rules for substitution for re.sub are the same.
- Regular expressions will only substitute on strings, meaning you cannot provide, for example, a regular expression matching floating point numbers and expect the columns in your frame that have a numeric dtype to be matched. However, if those floating point numbers *are* strings, then you can do this.
- This method has *a lot* of options. You are encouraged to experiment and play with this method to gain intuition about how it works.
- When dict is used as the *to\_replace* value, it is like key(s) in the dict are the to\_replace part and value(s) in the dict are the value parameter.

## **Examples**

## Scalar 'to\_replace' and 'value'

```
>>> df = pd.DataFrame({'A': [0, 1, 2, 3, 4],
                       'B': [5, 6, 7, 8, 9],
. . .
                       'C': ['a', 'b', 'c', 'd', 'e']})
. . .
>>> df.replace(0, 5)
  A B C
  5 5
0
        а
  1 6 b
1
2
  2 7
        С
3
  3
     8
        d
  4
     9
```

# List-like 'to\_replace'

```
>>> df.replace([0, 1, 2, 3], 4)

A B C

0 4 5 a

1 4 6 b

2 4 7 c

3 4 8 d

4 4 9 e
```

```
>>> df.replace([0, 1, 2, 3], [4, 3, 2, 1])

A B C

0 4 5 a

1 3 6 b

2 2 7 c

3 1 8 d

4 4 9 e
```

# dict-like 'to\_replace'

```
>>> df.replace({0: 10, 1: 100})

A B C

0 10 5 a

1 100 6 b

2 2 7 c

3 3 8 d

4 4 9 e
```

```
>>> df.replace({'A': 0, 'B': 5}, 100)
   A B C
0
 100 100
          а
      6
1
   1
          b
2
   2.
          С
3
   3
      8 d
4
  4 9 e
```

```
>>> df.replace({'A': {0: 100, 4: 400}})

A B C

0 100 5 a

1 1 6 b

2 2 7 c

3 3 8 d

4 400 9 e
```

# Regular expression 'to\_replace'

(continues on next page)

```
0 new abc
1 foo new
2 bait xyz
```

```
>>> df.replace(regex=r'^ba.$', value='new')

A B

0 new abc

1 foo new

2 bait xyz
```

```
>>> df.replace(regex={r'^ba.$':'new', 'foo':'xyz'})
        A    B
0    new abc
1    xyz    new
2    bait    xyz
```

```
>>> df.replace(regex=[r'^ba.$', 'foo'], value='new')

A B
0 new abc
1 new new
2 bait xyz
```

Note that when replacing multiple bool or datetime 64 objects, the data types in the *to\_replace* parameter must match the data type of the value being replaced:

This raises a TypeError because one of the dict keys is not of the correct type for replacement.

Compare the behavior of s.replace({'a': None}) and s.replace('a', None) to understand the pecularities of the *to\_replace* parameter:

```
>>> s = pd.Series([10, 'a', 'a', 'b', 'a'])
```

When one uses a dict as the *to\_replace* value, it is like the value(s) in the dict are equal to the *value* parameter. s.replace({'a': None}) is equivalent to s.replace(to\_replace={'a': None}, value=None, method=None):

(continues on next page)

```
4 None dtype: object
```

When value=None and to\_replace is a scalar, list or tuple, replace uses the method parameter (default 'pad') to do the replacement. So this is why the 'a' values are being replaced by 10 in rows 1 and 2 and 'b' in row 4 in this case. The command s.replace('a', None) is actually equivalent to s. replace(to\_replace='a', value=None, method='pad'):

```
>>> s.replace('a', None)
0    10
1    10
2    10
3    b
4    b
dtype: object
```

resample (rule, how=None, axis=0, fill\_method=None, closed=None, label=None, convention='start', kind=None, loffset=None, limit=None, base=0, on=None, level=None)

Convenience method for frequency conversion and resampling of time series. Object must have a datetime-like index (DatetimeIndex, PeriodIndex, or TimedeltaIndex), or pass datetime-like values to the on or level keyword.

#### **Parameters**

**rule** [string] the offset string or object representing target conversion

axis [int, optional, default 0]

**closed** [{'right', 'left'}] Which side of bin interval is closed. The default is 'left' for all frequency offsets except for 'M', 'A', 'Q', 'BM', 'BA', 'BQ', and 'W' which all have a default of 'right'.

label [{'right', 'left'}] Which bin edge label to label bucket with. The default is 'left' for all frequency offsets except for 'M', 'A', 'Q', 'BM', 'BA', 'BQ', and 'W' which all have a default of 'right'.

**convention** [{'start', 'end', 's', 'e'}] For PeriodIndex only, controls whether to use the start or end of *rule* 

kind: {'timestamp', 'period'}, optional Pass 'timestamp' to convert the resulting index to a DateTimeIndex or 'period' to convert it to a PeriodIndex. By default the input representation is retained.

**loffset** [timedelta] Adjust the resampled time labels

**base** [int, default 0] For frequencies that evenly subdivide 1 day, the "origin" of the aggregated intervals. For example, for '5min' frequency, base could range from 0 through 4. Defaults to 0

**on** [string, optional] For a DataFrame, column to use instead of index for resampling. Column must be datetime-like.

New in version 0.19.0.

**level** [string or int, optional] For a MultiIndex, level (name or number) to use for resampling. Level must be datetime-like.

New in version 0.19.0.

### Returns

# Resampler object

#### See also:

groupby Group by mapping, function, label, or list of labels.

### **Notes**

See the user guide for more.

To learn more about the offset strings, please see this link.

# **Examples**

Start by creating a series with 9 one minute timestamps.

```
>>> index = pd.date_range('1/1/2000', periods=9, freq='T')
>>> series = pd.Series(range(9), index=index)
>>> series
2000-01-01 00:00:00
2000-01-01 00:01:00
                       1
2000-01-01 00:02:00
                       2
2000-01-01 00:03:00
                       3
2000-01-01 00:04:00
                       4
2000-01-01 00:05:00
                       5
2000-01-01 00:06:00
                        6
2000-01-01 00:07:00
                       7
2000-01-01 00:08:00
                       8
Freq: T, dtype: int64
```

Downsample the series into 3 minute bins and sum the values of the timestamps falling into a bin.

Downsample the series into 3 minute bins as above, but label each bin using the right edge instead of the left. Please note that the value in the bucket used as the label is not included in the bucket, which it labels. For example, in the original series the bucket 2000-01-01 00:03:00 contains the value 3, but the summed value in the resampled bucket with the label 2000-01-01 00:03:00 does not include 3 (if it did, the summed value would be 6, not 3). To include this value close the right side of the bin interval as illustrated in the example below this one.

Downsample the series into 3 minute bins as above, but close the right side of the bin interval.

```
>>> series.resample('3T', label='right', closed='right').sum()
2000-01-01 00:00:00 0
```

(continues on next page)

```
2000-01-01 00:03:00 6

2000-01-01 00:06:00 15

2000-01-01 00:09:00 15

Freq: 3T, dtype: int64
```

Upsample the series into 30 second bins.

```
>>> series.resample('30S').asfreq()[0:5] #select first 5 rows
2000-01-01 00:00:00 0.0
2000-01-01 00:00:30 NaN
2000-01-01 00:01:00 1.0
2000-01-01 00:01:30 NaN
2000-01-01 00:02:00 2.0
Freq: 30S, dtype: float64
```

Upsample the series into 30 second bins and fill the NaN values using the pad method.

Upsample the series into 30 second bins and fill the NaN values using the bfill method.

Pass a custom function via apply

```
>>> def custom_resampler(array_like):
... return np.sum(array_like)+5
```

For a Series with a PeriodIndex, the keyword *convention* can be used to control whether to use the start or end of *rule*.

Resample by month using 'start' convention. Values are assigned to the first month of the period.

Resample by month using 'end' convention. Values are assigned to the last month of the period.

```
>>> s.resample('M', convention='end').asfreq()
2012-12
          1.0
2013-01
          NaN
2013-02
          NaN
2013-03
          NaN
2013-04
          NaN
2013-05
          NaN
2013-06
          NaN
2013-07
          NaN
2013-08
          NaN
2013-09
          NaN
2013-10
          NaN
2013-11
          NaN
2013-12
          2.0
Freq: M, dtype: float64
```

For DataFrame objects, the keyword on can be used to specify the column instead of the index for resampling.

For a DataFrame with MultiIndex, the keyword level can be used to specify on level the resampling needs to take place.

reset\_index (level=None, drop=False, name=None, inplace=False)

Generate a new DataFrame or Series with the index reset.

This is useful when the index needs to be treated as a column, or when the index is meaningless and needs to be reset to the default before another operation.

### **Parameters**

**level** [int, str, tuple, or list, default optional] For a Series with a MultiIndex, only remove the specified levels from the index. Removes all levels by default.

**drop** [bool, default False] Just reset the index, without inserting it as a column in the new DataFrame.

**name** [object, optional] The name to use for the column containing the original Series values. Uses self.name by default. This argument is ignored when *drop* is True.

inplace [bool, default False] Modify the Series in place (do not create a new object).

### **Returns**

**Series or DataFrame** When *drop* is False (the default), a DataFrame is returned. The newly created columns will come first in the DataFrame, followed by the original Series values. When *drop* is True, a *Series* is returned. In either case, if inplace=True, no value is returned.

### See also:

**DataFrame.reset\_index** Analogous function for DataFrame.

## **Examples**

```
>>> s = pd.Series([1, 2, 3, 4], name='foo',
... index=pd.Index(['a', 'b', 'c', 'd'], name='idx'))
```

Generate a DataFrame with default index.

```
>>> s.reset_index()
    idx foo
0 a 1
1 b 2
2 c 3
3 d 4
```

To specify the name of the new column use *name*.

```
>>> s.reset_index(name='values')
    idx values
0 a 1
1 b 2
2 c 3
3 d 4
```

To generate a new Series with the default set *drop* to True.

To update the Series in place, without generating a new one set *inplace* to True. Note that it also requires drop=True.

The level parameter is interesting for Series with a multi-level index.

To remove a specific level from the Index, use level.

If *level* is not set, all levels are removed from the Index.

```
>>> s2.reset_index()
    a    b    foo

0    bar    one    0
1    bar    two    1
2    baz    one    2
3    baz    two    3
```

# rfloordiv (other, level=None, fill\_value=None, axis=0)

Integer division of series and other, element-wise (binary operator rfloordiv).

Equivalent to other // series, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

other [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### Returns

result [Series]

# See also:

Series.floordiv

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
b
     1.0
С
     1.0
d
     NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
    1.0
b
     NaN
d
     1.0
е
     NaN
dtype: float64
>>> a.add(b, fill_value=0)
     2.0
b
     1.0
С
     1.0
d
     1.0
0
     NaN
dtype: float64
```

**rmod** (other, level=None, fill\_value=None, axis=0)

Modulo of series and other, element-wise (binary operator *rmod*).

Equivalent to other % series, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

**other** [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

# Returns

result [Series]

## See also:

Series.mod

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
a    1.0
b    1.0
c    1.0
d    NaN
```

(continues on next page)

```
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
а
    1.0
b
    NaN
d
     1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
а
    1.0
b
С
    1.0
d
    1.0
    NaN
dtype: float64
```

rmul (other, level=None, fill\_value=None, axis=0)

Multiplication of series and other, element-wise (binary operator *rmul*).

Equivalent to other \* series, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

other [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### **Returns**

result [Series]

# See also:

Series.mul

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
     1.0
b
С
    1.0
d
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
     1.0
а
b
     NaN
d
     1.0
е
     NaN
```

(continues on next page)

```
dtype: float64
>>> a.add(b, fill_value=0)
a    2.0
b    1.0
c    1.0
d    1.0
e    NaN
dtype: float64
```

Provides rolling window calculations.

New in version 0.18.0.

#### **Parameters**

window [int, or offset] Size of the moving window. This is the number of observations used for calculating the statistic. Each window will be a fixed size.

If its an offset then this will be the time period of each window. Each window will be a variable sized based on the observations included in the time-period. This is only valid for datetimelike indexes. This is new in 0.19.0

**min\_periods** [int, default None] Minimum number of observations in window required to have a value (otherwise result is NA). For a window that is specified by an offset, this will default to 1.

center [boolean, default False] Set the labels at the center of the window.

win\_type [string, default None] Provide a window type. If None, all points are evenly weighted. See the notes below for further information.

**on** [string, optional] For a DataFrame, column on which to calculate the rolling window, rather than the index

**closed** [string, default None] Make the interval closed on the 'right', 'left', 'both' or 'neither' endpoints. For offset-based windows, it defaults to 'right'. For fixed windows, defaults to 'both'. Remaining cases not implemented for fixed windows.

New in version 0.20.0.

axis [int or string, default 0]

### Returns

a Window or Rolling sub-classed for the particular operation

### See also:

expanding Provides expanding transformations.

ewm Provides exponential weighted functions

# **Notes**

By default, the result is set to the right edge of the window. This can be changed to the center of the window by setting center=True.

To learn more about the offsets & frequency strings, please see this link.

The recognized win\_types are:

- boxcar
- triang
- blackman
- hamming
- bartlett
- parzen
- bohman
- blackmanharris
- nuttall
- barthann
- kaiser (needs beta)
- gaussian (needs std)
- general\_gaussian (needs power, width)
- slepian (needs width).

If win\_type=None all points are evenly weighted. To learn more about different window types see scipy.signal window functions.

# **Examples**

Rolling sum with a window length of 2, using the 'triang' window type.

```
>>> df.rolling(2, win_type='triang').sum()

B
0 NaN
1 1.0
2 2.5
3 NaN
4 NaN
```

Rolling sum with a window length of 2, min\_periods defaults to the window length.

```
>>> df.rolling(2).sum()
B
0 NaN
1 1.0
2 3.0
```

(continues on next page)

```
3 NaN
4 NaN
```

Same as above, but explicitly set the min\_periods

```
>>> df.rolling(2, min_periods=1).sum()

B
0 0.0
1 1.0
2 3.0
3 2.0
4 4.0
```

A ragged (meaning not-a-regular frequency), time-indexed DataFrame

```
>>> df = pd.DataFrame({'B': [0, 1, 2, np.nan, 4]},
... index = [pd.Timestamp('20130101 09:00:00'),
pd.Timestamp('20130101 09:00:02'),
pd.Timestamp('20130101 09:00:03'),
pd.Timestamp('20130101 09:00:05'),
pd.Timestamp('20130101 09:00:06')])
```

```
>>> df

B

2013-01-01 09:00:00 0.0
2013-01-01 09:00:02 1.0
2013-01-01 09:00:03 2.0
2013-01-01 09:00:05 NaN
2013-01-01 09:00:06 4.0
```

Contrasting to an integer rolling window, this will roll a variable length window corresponding to the time period. The default for min\_periods is 1.

```
>>> df.rolling('2s').sum()

B

2013-01-01 09:00:00 0.0

2013-01-01 09:00:02 1.0

2013-01-01 09:00:03 3.0

2013-01-01 09:00:05 NaN

2013-01-01 09:00:06 4.0
```

round (decimals=0, \*args, \*\*kwargs)

Round each value in a Series to the given number of decimals.

### **Parameters**

**decimals** [int] Number of decimal places to round to (default: 0). If decimals is negative, it specifies the number of positions to the left of the decimal point.

# Returns

Series object

### See also:

```
numpy.around, DataFrame.round
```

```
rpow (other, level=None, fill_value=None, axis=0)
```

Exponential power of series and other, element-wise (binary operator *rpow*).

Equivalent to other \*\* series, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

other [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### **Returns**

result [Series]

### See also:

Series.pow

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
а
    1.0
b
    1.0
     1.0
С
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
     1.0
     NaN
     1.0
d
     NaN
е
dtype: float64
>>> a.add(b, fill_value=0)
     2.0
b
     1.0
     1.0
С
d
     1.0
     NaN
dtype: float64
```

rsub (other, level=None, fill\_value=None, axis=0)

Subtraction of series and other, element-wise (binary operator *rsub*).

Equivalent to other - series, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

**other** [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before

computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

## **Returns**

result [Series]

#### See also:

Series.sub

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
b
    1.0
     1.0
С
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
а
    1.0
     NaN
b
     1.0
d
    NaN
e
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
b
    1.0
    1.0
С
     1.0
    NaN
dtype: float64
```

## rtruediv (other, level=None, fill\_value=None, axis=0)

Floating division of series and other, element-wise (binary operator *rtruediv*).

Equivalent to other / series, but with support to substitute a fill\_value for missing data in one of the inputs.

### **Parameters**

**other** [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

## Returns

result [Series]

### See also:

Series.truediv

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
b
    1.0
     1.0
     NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
а
    1.0
b
     NaN
d
     1.0
0
     NaN
dtype: float64
>>> a.add(b, fill_value=0)
     2.0
а
b
     1.0
С
     1.0
d
     1.0
     NaN
e
dtype: float64
```

**sample** (n=None, frac=None, replace=False, weights=None, random\_state=None, axis=None) Return a random sample of items from an axis of object.

You can use *random\_state* for reproducibility.

## **Parameters**

**n** [int, optional] Number of items from axis to return. Cannot be used with frac. Default = 1 if frac = None.

**frac** [float, optional] Fraction of axis items to return. Cannot be used with n.

**replace** [boolean, optional] Sample with or without replacement. Default = False.

weights [str or ndarray-like, optional] Default 'None' results in equal probability weighting. If passed a Series, will align with target object on index. Index values in weights not found in sampled object will be ignored and index values in sampled object not in weights will be assigned weights of zero. If called on a DataFrame, will accept the name of a column when axis = 0. Unless weights are a Series, weights must be same length as axis being sampled. If weights do not sum to 1, they will be normalized to sum to 1. Missing values in the weights column will be treated as zero. inf and -inf values not allowed.

**random\_state** [int or numpy.random.RandomState, optional] Seed for the random number generator (if int), or numpy RandomState object.

**axis** [int or string, optional] Axis to sample. Accepts axis number or name. Default is stat axis for given data type (0 for Series and DataFrames, 1 for Panels).

# Returns

A new object of same type as caller.

# **Examples**

Generate an example Series and DataFrame:

```
>>> s = pd.Series(np.random.randn(50))
>>> s.head()
0
  -0.038497
1
    1.820773
2
  -0.972766
  -1.598270
  -1.095526
dtype: float64
>>> df = pd.DataFrame(np.random.randn(50, 4), columns=list('ABCD'))
>>> df.head()
                    В
                               С
0 0.016443 -2.318952 -0.566372 -1.028078
1 -1.051921 0.438836 0.658280 -0.175797
2 -1.243569 -0.364626 -0.215065 0.057736
3 \quad 1.768216 \quad 0.404512 \quad -0.385604 \quad -1.457834
4 1.072446 -1.137172 0.314194 -0.046661
```

Next extract a random sample from both of these objects...

3 random elements from the Series:

```
>>> s.sample(n=3)
27 -0.994689
55 -1.049016
67 -0.224565
dtype: float64
```

And a random 10% of the <code>DataFrame</code> with replacement:

```
>>> df.sample(frac=0.1, replace=True)

A B C D

35 1.981780 0.142106 1.817165 -0.290805
49 -1.336199 -0.448634 -0.789640 0.217116
40 0.823173 -0.078816 1.009536 1.015108
15 1.421154 -0.055301 -1.922594 -0.019696
6 -0.148339 0.832938 1.787600 -1.383767
```

You can use *random state* for reproducibility:

```
>>> df.sample(random_state=1)
A B C D
37 -2.027662 0.103611 0.237496 -0.165867
43 -0.259323 -0.583426 1.516140 -0.479118
12 -1.686325 -0.579510 0.985195 -0.460286
8 1.167946 0.429082 1.215742 -1.636041
9 1.197475 -0.864188 1.554031 -1.505264
```

## searchsorted(\*\*kwargs)

Find indices where elements should be inserted to maintain order.

Find the indices into a sorted Series *self* such that, if the corresponding elements in *value* were inserted before the indices, the order of *self* would be preserved.

# **Parameters**

**value** [array\_like] Values to insert into *self*.

**side** [{'left', 'right'}, optional] If 'left', the index of the first suitable location found is given. If 'right', return the last such index. If there is no suitable index, return either 0 or N (where N is the length of *self*).

**sorter** [1-D array\_like, optional] Optional array of integer indices that sort *self* into ascending order. They are typically the result of np.argsort.

## Returns

indices [array of ints] Array of insertion points with the same shape as value.

### See also:

```
numpy.searchsorted
```

### **Notes**

Binary search is used to find the required insertion points.

# **Examples**

```
>>> x.searchsorted(4)
array([3])
```

```
>>> x.searchsorted([0, 4])
array([0, 3])
```

```
>>> x.searchsorted([1, 3], side='left')
array([0, 2])
```

```
>>> x.searchsorted([1, 3], side='right')
array([1, 3])
```

```
>>> x.searchsorted('bread')
array([1])  # Note: an array, not a scalar
```

```
>>> x.searchsorted(['bread'], side='right')
array([3])
```

```
select (crit, axis=0)
```

Return data corresponding to axis labels matching criteria

Deprecated since version 0.21.0: Use df.loc[df.index.map(crit)] to select via labels

#### **Parameters**

crit [function] To be called on each index (label). Should return True or False
axis [int]

### Returns

**selection** [type of caller]

**sem** (axis=None, skipna=None, level=None, ddof=1, numeric\_only=None, \*\*kwargs)
Return unbiased standard error of the mean over requested axis.

Normalized by N-1 by default. This can be changed using the ddof argument

### **Parameters**

```
axis [\{index (0)\}]
```

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

**ddof** [int, default 1] Delta Degrees of Freedom. The divisor used in calculations is N - ddof, where N represents the number of elements.

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

sem [scalar or Series (if level specified)]

set\_axis (labels, axis=0, inplace=None)

Assign desired index to given axis.

Indexes for column or row labels can be changed by assigning a list-like or Index.

Changed in version 0.21.0: The signature is now *labels* and *axis*, consistent with the rest of pandas API. Previously, the *axis* and *labels* arguments were respectively the first and second positional arguments.

### **Parameters**

**labels** [list-like, Index] The values for the new index.

**axis** [{0 or 'index', 1 or 'columns'}, default 0] The axis to update. The value 0 identifies the rows, and 1 identifies the columns.

inplace [boolean, default None] Whether to return a new %(klass)s instance.

**Warning:** inplace=None currently falls back to to True, but in a future version, will default to False. Use inplace=True explicitly rather than relying on the default.

### Returns

**renamed** [%(klass)s or None] An object of same type as caller if inplace=False, None otherwise.

See also:

pandas.DataFrame.rename\_axis Alter the name of the index or columns.

# **Examples**

## **Series**

The original object is not modified.

# **DataFrame**

```
>>> df = pd.DataFrame({"A": [1, 2, 3], "B": [4, 5, 6]})
```

Change the row labels.

```
>>> df.set_axis(['a', 'b', 'c'], axis='index', inplace=False)

A B
a 1 4
b 2 5
c 3 6
```

Change the column labels.

Now, update the labels inplace.

```
>>> df.set_axis(['i', 'ii'], axis='columns', inplace=True)
>>> df
```

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```
i ii
0 1 4
1 2 5
2 3 6
```

### set\_value (label, value, takeable=False)

Quickly set single value at passed label. If label is not contained, a new object is created with the label placed at the end of the result index.

Deprecated since version 0.21.0: Please use .at[] or .iat[] accessors.

#### **Parameters**

```
label [object] Partial indexing with MultiIndex not allowedvalue [object] Scalar valuetakeable [interpret the index as indexers, default False]
```

### Returns

**series** [Series] If label is contained, will be reference to calling Series, otherwise a new object

### shape

return a tuple of the shape of the underlying data

```
shift (periods=1, freq=None, axis=0)
```

Shift index by desired number of periods with an optional time freq

### **Parameters**

```
periods [int] Number of periods to move, can be positive or negative
freq [DateOffset, timedelta, or time rule string, optional] Increment to use from the tseries module or time rule (e.g. 'EOM'). See Notes.
axis [{0 or 'index'}]
```

# Returns

shifted [Series]

# **Notes**

If freq is specified then the index values are shifted but the data is not realigned. That is, use freq if you would like to extend the index when shifting and preserve the original data.

### size

return the number of elements in the underlying data

```
skew (axis=None, skipna=None, level=None, numeric_only=None, **kwargs)
Return unbiased skew over requested axis Normalized by N-1
```

# **Parameters**

```
axis [{index (0)}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

**skew** [scalar or Series (if level specified)]

```
slice shift (periods=1, axis=0)
```

Equivalent to *shift* without copying data. The shifted data will not include the dropped periods and the shifted axis will be smaller than the original.

### **Parameters**

**periods** [int] Number of periods to move, can be positive or negative

#### Returns

shifted [same type as caller]

### **Notes**

While the *slice\_shift* is faster than *shift*, you may pay for it later during alignment.

Returns a new Series sorted by label if *inplace* argument is False, otherwise updates the original series and returns None.

#### **Parameters**

axis [int, default 0] Axis to direct sorting. This can only be 0 for Series.

**level** [int, optional] If not None, sort on values in specified index level(s).

ascending [bool, default true] Sort ascending vs. descending.

**inplace** [bool, default False] If True, perform operation in-place.

**kind** [{'quicksort', 'mergesort', 'heapsort'}, default 'quicksort'] Choice of sorting algorithm. See also numpy.sort() for more information. 'mergesort' is the only stable algorithm. For DataFrames, this option is only applied when sorting on a single column or label.

**na\_position** [{'first', 'last'}, default 'last'] If 'first' puts NaNs at the beginning, 'last' puts NaNs at the end. Not implemented for MultiIndex.

**sort\_remaining** [bool, default True] If true and sorting by level and index is multilevel, sort by other levels too (in order) after sorting by specified level.

## Returns

pandas.Series The original Series sorted by the labels

### See also:

```
DataFrame.sort_index Sort DataFrame by the index
DataFrame.sort_values Sort DataFrame by the value
Series.sort_values Sort Series by the value
```

## **Examples**

## Sort Descending

### Sort Inplace

# By default NaNs are put at the end, but use na\_position to place them at the beginning

# Specify index level to sort

```
>>> arrays = [np.array(['qux', 'qux', 'foo', 'foo',
                        'baz', 'baz', 'bar', 'bar']),
             np.array(['two', 'one', 'two', 'one',
. . .
                        'two', 'one', 'two', 'one'])]
>>> s = pd.Series([1, 2, 3, 4, 5, 6, 7, 8], index=arrays)
>>> s.sort_index(level=1)
bar one
           8
baz one
            6
foo one
           4
qux one
bar
    two
baz two
foo two
qux two
dtype: int64
```

Does not sort by remaining levels when sorting by levels

```
>>> s.sort_index(level=1, sort_remaining=False)
qux one 2
foo one 4
baz one 6
bar one 8
qux two 1
foo two 3
baz two 5
bar two 7
dtype: int64
```

**sort\_values** (axis=0, ascending=True, inplace=False, kind='quicksort', na\_position='last') Sort by the values.

Sort a Series in ascending or descending order by some criterion.

#### **Parameters**

**axis** [{0 or 'index'}, default 0] Axis to direct sorting. The value 'index' is accepted for compatibility with DataFrame.sort\_values.

**ascending** [bool, default True] If True, sort values in ascending order, otherwise descending.

inplace [bool, default False] If True, perform operation in-place.

**kind** [{'quicksort', 'mergesort' or 'heapsort'}, default 'quicksort'] Choice of sorting algorithm. See also numpy.sort() for more information. 'mergesort' is the only stable algorithm.

na\_position [{'first' or 'last'}, default 'last'] Argument 'first' puts NaNs at the beginning, 'last' puts NaNs at the end.

### Returns

Series Series ordered by values.

See also:

Series.sort\_index Sort by the Series indices.

**DataFrame.sort\_values** Sort DataFrame by the values along either axis.

DataFrame.sort\_index Sort DataFrame by indices.

# **Examples**

```
>>> s = pd.Series([np.nan, 1, 3, 10, 5])
>>> s

0    NaN
1    1.0
2    3.0
3    10.0
4    5.0
dtype: float64
```

Sort values ascending order (default behaviour)

## Sort values descending order

```
>>> s.sort_values(ascending=False)
3     10.0
4     5.0
2     3.0
1     1.0
0     NaN
dtype: float64
```

## Sort values inplace

```
>>> s.sort_values(ascending=False, inplace=True)
>>> s
3     10.0
4     5.0
2     3.0
1     1.0
0     NaN
dtype: float64
```

# Sort values putting NAs first

```
>>> s.sort_values(na_position='first')

0     NaN

1     1.0

2     3.0

4     5.0

3     10.0

dtype: float64
```

# Sort a series of strings

```
sortlevel (level=0, ascending=True, sort remaining=True)
```

Sort Series with MultiIndex by chosen level. Data will be lexicographically sorted by the chosen level followed by the other levels (in order),

Deprecated since version 0.20.0: Use Series.sort\_index()

### **Parameters**

**level** [int or level name, default None]

ascending [bool, default True]

Returns

sorted [Series]

#### See also:

Series.sort\_index

squeeze (axis=None)

Squeeze length 1 dimensions.

### **Parameters**

axis [None, integer or string axis name, optional] The axis to squeeze if 1-sized.

New in version 0.20.0.

### Returns

## scalar if 1-sized, else original object

std(axis=None, skipna=None, level=None, ddof=1, numeric\_only=None, \*\*kwargs)

Return sample standard deviation over requested axis.

Normalized by N-1 by default. This can be changed using the ddof argument

### **Parameters**

```
axis [\{index (0)\}]
```

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

**ddof** [int, default 1] Delta Degrees of Freedom. The divisor used in calculations is N - ddof, where N represents the number of elements.

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### Returns

**std** [scalar or Series (if level specified)]

str

 ${\bf alias} \ {\bf of} \ {\bf pandas.core.strings.String} \\ {\bf Methods}$ 

### strides

return the strides of the underlying data

```
sub (other, level=None, fill_value=None, axis=0)
```

Subtraction of series and other, element-wise (binary operator sub).

Equivalent to series - other, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

**other** [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### **Returns**

result [Series]

#### See also:

Series.rsub

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
    1.0
а
b
     1.0
     1.0
С
d
     NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
     1.0
а
b
     NaN
d
     1.0
     NaN
dtype: float64
>>> a.add(b, fill_value=0)
     2.0
b
     1.0
     1.0
С
d
     1.0
     NaN
dtype: float64
```

# subtract (other, level=None, fill\_value=None, axis=0)

Subtraction of series and other, element-wise (binary operator *sub*).

Equivalent to series - other, but with support to substitute a fill\_value for missing data in one of the inputs.

## **Parameters**

other [Series or scalar value]

**fill\_value** [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

#### Returns

```
result [Series]
```

### See also:

Series.rsub

## **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
а
    1.0
b
     1.0
     1.0
С
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
а
    1.0
    NaN
d
     1.0
    NaN
dtype: float64
>>> a.add(b, fill_value=0)
     2.0
     1.0
h
С
     1.0
d
     1.0
e
     NaN
dtype: float64
```

**sum** (axis=None, skipna=None, level=None, numeric\_only=None, min\_count=0, \*\*kwargs)
Return the sum of the values for the requested axis

#### **Parameters**

```
axis [\{index (0)\}]
```

skipna [boolean, default True] Exclude NA/null values when computing the result.

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

min\_count [int, default 0] The required number of valid values to perform the operation. If fewer than min\_count non-NA values are present the result will be NA.

New in version 0.22.0: Added with the default being 0. This means the sum of an all-NA or empty Series is 0, and the product of an all-NA or empty Series is 1.

#### Returns

sum [scalar or Series (if level specified)]

# **Examples**

By default, the sum of an empty or all-NA Series is 0.

```
>>> pd.Series([]).sum() # min_count=0 is the default 0.0
```

This can be controlled with the min\_count parameter. For example, if you'd like the sum of an empty series to be NaN, pass min count=1.

```
>>> pd.Series([]).sum(min_count=1)
nan
```

Thanks to the skipna parameter, min\_count handles all-NA and empty series identically.

```
>>> pd.Series([np.nan]).sum()
0.0
```

```
>>> pd.Series([np.nan]).sum(min_count=1)
nan
```

### swapaxes (axis1, axis2, copy=True)

Interchange axes and swap values axes appropriately

## Returns

y [same as input]

```
swaplevel(i=-2, j=-1, copy=True)
```

Swap levels i and j in a MultiIndex

#### **Parameters**

**i, j** [int, string (can be mixed)] Level of index to be swapped. Can pass level name as string.

### Returns

```
swapped [Series]
```

.. versionchanged:: 0.18.1 The indexes i and j are now optional, and default to the two innermost levels of the index.

## tail(n=5)

Return the last n rows.

This function returns last n rows from the object based on position. It is useful for quickly verifying data, for example, after sorting or appending rows.

#### **Parameters**

**n** [int, default 5] Number of rows to select.

## Returns

**type of caller** The last *n* rows of the caller object.

See also:

pandas.DataFrame.head The first n rows of the caller object.

# **Examples**

```
>>> df = pd.DataFrame({'animal':['alligator', 'bee', 'falcon', 'lion',
                        'monkey', 'parrot', 'shark', 'whale', 'zebra']})
>>> df
      animal
  alligator
1
         bee
2
      falcon
3
        lion
4
      monkey
5
      parrot
6
       shark
7
       whale
8
       zebra
```

### Viewing the last 5 lines

```
>>> df.tail()
animal
4 monkey
5 parrot
6 shark
7 whale
8 zebra
```

Viewing the last *n* lines (three in this case)

```
>>> df.tail(3)
animal
6 shark
7 whale
8 zebra
```

**take** (indices, axis=0, convert=None, is\_copy=True, \*\*kwargs)

Return the elements in the given positional indices along an axis.

This means that we are not indexing according to actual values in the index attribute of the object. We are indexing according to the actual position of the element in the object.

### **Parameters**

indices [array-like] An array of ints indicating which positions to take.

**axis** [{0 or 'index', 1 or 'columns', None}, default 0] The axis on which to select elements. 0 means that we are selecting rows, 1 means that we are selecting columns.

**convert** [bool, default True] Whether to convert negative indices into positive ones. For example, -1 would map to the len(axis) - 1. The conversions are similar to the behavior of indexing a regular Python list.

Deprecated since version 0.21.0: In the future, negative indices will always be converted.

is\_copy [bool, default True] Whether to return a copy of the original object or not.

\*\*kwargs For compatibility with numpy.take(). Has no effect on the output.

## **Returns**

taken [type of caller] An array-like containing the elements taken from the object.

See also:

DataFrame.loc Select a subset of a DataFrame by labels.

**DataFrame.iloc** Select a subset of a DataFrame by positions.

numpy.take Take elements from an array along an axis.

# **Examples**

```
>>> df = pd.DataFrame([('falcon', 'bird', 389.0),
                     ('parrot', 'bird',
                                          24.0),
                              'mammal', 80.5),
                     ('lion',
. . .
                     ('monkey', 'mammal', np.nan)],
. . .
                     columns=['name', 'class', 'max_speed'],
. . .
                     index=[0, 2, 3, 1])
. . .
>>> df
    name class max_speed
0 falcon bird 389.0
                    24.0
 parrot bird
3
  lion mammal
                     80.5
1 monkey mammal
                      NaN
```

Take elements at positions 0 and 3 along the axis 0 (default).

Note how the actual indices selected (0 and 1) do not correspond to our selected indices 0 and 3. That's because we are selecting the 0th and 3rd rows, not rows whose indices equal 0 and 3.

```
>>> df.take([0, 3])
    name class max_speed
0 falcon bird 389.0
1 monkey mammal NaN
```

Take elements at indices 1 and 2 along the axis 1 (column selection).

```
>>> df.take([1, 2], axis=1)
    class max_speed
0 bird 389.0
2 bird 24.0
3 mammal 80.5
1 mammal NaN
```

We may take elements using negative integers for positive indices, starting from the end of the object, just like with Python lists.

```
>>> df.take([-1, -2])
    name class max_speed
1 monkey mammal NaN
3 lion mammal 80.5
```

```
to_clipboard(excel=True, sep=None, **kwargs)
```

Copy object to the system clipboard.

Write a text representation of object to the system clipboard. This can be pasted into Excel, for example.

### **Parameters**

excel [bool, default True]

- True, use the provided separator, writing in a csv format for allowing easy pasting into excel.
- False, write a string representation of the object to the clipboard.

```
sep [str, default '\t'] Field delimiter.
```

\*\*kwargs These parameters will be passed to DataFrame.to\_csv.

See also:

DataFrame.to\_csv Write a DataFrame to a comma-separated values (csv) file.

read\_clipboard Read text from clipboard and pass to read\_table.

### **Notes**

Requirements for your platform.

- Linux : xclip, or xsel (with gtk or PyQt4 modules)
- Windows: none
- OS X : none

### **Examples**

Copy the contents of a DataFrame to the clipboard.

```
>>> df = pd.DataFrame([[1, 2, 3], [4, 5, 6]], columns=['A', 'B', 'C'])
>>> df.to_clipboard(sep=',')
... # Wrote the following to the system clipboard:
... # ,A,B,C
... # 0,1,2,3
... # 1,4,5,6
```

We can omit the the index by passing the keyword *index* and setting it to false.

```
>>> df.to_clipboard(sep=',', index=False)
... # Wrote the following to the system clipboard:
... # A,B,C
... # 1,2,3
... # 4,5,6
```

to\_csv (path=None, index=True, sep=', ', na\_rep=", float\_format=None, header=False, index\_label=None, mode='w', encoding=None, compression=None, date\_format=None, decimal='.')

Write Series to a comma-separated values (csv) file

### **Parameters**

**path** [string or file handle, default None] File path or object, if None is provided the result is returned as a string.

na\_rep [string, default ''] Missing data representation

float\_format [string, default None] Format string for floating point numbers

header [boolean, default False] Write out series name

index [boolean, default True] Write row names (index)

index\_label [string or sequence, default None] Column label for index column(s) if desired. If None is given, and *header* and *index* are True, then the index names are used. A sequence should be given if the DataFrame uses MultiIndex.

mode [Python write mode, default 'w']

sep [character, default ","] Field delimiter for the output file.

**encoding** [string, optional] a string representing the encoding to use if the contents are non-ascii, for python versions prior to 3

**compression** [string, optional] A string representing the compression to use in the output file. Allowed values are 'gzip', 'bz2', 'zip', 'xz'. This input is only used when the first argument is a filename.

date\_format: string, default None Format string for datetime objects.

**decimal: string, default '.'** Character recognized as decimal separator. E.g. use ',' for European data

#### to dense()

Return dense representation of NDFrame (as opposed to sparse)

```
to_dict (into=<type 'dict'>)
```

Convert Series to {label -> value} dict or dict-like object.

### **Parameters**

into [class, default dict] The collections. Mapping subclass to use as the return object. Can be the actual class or an empty instance of the mapping type you want. If you want a collections default dict, you must pass it initialized.

New in version 0.21.0.

### Returns

value\_dict [collections.Mapping]

# **Examples**

```
>>> s = pd.Series([1, 2, 3, 4])
>>> s.to_dict()
{0: 1, 1: 2, 2: 3, 3: 4}
>>> from collections import OrderedDict, defaultdict
>>> s.to_dict(OrderedDict)
OrderedDict([(0, 1), (1, 2), (2, 3), (3, 4)])
>>> dd = defaultdict(list)
>>> s.to_dict(dd)
defaultdict(<type 'list'>, {0: 1, 1: 2, 2: 3, 3: 4})
```

New in version 0.20.0.

### **Parameters**

**excel writer** [string or ExcelWriter object] File path or existing ExcelWriter

**sheet\_name** [string, default 'Sheet1'] Name of sheet which will contain DataFrame

na\_rep [string, default ''] Missing data representation

float\_format [string, default None] Format string for floating point numbers

columns [sequence, optional] Columns to write

**header** [boolean or list of string, default True] Write out the column names. If a list of strings is given it is assumed to be aliases for the column names

index [boolean, default True] Write row names (index)

index\_label [string or sequence, default None] Column label for index column(s) if desired. If None is given, and *header* and *index* are True, then the index names are used. A sequence should be given if the DataFrame uses MultiIndex.

**startrow:** upper left cell row to dump data frame

**startcol:** upper left cell column to dump data frame

engine [string, default None] write engine to use - you can also set this via the options io.excel.xlsx.writer, io.excel.xls.writer, and io.excel. xlsm.writer.

**merge\_cells** [boolean, default True] Write MultiIndex and Hierarchical Rows as merged cells.

**encoding: string, default None** encoding of the resulting excel file. Only necessary for xlwt, other writers support unicode natively.

inf\_rep [string, default 'inf'] Representation for infinity (there is no native representation
for infinity in Excel)

**freeze\_panes** [tuple of integer (length 2), default None] Specifies the one-based bottom-most row and rightmost column that is to be frozen

New in version 0.20.0.

### **Notes**

If passing an existing ExcelWriter object, then the sheet will be added to the existing workbook. This can be used to save different DataFrames to one workbook:

```
>>> writer = pd.ExcelWriter('output.xlsx')
>>> df1.to_excel(writer,'Sheet1')
>>> df2.to_excel(writer,'Sheet2')
>>> writer.save()
```

For compatibility with to\_csv, to\_excel serializes lists and dicts to strings before writing.

#### to frame (name=None)

Convert Series to DataFrame

#### **Parameters**

**name** [object, default None] The passed name should substitute for the series name (if it has one).

#### Returns

data frame [DataFrame]

## to\_hdf (path\_or\_buf, key, \*\*kwargs)

Write the contained data to an HDF5 file using HDFStore.

Hierarchical Data Format (HDF) is self-describing, allowing an application to interpret the structure and contents of a file with no outside information. One HDF file can hold a mix of related objects which can be accessed as a group or as individual objects.

In order to add another DataFrame or Series to an existing HDF file please use append mode and a different a key.

For more information see the user guide.

#### **Parameters**

path\_or\_buf [str or pandas.HDFStore] File path or HDFStore object.

**key** [str] Identifier for the group in the store.

**mode** [{'a', 'w', 'r+'}, default 'a'] Mode to open file:

- 'w': write, a new file is created (an existing file with the same name would be deleted).
- 'a': append, an existing file is opened for reading and writing, and if the file does not exist it is created.
- 'r+': similar to 'a', but the file must already exist.

**format** [{'fixed', 'table'}, default 'fixed'] Possible values:

- 'fixed': Fixed format. Fast writing/reading. Not-appendable, nor searchable.
- 'table': Table format. Write as a PyTables Table structure which may perform worse but allow more flexible operations like searching / selecting subsets of the data.

**append** [bool, default False] For Table formats, append the input data to the existing.

data\_columns [list of columns or True, optional] List of columns to create as indexed data columns for on-disk queries, or True to use all columns. By default only the axes of the object are indexed. See io.hdf5-query-data-columns. Applicable only to format='table'.

**complevel** [{0-9}, optional] Specifies a compression level for data. A value of 0 disables compression.

complib [{'zlib', 'lzo', 'bzip2', 'blosc'}, default 'zlib'] Specifies the compression library to be used. As of v0.20.2 these additional compressors for Blosc are supported (default if no compressor specified: 'blosc:blosclz'): {'blosc:blosclz', 'blosc:lz4', 'blosc:lz4hc', 'blosc:snappy', 'blosc:zlib', 'blosc:zstd'}. Specifying a compression library which is not available issues a ValueError.

**fletcher32** [bool, default False] If applying compression use the fletcher32 checksum.

dropna [bool, default False] If true, ALL nan rows will not be written to store.

**errors** [str, default 'strict'] Specifies how encoding and decoding errors are to be handled. See the errors argument for open () for a full list of options.

### See also:

DataFrame.read hdf Read from HDF file.

**DataFrame.to\_parquet** Write a DataFrame to the binary parquet format.

DataFrame.to\_sql Write to a sql table.

DataFrame.to\_feather Write out feather-format for DataFrames.

DataFrame.to\_csv Write out to a csv file.

## **Examples**

```
>>> df = pd.DataFrame({'A': [1, 2, 3], 'B': [4, 5, 6]},
... index=['a', 'b', 'c'])
>>> df.to_hdf('data.h5', key='df', mode='w')
```

We can add another object to the same file:

```
>>> s = pd.Series([1, 2, 3, 4])
>>> s.to_hdf('data.h5', key='s')
```

## Reading from HDF file:

```
>>> pd.read_hdf('data.h5', 'df')

A B
a 1 4
b 2 5
c 3 6
>>> pd.read_hdf('data.h5', 's')
0 1
1 2
2 3
3 4
dtype: int64
```

### Deleting file with data:

```
>>> import os
>>> os.remove('data.h5')
```

to\_json (path\_or\_buf=None, orient=None, date\_format=None, double\_precision=10, force\_ascii=True, date\_unit='ms', default\_handler=None, lines=False, compression=None, index=True)

Convert the object to a JSON string.

Note NaN's and None will be converted to null and datetime objects will be converted to UNIX timestamps.

# **Parameters**

**path\_or\_buf** [string or file handle, optional] File path or object. If not specified, the result is returned as a string.

**orient** [string] Indication of expected JSON string format.

- Series
  - default is 'index'
  - allowed values are: {'split','records','index'}
- DataFrame
  - default is 'columns'
  - allowed values are: {'split','records','index','columns','values'}
- The format of the JSON string
  - 'split': dict like { 'index' -> [index], 'columns' -> [columns], 'data' -> [values]}
  - 'records': list like [{column -> value}, ..., {column -> value}]
  - 'index' : dict like {index -> {column -> value}}
  - 'columns' : dict like {column -> {index -> value}}
  - 'values' : just the values array
  - 'table': dict like {'schema': {schema}, 'data': {data}} describing the data, and the data component is like orient='records'.

Changed in version 0.20.0.

- date\_format [{None, 'epoch', 'iso'}] Type of date conversion. 'epoch' = epoch milliseconds, 'iso' = ISO8601. The default depends on the orient. For orient='table',
  the default is 'iso'. For all other orients, the default is 'epoch'.
- **double\_precision** [int, default 10] The number of decimal places to use when encoding floating point values.
- force\_ascii [boolean, default True] Force encoded string to be ASCII.
- date\_unit [string, default 'ms' (milliseconds)] The time unit to encode to, governs timestamp and ISO8601 precision. One of 's', 'ms', 'us', 'ns' for second, millisecond, microsecond, and nanosecond respectively.
- **default\_handler** [callable, default None] Handler to call if object cannot otherwise be converted to a suitable format for JSON. Should receive a single argument which is the object to convert and return a serialisable object.
- **lines** [boolean, default False] If 'orient' is 'records' write out line delimited json format. Will throw ValueError if incorrect 'orient' since others are not list like.

New in version 0.19.0.

**compression** [{None, 'gzip', 'bz2', 'zip', 'xz'}] A string representing the compression to use in the output file, only used when the first argument is a filename.

New in version 0.21.0.

index [boolean, default True] Whether to include the index values in the JSON string. Not including the index (index=False) is only supported when orient is 'split' or 'table'.

New in version 0.23.0.

## See also:

pandas.read\_json

# **Examples**

Encoding/decoding a Dataframe using 'records' formatted JSON. Note that index labels are not preserved with this encoding.

```
>>> df.to_json(orient='records')
'[{"col 1":"a","col 2":"b"},{"col 1":"c","col 2":"d"}]'
```

Encoding/decoding a Dataframe using 'index' formatted JSON:

```
>>> df.to_json(orient='index')
'{"row 1":{"col 1":"a","col 2":"b"},"row 2":{"col 1":"c","col 2":"d"}}'
```

Encoding/decoding a Dataframe using 'columns' formatted JSON:

```
>>> df.to_json(orient='columns')
'{"col 1":{"row 1":"a","row 2":"c"},"col 2":{"row 1":"b","row 2":"d"}}'
```

Encoding/decoding a Dataframe using 'values' formatted JSON:

```
>>> df.to_json(orient='values')
'[["a","b"],["c","d"]]'
```

**Encoding with Table Schema** 

to\_latex (buf=None, columns=None, col\_space=None, header=True, index=True, na\_rep='NaN', formatters=None, float\_format=None, sparsify=None, index\_names=True, bold\_rows=False, column\_format=None, longtable=None, escape=None, encoding=None, decimal='.', multicolumn=None, multicolumn\_format=None, multirow=None)

Render an object to a tabular environment table. You can splice this into a LaTeX document. Requires \usepackage{booktabs}.

Changed in version 0.20.2: Added to Series

*to\_latex*-specific options:

**bold\_rows** [boolean, default False] Make the row labels bold in the output

column\_format [str, default None] The columns format as specified in LaTeX table format e.g 'rcl' for 3 columns

**longtable** [boolean, default will be read from the pandas config module] Default: False. Use a longtable environment instead of tabular. Requires adding a \usepackage{longtable} to your LaTeX preamble.

**escape** [boolean, default will be read from the pandas config module] Default: True. When set to False prevents from escaping latex special characters in column names.

**encoding** [str, default None] A string representing the encoding to use in the output file, defaults to 'ascii' on Python 2 and 'utf-8' on Python 3.

**decimal** [string, default '.'] Character recognized as decimal separator, e.g. ',' in Europe.

New in version 0.18.0.

**multicolumn** [boolean, default True] Use multicolumn to enhance MultiIndex columns. The default will be read from the config module.

New in version 0.20.0.

**multicolumn\_format** [str, default '1'] The alignment for multicolumns, similar to *column\_format* The default will be read from the config module.

New in version 0.20.0.

multirow [boolean, default False] Use multirow to enhance MultiIndex rows. Requires adding a \usepackage{multirow} to your LaTeX preamble. Will print centered labels (instead of top-aligned) across the contained rows, separating groups via clines. The default will be read from the pandas config module.

New in version 0.20.0.

to mol2 (filepath or buffer=None)

to\_msgpack (path\_or\_buf=None, encoding='utf-8', \*\*kwargs) msgpack (serialize) object to input file path

THIS IS AN EXPERIMENTAL LIBRARY and the storage format may not be stable until a future release.

### **Parameters**

path [string File path, buffer-like, or None] if None, return generated stringappend [boolean whether to append to an existing msgpack] (default is False)compress [type of compressor (zlib or blosc), default to None (no] compression)

to\_period (freq=None, copy=True)

Convert Series from DatetimeIndex to PeriodIndex with desired frequency (inferred from index if not passed)

### **Parameters**

**freq** [string, default]

### Returns

ts [Series with PeriodIndex]

to\_pickle (path, compression='infer', protocol=2)
Pickle (serialize) object to file.

## **Parameters**

path [str] File path where the pickled object will be stored.

**compression** [{'infer', 'gzip', 'bz2', 'zip', 'xz', None}, default 'infer'] A string representing the compression to use in the output file. By default, infers from the file extension in specified path.

New in version 0.20.0.

**protocol** [int] Int which indicates which protocol should be used by the pickler, default HIGHEST\_PROTOCOL (see [1] paragraph 12.1.2). The possible values for this parameter depend on the version of Python. For Python 2.x, possible values are 0, 1, 2. For Python>=3.0, 3 is a valid value. For Python>= 3.4, 4 is a valid value. A negative value for the protocol parameter is equivalent to setting its value to HIGH-EST PROTOCOL.

New in version 0.21.0.

### See also:

**read\_pickle** Load pickled pandas object (or any object) from file.

DataFrame.to\_hdf Write DataFrame to an HDF5 file.

DataFrame.to\_sql Write DataFrame to a SQL database.

**DataFrame.to\_parquet** Write a DataFrame to the binary parquet format.

# **Examples**

```
>>> original_df = pd.DataFrame({"foo": range(5), "bar": range(5, 10)})
>>> original_df
   foo bar
    0
          5
0
          6
     1
1
          7
2
     2
3
     3
          8
4
     4
>>> original_df.to_pickle("./dummy.pkl")
```

```
>>> import os
>>> os.remove("./dummy.pkl")
```

```
to_sdf (filepath_or_buffer=None)

to_smiles (filepath_or_buffer=None)

to_sparse (kind='block', fill_value=None)

Convert Series to SparseSeries

Parameters

kind [{'block', 'integer'}]

fill_value [float, defaults to NaN (missing)]

Returns
```

**sp** [SparseSeries]

Write records stored in a DataFrame to a SQL database.

Databases supported by SQLAlchemy [1] are supported. Tables can be newly created, appended to, or overwritten.

#### **Parameters**

name [string] Name of SQL table.

**con** [sqlalchemy.engine.Engine or sqlite3.Connection] Using SQLAlchemy makes it possible to use any DB supported by that library. Legacy support is provided for sqlite3.Connection objects.

**schema** [string, optional] Specify the schema (if database flavor supports this). If None, use default schema.

**if\_exists** [{'fail', 'replace', 'append'}, default 'fail'] How to behave if the table already exists.

- fail: Raise a ValueError.
- replace: Drop the table before inserting new values.
- append: Insert new values to the existing table.

**index** [boolean, default True] Write DataFrame index as a column. Uses *index\_label* as the column name in the table.

**index\_label** [string or sequence, default None] Column label for index column(s). If None is given (default) and *index* is True, then the index names are used. A sequence should be given if the DataFrame uses MultiIndex.

**chunksize** [int, optional] Rows will be written in batches of this size at a time. By default, all rows will be written at once.

**dtype** [dict, optional] Specifying the datatype for columns. The keys should be the column names and the values should be the SQLAlchemy types or strings for the sqlite3 legacy mode.

### Raises

**ValueError** When the table already exists and *if\_exists* is 'fail' (the default).

See also:

pandas.read\_sql read a DataFrame from a table

#### References

[1], [2]

## **Examples**

Create an in-memory SQLite database.

```
>>> from sqlalchemy import create_engine
>>> engine = create_engine('sqlite://', echo=False)
```

Create a table from scratch with 3 rows.

```
>>> df = pd.DataFrame({'name' : ['User 1', 'User 2', 'User 3']})
>>> df
    name
0    User 1
1    User 2
2    User 3
```

```
>>> df.to_sql('users', con=engine)
>>> engine.execute("SELECT * FROM users").fetchall()
[(0, 'User 1'), (1, 'User 2'), (2, 'User 3')]
```

Overwrite the table with just df1.

```
>>> df1.to_sql('users', con=engine, if_exists='replace',
... index_label='id')
>>> engine.execute("SELECT * FROM users").fetchall()
[(0, 'User 4'), (1, 'User 5')]
```

Specify the dtype (especially useful for integers with missing values). Notice that while pandas is forced to store the data as floating point, the database supports nullable integers. When fetching the data with Python, we get back integer scalars.

```
>>> from sqlalchemy.types import Integer
>>> df.to_sql('integers', con=engine, index=False,
... dtype={"A": Integer()})
```

```
>>> engine.execute("SELECT * FROM integers").fetchall()
[(1,), (None,), (2,)]
```

to\_string (buf=None, na\_rep='NaN', float\_format=None, header=True, index=True, length=False, dtype=False, name=False, max\_rows=None)

Render a string representation of the Series

# **Parameters**

**buf** [StringIO-like, optional] buffer to write to

na\_rep [string, optional] string representation of NAN to use, default 'NaN'

**float\_format** [one-parameter function, optional] formatter function to apply to columns' elements if they are floats default None

header: boolean, default True Add the Series header (index name)

```
index [bool, optional] Add index (row) labels, default True
```

length [boolean, default False] Add the Series length

dtype [boolean, default False] Add the Series dtype

name [boolean, default False] Add the Series name if not None

max\_rows [int, optional] Maximum number of rows to show before truncating. If None, show all.

### Returns

formatted [string (if not buffer passed)]

to\_timestamp (freq=None, how='start', copy=True)

Cast to datetimeindex of timestamps, at beginning of period

### **Parameters**

freq [string, default frequency of PeriodIndex] Desired frequency

**how** [{'s', 'e', 'start', 'end'}] Convention for converting period to timestamp; start of period vs. end

#### Returns

ts [Series with DatetimeIndex]

#### to xarray()

Return an xarray object from the pandas object.

#### Returns

- a DataArray for a Series
- a Dataset for a DataFrame
- a DataArray for higher dims

## **Notes**

See the xarray docs

## **Examples**

```
>>> df = pd.DataFrame({'A' : [1, 1, 2],
                       'B' : ['foo', 'bar', 'foo'],
                       'C' : np.arange(4.,7) })
>>> df
       В
            С
  A
     foo 4.0
  1
  1
     bar
          5.0
          6.0
2
  2
     foo
```

```
>>> df.to_xarray()
<xarray.Dataset>
Dimensions: (index: 3)
Coordinates:
  * index (index) int64 0 1 2
```

(continues on next page)

(continued from previous page)

```
Data variables:

A (index) int64 1 1 2

B (index) object 'foo' 'bar' 'foo'

C (index) float64 4.0 5.0 6.0
```

```
>>> p.to_xarray()
<xarray.DataArray (items: 4, major_axis: 3, minor_axis: 2)>
array([[[ 0, 1],
       [ 2, 3],
        [ 4,
             5]],
             7],
       [[6,
       [8, 9],
       [10, 11]],
       [[12, 13],
       [14, 15],
       [16, 17]],
       [[18, 19],
       [20, 21],
        [22, 23]])
Coordinates:
                (items) object 'A' 'B' 'C' 'D'
 * items
 * major_axis (major_axis) datetime64[ns] 2013-01-01 2013-01-02 2013-01-03,
→ # noqa
 * minor_axis (minor_axis) object 'first' 'second'
```

#### tolist()

Return a list of the values.

These are each a scalar type, which is a Python scalar (for str, int, float) or a pandas scalar (for Timestamp/Timedelta/Interval/Period)

### See also:

```
numpy.ndarray.tolist
```

## transform(func, \*args, \*\*kwargs)

Call function producing a like-indexed NDFrame and return a NDFrame with the transformed values

New in version 0.20.0.

#### **Parameters**

**func** [callable, string, dictionary, or list of string/callables] To apply to column

Accepted Combinations are:

- string function name
- function
- · list of functions
- dict of column names -> functions (or list of functions)

### Returns

## transformed [NDFrame]

#### See also:

```
pandas.NDFrame.aggregate, pandas.NDFrame.apply
```

### **Examples**

```
>>> df = pd.DataFrame(np.random.randn(10, 3), columns=['A', 'B', 'C'],
... index=pd.date_range('1/1/2000', periods=10))
df.iloc[3:7] = np.nan
```

```
>>> df.transform(lambda x: (x - x.mean()) / x.std())
                  A
                           В
2000-01-01 0.579457 1.236184 0.123424
2000-01-02 0.370357 -0.605875 -1.231325
2000-01-03 1.455756 -0.277446 0.288967
2000-01-04
              NaN
                        NaN
2000-01-05
              NaN
                        NaN
                                   NaN
2000-01-06 NaN NaN 2000-01-07 NaN NaN
                                  NaN
                                   NaN
2000-01-08 -0.498658 1.274522 1.642524
2000-01-09 -0.540524 -1.012676 -0.828968
2000-01-10 -1.366388 -0.614710 0.005378
```

## transpose (\*args, \*\*kwargs)

return the transpose, which is by definition self

```
truediv (other, level=None, fill_value=None, axis=0)
```

Floating division of series and other, element-wise (binary operator truediv).

Equivalent to series / other, but with support to substitute a fill\_value for missing data in one of the inputs.

#### **Parameters**

other [Series or scalar value]

fill\_value [None or float value, default None (NaN)] Fill existing missing (NaN) values, and any new element needed for successful Series alignment, with this value before computation. If data in both corresponding Series locations is missing the result will be missing

**level** [int or name] Broadcast across a level, matching Index values on the passed Multi-Index level

### **Returns**

result [Series]

### See also:

Series.rtruediv

# **Examples**

```
>>> a = pd.Series([1, 1, 1, np.nan], index=['a', 'b', 'c', 'd'])
>>> a
а
    1.0
b
    1.0
     1.0
С
    NaN
dtype: float64
>>> b = pd.Series([1, np.nan, 1, np.nan], index=['a', 'b', 'd', 'e'])
>>> b
     1.0
     NaN
h
d
     1.0
     NaN
е
dtype: float64
>>> a.add(b, fill_value=0)
    2.0
b
     1.0
     1.0
С
d
     1.0
     NaN
dtype: float64
```

truncate (before=None, after=None, axis=None, copy=True)

Truncate a Series or DataFrame before and after some index value.

This is a useful shorthand for boolean indexing based on index values above or below certain thresholds.

### **Parameters**

```
before [date, string, int] Truncate all rows before this index value.
```

after [date, string, int] Truncate all rows after this index value.

axis [{0 or 'index', 1 or 'columns'}, optional] Axis to truncate. Truncates the index (rows) by default.

**copy** [boolean, default is True,] Return a copy of the truncated section.

### Returns

type of caller The truncated Series or DataFrame.

See also:

**DataFrame.loc** Select a subset of a DataFrame by label.

**DataFrame.iloc** Select a subset of a DataFrame by position.

### **Notes**

If the index being truncated contains only datetime values, *before* and *after* may be specified as strings instead of Timestamps.

# **Examples**

```
>>> df = pd.DataFrame({'A': ['a', 'b', 'c', 'd', 'e'],
                       'B': ['f', 'g', 'h', 'i', 'j'],
. . .
                       'C': ['k', 'l', 'm', 'n', 'o']},
. . .
                       index=[1, 2, 3, 4, 5])
. . .
>>> df
  A B C
  a f
  b
     g
        1
3
  c h m
4
  d i
        n
5
  ејо
```

```
>>> df.truncate(before=2, after=4)

A B C

2 b g l

3 c h m

4 d i n
```

The columns of a DataFrame can be truncated.

```
>>> df.truncate(before="A", after="B", axis="columns")

A B

1 a f

2 b g

3 c h

4 d i

5 e j
```

For Series, only rows can be truncated.

The index values in truncate can be datetimes or string dates.

```
>>> dates = pd.date_range('2016-01-01', '2016-02-01', freq='s')
>>> df = pd.DataFrame(index=dates, data={'A': 1})
>>> df.tail()

A
2016-01-31 23:59:56 1
2016-01-31 23:59:57 1
2016-01-31 23:59:58 1
2016-01-31 23:59:59 1
2016-02-01 00:00:00 1
```

```
>>> df.truncate(before=pd.Timestamp('2016-01-05'),
... after=pd.Timestamp('2016-01-10')).tail()

A
2016-01-09 23:59:56 1
2016-01-09 23:59:57 1
2016-01-09 23:59:58 1
2016-01-09 23:59:59 1
2016-01-10 00:00:00 1
```

Because the index is a DatetimeIndex containing only dates, we can specify *before* and *after* as strings. They will be coerced to Timestamps before truncation.

```
>>> df.truncate('2016-01-05', '2016-01-10').tail()

A
2016-01-09 23:59:56 1
2016-01-09 23:59:57 1
2016-01-09 23:59:58 1
2016-01-09 23:59:59 1
2016-01-10 00:00:00 1
```

Note that truncate assumes a 0 value for any unspecified time component (midnight). This differs from partial string slicing, which returns any partially matching dates.

```
>>> df.loc['2016-01-05':'2016-01-10', :].tail()

A
2016-01-10 23:59:55 1
2016-01-10 23:59:56 1
2016-01-10 23:59:57 1
2016-01-10 23:59:58 1
2016-01-10 23:59:59 1
```

## tshift (periods=1, freq=None, axis=0)

Shift the time index, using the index's frequency if available.

### **Parameters**

periods [int] Number of periods to move, can be positive or negative

**freq** [DateOffset, timedelta, or time rule string, default None] Increment to use from the tseries module or time rule (e.g. 'EOM')

axis [int or basestring] Corresponds to the axis that contains the Index

## Returns

shifted [NDFrame]

### **Notes**

If freq is not specified then tries to use the freq or inferred\_freq attributes of the index. If neither of those attributes exist, a ValueError is thrown

tz\_convert (tz, axis=0, level=None, copy=True)

Convert tz-aware axis to target time zone.

### **Parameters**

tz [string or pytz.timezone object]

axis [the axis to convert]

**level** [int, str, default None] If axis ia a MultiIndex, convert a specific level. Otherwise must be None

**copy** [boolean, default True] Also make a copy of the underlying data

### **Raises**

**TypeError** If the axis is tz-naive.

tz\_localize(tz, axis=0, level=None, copy=True, ambiguous='raise')

Localize tz-naive TimeSeries to target time zone.

#### **Parameters**

tz [string or pytz.timezone object]

axis [the axis to localize]

**level** [int, str, default None] If axis ia a MultiIndex, localize a specific level. Otherwise must be None

copy [boolean, default True] Also make a copy of the underlying data

ambiguous ['infer', bool-ndarray, 'NaT', default 'raise']

- 'infer' will attempt to infer fall dst-transition hours based on order
- bool-ndarray where True signifies a DST time, False designates a non-DST time (note that this flag is only applicable for ambiguous times)
- 'NaT' will return NaT where there are ambiguous times
- 'raise' will raise an Ambiguous TimeError if there are ambiguous times

## Raises

**TypeError** If the TimeSeries is tz-aware and tz is not None.

## unique()

Return unique values of Series object.

Uniques are returned in order of appearance. Hash table-based unique, therefore does NOT sort.

### Returns

**ndarray or Categorical** The unique values returned as a NumPy array. In case of categorical data type, returned as a Categorical.

### See also:

pandas.unique top-level unique method for any 1-d array-like object.

**Index.unique** return Index with unique values from an Index object.

# **Examples**

```
>>> pd.Series([2, 1, 3, 3], name='A').unique()
array([2, 1, 3])
```

```
>>> pd.Series([pd.Timestamp('2016-01-01') for _ in range(3)]).unique() array(['2016-01-01T00:00:00.000000000'], dtype='datetime64[ns]')
```

An unordered Categorical will return categories in the order of appearance.

```
>>> pd.Series(pd.Categorical(list('baabc'))).unique()
[b, a, c]
Categories (3, object): [b, a, c]
```

An ordered Categorical preserves the category ordering.

### unstack (level=-1, fill\_value=None)

Unstack, a.k.a. pivot, Series with MultiIndex to produce DataFrame. The level involved will automatically get sorted.

### **Parameters**

**level** [int, string, or list of these, default last level] Level(s) to unstack, can pass level name

**fill\_value** [replace NaN with this value if the unstack produces] missing values New in version 0.18.0.

### **Returns**

unstacked [DataFrame]

# **Examples**

```
>>> s = pd.Series([1, 2, 3, 4],
... index=pd.MultiIndex.from_product([['one', 'two'], ['a', 'b']]))
>>> s
one a  1
        b   2
two a   3
        b   4
dtype: int64
```

```
>>> s.unstack(level=-1)
    a    b
    one    1    2
    two    3    4
```

```
>>> s.unstack(level=0)
one two
a 1 3
b 2 4
```

## update(other)

Modify Series in place using non-NA values from passed Series. Aligns on index

## **Parameters**

other [Series]

# **Examples**

```
>>> s = pd.Series(['a', 'b', 'c'])
>>> s.update(pd.Series(['d', 'e'], index=[0, 2]))
>>> s
0 d
1 b
2 e
dtype: object
```

If other contains NaNs the corresponding values are not updated in the original Series.

```
>>> s = pd.Series([1, 2, 3])
>>> s.update(pd.Series([4, np.nan, 6]))
>>> s
0     4
1     2
2     6
dtype: int64
```

## valid (inplace=False, \*\*kwargs)

Return Series without null values.

Deprecated since version 0.23.0: Use Series.dropna() instead.

**value\_counts** (normalize=False, sort=True, ascending=False, bins=None, dropna=True)
Returns object containing counts of unique values.

The resulting object will be in descending order so that the first element is the most frequently-occurring element. Excludes NA values by default.

#### **Parameters**

**normalize** [boolean, default False] If True then the object returned will contain the relative frequencies of the unique values.

sort [boolean, default True] Sort by values

ascending [boolean, default False] Sort in ascending order

**bins** [integer, optional] Rather than count values, group them into half-open bins, a convenience for pd.cut, only works with numeric data

dropna [boolean, default True] Don't include counts of NaN.

#### Returns

counts [Series]

#### values

Return Series as ndarray or ndarray-like depending on the dtype

#### Returns

arr [numpy.ndarray or ndarray-like]

### **Examples**

```
>>> pd.Series([1, 2, 3]).values
array([1, 2, 3])
```

```
>>> pd.Series(list('aabc')).values
array(['a', 'a', 'b', 'c'], dtype=object)
```

```
>>> pd.Series(list('aabc')).astype('category').values
[a, a, b, c]
Categories (3, object): [a, b, c]
```

Timezone aware datetime data is converted to UTC:

var (axis=None, skipna=None, level=None, ddof=1, numeric\_only=None, \*\*kwargs)
Return unbiased variance over requested axis.

Normalized by N-1 by default. This can be changed using the ddof argument

### **Parameters**

axis  $[\{index (0)\}]$ 

**skipna** [boolean, default True] Exclude NA/null values. If an entire row/column is NA, the result will be NA

**level** [int or level name, default None] If the axis is a MultiIndex (hierarchical), count along a particular level, collapsing into a scalar

**ddof** [int, default 1] Delta Degrees of Freedom. The divisor used in calculations is N - ddof, where N represents the number of elements.

numeric\_only [boolean, default None] Include only float, int, boolean columns. If None, will attempt to use everything, then use only numeric data. Not implemented for Series.

### **Returns**

var [scalar or Series (if level specified)]

## view(dtype=None)

Create a new view of the Series.

This function will return a new Series with a view of the same underlying values in memory, optionally reinterpreted with a new data type. The new data type must preserve the same size in bytes as to not cause index misalignment.

#### **Parameters**

**dtype** [data type] Data type object or one of their string representations.

#### Returns

**Series** A new Series object as a view of the same data in memory.

### See also:

numpy .ndarray .view Equivalent numpy function to create a new view of the same data in memory.

# **Notes**

Series are instantiated with dtype=float64 by default. While numpy.ndarray.view() will return a view with the same data type as the original array, Series.view() (without specified dtype) will try using float64 and may fail if the original data type size in bytes is not the same.

## **Examples**

```
>>> s = pd.Series([-2, -1, 0, 1, 2], dtype='int8')
>>> s
0    -2
1    -1
2    0
3    1
4    2
dtype: int8
```

The 8 bit signed integer representation of -1 is 0b111111111, but the same bytes represent 255 if read as an 8 bit unsigned integer:

The views share the same underlying values:

```
>>> us[0] = 128

>>> s

0 -128

1 -1

2 0

3 1

4 2

dtype: int8
```

**where** (cond, other=nan, inplace=False, axis=None, level=None, errors='raise', try\_cast=False, raise\_on\_error=None)

Return an object of same shape as self and whose corresponding entries are from self where *cond* is True and otherwise are from *other*.

#### **Parameters**

**cond** [boolean NDFrame, array-like, or callable] Where *cond* is True, keep the original value. Where False, replace with corresponding value from *other*. If *cond* is callable, it is computed on the NDFrame and should return boolean NDFrame or array. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as cond.

**other** [scalar, NDFrame, or callable] Entries where *cond* is False are replaced with corresponding value from *other*. If other is callable, it is computed on the NDFrame and should return scalar or NDFrame. The callable must not change input NDFrame (though pandas doesn't check it).

New in version 0.18.1: A callable can be used as other.

inplace [boolean, default False] Whether to perform the operation in place on the data

axis [alignment axis if needed, default None]

level [alignment level if needed, default None]

**errors** [str, {'raise', 'ignore'}, default 'raise']

- raise: allow exceptions to be raised
- ignore: suppress exceptions. On error return original object

Note that currently this parameter won't affect the results and will always coerce to a suitable dtype.

try\_cast [boolean, default False] try to cast the result back to the input type (if possible),

raise\_on\_error [boolean, default True] Whether to raise on invalid data types (e.g. trying to where on strings)

Deprecated since version 0.21.0.

### **Returns**

**wh** [same type as caller]

### See also:

```
DataFrame.mask()
```

## **Notes**

The where method is an application of the if-then idiom. For each element in the calling DataFrame, if cond is True the element is used; otherwise the corresponding element from the DataFrame other is used.

The signature for DataFrame.where() differs from numpy.where(). Roughly df1.where(m, df2) is equivalent to np.where(m, df1, df2).

For further details and examples see the where documentation in indexing.

# **Examples**

```
>>> s.where(s > 1, 10)
0     10.0
1     10.0
2     2.0
3     3.0
4     4.0
```

(continues on next page)

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```
True
        True
2
  True
        True
  True
3
        True
  True
        True
>>> df.where(m, -df) == df.mask(~m, -df)
  True
        True
  True
        True
  True True
3
  True True
  True True
```

**xs** (key, axis=0, level=None, drop\_level=True)

Returns a cross-section (row(s) or column(s)) from the Series/DataFrame. Defaults to cross-section on the rows (axis=0).

## **Parameters**

key [object] Some label contained in the index, or partially in a MultiIndex

axis [int, default 0] Axis to retrieve cross-section on

**level** [object, defaults to first n levels (n=1 or len(key))] In case of a key partially contained in a MultiIndex, indicate which levels are used. Levels can be referred by label or position.

**drop\_level** [boolean, default True] If False, returns object with same levels as self.

## Returns

xs [Series or DataFrame]

## **Notes**

xs is only for getting, not setting values.

MultiIndex Slicers is a generic way to get/set values on any level or levels. It is a superset of xs functionality, see MultiIndex Slicers

## **Examples**

```
>>> df
  A B
  4 5
         2
а
  4 0
  9 7
        3
>>> df.xs('a')
    4
В
     5
С
     2
Name: a
>>> df.xs('C', axis=1)
    2
а
     9
b
С
     3
Name: C
```

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```
>>> df
                  A B C D
first second third
                          9
                  4 1 8
    one 1
           1
                  7
                    5 5 0
     two
                  6 6 8 0
           1
baz
    one
     three 2
                  5 3 5
>>> df.xs(('baz', 'three'))
      A B C D
third
      5 3 5 3
2
>>> df.xs('one', level=1)
           A B C D
first third
           4
             1 8
bar
     1
     1
           6 6 8
                   0
baz
>>> df.xs(('baz', 2), level=[0, 'third'])
       A B C D
second
three
       5 3 5 3
```

oddt.pandas.read\_csv(\*args, \*\*kwargs)

**TODO: Support Chunks** 

```
oddt.pandas.read_mol2 (filepath_or_buffer=None, usecols=None, molecule_column='mol', molecule_name_column='mol_name', smiles_column=None, skip bad mols=False, chunksize=None, **kwargs)
```

Read Mol2 multi molecular file to ChemDataFrame. UCSF Dock 6 comments style is supported, i.e. #### var\_name: value before molecular block.

New in version 0.3.

## **Parameters**

filepath\_or\_buffer [string or None] File path

**usecols** [list or None, optional (default=None)] A list of columns to read from file. If None then all available fields are read.

**molecule\_column** [string or None, optional (default='mol')] Name of molecule column. If None the molecules will be skipped and the reading will be speed up significantly.

**molecule\_name\_column** [string or None, optional (default='mol\_name')] Column name which will contain molecules' title/name. Column is skipped when set to None.

**smiles\_column** [string or None, optional (default=None)] Column name containg molecules' SMILES, by default it is disabled.

**skip\_bad\_mols** [bool, optional (default=False)] Switch to skip empty (bad) molecules. Useful for RDKit, which Returns None if molecule can not sanitize.

**chunksize** [int or None, optional (default=None)] Size of chunk to return. If set to None whole set is returned.

## Returns

**result :** A *ChemDataFrame* containg all molecules if *chunksize* is None or genrerator of *ChemDataFrame* with *chunksize* molecules.

```
oddt.pandas.read_sdf (filepath_or_buffer=None, usecols=None, molecule_column='mol', molecule_name_column='mol_name', smiles_column=None, skip_bad_mols=False, chunksize=None, **kwargs)
```

Read SDF/MDL multi molecular file to ChemDataFrame

New in version 0.3.

## **Parameters**

- filepath\_or\_buffer [string or None] File path
- **usecols** [list or None, optional (default=None)] A list of columns to read from file. If None then all available fields are read.
- **molecule\_column** [string or None, optional (default='mol')] Name of molecule column. If None the molecules will be skipped and the reading will be speed up significantly.
- **molecule\_name\_column** [string or None, optional (default='mol\_name')] Column name which will contain molecules' title/name. Column is skipped when set to None.
- smiles\_column [string or None, optional (default=None)] Column name containg molecules' SMILES, by default it is disabled.
- **skip\_bad\_mols** [bool, optional (default=False)] Switch to skip empty (bad) molecules. Useful for RDKit, which Returns None if molecule can not sanitize.
- **chunksize** [int or None, optional (default=None)] Size of chunk to return. If set to None whole set is returned.

## **Returns**

**result :** A *ChemDataFrame* containg all molecules if *chunksize* is None or genrerator of *ChemDataFrame* with *chunksize* molecules.

## 5.1.8 oddt.shape module

oddt.shape.common\_usr(molecule, ctd=None, cst=None, fct=None, ftf=None, atoms\_type=None)
Function used in USR and USRCAT function

## **Parameters**

- molecule [oddt.toolkit.Molecule] Molecule to compute USR shape descriptor
- **ctd** [numpy array or None (default = None)] Coordinates of the molecular centroid If 'None', the point is calculated
- **cst** [numpy array or None (default = None)] Coordinates of the closest atom to the molecular centroid If 'None', the point is calculated
- **fct** [numpy array or None (default = None)] Coordinates of the farthest atom to the molecular centroid If 'None', the point is calculated
- **ftf** [numpy array or None (default = None)] Coordinates of the farthest atom to the molecular centroid If 'None', the point is calculated
- **atoms\_type** [str or None (default None)] Type of atoms to be selected from atom\_dict If 'None', all atoms are used to calculate shape descriptor

## Returns

**shape\_descriptor** [numpy array, shape = (12)] Array describing shape of molecule

oddt.shape.electroshape(mol)

Computes shape descriptor based on Armstrong, M. S. et al. ElectroShape: fast molecular similarity calculations incorporating shape, chirality and electrostatics. J Comput Aided Mol Des 24, 789-801 (2010). http://dx.doi.org/doi:10.1007/s10822-010-9374-0

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Aside from spatial coordinates, atoms' charges are also used as the fourth dimension to describe shape of the molecule.

#### **Parameters**

mol [oddt.toolkit.Molecule] Molecule to compute Electroshape descriptor

## Returns

**shape descriptor** [numpy array, shape = (15)] Array describing shape of molecule

oddt.shape.usr(molecule)

Computes USR shape descriptor based on Ballester PJ, Richards WG (2007). Ultrafast shape recognition to search compound databases for similar molecular shapes. Journal of computational chemistry, 28(10):1711-23. http://dx.doi.org/10.1002/jcc.20681

#### **Parameters**

molecule [oddt.toolkit.Molecule] Molecule to compute USR shape descriptor

## **Returns**

**shape\_descriptor** [numpy array, shape = (12)] Array describing shape of molecule

oddt.shape.usr\_cat(molecule)

Computes USRCAT shape descriptor based on Adrian M Schreyer, Tom Blundell (2012). USRCAT: real-time ultrafast shape recognition with pharmacophoric constraints. Journal of Cheminformatics, 2012 4:27. http://dx.doi.org/10.1186/1758-2946-4-27

#### **Parameters**

molecule [oddt.toolkit.Molecule] Molecule to compute USRCAT shape descriptor

#### Returns

**shape\_descriptor** [numpy array, shape = (60)] Array describing shape of molecule

oddt.shape.usr\_similarity ( $mol1\_shape$ ,  $mol2\_shape$ , ow=1.0, hw=1.0, rw=1.0, aw=1.0, dw=1.0) Computes similarity between molecules

## **Parameters**

mol1\_shape [numpy array] USR shape descriptor

mol2\_shape [numpy array] USR shape descriptor

- **ow** [float (default = 1.)] Scaling factor for all atoms Only used for USRCAT, ignored for other types
- **hw** [float (default = 1.)] Scaling factor for hydrophobic atoms Only used for USRCAT, ignored for other types
- **rw** [float (default = 1.)] Scaling factor for aromatic atoms Only used for USRCAT, ignored for other types
- aw [float (default = 1.)] Scaling factor for acceptors Only used for USRCAT, ignored for other types
- **dw** [float (default = 1.)] Scaling factor for donors Only used for USRCAT, ignored for other types

## Returns

**similarity** [float from 0 to 1] Similarity between shapes of molecules, 1 indicates identical molecules

## 5.1.9 oddt.spatial module

Spatial functions included in ODDT Mainly used by other modules, but can be accessed directly.

```
oddt.spatial.angle (p1, p2, p3)
```

Returns an angle from a series of 3 points (point #2 is centroid). Angle is returned in degrees.

#### **Parameters**

**p1,p2,p3** [numpy arrays, shape = [n\_points, n\_dimensions]] Triplets of points in n-dimensional space, aligned in rows.

#### Returns

**angles** [numpy array, shape =  $[n_points]$ ] Series of angles in degrees

```
oddt.spatial.angle_2\mathbf{v}(v1, v2)
```

Returns an angle between two vecors. Angle is returned in degrees.

## **Parameters**

**v1,v2** [numpy arrays, shape = [n\_vectors, n\_dimensions]] Pairs of vectors in n-dimensional space, aligned in rows.

#### Returns

**angles** [numpy array, shape = [n\_vectors]] Series of angles in degrees

```
oddt.spatial.dihedral (p1, p2, p3, p4)
```

Returns an dihedral angle from a series of 4 points. Dihedral is returned in degrees. Function distingishes clockwise and antyclockwise dihedrals.

#### **Parameters**

**p1, p2, p3, p4** [numpy arrays, shape = [n\_points, n\_dimensions]] Quadruplets of points in n-dimensional space, aligned in rows.

## Returns

**angles** [numpy array, shape = [n\_points]] Series of angles in degrees

```
oddt.spatial.distance (x, y)
```

Computes distance between each pair of points from x and y.

## **Parameters**

- $\mathbf{x}$  [numpy arrays, shape = [n\_x, 3]] Array of poinds in 3D
- y [numpy arrays, shape = [n\_y, 3]] Array of poinds in 3D

## Returns

```
dist_matrix [numpy arrays, shape = [n_x, n_y]] Distance matrix
```

oddt.spatial.rmsd(ref, mol, ignore\_h=True, method=None, normalize=False)

Computes root mean square deviation (RMSD) between two molecules (including or excluding Hydrogens). No symmetry checks are performed.

#### **Parameters**

ref [oddt.toolkit.Molecule object] Reference molecule for the RMSD calculation

mol [oddt.toolkit.Molecule object] Query molecule for RMSD calculation

**ignore\_h** [bool (default=False)] Flag indicating to ignore Hydrogen atoms while performing RMSD calculation. This toggle works only with 'hungarian' method and without sorting (method=None).

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**method** [str (default=None)] The method to be used for atom asignment between ref and mol. None means that direct matching is applied, which is the default behavior. Available methods:

• canonize - match heavy atoms using canonical ordering (it forces

ignoring H's) - hungarian - minimize RMSD using Hungarian algorithm - min\_symmetry - makes multiple molecule-molecule matches and finds minimal RMSD (the slowest). Hydrogens are ignored.

normalize [bool (default=False)] Normalize RMSD by square root of rot. bonds

## Returns

rmsd [float] RMSD between two molecules

oddt.spatial.rotate(coords, alpha, beta, gamma)

Rotate coords by cerain angle in X, Y, Z. Angles are specified in radians.

## **Parameters**

**coords** [numpy arrays, shape = [n\_points, 3]] Coordinates in 3-dimensional space.

**alpha, beta, gamma: float** Angles to rotate the coordinates along X, Y and Z axis. Angles are specified in radians.

## Returns

**new\_coords** [numpy arrays, shape = [n\_points, 3]] Rorated coordinates in 3-dimensional space.

## 5.1.10 oddt.surface module

This module generates and does computation with molecular surfaces.

oddt.surface.find\_surface\_residues(molecule, max\_dist=None, scaling=1.0)

Finds residues close to the molecular surface using generate\_surface\_marching\_cubes. Ignores hydrogens and waters present in the molecule.

#### **Parameters**

molecule [oddt.toolkit.Molecule] Molecule to find surface residues in.

max\_dist [array\_like, numeric or None (default = None)] Maximum distance from the surface where residues would still be considered close. If None, compares distances to radii of respective atoms.

**scaling** [float (default = 1.0)] Expands the grid in which computation is done by generate\_surface\_marching\_cubes by a factor of scaling. Results in a more accurate representation of the surface, and therefore more accurate computation of distances but increases computation time.

## Returns

**atom\_dict** [numpy array] An atom\_dict containing only the surface residues from the original molecule.

oddt.surface.generate\_surface\_marching\_cubes (molecule, remove\_hoh=False, scaling=1.0, probe radius=1.4)

Generates a molecular surface mesh using the marching\_cubes method from scikit-image. Ignores hydrogens present in the molecule.

## **Parameters**

molecule [oddt.toolkit.Molecule object] Molecule for which the surface will be generated

**remove\_hoh** [bool (default = False)] If True, remove waters from the molecule before generating the surface. Requires molecule.protein to be set to True.

**scaling** [float (default = 1.0)] Expands the grid in which computation is done by a factor of scaling. Results in a more accurate representation of the surface, but increases computation time.

**probe\_radius** [float (default = 1.4)] Radius of a ball used to patch up holes inside the molecule resulting from some molecular distances being larger (usually in protein). Basically reduces the surface to one accesible by other molecules of radius smaller than probe\_radius.

#### Returns

verts [numpy array] Spatial coordinates for mesh vertices.

faces [numpy array] Faces are defined by referencing vertices from verts.

## 5.1.11 oddt.utils module

Common utilities for ODDT

```
oddt.utils.check_molecule(mol, force\_protein=False, force\_coords=False, non\_zero\_atoms=False)
```

Universal validator of molecule objects. Usage of positional arguments is allowed only for molecule object, otherwise it is prohibitted (i.e. the order of arguments **will** change). Desired properties of molecule are validated based on specified arguments. By default only the object type is checked. In case of discrepancy to the specification a *ValueError* is raised with appropriate message.

New in version 0.6.

## **Parameters**

mol: oddt.toolkit.Molecule object Object to verify

**force\_protein: bool** (**default=False**) Force the molecule to be marked as protein (mol.protein).

force coords: bool (default=False) Force the molecule to have non-zero coordinates.

non\_zero\_atoms: bool (default=False) Check if molecule has at least one atom.

```
oddt.utils.chunker(iterable, chunksize=100)
```

Generate chunks from a generator object. If iterable is passed which is not a generator it will be converted to one with *iter()*.

New in version 0.6.

```
oddt.utils.compose iter(iterable, funcs)
```

Chain functions and apply them to iterable, by exhausting the iterable. Functions are executed in the order from funcs.

New in version 0.6.

```
oddt.utils.is_molecule(obj)
```

Check whether an object is an *oddt.toolkits.{rdk,ob}.Molecule* instance.

New in version 0.6.

```
oddt.utils.is_openbabel_molecule(obj)
```

Check whether an object is an *oddt.toolkits.ob.Molecule* instance.

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New in version 0.6.

## oddt.utils.is\_rdkit\_molecule(obj)

Check whether an object is an *oddt.toolkits.rdk.Molecule* instance.

New in version 0.6.

oddt.utils.method\_caller(obj, methodname, \*args, \*\*kwargs)

Helper function to workaround Python 2 pickle limitations to parallelize methods and generator objects

## 5.1.12 oddt.virtualscreening module

ODDT pipeline framework for virtual screening

 $\textbf{class} \ \, \textbf{oddt.virtualscreening.virtualscreening} \, (\textit{n\_cpu=-1, verbose=False, chunksize=100}) \\ \text{Virtual Screening pipeline stack}$ 

## **Parameters**

 $n\_cpu: int (default=-1)$  The number of parallel procesors to use

verbose: bool (default=False) Verbosity flag for some methods

## Methods

<pre>apply_filter(expression[, soft_fail])</pre>	Filtering method, can use raw expressions
	(strings to be evaled in if statement, can use
	oddt.toolkit.Molecule methods, eg.
dock(engine, protein, *args, **kwargs)	Docking procedure.
fetch()	A method to exhaust the pipeline.
load_ligands(fmt, ligands_file, **kwargs)	Loads file with ligands.
score(function[, protein])	Scoring procedure compatible with any scoring func-
	tion implemented in ODDT and other pickled SFs
	which are subclasses of oddt.scoring.scorer.
<pre>similarity(method, query[, cutoff, protein])</pre>	Similarity filter.
write(fmt, filename[, csv_filename])	Outputs molecules to a file
write_csv(csv_filename[, fields, keep_pipe])	Outputs molecules to a csv file

## apply\_filter(expression, soft\_fail=0)

Filtering method, can use raw expressions (strings to be evaled in if statement, can use oddt.toolkit.Molecule methods, eg. mol.molwt < 500) Currently supported presets:

- Lipinski Rule of 5 ('ro5' or '15')
- Fragment Rule of 3 ('ro3')
- PAINS filter ('pains')

## **Parameters**

**expression: string or list of strings** Expresion(s) to be used while filtering.

soft\_fail: int (default=0) The number of faulures molecule can have to pass filter, aka. soft-fails.

**dock** (*engine*, *protein*, \**args*, \*\**kwargs*)

Docking procedure.

## **Parameters**

engine: string Which docking engine to use.

## **Notes**

Additional parameters are passed directly to the engine. Following docking engines are supported:

1. Audodock Vina (`engine="autodock\_vina"`), see oddt.docking.autodock\_vina.

#### fetch()

A method to exhaust the pipeline. Itself it is lazy (a generator)

load\_ligands (fmt, ligands\_file, \*\*kwargs)

Loads file with ligands.

## **Parameters**

file\_type: string Type of molecular file

**ligands file: string** Path to a file, which is loaded to pipeline

score (function, protein=None, \*args, \*\*kwargs)

Scoring procedure compatible with any scoring function implemented in ODDT and other pickled SFs which are subclasses of *oddt.scoring.scorer*.

## **Parameters**

**function: string** Which scoring function to use.

protein: oddt.toolkit.Molecule Default protein to use as reference

## **Notes**

Additional parameters are passed directly to the scoring function.

similarity (method, query, cutoff=0.9, protein=None)

## Similarity filter. Supported structural methods:

- ift: interaction fingerprints
- sift: simple interaction fingerprints
- usr: Ultrafast Shape recognition
- usr cat: Ultrafast Shape recognition, Credo Atom Types
- electroshape: Electroshape, an USR method including partial charges

## **Parameters**

**method: string** Similarity method used to compare molecules. Avaiale methods: \* *ifp* - interaction fingerprint (requires a receptor) \* *sifp* - simple interaction fingerprint (requires a receptor) \* *usr* - Ultrafast Shape Reckognition \* *usr\_cat* - USR, with CREDO atom types \* *electroshape* - Electroshape, USR with moments representing partial charge

**query: oddt.toolkit.Molecule or list of oddt.toolkit.Molecule** Query molecules to compare the pipeline to.

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cutoff: float Similarity cutoff for filtering molecules. Any similarity lower than it will be filtered out.

**protein: oddt.toolkit.Molecule (default = None)** Protein for underling method. By default it's empty, but sturctural fingerprints need one.

```
write (fmt, filename, csv_filename=None, **kwargs)
Outputs molecules to a file
```

#### **Parameters**

file\_type: string Type of molecular file
ligands\_file: string Path to a output file
csv\_filename: string Optional path to a CSV file

write\_csv (csv\_filename, fields=None, keep\_pipe=False, \*\*kwargs)
Outputs molecules to a csv file

## **Parameters**

csv\_filename: string Optional path to a CSV file

fields: list (default None) List of fields to save in CSV file

**keep\_pipe: bool (default=False)** If set to True, the ligand pipe is sustained.

## 5.1.13 Module contents

## **Open Drug Discovery Toolkit**

Universal and easy to use resource for various drug discovery tasks, ie docking, virutal screening, rescoring.

## **Attributes**

toolkit [module,] Toolkits backend module, currenlty OpenBabel [ob] and RDKit [rdk]. This setting is toolkit-wide, and sets given toolkit as default

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To be announced.

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# Documentation Indices and tables

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