

Twister: Construction and structural relaxation of commensurate moiré superlattices ^{☆,☆☆}

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ABSTRACT

Introduction of a twist between layers of two-dimensional materials which leads to the formation of a moiré pattern is an emerging pathway to tune the electronic, vibrational and optical properties. The fascinating properties of these systems are often linked to large-scale structural reconstruction of the moiré pattern. Hence, an essential first step in the theoretical study of these systems is the construction and structural relaxation of the atoms in the moiré superlattice. We present the Twister package, a collection of tools that constructs commensurate superlattices for any combination of 2D materials and also helps perform structural relaxations of the moiré superlattice. Twister constructs commensurate moiré superlattices using the coincidence lattice method and provides an interface to perform structural relaxations using classical forcefields.

Program summary

Program Title: Twister

CPC Library link to program files: <https://doi.org/10.17632/frfrj8wpbk.1>

Developer's repository link: <https://github.com/qtm-iisc/Twister>

Licensing provisions: BSD 3-clause

Programming language: Python

External routines/libraries: numpy, scipy, mpi4py, matplotlib

Nature of problem: Moiré patterns can be constructed from 2D materials by twisting the layers with respect to each other. Simulation of these patterns using forcefields or density functional theory requires the construction of periodically repeating moiré superlattices.

Solution method: Moiré superlattices are constructed using the coincidence site lattice theory while also allowing the user to accommodate strains in the layer. Twister also helps perform structural relaxation of the moiré superlattice using forcefields by interfacing to the LAMMPS package.

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1. Introduction

Two-dimensional (2D) materials, with their extensive spectrum of physical properties and long reaching applicability in technology, have triggered a wave of innovation in the realm of material science. Twisting the layers of 2D materials to form moiré patterns is an emerging direction of research [1–10] which was triggered by the recent discovery of unconventional supercon-

ductivity and Mott-insulating phases in twisted bilayer graphene [11,12]. An important feature of the moiré patterns is that they undergo structural reconstructions as predicted by theoretical simulations [13–25] and corroborated by experiments [26–32]. The atomic reconstructions can strongly modify the electronic structure [28,16,33,30,29,34,35,17,36–39] and vibrational modes [40–42] of the moiré superlattice.

Most theoretical methods, particularly those based on classical forcefields or first-principles density functional theory [43] (DFT), require the creation of a periodic simulation cell. Application of an arbitrary twist between the layers of a 2D material leads to an incommensurate moiré which is intractable for these methods. It is thus essential to find special twist angles [44,45] which lead to moiré superlattices that are commensurate with periodic boundary conditions. A common approach to find these twist angles is

[☆] The review of this paper was arranged by Prof. Blum Volker.

^{☆☆} This paper and its associated computer program are available via the Computer Physics Communications homepage on ScienceDirect (<http://www.sciencedirect.com/science/journal/00104655>).

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the coincidence lattice method [44,46]. In this method, the commensurate twist angle is determined as the one at which perfect coincidence occurs for a set of lattice points in the top and bottom layers. This method has been used to derive an analytical expression for the commensurate twist-angles in systems which have a hexagonal lattice and exactly same lattice constant in the top and bottom layer. To go beyond this restricted class to 2D materials, a general implementation of this method is necessary. Furthermore, it is important to note that perfect coincidence of lattice sites is often not possible or only possible at very large length scales. In this case a small strain may be applied to each layer to obtain a commensurate superlattice of reasonable size [46].

At large moiré length scales, all moiré superlattices undergo structural reconstructions arising from in-plane and out-of-plane displacement of atoms from the rigidly-twisted structure. In-plane displacements lead to a change in the local stackings giving rise to strain-soliton networks [17,16,13,14,26,30] and out-of-plane displacements lead to a varying interlayer spacing and/or buckling of the layers [29,47–50] in the moiré. Hence, once the commensurate moiré superlattice is constructed, performing structural relaxation is the essential next step before exploring its electronic or optical properties. Despite evidence of structural reconstructions, many theoretical studies of moiré superlattices do not take this into account [10,51–55]. This is because lattice relaxations using van der Waals corrected DFT is a major computational bottleneck due to the large number of atoms that constitute the moiré superlattice. An alternative computationally cheaper approach to perform the structural relaxation of the moiré superlattices is based on classical forcefields. This approach eliminates the electronic degrees of freedom and employs a theoretical description involving only nuclei. The interactions between nuclei are represented with simple inter-atomic potentials, whose parameters are obtained by fitting to accurate DFT calculations. For instance, the structural relaxations of the moiré superlattices of twisted bilayer of transition metal dichalcogenides (TMDs) are performed with a combination of Stillinger-Weber [56] (SW) and registry-dependent Kolmogorov-Crespi [57] (KC) potentials. The SW and KC potential describes intralayer and interlayer interaction in these systems. The SW potential parameters were obtained by fitting to DFT-derived phonon dispersion of single layer TMD [58], whereas the KC potential parameters were obtained by fitting interlayer binding energy landscape to van der Waals corrected DFT results [59–63]. The fit forcefields can be used to obtain moiré reconstructions in good agreement with DFT relaxations [59] at a fraction of the computational cost. However, the technical details of classical force-field based simulation are non-trivial. An all-inclusive package that constructs and automates the performance of lattice relaxations of the moiré superlattice can bridge this gap for non-experts.

We present the Twister package to construct commensurate moiré superlattices for any 2D homobilayer or heterobilayer material. Twister helps find minimum-area commensurate superlattices using the coincidence lattice method for a given range of twist angles and user-defined strain tolerance. The superlattice vectors and atoms within the superlattice are written to a file for further study of its structural and electronic properties. The construction of commensurate superlattice has been tested for various homobilayer and heterobilayer cases: graphene/hBN, MoS₂/MoSe₂, hBN/MoSe₂, twisted bilayer phosphorene and MoS₂. We use the package to find a range of commensurate twist angles for the case of MoS₂/MoSe₂ and twisted bilayer phosphorene. Twister further also streamlines the process of studying moiré reconstruction using classical force-field calculations as implemented in the LAMMPS [64] package. The streamlined process is demonstrated for twisted bilayer MoS₂, twisted bilayer graphene and twisted MoS₂/MoSe₂ moiré patterns.

2. Construction of commensurate superlattices

The method used to construct the commensurate superlattices depends on the type of lattice vectors in the bottom and top layer forming the moiré. If the two lattices are hexagonal and with identical lattice constants, we use an analytical expression [45] to obtain the commensurate twist-angles (θ_i).

$$\cos(\theta_i) = \frac{3i^2 + 3i + 0.5}{3i^2 + 3i + 1}, \quad (1)$$

where i is a whole number. The corresponding superlattice vectors are given by: $\mathbf{A}_1 = i\mathbf{a}_1 + (i+1)\mathbf{a}_2$, $\mathbf{A}_2 = -(i+1)\mathbf{a}_1 + (2i+1)\mathbf{a}_2$ where \mathbf{a}_1 and \mathbf{a}_2 are the unit-cell lattice vectors. In the superlattices generated using this equation, the periodicity of the stackings matches the moiré unit-cell periodicity. I.e. each type of stacking (say AA) is present once per moiré unit-cell. For commensurate twist-angles in addition to θ_i , a general form for commensurate twist-angles is given by [44]:

$$\cos(\theta_{m,n}) = \frac{n^2 + 4nm + m^2}{2(n^2 + mn + m^2)}, \quad (2)$$

where m, n are whole numbers. The superlattice vectors are: $\mathbf{A}_1 = m\mathbf{a}_1 + n\mathbf{a}_2$, $\mathbf{A}_2 = -n\mathbf{a}_1 + (m+n)\mathbf{a}_2$. This general equation yields all possible commensurate superlattices and not just ones with minimum area. The superlattice can have multiple AA stackings per moiré unit-cell.

For a combination of layers with dissimilar lattices we use the coincidence site lattice theory to generate the superlattice. Consider the unit-cell lattice vectors of bottom layer to be $\mathbf{a}_1, \mathbf{a}_2, \mathbf{a}_3$ and of top layer to be $\mathbf{b}_1, \mathbf{b}_2, \mathbf{b}_3$. Let \mathbf{R}_θ be the rotation matrix corresponding to a rotation angle θ in the xy plane. Top layer is rotated about the origin keeping the bottom layer fixed. The rotated lattice vectors, \mathbf{b}_i^r are given by:

$$\mathbf{b}_i^r = \mathbf{b}_i \mathbf{R}_\theta^T \quad (3)$$

Coincidence of a subset of lattice sites occurs when a lattice vector in bottom layer, $\mathbf{v}_1 = n_1\mathbf{a}_1 + n_2\mathbf{a}_2$, superposes with a lattice vector in the rotated top layer, $\mathbf{v}_2^r = m_1\mathbf{b}_1^r + m_2\mathbf{b}_2^r$, i.e.

$$n_1\mathbf{a}_1 + n_2\mathbf{a}_2 = m_1\mathbf{b}_1^r + m_2\mathbf{b}_2^r \quad (4)$$

Fig. 1 graphically represents the above steps for 27.8° twisted bilayer MoS₂. The shortest coinciding vectors have been plotted and any two non-parallel vectors can be equivalently used as superlattice vectors for the given twist angle.

For heterobilayer materials, it is often difficult to obtain exact coincidence since lattice parameters of the constituent layers are different. Hence, we introduce a user-defined upper limit of mismatch, Δ , between the vectors in Å units,

$$|\mathbf{v}_1 - \mathbf{v}_2^r| \leq \Delta \quad (5)$$

It is important to find superlattices with as small a mismatch as possible to ensure that the superlattice is commensurate with the twist-angle. A large value of Δ can introduce spurious strains in the superlattice. For an arbitrary combination of layers, the value of mismatch or the size of superlattice is often too large despite sweeping a large range of twist-angles. In these cases, more effort is needed to find perfectly commensurate superlattices. In the following subsections, we describe two methods to improve coincidence.

After the computation of all vectors co-incident within the mismatch, pairs of solutions ($\{n_1\mathbf{a}_1 + n_2\mathbf{a}_2, n'_1\mathbf{a}_1 + n'_2\mathbf{a}_2\}$) are checked to be non-parallel by ensuring their vector cross products is non-zero. The solutions can also be further restricted to have a specific

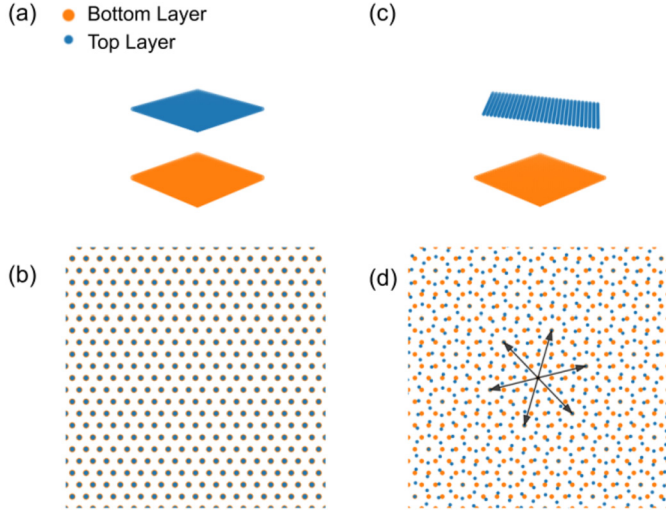


Fig. 1. a) and c) 3-D representations of untwisted and twisted MoS₂ homobilayer respectively. Twist angle is 27.8° in c). b) and d) Top view of a) and c), respectively. Six solution vectors of Eqn. (4) (with minimum superlattice area) are shown in d).

angle between the vectors. In the final step, we choose superlattices with minimum area with a variation of 1% allowed around the minima. The final output is a list of appropriate superlattice vectors and the corresponding strained lattice parameters.

2.1. Straining lattice parameters

The lattice parameters of each layer can be uniformly strained as demonstrated for the case of hBN/MoSe₂ in Fig. 2. A biaxial strain (β) is applied to each layer such that the following equation is satisfied [46]:

$$(1 + \beta)|\mathbf{v}_1| \approx (1 - \beta)|\mathbf{v}_2^r| \quad (6)$$

Lattice parameters of one layer are strained by β and those of the lower layer by $-\beta$ to improve coincidence. In certain situations, it is desirable to allow for only one layer to be strained. If only lattice parameters of the top layer are to be strained, β is applied to the top layer such that:

$$|\mathbf{v}_1| \approx (1 + \beta)|\mathbf{v}_2^r| \quad (7)$$

If only lattice parameters of the bottom layer are to be strained, β is applied to the bottom layer such that:

$$(1 + \beta)|\mathbf{v}_1| \approx |\mathbf{v}_2^r| \quad (8)$$

Two values of β are generated for each pair of solutions ($\{n_1\mathbf{a}_1 + n_2\mathbf{a}_2, n'_1\mathbf{a}_1 + n'_2\mathbf{a}_2\}$) and the averaged value is used to form the superlattice vectors. The mismatch after straining the superlattice vectors is computed as demonstrated in Eqn. (5).

2.2. Applying a strain tensor

When straining the lattice parameters alone does not yield small mismatch, the unit-vectors of the top layer can be deformed to ensure perfect coincidence. However, this may change the angle between the lattice vectors of the strained layer leading to breaking of symmetries within that layer. The deformation tensor (\mathbf{D}) is computed such that:

$$\begin{bmatrix} n_1 & n_2 \\ n'_1 & n'_2 \end{bmatrix} \begin{bmatrix} \mathbf{a}_1 \\ \mathbf{a}_2 \end{bmatrix} = \begin{bmatrix} m_1 & m_2 \\ m'_1 & m'_2 \end{bmatrix} \begin{bmatrix} \mathbf{b}_1 \\ \mathbf{b}_2 \end{bmatrix} \mathbf{D} \mathbf{R}_\theta^T \quad (9)$$

Table 1
Lattice parameters of each material.

Materials	Lattice parameter (Å)
MoS ₂	3.164
MoSe ₂	3.301
Graphene	2.467
hBN	2.512
Phosphorene	3.321, 4.62

Table 2

Results obtained from Twister for the test systems to compare with the commensurate twist-angles reported in Ref. [46].

Materials	Twist angle	Superlattice vectors (n_1, n_2, n'_1, n'_2)	Atoms	Biaxial strain
MoS ₂ /MoS ₂	27.8°	(-4, 3), (-3, -1)	78	±0.0
MoS ₂ /MoSe ₂	16.1°	(-1, -3), (3, -4)	75	±0.119
Graphene/hBN	0.0°	(-37, 0), (0, 37)	5330	±0.466
hBN/MoSe ₂	19.1°	(1, -3), (3, -2)	26	±0.333

Any deformation of a lattice can be expressed in terms of a pure rotation of the lattice and a symmetric strain tensor. Through a polar decomposition [65], we have $\mathbf{D} = \epsilon \mathbf{R}_\phi$:

$$\begin{bmatrix} n_1 & n_2 \\ n'_1 & n'_2 \end{bmatrix} \begin{bmatrix} \mathbf{a}_1 \\ \mathbf{a}_2 \end{bmatrix} = \begin{bmatrix} m_1 & m_2 \\ m'_1 & m'_2 \end{bmatrix} \begin{bmatrix} \mathbf{b}_1 \\ \mathbf{b}_2 \end{bmatrix} \epsilon \mathbf{R}_\phi \mathbf{R}_\theta^T \quad (10)$$

The pure rotation leads to a further change in the twist-angle by ϕ . The change in the twist-angle along with a strain in the unit-vectors (by ϵ) leads to a perfectly commensurate superlattice.

3. Test systems

The construction of commensurate superlattices using our implementation (Eqn (4)) is verified for five test systems – twisted bilayer MoS₂, MoS₂/MoSe₂ heterobilayer, Graphene/hBN heterobilayer and hBN/MoSe₂ heterobilayer to compare with Ref. [46]. To find the commensurate superlattices (using the procedure described in section 2) we use the following constraints. For MoS₂/MoSe₂, MoS₂ / MoS₂ and hBN/MoSe₂, the angle between superlattice vectors is fixed at 60°. For Graphene/hBN, the angle between superlattice vectors is fixed to be 120°. Maximum strain, β , allowed in the lattice parameters is 1%. Mismatch (Å) threshold, Δ , is set to be less than 0.1 Å. The commensurate superlattices, number of atoms and strains are in agreement with the reported results [46,33]. The lattice parameters used in our simulation and the results are shown in Table 1 and Table 2, respectively. The input files along with instructions on how to use the code to generate the commensurate superlattices for these test systems is included in the Twister package.

4. Examples

4.1. Commensurate twist-angles

Apart from the above test examples, we use Twister to compute commensurate twist-angles for the heterobilayer systems of MoS₂/MoSe₂ and bilayer phosphorene. For systems of MoS₂/MoSe₂, the angle between superlattice vectors is fixed at 60°. Strain for lattice parameters is kept below 1% and distributed over both layers. Mismatch threshold is 10^{-5} Å. The size of the moiré superlattice is expected to increase as a function of reducing twist-angle. We accordingly increase the search range of n_1, n_2, m_1, m_2 (Eqn. (4)) as a function of decreasing twist-angle. The commensurate twist-angles for MoS₂/MoSe₂ can be found in Table 3. The commensurate twist-angles for twisted bilayer phosphorene are

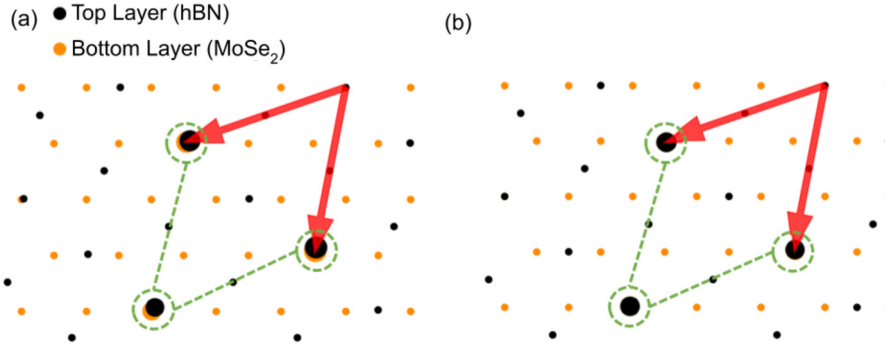


Fig. 2. Red arrows are the superlattice vectors of hBN/ MoSe₂ bilayer twisted at 19.1°. The corner lattice points are encircled. a) Supercell with unstrained lattice parameters. b) Supercell with strained parameters with improved overlap. (For interpretation of the colors in the figure(s), the reader is referred to the web version of this article.)

Table 3

Commensurate twist-angles in the range 0-30° generated for MoS₂/MoSe₂ using Twister. Lattice parameters were strained to improve mismatch.

Twist angle	Atoms	Superlattice vectors (n_1, n_2), (n'_1, n'_2)	Strain (%)	Mismatch (Å)
0.0°	3315	(24, 0), (0, 24)	0.00856	10 ⁻⁹
1.56654°	2283	(11, -23), (-12, -11)	0.0501	5 × 10 ⁻⁶
2.84558°	1362	(4, -17), (17, -13)	0.0847	2 × 10 ⁻⁶
3.10565°	1248	(-16, 13), (-13, -3)	0.0454	2 × 10 ⁻⁶
4.715°	696	(0, -11), (11, -11)	0.0371	2 × 10 ⁻⁶
7.3111°	327	(-8, 1), (-7, 8)	0.176	4 × 10 ⁻⁶
11.17856°	456	(3, -10), (10, -7)	0.145	1 × 10 ⁻⁶
14.11315°	384	(7, -9), (-2, -7)	0.226	2 × 10 ⁻⁶
16.1021°	75	(-4, 1), (-3, 4)	0.118	3 × 10 ⁻⁶
21.45415°	384	(7, -9), (-2, -7)	0.226	10 ⁻⁶

generated using the deformation tensor method described above. One of the layers is allowed to deform to form a perfectly commensurate superlattice. A fine tuning of the twist-angles is not necessary for this method. For example, starting with $\theta = 6.9^\circ$ leads to the following strain and rotation matrices (Eqn. (10)):

$$\epsilon = \begin{bmatrix} 0.992893 & -0.001897 \\ -0.001897 & 0.993364 \end{bmatrix};$$

$$\mathbf{R}_\phi = \begin{bmatrix} 0.999995 & -0.003102 \\ 0.003102 & 0.999995 \end{bmatrix} \quad (11)$$

The ϵ tensor is applied to the lattice vectors and the \mathbf{R}_ϕ leads to a change in the twist-angle between the layers by $\phi = -0.17773^\circ$. The commensurate twist-angle is thus, $\theta + \phi = 6.72227^\circ$. To test the implementation, we also start with $\theta = 6.5^\circ$ and find a solution:

$$\epsilon = \begin{bmatrix} 0.992919 & -0.001900 \\ -0.001900 & 0.993338 \end{bmatrix};$$

$$\mathbf{R}_\phi = \begin{bmatrix} 0.999992 & 0.003879 \\ -0.003879 & 0.999992 \end{bmatrix} \quad (12)$$

The strain tensor is nearly the same, while the change in twist-angle, $\phi = 0.2222699^\circ$. As expected, the final commensurate twist-angle ($\theta + \phi$) remains identical for different starting values of θ . The commensurate twist-angles for twisted bilayer phosphorene generated using this method are provided in Table 4. The input files, along with instructions on how to run the code, are provided with the Twister package. More details on using the code and the workflow is described in the next section.

4.2. Structural relaxations

We demonstrate the automation of construction and relaxation of moiré superlattices (using the LAMMPS package) for the case

of twisted bilayer MoS₂. Using a single script, *homo.sh*, a range of commensurate superlattices can be generated and relaxed. Using this script we perform the following:

1. A file, *hex.table*, which contains a series of twist angles and superlattice vectors is created.
2. The moiré superlattice and atom positions are generated for each twist-angle in this file.
3. To perform the relaxation using LAMMPS, the superlattice vectors must be reoriented so that one of the superlattice vectors is along the (100) direction. Consider three superlattice vectors of the moiré superlattice generated using Eqn. (1): $\mathbf{v}_1, \mathbf{v}_2$ and \mathbf{v}_3 . $\hat{\mathbf{v}}_3$ is along the out-of-plane, z direction. The transformation is thus computed in the following manner,

$$\bar{\mathbf{A}}_1 = \|\bar{\mathbf{v}}_1\| \hat{\mathbf{x}}$$

$$\bar{\mathbf{A}}_2 = (\bar{\mathbf{v}}_2 \cdot \hat{\mathbf{v}}_1) \hat{\mathbf{x}} + (|\hat{\mathbf{v}}_1 \times \bar{\mathbf{v}}_2|) \hat{\mathbf{y}}$$

$$\bar{\mathbf{A}}_3 = \bar{\mathbf{v}}_3$$

4. With the user-provided force-fields, the atomic relaxations are performed using the LAMMPS package for all the twist-angles in *hex.table*.

The range of twist-angles in *hex.table* can be generated using both Eqn. (1) or (2) according to the user's preference. The relaxed structures can be visualized with the OVITO [66] software directly. The order-parameter and interlayer spacing distribution for relaxed 1.61° and 58.39° twisted bilayer MoS₂ is shown in Fig. 3. The order-parameter is defined as the shortest displacement of the top layer that transforms a given local stacking to the highest energy stacking in the moiré – AA for 1.61° twist and A'B for 58.39° twist. The relaxation leads to increase in the area of low-energy stackings (AB, BA, AA', AB') and formation of shear-strain solitons at the boundary between adjacent low-energy stackings [41,17,36,13]. The strain distribution in the bottom layer of the relaxed moiré pattern is shown in Fig. 4. The strains are localized along the solitonic networks.

5. Workflow

In this section the sequential steps for operating the code are described. The code can be divided into three sections.

5.1. Finding commensurate angle

This part of the package is used to generate the commensurate twist-angles and the corresponding lattice vectors using the method described in section 1. The code, *get_ang.py*, is used for this purpose and uses an input file: *get_ang.inp*.

Table 4
Commensurate twist-angles of bilayer black phosphorous generated using Twister. A strain tensor is applied to one of the layers to improve mismatch.

Twist angle	Atoms	Superlattice vectors (n_1, n_2), (n'_1, n'_2)	Strain tensor	Mismatch (Å)
1.595323°	10444	(45, 0), (0, 29)	$\begin{bmatrix} 1.00047 & 0.00306 \\ 0.00306 & 1.00030 \end{bmatrix}$	0.0
2.793987°	3364	(0, -15), (28, 0)	$\begin{bmatrix} 0.999 & 0.000877 \\ 0.000877 & 0.999 \end{bmatrix}$	0.0
3.444685°	2212	(-12, 0), (0, -23)	$\begin{bmatrix} 0.998 & 0.00289 \\ 0.00289 & 0.998 \end{bmatrix}$	0.0
4.62238°	1228	(0, -9), (-17, 0)	$\begin{bmatrix} 0.997 & 0.000972 \\ 0.000972 & 0.997 \end{bmatrix}$	0.0
5.768955°	788	(0, -7), (14, 0)	$\begin{bmatrix} 0.996 & -0.00163 \\ -0.00163 & 0.995 \end{bmatrix}$	0.0
6.722269°	580	(0, -6), (12, 0)	$\begin{bmatrix} 0.993 & -0.00189 \\ -0.00189 & 0.993 \end{bmatrix}$	0.0

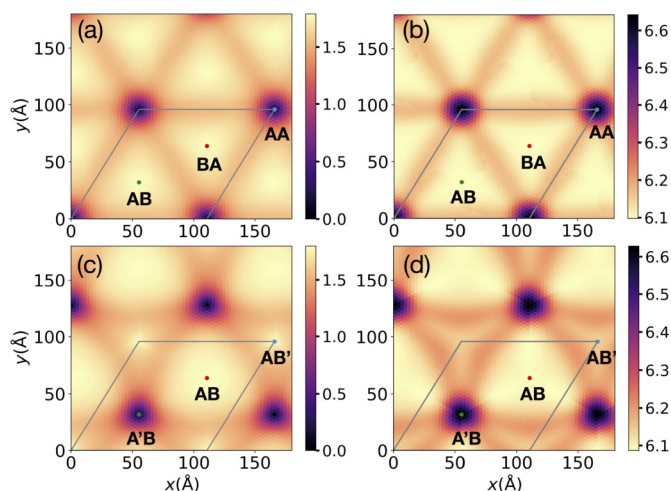


Fig. 3. a and b (c and d) Order-parameter distribution and interlayer spacing distribution in 1.61° (58.39°) twisted bilayer MoS₂, respectively.

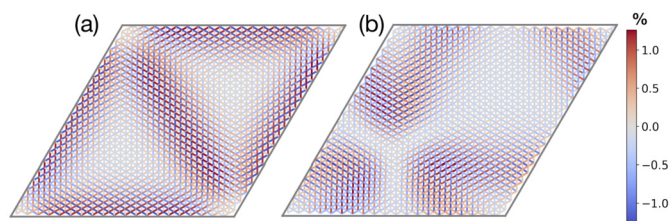


Fig. 4. a and b Distribution of strain in the bottom layer of 1.61° and 58.39° twisted bilayer MoS₂. Lines are drawn between every Mo atom and its 6 nearest neighbors. The color of the line represents the strain in that direction.

1. Prepare input file, *get_ang.inp*. The lattice vectors for the top and bottom layer are provided by the user in this file. The basis atom positions are not necessary at this stage. The input file is also designed to contain an exhaustive range of constraints for the superlattice to be found. The user can fix the angle between the superlattice vectors, the type and amount of strain applied on the layer parameters and the final and initial mismatch thresholds, Δ . The user also provides a range of twist-angles to be searched through this input file. The initial mismatch threshold is used to specify the maximum mismatch that will be tolerated for the twist angles provided in the input file. The final mismatch threshold is employed only when *DeepSearch* function is used, as the twist-angle from

the input range with the minimum mismatch is automatically refined further by searching in a narrower angle range around the chosen twist-angle until the mismatch is smaller than the specified final mismatch threshold. If *DeepSearch* is 'False', only the angles provided in the range in the input file are searched for solutions. The output file is generated for the twist-angle having the minimum mismatch with the solution found by solving Eqn. (5). This solution may not have mismatch less than the final mismatch threshold provided. Output files for other angles in the input range are stored in a folder named 'solutions'. If *DeepSearch* is 'True', the output file for the refined twist-angle is generated. As mentioned above, the refined twist-angle will have mismatch lesser than final mismatch threshold. Output files for rest of the twist-angles searched are stored in a folder named 'solutions'. On running the code, a plot is generated with the lattice points and the superlattice vectors. Vectors of the same color show super-cells of the same area. An output file is generated for every twist-angle in the range provided by the user and contains all the solutions found by solving Eqn. (5).

2. If too many solutions are found, the parameters can be changed to reduce the number of vectors and plot them for confirmation. Usually, it involves reducing ranges of n_1, n_2, m_1, m_2 to look for smaller superlattices or adjusting the strain percent and mismatch allowed. If no superlattice vectors are found, changes to the input are suggested.
3. Pick the appropriate super-cell vectors from the output and prepare the second input file, *twist.inp*, to generate the atom positions.

5.2. Atoms in superlattice

This part of the package, *twister.py*, generates the atom coordinates in the moiré superlattice. The input files needed are *twist.inp* and basis atom positions of the unit-cell for top and bottom layer. The code also verifies that the atoms lying on the boundary of the supercell are not missed or counted twice. The code verifies that the number of atoms computed matches the expected value from area ratios. The expected number of atoms in top layer is given by: $n_{top} = n_b * Ar_{sl} / Ar_{uc}$, where n_b is the number of basis atoms in top layer, Ar_{sl} is the area of the superlattice and Ar_{uc} is the area of the unit-cell. The output file, *superlattice.dat*, contains the superlattice vectors and the atom coordinates.

A flowchart summary of all the above steps can be found in Fig. 5.

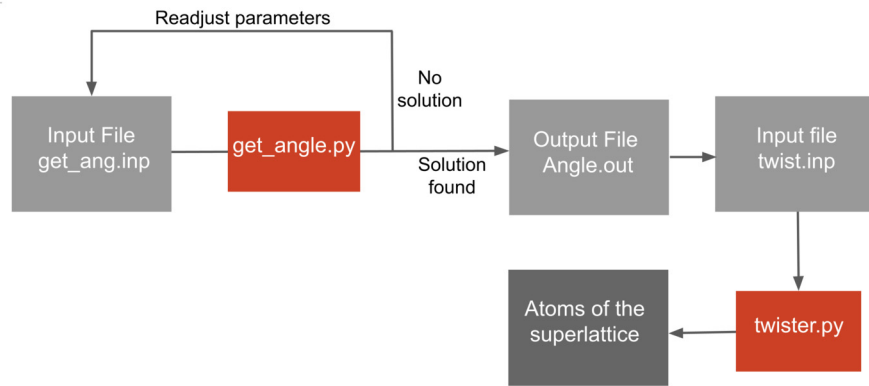


Fig. 5. The program flow for generating the supercell lattice and number of atoms for given material and twist angle.

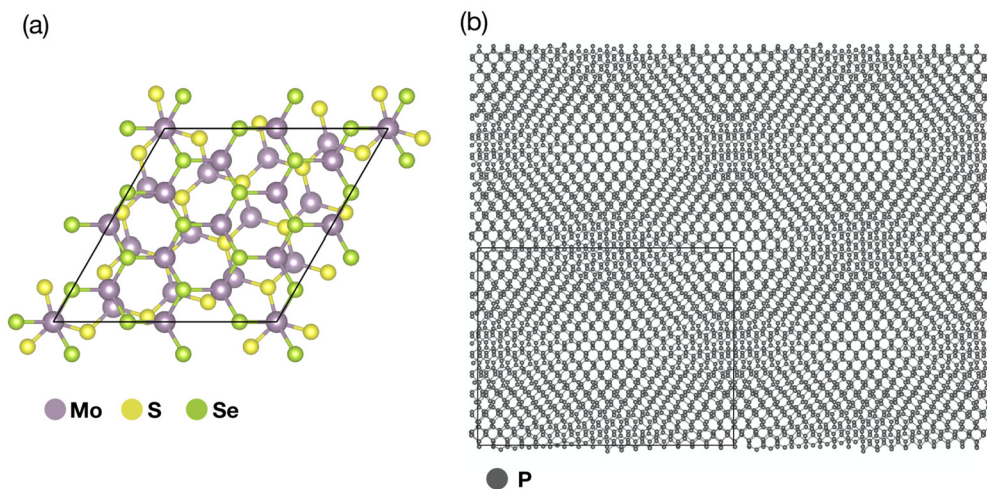


Fig. 6. Superlattices of a) 16.1° twisted MoS₂/MoSe₂ and b) 4.265° twisted bilayer black phosphorous.

6. Twister code framework

6.1. Input file

We provide an example input file, *get_ang.inp* for the code *get_ang.py* described in the previous section to showcase the various options available to the user (Fig. 6).

```

range_nm:
-20 21
celldm_a:
3.164 3.164 25.0
a1:
0.5 0.8660254 0.0
a2:
-0.5 0.8660254 0.0
a3:
0.0 0.0 1.0
Number_basis_atoms_a:
3
celldm_b:
3.301 3.301 25.0
b1:
0.5 0.8660254 0.0
b2:
-0.5 0.8660254 0.0
b3:
0.0 0.0 1.0
Number_basis_atoms_b:

```

```

3
theta_range:
16.05 16.15 0.01
initial_mismatch_threshold (Angstrom):
0.0271
final_mismatch_threshold (Angstrom):
0.00001
strain_per:
1.0
strain_tensor_vector:
'False'
strain_layer:
'Both'
DeepSearch:
'True'
fix_ang:
'True'
f_ang:
60
plot:
'True'

```

The unit of length is Angstrom (Å). Lattice parameters, unit-cell vectors and basis atoms for the top layer are provided in the line following the key-word: *celldm_a*, (*a1,a2,a3*) and *Number_basis_atoms_a* respectively and similarly for the bottom layer. Lower bound, upper bound and step size for the twist-angle range is given by the string below *theta_range* in the same order. Threshold for mismatch between lattice vectors of top and bottom layer

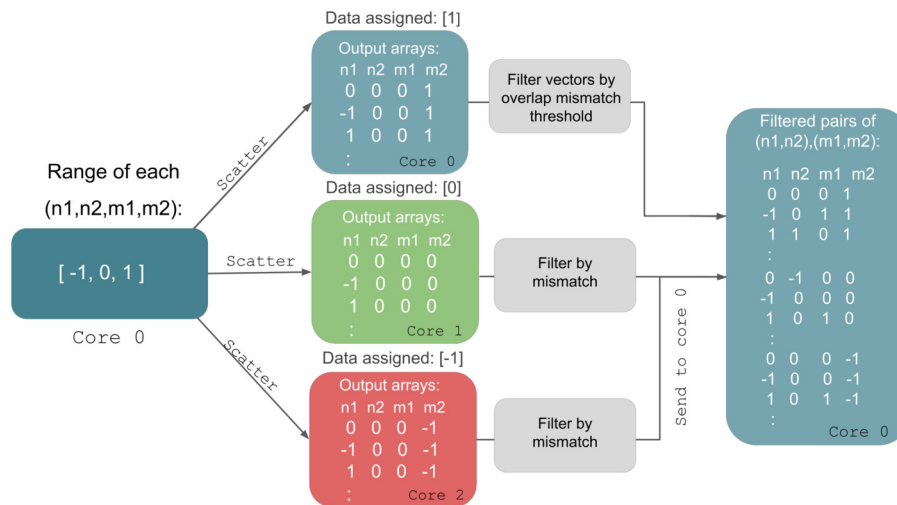


Fig. 7. Schematic of the parallelization used in Twister to find commensurate twist angles.

for the twist angles provided in the input file can be specified in the string below *initial_mismatch_threshold* (Angstroms). If the *DeepSearch* function is being used, threshold mismatch for the refined solutions can be specified in the string below *final_mismatch_threshold* (Angstroms). The user can apply a strain tensor by specifying 'True' below *strain_tensor_vector* otherwise, if specified 'False', only the lattice parameters will be strained. Threshold for percentage strain allowed to improve mismatch is given below *strain_per*. When entry under *strain_tensor_vector* is 'False', uniaxial strain in the top or bottom layer, or biaxial strain in both layer can be specified below *strain_layer*. When entry under *strain_tensor_vector* is 'True', only uniaxial strain in the top or bottom layer can be applied. When *DeepSearch* is 'True', the twist-angle is refined and the solution is printed if one with mismatch less than the final threshold is found. If *DeepSearch* is 'False', solution for the twist-angle from the range given in the input file with the minimum mismatch will be printed. Fixing angle between superlattice vectors and specifying the angle between superlattice vectors are allowed by *fix_ang* and *f_ang* respectively. The user can plot the resultant superlattice vectors for each angle by specifying 'True' below *plot*. Exact format of input can be found in the README files provided with each example included with the package.

7. Parallelization and optimization

The primary function for finding superlattice vectors in *get_ang.py* is the function *clt()* and scales as $O(N^4)$. We therefore optimize and parallelize this part of the code. Consider the unit-cell lattice vectors of bottom layer to be $\mathbf{a}_1, \mathbf{a}_2, \mathbf{a}_3$ and of top layer to be $\mathbf{b}_1, \mathbf{b}_2, \mathbf{b}_3$. Let superlattice in bottom layer be given as $\mathbf{v}_1 = n_1\mathbf{a}_1 + n_2\mathbf{a}_2$, and top layer as $\mathbf{v}_2 = m_1\mathbf{b}_1 + m_2\mathbf{b}_2$. The function *clt()* utilizes user defined unit-cell lattice vectors and range of n_1, n_2, m_1, m_2 to checks all possible combinations of superlattice vectors for solutions. To parallelize computing the permutations, the range is scattered across the cores via MPI, using package *mpi4py*. Each core gets N_{nm}/N_c numbers from the range, where N_{nm} is the length of range of n_1, n_2, m_1, m_2 and N_c is the user-defined number of cores. For each element in its assigned data, the core generates a four-columned *NumPy* array, each column corresponding to n_1, n_2, m_1, m_2 . The column for m_2 is uni-valued and only contains corresponding element in the data assigned to the core. The rows for n_1, n_2, m_1 contain all possible ordered triplets from the range of n_1, n_2, m_1, m_2 (with repeats allowed). These triplets were made using the *product()* iterator from the *Itertools* module

of Python. A schematic for the parallelized permutation calculations can be found in Fig. 7.

Using *NumPy* operations, the columns for n_1, n_2 are vector multiplied with $\mathbf{a}_1, \mathbf{a}_2$ and columns for m_1, m_2 are vector multiplied with rotated top layer vectors, $\mathbf{b}'_1, \mathbf{b}'_2$. The resulting *NumPy* arrays, \mathbf{v}_1 and \mathbf{v}'_2 , are used to calculate mismatch between them and the arrays are filtered according to the user-defined threshold (Eqn. (5)). *NumPy* arrays for $\mathbf{v}_1, \mathbf{v}'_2, (n_1, n_2), (m_1, m_2)$ and mismatch between the pairs, are sent by different cores and received by the core with *rank* = 0. All possible pairs of superlattice vectors are then made using *combinations()* iterator from the *Itertools* module is used to make all possible pairs of superlattice vectors. They are filtered for non-parallel pairs. No parallelization is applied in the remaining steps. *NumPy* operations and arrays are used for further filtering depending on angle between superlattice vectors, strain in lattice parameters, and area of supercell.

8. Conclusion

We present the Twister package to construct commensurate moiré superlattice for any combination of 2D material. The package also streamlines atomic relaxation of the moiré superlattices by interfacing with classical forcefield calculations implemented in the LAMMPS package. We demonstrate the construction of the superlattices for five test systems, twisted bilayer MoS_2 , $\text{MoS}_2/\text{MoSe}_2$, graphene/hBN, hBN/ MoSe_2 and twisted bilayer phosphorene. We also use the package to find commensurate twist-angles for the phosphorene bilayer and $\text{MoS}_2/\text{MoSe}_2$ heterobilayer. The package is written in Python and uses MPI parallelization for the computationally demanding steps.

Declaration of competing interest

The authors declare that they have no known competing financial interests or personal relationships that could have appeared to influence the work reported in this paper.

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