Magnetic Response of Nematic Superconductors: Skyrmion Stripes and Their Signatures in Muon Spin Relaxation Experiments

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We investigate the magnetic response of nematic superconductors, presenting a new approach to find vortex and skyrmion structures beyond symmetry-constraining Ansätze. Using this approach we show that nematic superconductors form distinctive skyrmion stripes. Our approach lends itself to accurate determination of the field distribution for muon spin rotation probes. We use this to show that the skyrmion structure manifests as a double peak in the field distribution, markedly different from the signal of standard vortex lattices.

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Understanding nematicity in superconducting states is currently one of the central questions of the field [\[1](#page-6-0)–[13](#page-6-1)]. Various indications of nematicity have been observed in a broad range of materials including cuprates [[14](#page-6-2),[15](#page-6-3)], heavy fermion materials [[16](#page-6-4)[,17\]](#page-6-5), and iron based superconductors [\[18](#page-6-6)–[22\]](#page-6-7). One particular example of a material that has been the target of intense experimental study is the doped topological insulator $Bi₂Se₃$ [\[23](#page-6-8)–[25\]](#page-7-0). Superconducting candidate materials of interest include $Cu_xBi₂Se₃$, $Sr_xBi_2Se_3$ [[26](#page-7-1)–[31](#page-7-2)], and $Nb_xBi_2Se_3$ [[32](#page-7-3)]. One of the interesting properties of nematic superconductors is that the flux carrying objects could be skyrmions [\[5](#page-6-9),[33](#page-7-4)[,34\]](#page-7-5), which are fundamentally distinct from the vortices that normally appear in superconductivity.

The appearance of skyrmions raises two questions: how do the skyrmions affect the magnetic response of the material, and what unique experimental signatures does this produce? The first question is technically nontrivial, since the interskyrmion interactions are complicated and anisotropic, driven by mode mixing in general multicomponent anisotropic models [[33](#page-7-4)[,35](#page-7-6)–[38](#page-7-7)], which dictates the appearance of multiple scales in the interskyrmion forces. This anisotropic behavior cannot be captured by a simple Ansatz, thus it is challenging to approximate the resulting lattice with any certainty. Using brute-force numerics, on the other hand, would require simulating complicated Ginzburg-Landau (GL) models at large system sizes, with boundaries that may affect structure formation.

In this Letter and in the companion extended paper [\[33\]](#page-7-4) we study this magnetic response by developing and applying a new method that can determine vortex lattice structure for any model. We find that nematic superconductors exhibit an unusual magnetic response in the form of skyrmion stripes.

Since surface probes visualize vortex structures at limited scales and structure formation may be affected by pinning, strain, inhomogeneities, and surface effects, the most reliable method to detect these structures is muon spin rotation (μSR) experiments. These fire muons into the sample, detecting the orientation of positrons emitted when the muons decay. These data can then be used to calculate the statistical distribution of the magnetic field in the bulk of a superconductor [[39](#page-7-8)].

We will demonstrate that for nematic superconductors, this magnetic field distribution exhibits a double peak structure. We posit that the number of peaks in the distribution correlates with the number of length scales in the lattice. For a nematic superconductor this is two, the separation between chains and the interskyrmion separation along the chain, which we will call the chain link length.

The model we consider in this Letter is derived from the proposal that certain topological insulators, such as electron-doped nematic superconductors (e.g., $Cu_xBi₂Se₃$) are described by a two-component order parameter [[1](#page-6-0),[23](#page-6-8)–[25](#page-7-0)]. This order parameter breaks rotational symmetry in the basal plane to a mixed symmetry, combining the phase difference of the order parameter and spatial rotations [[33](#page-7-4)]. This matches the observed twofold rotational symmetry of specific heat and upper critical field [\[40\]](#page-7-9). The associated Ginzburg-Landau Gibbs free energy for such a nematic

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superconductor was microscopically derived in [\[5](#page-6-9)]. In dimensionless units, it reads

$$
G = \int_{\Omega} \left\{ \frac{1}{2} Q_{ij}^{\alpha \beta} \overline{D_i \psi_{\alpha}} D_j \psi_{\beta} + \frac{1}{2} |B - H|^2 + F_p(|\psi_{\alpha}|, \theta_{\alpha \beta}) \right\},\tag{1}
$$

where $\psi_{\alpha} = \rho_{\alpha} e^{i\varphi_{\alpha}}$ are two complex superconducting order parameters $\alpha \in \{1, 2\}$. Note that Greek indices will always be used for superconducting components, Latin indices will be used for spatial directions and repeated indices imply summation. F_p denotes the collected potential terms and varies between models, but we will use the standard form,

$$
F_p = \frac{\eta^2}{2} \left(-|\psi_1|^2 - |\psi_2|^2 + \frac{1}{2} (|\psi_1|^4 + |\psi_2|^4) + \gamma |\psi_1|^2 |\psi_2|^2 \right),
$$
\n(2)

where γ is a microscopically derived parameter [\[5](#page-6-9)] related to the nematic nature of the system. It directly determines the ground state value for the condensates and is set to $\gamma = 1/3$ according to [[5\]](#page-6-9). The parameter η scales the potential and hence the ratio of critical fields H_{c_2}/H_{c_1} . Note that for $\eta \ll 1$ the system starts to act like a type-I superconductor and skyrmions do not form.

The constant matrices $Q^{\alpha\beta}$ define the anisotropy in the system and must obey the symmetry $Q_{ij}^{\alpha\beta} = \overline{Q}_{ji}^{\beta\alpha}$, ensuring that the energy is real. $\overline{Q}_{ij} = \overline{d}_{ji}^{\beta\alpha}$, the energy that the energy is real. $D_i = \partial_i - iA_i$ is the covariant derivative, where A is the vector potential corresponding to a $U(1)$ gauge field, with corresponding magnetic field $B = \nabla \times A$. For a nematic superconductor the anisotropy matrices take the form

$$
Q^{11} = Q^{22} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & \beta_z \end{pmatrix}, \quad Q^{12} = \beta_{\perp} \begin{pmatrix} 1 & i & 0 \\ i & -1 & 0 \\ 0 & 0 & 0 \end{pmatrix}, \quad (3)
$$

where β_z and β_\perp are positive, we will follow [\[5\]](#page-6-9) with the choice $\beta_z = 4/3$ and $\beta_{\perp} = 1/3$. The β_{\perp} term drives skyrmion formation or fractional vortex splitting. This can be seen by considering co-centered fractional vortices where the term integrates to zero. Hence, by splitting the fractional vortices this term can become negative, thus if β_{\perp} is sufficiently large it is energetically favorable to have fractional vortex splitting. The field equations in the bulk of the superconductor can then be found by varying the Gibbs free energy with respect to the fields ψ_a and A_i . Note that the resulting equations are independent of β_z which will have no effect on our results.

Solutions for isolated flux-carrying objects show that in nematic superconductors these are skyrmions, that have the form of spatially separated half-quantum vortices [[5](#page-6-9)[,33,](#page-7-4)[34](#page-7-5)]. These solutions were found by simulating the field

FIG. 1. Diagram of the geometry of a general unit cell for a vortex lattice, defined by the two vectors V_1 and V_2 with angle α .

equations on a plane ($\Omega = \mathbb{R}^2$), however, to find a lattice solution, we must solve them over the unit cell of a lattice with periodic boundary conditions (up to gauge). While it is often assumed that the unit cell has hexagonal symmetry, there is no reason to expect this for skyrmions in an anisotropic model. Hence, we take a new approach, considering a general unit cell, which takes the form of a parallelogram that is periodic (up to gauge), as shown in Fig. [1.](#page-1-0) A lattice solution is then given by the unit cell and field configuration that minimizes G/A , where A is the area of the unit cell. Hence, numerically, we seek minimizers of G/A with respect to the fields and with respect to the geometry of the unit cell itself, represented as the two vectors V_1 V_1 and V_2 in Fig. 1 (note that this will also control the area A). The full technical detail of the procedure we propose is given in a companion paper [[33](#page-7-4)] (along with a generalization that is not of use in this Letter). A brief account is also presented in Appendix A.

By tessellating the solution on the periodic unit cell, we find a solution that covers the plane with a vortex lattice. Hence, in applying the above method we find the field configuration that minimizes G/A in the bulk of a superconductor with no boundary effects. In addition, as we know the shape of the unit cell, we know the symmetry of the lattice.

Applying the above approach for the basal plane of a nematic system gives the results in Fig. [2,](#page-2-0) where we see that for low and medium strength external field H , the unit cell is rectangular ($\alpha = \pi/2$), containing four half-quantum vortices (totalling winding two) in each cell, forming a skyrmion chain. The lattice vectors lie parallel to the chain and perpendicular to the chain. Hence, we can understand such a nematic unit cell by considering two lengths: (i) $|V_1|$, chain link length and (ii) $|V_2|$, the distance between the chains.

We can see how these values change in relation to an increasing external field strength in Fig. [2.](#page-2-0) As the strength of the external field H is initially increased, the density of chains increases and $|V_2|$ decreases. Then we see that $|V_1|$ also decreases as the chain segments are squashed. Eventually when the external field H becomes large and approaches H_{c_2} , the vortices are forced together and form the familiar tightly packed triangular lattice. Note that V_1 is related to the fractional vortex separation of the model which is driven by β_{\perp} , and both V_1 and V_2 will

FIG. 2. Contour plots of the lattice solutions (skyrmion chains) for the model Eq. [\(1\)](#page-1-1) in the basal plane, with $\eta = 3$ and increasing external field H. The fields plotted are the order parameter magnitudes $|\psi_1|$ and $|\psi_2|$, the phase difference between the two complex order parameters $\theta_{12} = \varphi_1 - \varphi_2$ and the local magnetic field strength B. The lattice consists of skyrmion chains, formed by rectangular unit cells (marked with black lines) containing four half-quantum vortices or winding two. As the external field increases the skyrmion chains squash together. Then near H_c , a transition occurs and the vortices are forced into the conventional triangular lattice with unit cells of winding one. Note that there is no simple symmetry between $|\psi_1|$ and $|\psi_2|$ as the symmetry is broken to a mixed symmetry with spatial rotations [\[33\]](#page-7-4).

approximately scale inversely with η (the exact relation is nonlinear). We have also shown this behaviour is similar for other $n \gg 1$ [[31](#page-7-2)].

Theoretically in clean samples these skyrmion structures should be visible on the surface using scanning SQUID or scanning Hall probes. However, flux structure near the surface can be affected by surface physics and it is difficult to probe large areas. Hence, one of the most powerful tools to identify skyrmions is the muon spin relaxation technique, which gives statistical information regarding the flux distribution in the bulk of the system. To compare the lattice configurations in Fig. [2](#page-2-0) with the μ SR experiment, we must convert the numerical field configuration for a given unit cell to a probability density distribution $p(B)$. Our method is ideal for this as it produces the field configuration on a single unit cell, with no boundary effects and no deformations. We will use a standard kernel density estimation [\[41\]](#page-7-10) to approximate a continuous distribution from our discrete field configuration. Namely, our numerical algorithm produces a discrete configuration, where the field distribution is a set of normalized delta functions. We choose to smooth these delta functions by replacing them with localized Gaussians and taking the normalized sum. Hence, the probability density of field strength B in the unit cell is

$$
p(B) = \frac{1}{n_1 n_2 h} \sum_{i=1}^{n_1} \sum_{j=1}^{n_2} \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2} \left(\frac{B - B_{ij}}{h}\right)^2},
$$
(4)

FIG. 3. Magnetic field distribution for a skyrmion state in nematic superconductor (top) and vortex lattice in single component superconductor (bottom) for external fields closer to H_{c1} than H_{c2} . The nematic superconductor is described in Eq. [\(1\)](#page-1-1) with $\eta = 3$. The structure of the field distribution is markedly different for the skyrmion state in the nematic superconductor with a double peak, rather than the familiar single peak for the standard GL model.

where B_{ij} is the local magnetic field at lattice site (i, j) with $i \in [1, n_1]$ and $j \in [1, n_2]$. The parameter h determines the amount of smoothing and should be set relative to the amount of smoothing and should be set relative to the number of grid points used to simulate the lattice. We used $h = 0.01$ throughout.

We used the above method to construct the field distribution $p(B)$ for a nematic superconductor and a standard single component superconductor for comparison. The latter's parameters were chosen so that H_{c_1} and H_{c_2} match those of the nematic model for $\eta = 3$. This process and the parameters are described in Appendix B. We have plotted the nematic field distribution for field strengths nearer H_{c1} than H_{c2} in Fig. [3](#page-3-0), and also compared the results to the standard GL model. There is a clear difference, with the nematic superconductor exhibiting a double peak structure as opposed to the well known single peak of the standard GL model. It is also different from the μ SR signature of a system that phase separates into vortex clusters and domains of Meissner states [[42](#page-7-11),[43](#page-7-12)].

Consider the single component Ginzburg-Landau model. Here, the neighboring vortex separation is determined by a single parameter, the Ginzburg-Landau parameter κ . While this parameter is all that is needed to describe the lattice solution, it is actually the ratio of two length scales. The form of the magnetic field distribution is determined by how the magnetic field decays between neighboring vortices, hence as there is a single defining parameter, there is a single corresponding peak in the field distribution. In contrast, multicomponent anisotropic models have a continuous set of coupled length scales that are directionally

FIG. 4. Comparing the field distribution $p(B)$ of a nematic superconductor in the lower plot with the areas of the local magnetic field configuration $B(x)$ each peak corresponds to for external field strength $H = 2$.

dependent [[33](#page-7-4),[38](#page-7-7)]. However, given the rectangular nematic unit cell shown in Fig. [2](#page-2-0), we can describe the vortex separation in the unit cell using just two length scales, $|V_1|$ and $|V_2|$. Hence, the decay of the magnetic field between two neighboring vortices depends on whether they are in the same chain or neighboring chains. If the vortices are in the same chain the decay is determined by $|V_1|$, if they are in different chains the decay is determined by $|V_2|$. Hence, each of these neighboring decay possibilities, or length scales, contributes a different peak to the field distribution $p(B)$. This can be seen in Fig. [4](#page-3-1), where we have compared the probability density distribution $p(B)$ with the corresponding magnetic field plot $B(x)$ for $H = 2$. It can be seen in this figure that the first sharp peak corresponds to chain separation, and the second peak corresponds to the chain link length.

The two peaks for low fields start far apart, due to the large chain separation $|V_2|$. Then as the field strength increases, the chains are packed in much tighter and $|V_2|$ decreases, causing the two peaks to get closer. Then as the external field strength is increased further, both $|V_1|$ and $|V_2|$ shrink. This explains why the peaks begin far apart for particularly low fields. Hence, we ultimately end up with a very distinctive double peak structure for nematic skyrmion lattices. We expect that in the presence of weak pinning the statistics of the magnetic fields in the bulk will still retain these features.

While the above analysis holds for low and medium strength external fields, once the external field is increased to be near H_{c2} , the vortices are forced into a tightly packed triangular lattice. This triangular structure leads to the most efficient packing for the vortices, hence it is natural that the increased pressure will eventually force a triangular lattice to be lower in energy. The vortex separation for the nematic superconductor is now described by a single length scale and forms a single peak, similar to that of the single component model. This can be seen for strong external

FIG. 5. Magnetic field distribution for a nematic superconductor (top) and single component superconductor (bottom) for external fields $H_{c2} \approx H \gg H_{c1}$. The nematic superconductor is described in Eq. [\(1\)](#page-1-1) with $\eta = 3$. As one approaches H_{c2} , the structure of the field distribution for the nematic superconductor becomes similar to the single peak structure of the familiar single peak for the standard GL model. This corresponds with the lattice becoming a triangular lattice for external fields near H_{c2} .

fields in Fig. [5,](#page-4-0) where we have again compared the nematic and single component results.

Note that Fig. [3](#page-3-0) has a low field peak separation of about 0.5 which is approximately 0.13 $H_{c₂}$ (where $H_{c₂}$ is calculated in [[33](#page-7-4)]). Modern muon spectroscopy techniques are capable of resolving peak separations far lower than this in various materials. Experiments often exhibit a background peak at H_0 from muons that overshoot the sample. This can be resolved from the single type-2 peak despite being far closer than 10% of H_c [[39](#page-7-8)].

In conclusion, we have shown that the magnetic response of nematic superconductors takes the form of skyrmion stripes, which is markedly different from the standard vortex lattice. As the field is increased, these skyrmion stripes decrease their separation. Eventually the magnetization process compresses the skyrmion stripes themselves, forming the conventional triangular vortex lattice close to upper critical field H_{c2} . We also find that skyrmion stripes give qualitatively distinct signatures in muon spin rotation experiments. In particular, we observe a distinctive two peak structure, where the second peak changes relative to the first as the external field strength is increased. These findings can be used to experimentally confirm (i) nematic superconducting states and (ii) that flux-carrying objects in superconductors take the form of skyrmions. We have also presented new numerical tools, to find unit cells of any symmetry, that can be used to analyse the signals of other unconventional materials.

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Appendix A: Finding vortex lattice solutions.—In this Appendix we give a brief account of the new method used to find vortex lattice solutions of arbitrary complexity in the presence of an external field H. Detailed discussion of the method is presented in the companion paper [[33](#page-7-4)]. Namely, we find the unit cell of the periodic vortex solution in the bulk, with external field strength $H_{c_1} < H_0 < H_{c_2}$. The standard approach for a Ginzburg-Landau model is to consider a unit cell of degree $n = 1$ with either triangular $(\alpha = \pi/3)$ or square $(\alpha = \pi/2)$ symmetry. However, as the nematic system we are considering is anisotropic, there is no reason to expect that a lattice with such high symmetry will be the global minimizer. Hence, the correct approach is to minimize energy, not just with respect to the periodic fields, but also with respect to the geometry of the unit cell itself. We present here a general method to find the optimal unit cell, without assuming the symmetry of the underlying lattice, based upon the method first presented in [\[33\]](#page-7-4).

We first assume that far from the boundary of the system (deep in the bulk), the gauge invariant quantities $(\rho_{\alpha}, \theta_{12}, B)$ are periodic on the unit cell. We also note that we can represent a general periodic structure as a tessellation of a general parallelogram (unit cell), as seen in Fig. [1](#page-1-0), formed by two vectors V_1 and V_2 with angle α .

We then seek local minimizers of the Gibbs free energy in Eq. [\(1\)](#page-1-1) with respect to the fields $\psi_{\alpha}(x)$, $A_i(x)$, now defined on a general flat periodic 2-torus $T_{\Lambda}^2 = \mathbb{R}^2/\Lambda$,
where the period lattice Λ is spanned by vectors V_{Λ} . V_{Λ} in where the period lattice Λ is spanned by vectors V_1 , V_2 in Fig. [1](#page-1-0),

$$
\Lambda = \{ n_1 V_1 + n_2 V_2 | n_1, n_2 \in \mathbb{Z}, V_1, V_2 \in \mathbb{R}^2 \}. \tag{A1}
$$

The unit cell has periodic boundary conditions, up to gauge, so that the fields have boundary conditions,

$$
\psi_{\alpha}(x + V_i) = \psi_{\alpha}(x)e^{if_i},
$$

\n
$$
A_j(x + V_i) = A_j(x) + \partial_j f_i.
$$
 (A2)

Note that while the fields are not purely periodic, the above boundary conditions do leave all physical quantities periodic as desired,

$$
B(x + V_i) = B(x), \qquad J(x + V_i) = J(x),
$$

$$
|\psi(x + V_i)| = |\psi(x)|.
$$
 (A3)

To simulate the fields on the unit cell, we will simplify the above formulation by transforming to a more convenient coordinate system in the x, y plane. Let us define X_1 , X_2 so that $(x, y) = X_1V_1 + X_2V_2$. The unit cell spanned by V_1, V_2 is now covered by $(X_1, X_2) \in [0, 1] \times [0, 1]$. Let L
be the matrix with columns V, V, Note that det $I - 4$ be the matrix with columns V_1 , V_2 . Note that $\det L = A$, the area of the unit cell. Now define the unimodular matrix $M = \sqrt{A}L^{-1}$ such that $X_i = (M_{ij}/\sqrt{A})x_j$. Then

$$
F(M, \mathcal{A}, \psi_{\alpha}, A) = \int_{[0,1]^2} \left\{ \frac{1}{2} (MQ^{\alpha\beta} M^T)_{ij} \overline{D_i \psi_{\alpha}} D_j \psi_{\beta} + \frac{1}{2\mathcal{A}} [(\partial_1 A_2 - \partial_2 A_1)^2] + \mathcal{A} F_p(\psi_{\alpha}) \right\} \times dX_1 dX_2,
$$

leading in turn to the Gibbs free energy,

$$
G(M, A, \psi_{\alpha}, A) = F - \int_{[0,1]^2} H_0(\partial_1 A_2 - \partial_2 A_1) dX_1 dX_2
$$

= $F - 2n\pi H_0,$ (A4)

where n is the winding number of the field configuration. Note that this is only valid in the basal plane, as otherwise one must consider nonzero A_3 in the free energy. This would allow in-plane spontaneous fields and magnetic field twisting, which occurs away from the basal plane due to the coupled length scales [\[33\]](#page-7-4). Models that exhibit such fields have been studied in [\[33](#page-7-4)[,44](#page-7-13)[,45\]](#page-7-14).

If we exclude boundary effects, the optimal lattice solution is the one that minimizes the total Gibbs free energy of the system. If the system area is $A_{\rm sys}$, this can be calculated from the unit cell as $G_{\text{sys}} = A_{\text{sys}}G/A$. Hence, we seek minimizers of G/A or Gibbs free energy per unit area, with respect to the fields, geometry of the unit cell M and area A. Note that the degree or winding number of a given unit cell n is fixed and strictly speaking we must the find the global minimizer of all G_n/\mathcal{A} , where G_n is the Gibbs free energy of a unit cell of degree n and then find the infimum of this set.

In practice, it is sufficient to find minima of G_n/\mathcal{A} until we get a repeated minimizer, that is, until we find a minimizer of G_n/\mathcal{A} whose cell and field configuration is two cells of the $G_{n/2}/\mathcal{A}$ minimizer joined together.

Numerical method: To find minimizers of $G(M, \mathcal{A}, \psi_{\alpha}, A)$ we discretize the standard square unit cell on an $n_1 \times n_2$ regular grid. We utilize an arrested Newton flow algorithm subject to the periodic boundary conditions,

$$
\psi_{\alpha}[x + (1, 0)] = \psi_{\alpha}(x)e^{in2\pi nx_2},
$$

\n
$$
\psi_{\alpha}[x + (0, 1)] = \psi_{\alpha}(x),
$$

\n
$$
A_2[x + (1, 0)] = A_1(x) + 2\pi n,
$$

\n
$$
A_i[x + (1, 0)] = A_2(x) \quad i \neq 2,
$$

\n
$$
A_i[x + (0, 1)] = A_i(x), \quad (A5)
$$

where n is the winding number of the unit cell. We set the fields $\phi = (\psi, A)$ on the unit square torus and the geometry of the unit cell (M, \mathcal{A}) to be some initial condition. Fixing the unit cell (M, A) , we find a local minimum with respect to the collected fields ϕ , using arrested Newton flow for a notional particle subject to the potential $\ddot{\phi} =$ $-\text{grad}G_{\text{dis}}(M, \mathcal{A}, \phi)$. A more detailed description of arrested Newton flow is given in [[46](#page-7-15)]. This is continued until the discretized function is minimized with respect to some tolerance. We then fix the field configuration ϕ and area A, and minimize $G_{dis}(M, \mathcal{A}, \phi)/\mathcal{A}$ with respect to $M \in SL(2,\mathbb{R})$, where we note that only the gradient term of the energy is dependent on this matrix. This can be done exactly using linear algebra, described below in the section on finding the minimal geometry. Finally, we fix the fields ϕ and the lattice shape M and minimize the free energy with respect to A , by solving the algebraic equation in A with coefficients given by the integrals of the different energy terms.

Finding the minimal geometry: We can find the minimal geometry of a unit cell with a given configuration (ψ_{α}, A) explicitly. We first note that the only term that is dependent on the geometry M is the gradient term of the free energy. Hence, we write the gradient energy as a 4×4 complex matrix,

$$
P_{ik,jl}^c := \int_{[0,1]^2} Q_{kl}^{\alpha\beta} \overline{D_i \psi_\alpha} D_j \psi_\beta, \tag{A6}
$$

where we take each index pair $(i, k), (j, l)$ to range over the values $\{(1, 1), (1, 2), (2, 1), (2, 2)\}.$ Because of the symmetry $Q_{ij}^{\alpha\beta} = Q_{ji}^{\beta\alpha}$ we know P^c is Hermitian. Hence, if we split it into its real and imaginary parts $P_c^c = P + iP$, we split it into its real and imaginary parts $P^c = P + iP_I$ we know that the real part P is symmetric and the imaginary part P_I is skew. If we then consider the vector,

$$
X := (M_{11}M_{12}M_{21}M_{22})^T, \tag{A7}
$$

we can rewrite the gradient part of the free energy as

$$
F_{\text{grad}}(X) = \frac{1}{2} X^T P^c X = \frac{1}{2} X^T P X. \tag{A8}
$$

Hence, we must minimize the quadratic form X^T PX subject to the constraint that det $M = 1$ [or $M \in SL(2, \mathbb{R})$], which can be written

$$
\frac{1}{2}X^T J X = 1, \qquad J := \begin{pmatrix} 0 & 0 & 0 & 1 \\ 0 & 0 & -1 & 0 \\ 0 & -1 & 0 & 0 \\ 1 & 0 & 0 & 0 \end{pmatrix}.
$$
 (A9)

Hence, we seek critical points of F_{grad} : $SL(2,\mathbb{R}) \to \mathbb{R}$, which are $X \in \mathbb{R}^4$ with $X^T J X = 1$ and,

$$
PX = \lambda (X_4, -X_3, -X_2, X_1)^T = \lambda JX, \quad (A10)
$$

for some $\lambda \in \mathbb{R}$. We can see that $J^2 = \mathbb{I}_4$ is the identity matrix, hence Eq. [\(A10\)](#page-6-10) becomes

$$
JPX = \lambda X. \tag{A11}
$$

In other words, X is an eigenvector of JP .

This allows us to minimize F_{grad} with respect to M explicitly by constructing JP and finding its 4 eigenvectors Y. We discard any complex solutions and any such that $Y^T J Y \leq 0$. We compute the solution we seek $X =$ $Y/\sqrt{Y^T J Y}$ such that

$$
F_{\text{grad}} = \frac{1}{2} X^T P X = \frac{1}{2} X^T \lambda J X = \lambda.
$$
 (A12)

We can then read off the new period lattice as

$$
L = \sqrt{\mathcal{A}}M^{-1}, \qquad M = \begin{pmatrix} X_1 & X_2 \\ X_3 & X_4 \end{pmatrix}.
$$
 (A13)

The period lattice is spanned by the columns of L , named $V_1, V_2.$

Appendix B: Single component field distribution.—It is instructive to compare the results for nematic systems to a standard single component Ginzburg-Landau model. We will consider the following model,

$$
G = \int_{\mathbb{R}^2} \left\{ \frac{1}{2} \overline{D_i \psi} D_i \psi + \frac{1}{2} |B - H|^2 + F_p(|\psi|) \right\}, \quad (B1)
$$

where there is a single order parameter ψ , and

$$
F_p(|\psi|) = -a|\psi|^2 + \frac{b}{2}|\psi|^4.
$$
 (B2)

To effectively compare this model to the results for the nematic model, we will select the parameters a, b so that $H_{c₁}$ and $H_{c₂}$ equal those of the nematic model. By a standard computation, for example, see Tinkham [\[47\]](#page-7-16),

$$
H_{c_2} = 2a,\t\t(B3)
$$

while, by definition, H_{c1} is the value of external field strength H for which the Gibbs energy of a vortex solution equals that of the superconducting ground state, yielding

$$
H_{c1} = \frac{F_{\text{vortex}} - F_{\text{ground}}}{2\pi}.
$$
 (B4)

The right hand side of Eq. $(B4)$ depends implicitly on a, b. We find that $a = 5.70$, $b = 2.21$ produces $H_{c1} = 1.18$ and $H_{c2} = 11.41$, matching the nematic model with $\eta = 3$.

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