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High critical temperature and field superconductivity in Nb_{0.85}X_{0.15}, (X = Ti, Zr, Hf) alloys: promising candidates for superconducting devices

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Niobium and its alloys with early transition metals have been extensively studied for their excellent superconducting properties. They have high transition temperatures, strong upper critical fields, and high critical current densities, making them ideal for superconducting applications such as SQUIDs, MRI, NMR, particle accelerators, and Qubits. Here, we report a systematic investigation of as-cast Nb-rich alloys, Nb_{0.85}X_{0.15} (X = Ti, Zr, Hf), using magnetization, electrical transport, and specific heat measurements. They exhibit strong type-II bulk superconductivity with moderate superconducting transition temperatures and upper critical fields. The estimated magnetic field-dependent critical current density lies in the range of 10⁵–10⁶ A cm⁻² across various temperatures, while the corresponding flux-pinning force density is on the order of GNm⁻³, suggesting the potential of these materials for practical applications. Electronic-specific heat data reveal a strongly coupled, single, isotropic, nodeless superconducting gap. These Nb-rich alloys, characterized by robust superconducting properties, hold significant potential for applications in superconducting device technologies.

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1. Introduction

Superconductivity is one of the most fascinating phenomena in condensed matter physics, offering both deep insights into quantum states of matter and broad technological potential. Among various superconducting materials, niobium (Nb)-based superconductors have emerged as highly sought-after materials in research due to their potential applications, characterized by high superconducting transition temperatures (T_c) and upper critical fields (H_{c2}).^{1–4} Extensive studies have been conducted on Nb-based A-15^{5–9} and Nb-T (T: early transition metals)¹⁰ superconductors.¹¹ These materials show a wide array of physical properties, including large magnetoresistance,¹² non-trivial band topology,¹³ and the ability to form wires^{6,8} and superconducting joints.^{14,15} Such characteristics make them promising candidates for advanced technologies.^{16–18} In addition to industrial applications, Nb and Ta are widely recognized as suitable materials for superconducting radio frequency (SRF) cavities¹⁹ and Josephson junctions,^{20,21} which facilitate the development of superconducting qubits.^{22,23} Recent studies have shown that Ta–Zr²⁴ and Ta–Hf²⁵ alloys also exhibit promising superconducting properties. Owing to the high electropositivity of Hf and Zr, their alloys are considered strong candidates for meeting theoretical benchmarks

aimed at enhancing qubit performance. Nb and Ta-based alloys, characterized by diverse crystal structures and physical properties, can provide a foundation for exploring similar compounds.

Among Nb–T alloys, Nb–Ti alloys^{26–29} have achieved widespread commercial success in superconducting magnets and have also recently been studied under high pressure, setting new records for both T_c (= 19.1 K) and H_{c2} (= 19 T) among all alloys composed only of transition metals.³⁰ Additionally, Nb–Ti alloys with an Nb content of 50% or more are applicable in osseous implant devices, such as orthopedic and orthodontic implants.³¹ Furthermore, Nb–Zr^{32–37} and Nb–Hf^{38–40} alloys are also decisive in the same context. Nb–Zr alloys also present significant advantages for SRF cavities, including higher transition temperature, which facilitates reduced energy dissipation and cryogenic costs. Their simplified phase diagrams enable easier processing and optimization, while a stable ZrO₂ native oxide enhances RF performance by minimizing quench-inducing defects.⁴¹ In addition, the combination of high values of T_c and H_{c2} with low surface resistance, the absence of nodes in the gap symmetry, metallic behavior, high thermodynamic critical field and superheating field, and material morphology⁴² make these alloys excellent candidates for high-performance SRF cavities. However, most of the research on these Nb-based alloys has focused on the A15 phase; a detailed study of the cubic α -W structure, which is the same as pure niobium, is lacking. These alloys are expected to exhibit superior superconducting and

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physical properties, making them suitable for superconducting device applications. Therefore, a detailed study of niobium-rich alloys in the α -W structure is crucial to identify promising candidates for practical applications.

In this paper, we have investigated the superconducting and normal-state properties of the Nb-rich family of alloys, $\text{Nb}_{0.85}\text{X}_{0.15}$ (nominal composition: Nb_6X , X = Ti, Zr, Hf). We substitute elements Ti (3d), Zr (4d), and Hf (5d) at the X site with a fixed Nb content. We perform a comparative analysis of superconducting and normal-state parameters derived from electrical resistivity, magnetization, and specific heat measurements, which remain relatively unexplored in the literature. Temperature-dependent magnetization loops yield critical current densities in the range of 10^5 – 10^6 A cm⁻², while the corresponding flux-pinning force density reaches several GNm^{-3} , underscoring their viability in technological applications. All alloys crystallized in the α -W bcc structure and show type-II strongly coupled superconductivity with an isotropic gap. The results obtained from this study on alloys, combined with a malleable morphology required for surface preparation, highlight their potential for further investigation in a thin film form for future applications in superconducting devices, such as single-photon detectors and superconducting qubits.

2. Materials and methods

Polycrystalline samples of the $\text{Nb}_{0.85}\text{X}_{0.15}$ (X = Ti, Zr, Hf) series were synthesized by arc melting in a pure argon atmosphere on a water-cooled copper hearth. The chamber was first evacuated to a base pressure and then backfilled to a partial pressure. This evacuation and backfilling cycle was performed three times to ensure a highly inert atmosphere. To ensure an oxygen-free environment, a Zr getter was employed to absorb residual gaseous impurities. Stoichiometric amounts of the high-purity constituent elements (purity $\geq 99.9\%$) were melted together on a water-cooled copper hearth using a tungsten electrode. To achieve optimal homogeneity, the resulting ingot/button was flipped and remelted at least four to six times. The buttons were subsequently cooled naturally on the copper hearth. The weight loss of the buttons after melting was observed to be negligible. The final samples used for various measurements were prepared by mechanically sectioning the arc-melted buttons using a low-speed diamond saw with a cutting fluid to prevent localized heating and structural damage, yielding pieces of the required dimensions. Phase purity and crystal structure were analyzed by room-temperature powder X-ray diffraction (XRD) using $\text{K}\alpha$ ($\lambda = 1.5406 \text{ \AA}$) radiation using a PANalytical powder X-ray diffractometer. XRD refinement was performed using Fullprof Suite software.⁴³ Elemental composition and homogeneity were assessed by energy-dispersive X-ray analysis (EDXA) on a scanning electron microscope (SEM). Magnetization and AC susceptibility were measured using the magnetic property measurement system (MPMS3). Full range magnetization loops at different temperatures were measured to calculate the critical current density on the rectangular slab of all alloys. Electrical resistivity was measured using the standard four-probe method

at 1.9–300 K, while specific heat measurement was performed using the two-tau relaxation technique using a physical property measurement system (PPMS).

3. Results and discussion

3.1. Sample characterization

All alloys are crystallized in the cubic structure α -W, and the crystal structure of $\text{Nb}_{0.85}\text{X}_{0.15}$ (X = Ti, Zr, Hf) is shown in Fig. 1(a). The Rietveld refinement of the $\text{Nb}_{0.85}\text{Ti}_{0.15}$ alloy is shown in Fig. 1(b). The elemental mapping of all alloys is shown in Fig. 1(c). The XRD patterns of the polycrystalline samples of $\text{Nb}_{0.85}\text{X}_{0.15}$ (X = Ti, Zr, Hf) are shown in Fig. 1(d) and the lattice parameters determined from the refinement are summarized in Table 1.

3.2. Electrical resistivity

The temperature-dependent resistivity ($\rho(T)$) of Nb_6X (X = Ti, Zr, Hf) measured at zero field confirms the transition temperatures T_c of 9.68(2), 11.06(4), and 9.78(3) K, as shown in Fig. 2(a)–(c), and the insets show the corresponding zero drops in resistivity, respectively. The low residual resistivity ratio (RRR) is approximately 2, consistent with the RRR for polycrystalline samples of some Heusler-type superconducting alloys.⁴⁴

3.3. Magnetization

The superconducting transition temperature of all alloys was confirmed through temperature-dependent magnetization measurements performed in zero field-cooled warming (ZFCW) and field-cooled cooling (FCC) modes under an applied magnetic field of 1 mT, where the quantity $4\pi\chi_v$ is corrected for the demagnetization factor. The results, presented in Fig. 2(d)–(f), indicate T_c values of 9.65 (2) K, 11.05 (2) K and 9.67 (3) K for Nb_6X (X = Ti, Zr, Hf), respectively. Temperature and field-dependent magnetization measurements were performed to determine the lower critical field ($H_{c1}(0)$) and the upper critical field ($H_{c2}(0)$). The H_{c1} values at each temperature were identified as the point where the magnetization curve deviates from the Meissner line (solid black line). The temperature dependence of H_{c1} was analyzed using the Ginzburg–Landau (GL) relation given $H_{c1}(T) = H_{c1}(0) \left[1 - \left(\frac{T}{T_c} \right)^2 \right]$, which yields $H_{c1}(0) = 39.4$ (2) mT, 107.4 (2) mT and 46.5 (3) mT for Nb_6X (X = Ti, Zr, Hf), respectively. The temperature dependence of H_{c1} is presented in Fig. 3(a)–(c), where the insets display magnetization magnetic field curves at various temperatures for each sample. In addition, we estimate $H_{c2}(0)$ through magnetic field and temperature-dependent magnetization and resistivity measurements. The temperature dependence of H_{c2} was plotted and well fitted by the GL equation (eqn (1)), as shown in Fig. 3(d)–(f) for Nb_6X , (X = Ti, Zr, Hf), respectively.

$$H_{c2}(T) = H_{c2}(0) \left[\frac{1 - (T/T_c)^2}{1 + (T/T_c)^2} \right] \quad (1)$$

Eqn (1) yields estimated values of $H_{c2}(0) = 5.57(6)$, $9.51(1)$, and $5.81(5)$ T from magnetization and $7.57(7)$, $9.67(3)$ and $7.83(3)$ T



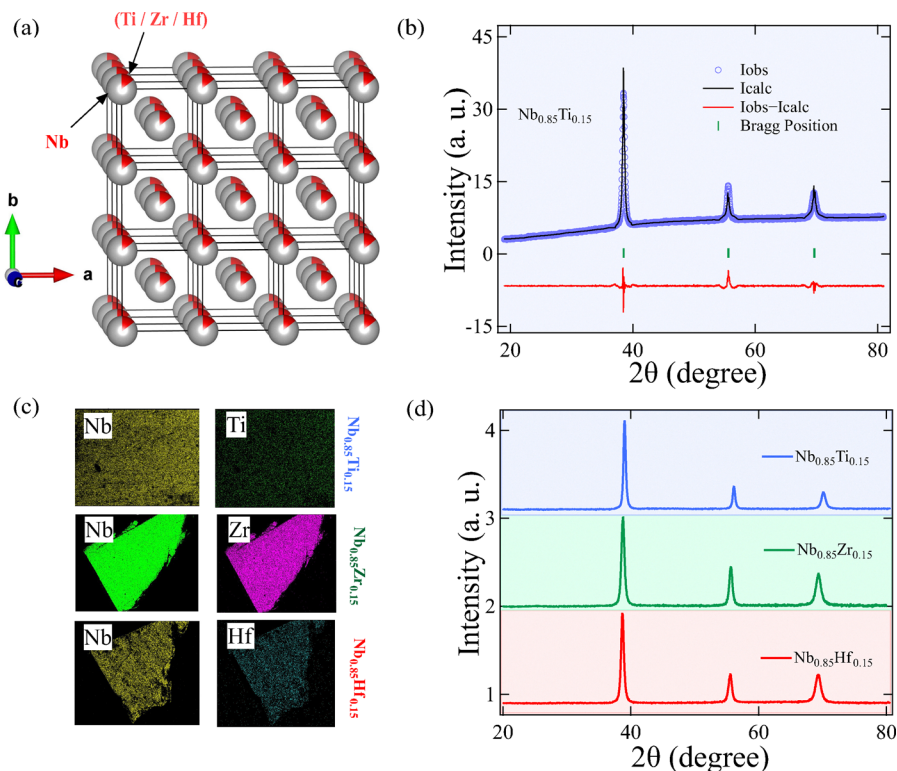


Fig. 1 (a) The bcc crystal structure of $\text{Nb}_{0.85}\text{X}_{0.15}$ ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$). (b) Rietveld refinement of the powder XRD pattern of Nb_6Ti alloy. (c) Elemental mapping of $\text{Nb}_{0.85}\text{X}_{0.15}$ ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$) alloys. (d) Room temperature XRD patterns of $\text{Nb}_{0.85}\text{X}_{0.15}$ ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$).

Table 1 Lattice parameters for $\text{Nb}_{0.85}\text{X}_{0.15}$ ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$) obtained from XRD refinement

Parameters	$\text{Nb}_{0.85}\text{Ti}_{0.15}$	$\text{Nb}_{0.85}\text{Zr}_{0.15}$	$\text{Nb}_{0.85}\text{Hf}_{0.15}$
$a = b = c$ (Å)	3.306(8)	3.350(1)	3.349(5)
V_{cell} (Å ³)	36.16(6)	37.59(3)	37.56(4)

from resistivity measurements for Nb_6X ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$), respectively, much higher than bulk Nb. The discrepancy between the $H_{c2}(0)$ values determined from the resistivity and magnetization measurements likely arises from the inherent limitations of the techniques, as well as their sensitivity to phenomena such as surface superconductivity and flux creep; however, for practical applications, the magnetization-derived value is more reliable, as it better reflects the bulk response of the sample.

The mechanisms behind Cooper pair-breaking when an external magnetic field is applied include the orbital pair-breaking effect and the Pauli limiting field effect. The orbital limiting field, $H_2^{\text{orb}}(0)$, can be described by the Werthamer–Helfand–Hohenberg (WHH) model,^{45,46} which neglects the spin–orbit interaction and Pauli paramagnetism. According to the WHH model, $H_2^{\text{orb}}(0)$ is expressed as

$$H_2^{\text{orb}}(0) = -0.693T_c \left. \frac{dH_{c2}(T)}{dT} \right|_{T=T_c} \quad (2)$$

By estimating the initial slopes of the H_{c2} vs. T curves, $-\frac{dH_{c2}(T)}{dT}$ at $T = T_c$, we obtain the orbital limiting fields using

eqn (2) as $H_2^{\text{orb}}(0) = 4.13(6)$, $7.11(5)$, and $4.43(3)$ T for Nb_6X ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$), respectively. For conventional superconductors, the Pauli limiting field is given by the expression $H_2^{\text{p}} = 1.86T_c$.^{47,48} Using the T_c values obtained from magnetization measurements, we estimate the Pauli limiting fields H_2^{p} as $17.94(3)$, $20.5(5)$, and $18.54(2)$ T for Nb_6X ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$), respectively, which are significantly higher than the corresponding estimated values of $H_{c2}(0)$. The coherence length $\xi_{\text{GL}}(0)$ ⁴⁹ can be estimated from $H_{c2}(0)$ using the relation $H_{c2}(0) = \frac{\Phi_0}{2\pi\xi_{\text{GL}}^2}$ where the magnetic flux quantum $\Phi_0 = 2.07 \times 10^{-15}$ T m².⁴⁹ The penetration depth $\lambda_{\text{GL}}(0)$ ⁵⁰ is obtained from the equation $H_{c1}(0) = \frac{\Phi_0}{4\pi\lambda_{\text{GL}}^2(0)} \left(\ln \frac{\lambda_{\text{GL}}(0)}{\xi_{\text{GL}}(0)} + 0.12 \right)$. The estimated values of $H_{c1}(0)$ and $H_{c2}(0)$ yield coherence lengths $\xi_{\text{GL}}(0)$ and penetration depths $\lambda_{\text{GL}}(0)$ as $76.9(1)$, $58.8(5)$ and $75.6(4)$ Å for $\xi_{\text{GL}}(0)$ and $1073(3)$, $615(2)$ and $975(4)$ Å for $\lambda_{\text{GL}}(0)$, for Nb_6X ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$), respectively. The GL parameter defined as $k_{\text{GL}} = \frac{\lambda_{\text{GL}}(0)}{\xi_{\text{GL}}(0)}$

was found to be much higher than $\frac{1}{\sqrt{2}}$ for the three alloys, indicating strong type II superconductivity. In addition, the thermodynamic critical field ($H_c(0)$) is defined as $H_{c1}(0)H_{c2}(0) = H_c^2(0) \ln(k_{\text{GL}})$. This yields the values of $H_c(0) = 288.6(2)$, $659.7(5)$, and $325.6(8)$ mT for Nb_6X ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$), respectively. However, the RF performance in the SRF cavities using bulk Nb reaches the theoretical limit of $H_c(0) \sim 200$ mT, and our results indicate that the Nb_6X series has a higher thermodynamic



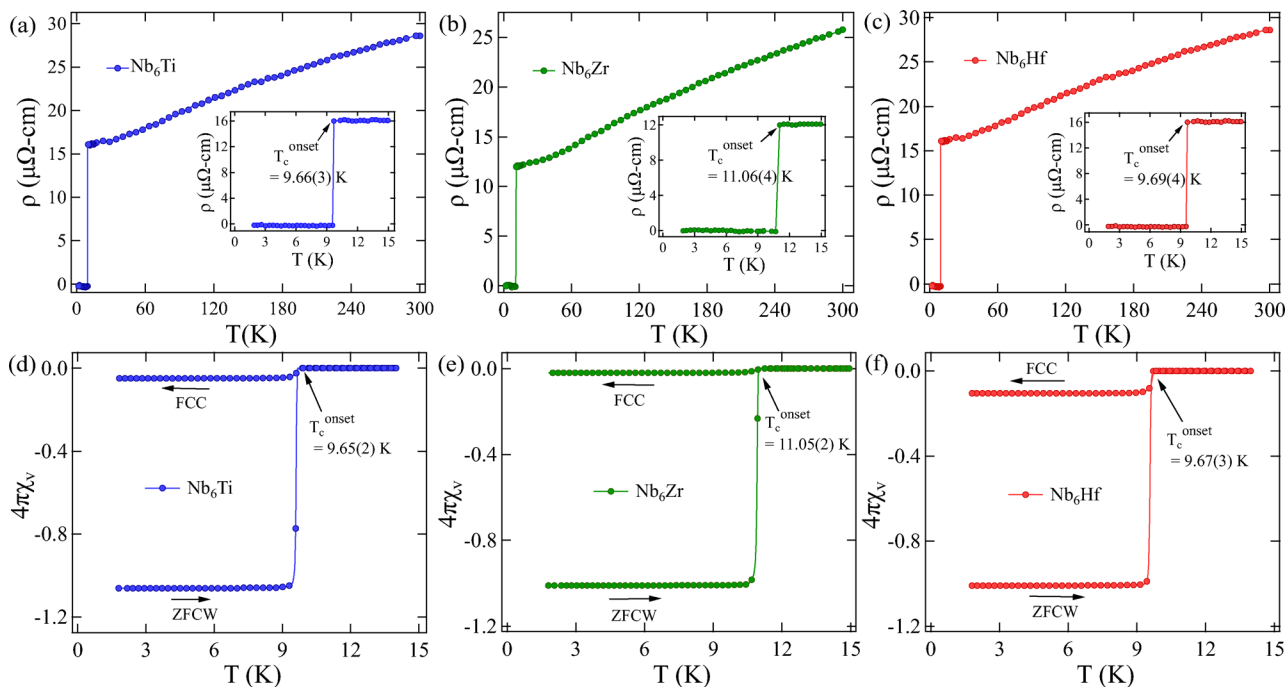


Fig. 2 Temperature-dependent electrical resistivity with the insets showing zero drops in resistivity for (a) Nb₆Ti, (b) Nb₆Zr, and (c) Nb₆Hf. Magnetization (corrected with demagnetization factor) in ZFCW and FCC mode at an applied field of 1 mT for (d) Nb₆Ti, (e) Nb₆Zr, and (f) Nb₆Hf, respectively.

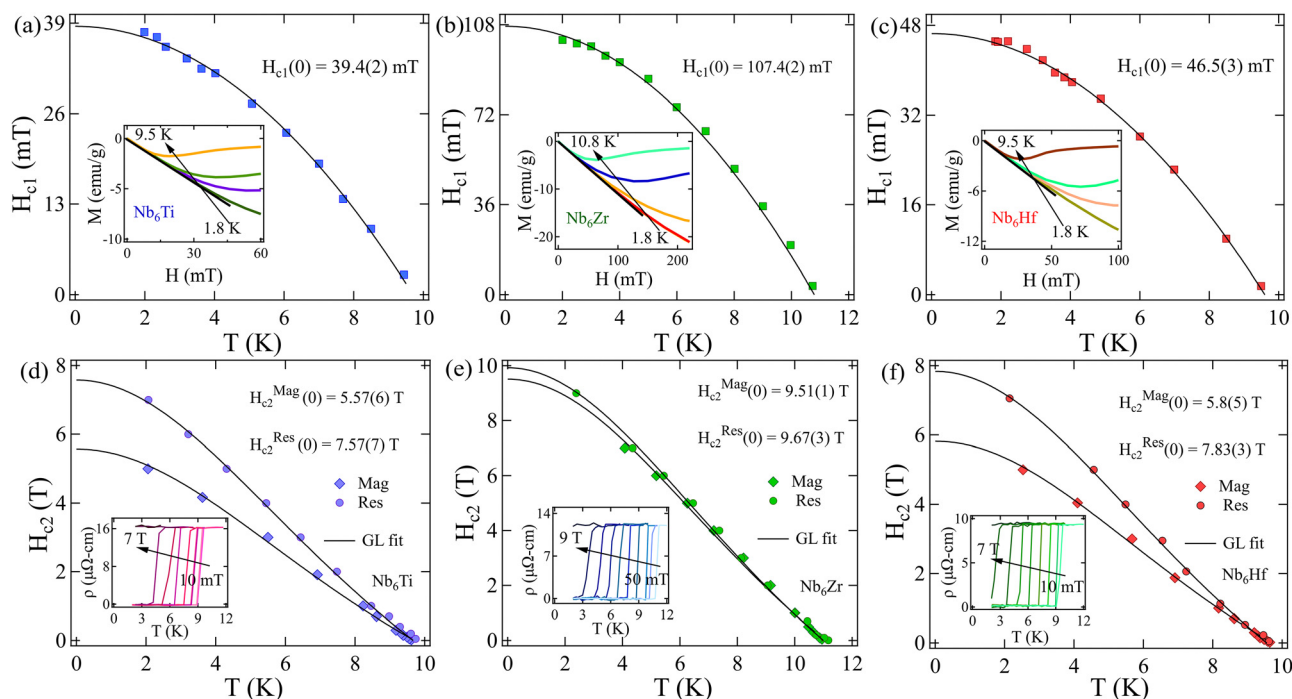


Fig. 3 Temperature-dependent lower critical fields and the insets show the field-dependent magnetization curves for (a) Nb₆Ti, (b) Nb₆Zr, and (c) Nb₆Hf, respectively. Temperature-dependent profiles of the upper critical field, estimated from magnetization and resistivity, where the insets show magnetic field-dependent resistivity data for (d) Nb₆Ti, (e) Nb₆Zr, and (f) Nb₆Hf, respectively.

critical field (>200 mT), hence the higher superheating field

$$\left(H_{sh} \approx \frac{0.89}{\sqrt{k_{GL}}} H_c\right)^{51}$$

The Ginzburg–Levanyuk number (Gi) quantifies the strength of thermal fluctuations relative to vortex unpinning in type-II

superconductors,⁵² which is described by the following expression

$$Gi = \frac{1}{2} \left(\frac{k_B \mu_0 \tau T_c}{4\pi \xi(0)^3 H_c^2(0)} \right)^2, \quad (3)$$



where τ is the anisotropy factor that is 1 for cubic Nb_6X ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$). Using the estimated T_c , $\xi(0)$ and $H_c(0)$ of magnetization measurement, we obtain $G_i = 6.18(2) \times 10^{-8}$, $1.48(4) \times 10^{-8}$, $4.43(8) \times 10^{-8}$ for Nb_6X ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$), respectively. These values are comparable to typical values of G_i observed for low T_c superconductors ($\sim 10^{-8}$). This suggests that thermal fluctuations do not contribute significantly to vortex unpinning in the Nb_6X series.⁵³

3.4. Critical current density and flux pinning force

Magnetic hysteresis loops (see Fig. 4(a)–(c)) of the Nb_6X ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$) series were measured at different temperatures to assess the critical current density (J_c) and its variation with magnetic field. Both Nb_6Zr and Nb_6Hf alloys show several low-field flux jumps in the M – H curves, as shown in Fig. 4(b) and (c). These flux jumps weaken and disappear when the temperature rises above 4 K or the magnetic field exceeds 3 T. Usually, flux jump phenomena are common in these alloys because of the strong magnetic flux pinning, while Nb_6Ti does not have a flux

pinning force at all, suggesting a relatively weak flux pinning force. The magnetic flux pinning force inside the Nb_6Zr sample is strongest, reflecting the high-low field flux jumps. The irreversible field (H_{irr}) of all samples is evaluated from the corresponding M – H loops at different temperatures. H_{irr} is the magnetic field above which vortices unpinning begins, and as the temperature increases, the irreversible fields of all samples decrease to various degrees. J_c is obtained by the Bean model⁵⁴ as described for J_c in the following equation:

$$J_c = \frac{20\Delta M}{w\left(1 - \frac{w}{3l}\right)} \quad (4)$$

where w and l are the width and length of the sample ($l \gg w$) perpendicular to the direction of the applied magnetic field, and ΔM is the width of the magnetization at the same magnetic field and Fig. 4(d)–(f) show the magnetic field dependence of J_c for Nb_6X ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$), respectively. The $J_c(0)$ at $T = 2$ K is extracted as $J_c(0) = 2.48 \times 10^5$ A cm⁻², 1.02×10^6 A cm⁻² and 4.87×10^5 A cm⁻² for Nb_6X ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$), respectively. In

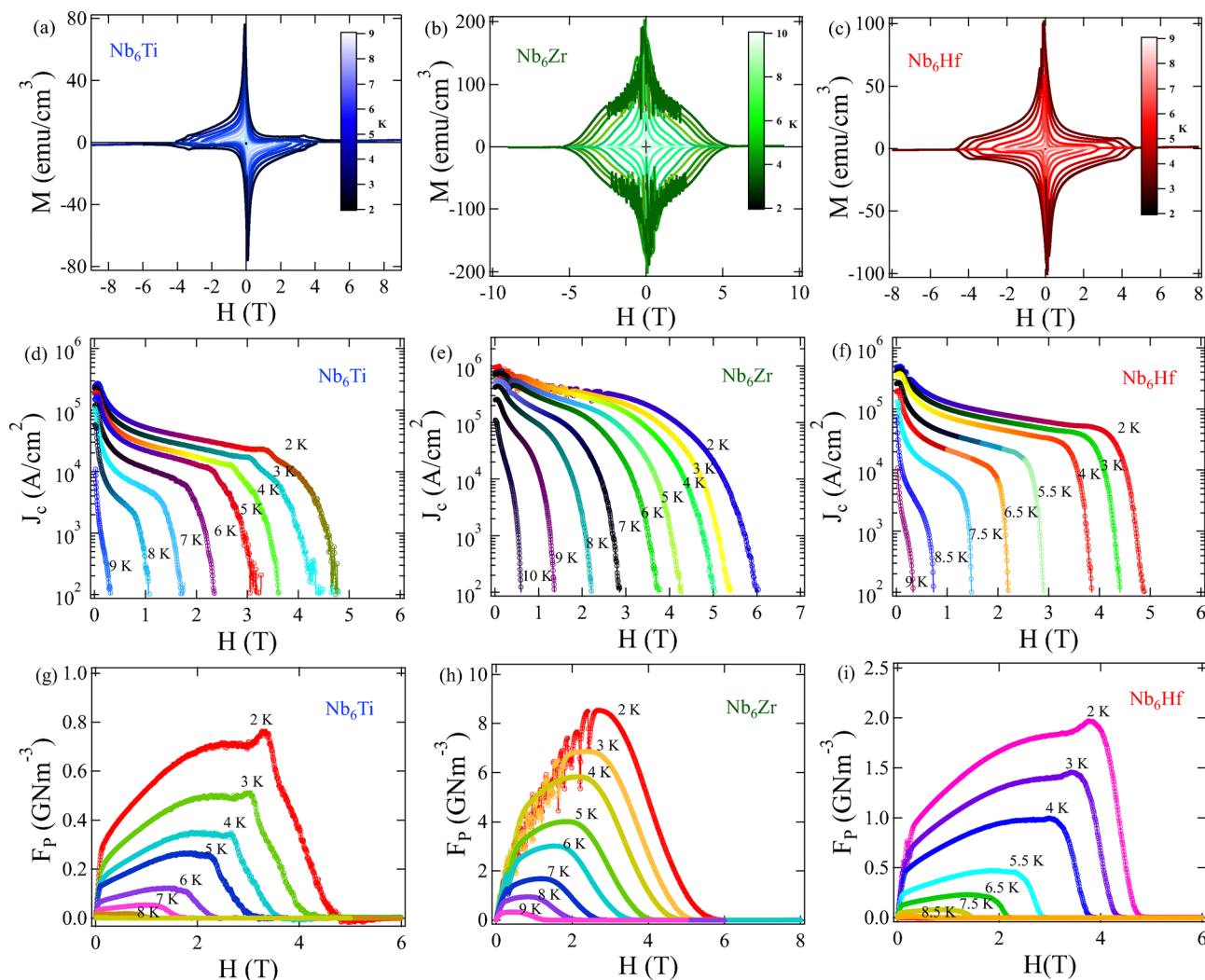


Fig. 4 (a)–(c) M – H loops at different temperatures, whereas (d)–(f) magnetic field dependent J_c variation at different temperatures and (g)–(i) variation of flux pinning force with magnetic field for Nb_6X , ($\text{X} = \text{Ti}, \text{Zr}, \text{Hf}$), respectively.



addition, the pinning force density was calculated using the relation $F_p = \mu_0 H \times J_c$.⁵⁵ The variations of the flux pinning force density with magnetic field for Nb₆X, (X = Ti, Zr, Hf) are shown in Fig. 4(g)–(i), respectively. The density of the pinning force is obtained in the order of Gnm⁻³ for this series, with the highest value obtained for Nb₆Zr due to the presence of low-field flux jumps, which is relatively low in Nb₆Hf and negligible in Nb₆Ti, suggesting the lowest value of the pinning force in Nb₆Ti. While flux jumps are commonly observed in Nb-based superconductors due to thermomagnetic instabilities, their absence in certain Nb alloys can be attributed to enhanced thermal conductivity and homogeneous flux pinning, which suppress abrupt vortex avalanches.⁵⁶ The critical current density and flux pinning force values for all alloys exceed the practical threshold value.

3.5. Specific heat

Temperature-dependent specific heat measurements were performed for all samples without an external magnetic field to study the thermal properties of Nb₆X (X = Ti, Zr, Hf). Normal state C/T vs. T^2 data were fitted using the Debye–Sommerfeld model represented by eqn (5) and shown in Fig. 5(a)–(c) with solid black curves.

$$C/T = \gamma_n + \beta T^2 \quad (5)$$

where γ_n is the Sommerfeld coefficient, and β holds the phononic contributions to the specific heat. Fitting normal state specific heat data using eqn (5) gives $\gamma_n = 51.17(6)$, 61.45(8), 51.93(7) mJ mol⁻¹ K⁻² and $\beta = 0.63(1)$, 0.79(2),

0.99(3) mJ mol⁻¹ K⁻⁴ for Nb₆X (X = Ti, Zr, Hf), respectively. γ_n is related to the density of states on the Fermi surface $D_c(E_f)$ by the relation $\gamma_n = \left(\frac{\pi^2 k_B^2}{3}\right) D_c(E_f)$, where $k_B \approx 1.38 \times 10^{-23}$ J K⁻¹. $D_c(E_f)$ is estimated to be 21.7(1), 26.1(2), and 22.1(1) states per eV f.u. for Nb₆X, (X = Ti, Zr, Hf), respectively. The Debye temperature (θ_D) is related to β as $\theta_D = \left(\frac{12\pi^4 RN}{5\beta_3}\right)^{1/3}$, where N is the number of atoms per formula unit and R is the molar gas constant (8.314 J mol⁻¹ K⁻¹), which calculates $\theta_D = 277.8(9)$, 258.2(4), and 239.5(3) K for Nb₆X, (X = Ti, Zr, Hf), respectively.

McMillan's model estimates the electron–phonon coupling strength from a dimensionless quantity λ_{e-ph} ,⁵⁷ which depends on the estimated values of θ_D and T_c as

$$\lambda_{e-ph} = \left[\frac{1.04 + \mu^* \ln(\theta_D/1.45 T_c)}{(1 - 0.62\mu^*) \ln(\theta_D/1.45 T_c) - 1.04} \right] \quad (6)$$

where μ^* is screened Coulomb repulsion, considering $\mu^* = 0.13$ as described for transition metals,⁵⁷ and estimated $\lambda_{e-ph} = 0.83(6)$, 0.92(4), and 0.89(7) for Nb₆X, (X = Ti, Zr, Hf), respectively. This classifies them as strongly coupled superconductors. After subtracting the term βT^3 from total specific heat (C), the temperature-dependent electronic specific heat (C_{el}) was extracted for all samples and fitted with the isotropic single gap model, represented as solid black curves in Fig. 5(d)–(f) and the quantity $\frac{\Delta C_{el}}{\gamma_n T_c} = 1.87(2)$, 1.79(5), and 1.85(3) for Nb₆X, (X = Ti, Zr, Hf), respectively.

Furthermore, C_{el} vs. T data can provide information about the symmetry of the superconducting gap around the Fermi

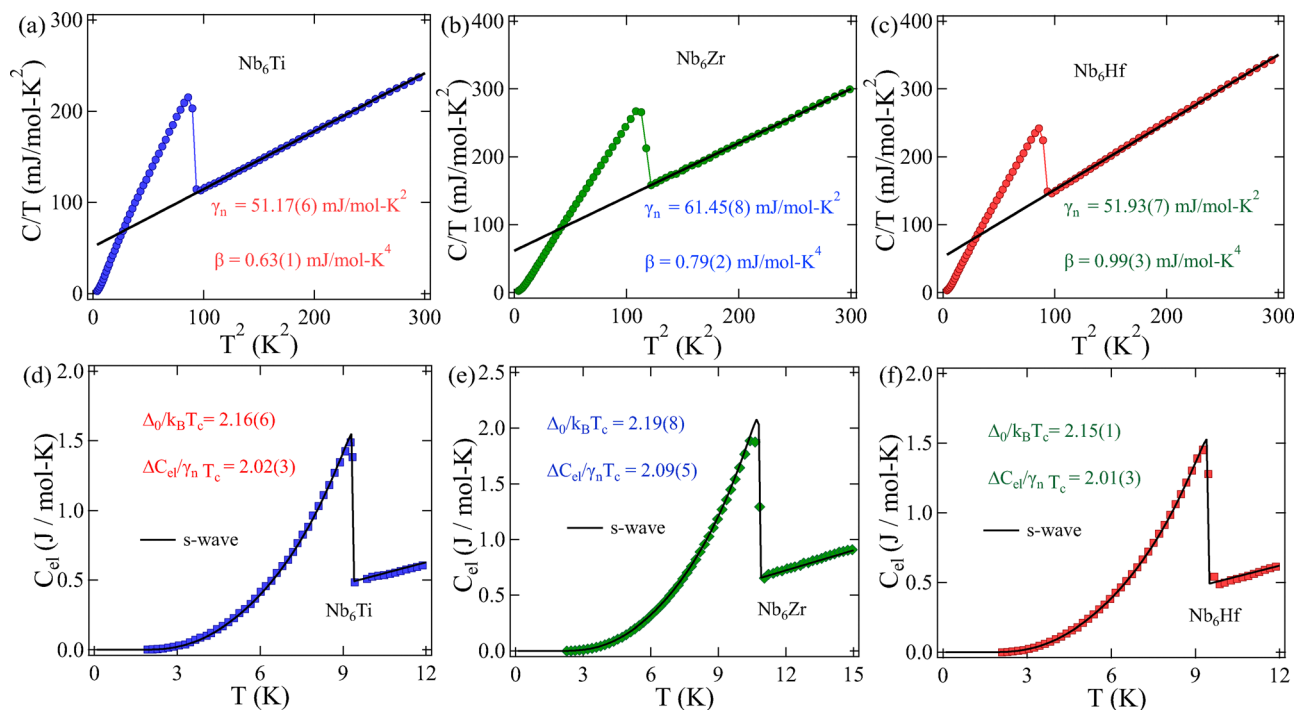


Fig. 5 (a)–(c) C/T vs. T^2 plot for Nb₆X, (X = Ti, Zr, Hf), respectively, where the solid black lines represent the Debye–Sommerfeld fitting represented by (eqn (5)). (d)–(f) C_{el} vs. T plot for Nb₆X (X = Ti, Zr, Hf), respectively, where solid black curves represent the s-wave model fit with BCS type single gap function.



surface, revealing the intricate pairing mechanism. Normalized entropy (S) in the superconducting region and C_{el} can be related as

$$\frac{C_{el}}{\gamma_n T_c} = t \frac{d(S/\gamma_n T_c)}{dt} \quad (7)$$

where $t = T/T_c$, the reduced temperature. Within the framework of the BCS approximation,^{58,59} the normalized entropy for a single BCS-like gap is defined by the following relation.

$$\frac{S}{\gamma_n T_c} = -\frac{6}{\pi^2} A_T^s \int_0^\infty [f \ln(f) + (1-f) \ln(1-f)] dy \quad (8)$$

where, $A_T^s = \frac{\Delta(0)}{k_B T_c}$, $f(\xi) = [\exp(E(\xi)/k_B T) + 1]^{-1}$ is the Fermi function, $E(\xi) = \sqrt{\xi^2 + \Delta^2(t)}$, where $E(\xi)$ is the energy of normal electrons measured relative to the Fermi energy, $y = \xi/\Delta(0)$, and $\Delta(t) = \tanh[1.82(1.018((1/t) - 1))^{0.51}]$ represents the temperature-dependent superconducting energy gap. The solid black curves in Fig. 5(d)–(f) represent the fit to electronic specific heat data using a single isotropic nodeless gap model as described by eqn (8), which provides the normalized superconducting gap values $\Delta(0)/k_B T_c = 2.16(1)$, $2.19(6)$, and $2.15(3)$ for Nb_6X ($X = Ti, Zr, Hf$), respectively, which confirms the strong electron–phonon coupled superconductivity.

4. Normal state properties and Uemura classification

In this section, we have analytically solved a set of equations simultaneously to investigate electronic properties and discussed the Uemura classification of Nb_6X ($X = Ti, Zr, Hf$).⁶⁰ The Sommerfeld coefficient (γ_n) is related to the quasiparticle number-density (n), defined as

$$\gamma_n = \left(\frac{\pi}{3}\right)^{2/3} \frac{k_B^2 m^* V_{f.u.} n^{1/3}}{\hbar^2 N_A} \quad (9)$$

and Fermi velocity (v_F) is related to the electronic mean free path (l_e) and n by the expressions

$$l_e = \frac{3\pi^2 \hbar^2}{e^2 \rho_0 m^{*2} v_F^2} \text{ and } n = \frac{1}{3\pi^2} \left(\frac{m^* v_F}{\hbar}\right)^3 \quad (10)$$

respectively, where k_B is the Boltzmann constant, m^* is the effective mass of the quasiparticles, $V_{f.u.}$ is the volume of the formula unit, N_A is the Avogadro number and ρ_0 is the residual resistivity.

Within the dirty limit superconductivity, the BCS coherence length of a superconductor is much larger than the mean free path ($\frac{\xi_0}{l_e} \gg 1$), and the scattering of electrons with impurities and defects may affect the superconducting properties. The GL penetration depth (λ_{GL}) is related to the London penetration depth (λ_L) at absolute zero temperature followed by the expression

$$\lambda_{GL}(0) = \lambda_L \left(1 + \frac{\xi_0}{l_e}\right)^{1/2}, \text{ where } \lambda_L = \left[\frac{m^*}{\mu_0 n e^2}\right]^{1/2} \quad (11)$$

and BCS coherence length relates to GL coherence length and is expressed by

$$\frac{\xi_{GL}(0)}{\xi_0} = \frac{\pi}{2\sqrt{3}} \left(1 + \frac{\xi_0}{l_e}\right)^{-1/2}. \quad (12)$$

We have solved the above equations simultaneously and estimated the values of m^* , n , v_F , ξ_0 , and l_e , using the previously obtained values of γ_n , ρ_0 , $\xi_{GL}(0)$ and $\lambda_{GL}(0)$. The Fermi temperature for an isotropic spherical Fermi surface is defined as

$$T_F = \frac{\hbar^2 k_F^2/3}{2m^* k_B} \quad (13)$$

where $k_F = 3\pi^2 n$, the Fermi wave vector. The ratio T_c/T_F classifies superconductors into the conventional or unconventional category. According to Uemura *et al.*,⁶⁰ the unconventional range $0.01 \leq \frac{T_c}{T_F} \leq 0.1$, shown by the green shaded band in Fig. 6, whereas the T_c/T_F values for Nb_6X ($X = Ti, Zr, Hf$) lie outside this unconventional region. All superconducting and normal state parameters of Nb_6X ($X = Ti, Zr, Hf$), estimated from various techniques, are listed in Table 2. This comparison highlights that, despite variations in microstructure and alloying, Nb-based systems generally exhibit weak to moderate electronic correlations and conventional electron–phonon-mediated pairing. The similarity of T_c/T_F values across these families underscores that superconductivity in Nb_6X alloys is governed by the same fundamental mechanism, situating them firmly within the conventional regime while still offering competitive superconducting parameters for practical applications.

5. Discussion

Fig. 7 presents a comparison of key superconducting parameters (T_c , H_{c2} and J_c) for Nb–T alloys α -W type ($T = Ti, Zr, Hf$) and other well-known compounds based on niobium. The data provide insight into the variation in superconducting behavior between different material classes. Niobium-based

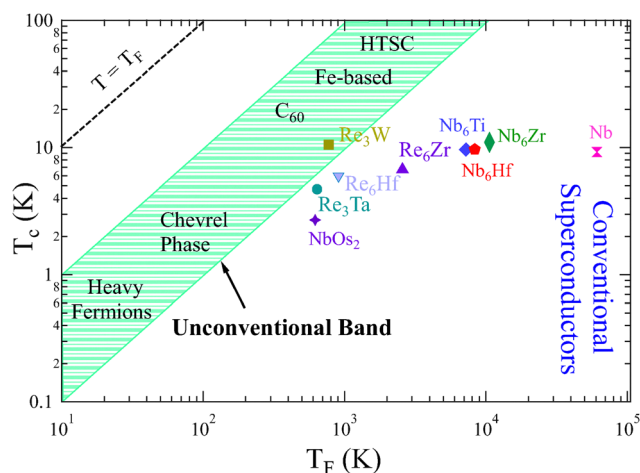


Fig. 6 Uemura plot defined by T_c and T_F for Nb_6X ($X = Ti, Zr, Hf$), where the green shaded region represents the unconventional band.



Table 2 Superconducting and normal state parameters of Nb₆X, (X = Ti, Zr, Hf)

Parameters	Unit	Nb ₆ Ti	Nb ₆ Zr	Nb ₆ Hf
T_c	K	9.65(2)	11.05(2)	9.67(3)
$H_{c1}(0)$	mT	39.4(2)	107.4(2)	46.5(3)
$H_{c2}(0)$	T	5.57(6)	9.51(1)	5.81(5)
$H_{c2}(0)^{\text{Pauli}}$	T	17.94(3)	20.5(5)	18.54(2)
$H_{c2}(0)^{\text{orb}}$	T	4.13(6)	7.11(5)	4.43(3)
$\xi_{\text{GL}}(0)$	Å	76.9(1)	58.8(5)	75.6(4)
$\lambda_{\text{GL}}(0)$	Å	1073(3)	615(2)	975(4)
$H_c(0)$	mT	288.6(2)	659.7(5)	325.6(8)
k_{GL}		13.9(5)	10.4(5)	12.9(4)
$\Delta C_{\text{el}}/\gamma_n T_c$		2.02(3)	2.09(5)	2.01(3)
$\Delta(0)/k_B T_c$		2.16(6)	2.19(8)	2.15(1)
θ_D	K	277.8(9)	258.2(4)	239.5(3)
$\lambda_{\text{e-ph}}$		0.83(6)	0.92(4)	0.89(7)
v_F	10^5 ms^{-1}	1.2(4)	1.4(6)	1.3(2)
n	10^{28} m^{-3}	13.4(2)	23.5(4)	15.6(3)
T_F	K	7234(4)	10 595(25)	8346(8)
T_c/T_F		0.0013(3)	0.0010(3)	0.0011(5)
m^*/m_e		15.2(3)	15.1(2)	14.6(2)

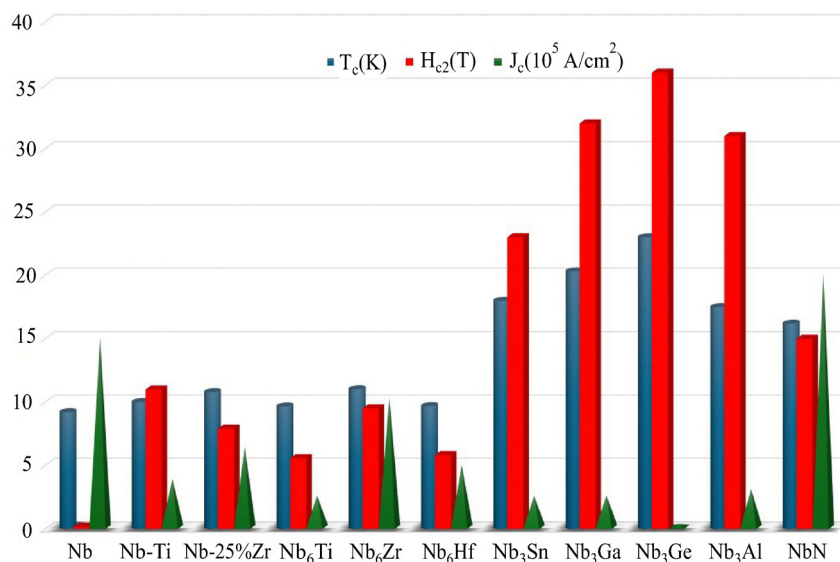
superconductors, especially binary alloys of Nb–T (where T = Ti, Zr, Hf) and Nb₃X compounds of type A15, represent two technologically significant families of low-temperature superconductors, each offering unique advantages tailored to specific operational demands. Nb–Ti alloys are widely used in commercial superconducting technologies due to their good thermal performance and suitability for wire manufacturing on a large scale.^{26,27,30,31} Introducing alloying elements such as Ti or Zr into niobium enhances flux pinning capabilities while maintaining critical superconducting parameters, with transition temperatures ranging from 9 to 11 K and upper critical fields in the range of 6–14 T⁶⁸ at ambient pressure. Specifically, incorporating zirconium into niobium has been shown to enhance surface superconductivity and radio-frequency characteristics, aided by the formation of a stable ZrO₂ surface layer and improved electron–phonon interactions.⁴¹ In addition, alloys such as

Nb–Zr and Nb–Hf exhibit strong resistance to radiation damage, making them particularly suitable for nuclear environments.⁶⁹ Recent computational studies suggest that irradiation promotes the formation of Nb-rich precipitates within Zr matrices, which adopt energetically favorable platelet morphologies. These structures contribute to enhanced microstructural integrity in nuclear fuel cladding materials under irradiation conditions.^{70–72}

In contrast, A15-type intermetallics such as Nb₃Sn, Nb₃Al, and Nb₃Ge demonstrate significantly enhanced superconducting characteristics, with critical temperatures (T_c) reaching up to 23 K and upper critical magnetic fields (H_{c2}) exceeding 30 T. These remarkable properties are mainly attributed to their distinctive crystal structure, characterized by quasi-one-dimensional chains of Nb atoms and a densely packed A₃B lattice.¹³ This arrangement strengthens electron–phonon interactions near the Fermi level, thereby promoting superconductivity. However, these materials are inherently brittle and exhibit a high sensitivity to slight variations in composition, which presents challenges during processing. Beyond the superconducting parameters summarized in Fig. 7, prior studies have shown that microstructural refinement (*e.g.*, grain size control in Nb₃Sn⁶¹ and stacking fault regulation by nano-oxide particles in Nb₃Al⁹), defect density introduced by irradiation,⁷³ and processing history such as hot rolling or equal channel angular pressing (ECAP) in NbTi⁷⁴ critically determine flux pinning and current-carrying capacity. These insights also contextualize the potential of Nb₆X alloys. The bulk properties observed here motivate future efforts in thin-film growth and characterization, where enhanced superconducting performance and device integration could be realized.

6. Conclusion

In summary, we have synthesized and characterized the Nb-rich series, Nb_{0.85}X_{0.15} (nominal composition: Nb₆X, X = Ti, Zr, Hf),

**Fig. 7** Comparison of T_c , H_{c2} , and J_c of Nb₆X (X = Ti, Zr, Hf) with niobium⁵⁴ and its alloys.^{9,32–37,61–67}

which crystallizes in a α -W bcc structure. We have thoroughly investigated their superconducting and normal state properties through XRD, electrical resistivity, magnetization, and specific heat measurements. All normal and superconducting state parameters follow nonmonotonic trends in the Nb₆X series, since the X component is substituted by elements of Ti (3d), Zr (4d) and Hf (5d). The calculated critical current density from temperature-dependent magnetization loops has values in the range 10⁵–10⁶ A cm⁻² with a flux pinning force on the order of GNm⁻³, which are good values for practical applications. All alloys have moderate transition temperatures and upper critical fields following isotropic nodeless superconducting gaps with strong electron–phonon coupling strength. The variation in spin–orbit coupling strength from 3d to 5d elements, while maintaining a fixed Nb content, shows only negligible influence on superconducting properties and gap symmetry, but it can provide further insight into the superconducting ground state and its relation to mechanical properties. Due to their morphological malleability and superior superconducting properties, this family of binary alloys holds significant promise and dedicated thin-film studies will be essential to fully explore their technological potential in superconducting devices.

Conflicts of interest

There are no conflicts to declare.

Data availability

The data supporting the findings of this study can be made available from the corresponding author upon reasonable request.

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