Polaritonic Behaviors of SiC, cBN and GaN for Spectrally-Selective Nano-Optics Applications

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1. Introduction
With the improved manipulation, handling, and fabrication of materials at the micro and nanoscale, the distinctive properties and functions of micro- and nanomaterials have become increasingly important for available and potential technological applications. Energy use with lowest possible waste and highest efficiency has reached to the level necessitating an urgent action to be taken due to environmental consciousness and limited sources. Thus far, this has motivated researchers to take a look at the bottom as Richard P. Feynmann suggested in his paper (Feynmann 1960) to see potential ways of employing nanomaterials, to improve efficiencies and sensitivities of processes. Thermal radiation becomes dependent on the...
distance between heat-exchanging objects when they are separated by a distance smaller than the dominant emitted wavelength, i.e., \( d < \lambda \) (Howell et al. 2021). In this case, the regime of radiative transfer is named the near-field (NF), for which the local density of electromagnetic states (LDOS) and consequently the amount of the transferred heat significantly exceeds the blackbody limit (Didari and Mengüç 2017). This increase depends on the material type of the objects, the distance between them \( (d) \), their temperatures, the environment they are held in, their shapes (Didari and Mengüç 2015) and surfaces (Chen and Xuan 2015), to name a few variables. The radiative transfer in the NF may find applications in areas including energy harvesting such as NF-thermophotovoltaics (Elçioğlu 2018), tip-based applications such as nano-manufacturing (e.g., as described by (Dönmez 2009, Dönmez et al. 2010), material characterization (e.g., with tunneling microscopy), and radiation-based thermal rectification (Park and Zhang 2013). So far, the near-field radiative transfer (NFRT) has been investigated experimentally and theoretically to unfold the mechanisms involved in this increased heat transfer. The micro-nano fabrication techniques also enabled optimizing materials to end up with higher LDOS, higher NFRT, and higher performance (e.g., as shown by (Didari and Mengüç 2017, Didari et al. 2018a)).

The main mechanism that makes NFRT different from that at the conventional scale (far-field, FF) is the contribution of surface modes, whose impact is negligible in the FF, as these surface (evanescent) waves having a decay length of around a wavelength (Francoeur 2010) making their contribution to FF radiation (e.g., for the objects separated by a distance much larger than the dominant emitted wavelength) negligible. In the NF, these evanescent modes can dominantly be harvested upon proper system configuration and setting. Figure 1 shows the generic NF – FF distinction based on a plane surface of a smooth slab.

Surface polaritons as excitations with the incident electromagnetic field, also propagate along the interface of two objects (depending on the type of materials) with decaying evanescent field in both (Francoeur 2010). The type of surface polaritons that contribute to the NFRT is surface plasmon polaritons for metals and semiconductors, and surface phonon polaritons in polar materials.

The main configurations studied in terms of the NFRT are plate-plate, sphere-plate, and sphere-sphere, all of which includes objects with a finite temperature gradient separated by a sub-wavelength distance. The most commonly investigated system setting has been parallel plates which holds potential for NF energy conversion and harvesting. The LDOS and NFRT have been investigated by covering a wide range of materials and system settings (e.g., separation, temperatures, etc.) The formulations of NFRT (as presented by (Francoeur and Mengüç 2008, Basu 2016), among others) reveal that once the separation between objects and their temperature are set, the material selection plays the major role in enhancing, or suppressing (Zhao et al. 2012) the NFRT. The parallel-plates configuration consists of two materials facing one another and separated by a vacuum gap. In this regard, understanding the importance of the material from its polaritonic performance at the material-vacuum interface is important.

SiC, cBN, and GaN support surface phonon polaritons in the IR, and as oppose to their plasmonic counterparts, they exhibit lower optical losses (Caldwell et al. 2015). This phenomenon is expected to cause a quasi-monochromatic spectral
radiative heat flux in the NF, potentially leading to higher performance (Francoeur 2010). It has been shown that the spectral NFRT flux between pairs of SiC, cBN, and GaN in parallel plates configuration exceeds the blackbody limit by orders of magnitudes (e.g., as mentioned by (Elçioğlu et al. 2017, Francoeur 2015) among others).

From a more materials-based point of view, performance measures for systems utilizing surface polaritons have been considered useful in conveniently assessing performance. The term figure of merit stands for a quantitative performance indicator (for a material or a system) that can be obtained more easily and conveniently than performing a whole system treatment. This applies to systems utilizing surface polaritons as well, generally by considering energy carried by polaritons, as well as their propagation and confinement. Maier (2006) mentioned a quality factor, Q, which was defined as the ratio of the energy stored to the loss per cycle; and for metals, Q was limited by the dissipative losses inside the material. It was shown that electromagnetic energy could be confined inside plasmonic cavities far below the diffraction limit, and a case study for Au half spaces was presented. More detailed figures of merit were provided for both surface plasmon polaritons and surface phonon polaritons by Caldwell et al. (2015) and comparisons were carried out to assess phononic and plasmonic materials performances by accounting for the quality factor values, locations of electrostatic resonances, and damping times. They presented figure of merit values for certain polar dielectrics, metals, metal alloys, and doped semiconductors. Berini (2006) proposed three figures of merit to quantify the spectral response of plasmon waveguides. The definition was based on a benefit-to-cost ratio where benefit was the confinement of surface plasmons, and the cost was the attenuation of them. Defining a figure of merit was considered important, since the confinement and attenuation are correlated, and there was a trade-off between them. Three figures of merit were defined for surface plasmons for a single-interface, metal slab bounded by dielectric, and dielectric slab bounded by metal cases. Configurations with Ag and SiO2 were considered, and their figures of merit were calculated. Buckley and Berini (2007) studied three figures of merit for 2D surface plasmon waveguides and considered different shaped (narrow, wide, thin, thick) metal (Au, Ag, Al) stripes (as one, two, three, and cladded) as the configuration. Each metal considered exhibited their highest performance at a wavelength region specific to them. Ordonez-Miranda et al. (2017) developed a polaritonic figure of merit expression and provided explicit and analytical expressions for (i) single interface, and (ii) suspended thin film cases. These polaritonic figure of merits are functions of the configuration (i or ii) and the material permittivity. They also analyzed the three polaritonic figures of merit they derived considering a single interface between (i) a SiC thin film-dielectric interface, and (ii) suspended SiC thin film cases, considering two different temperatures, 295 K and 686 K, for SiC’s Drude-Lorentz permittivity terms. They concluded that for SiC-dielectric interface case, the surface phonon polaritons carried more energy at around transverse optical phonon frequency (ωTO) of SiC.

In order to investigate polaritonic performance from materials point of view, the fundamental material-vacuum interface case has been considered in this work. The calculation performed by Ordonez-Miranda et al. (2017) for SiC (at 295 K) has been repeated in this work for comparison purposes against the polaritonic figure of merits of cBN and GaN calculated in this current work. In this way, it is aimed to present an assessment on the polaritonic behaviors of these three very common NFRT system component, and to highlight these materials performance in spectrally selective potential nano-optics applications. For this purpose, the polaritonic figure of merit expressions as developed by (Ordonez-Miranda et al. 2017) have been used. By analytically solving these relations, both the energy (phonon thermal conductivity) and the phonon propagation and confinement aspects were highlighted along with the phonon propagation length and penetration depth results. To the best of the authors literature review, this paper is to first to
calculate and assess the polaritonic figures of merit of cBN and GaN, and compare to those of SiC.

2. Material and Methodology

2.1. Methodology

This work has an analytical approach in determining the polaritonic figure of merit of cBN, and GaN and comparing them with those of SiC (as previously presented by Ordonez-Miranda et al. (2017)) and recalculated here for comparison purposes. Ordonez-Miranda et al. (2017) derived explicit analytical formulations for polaritonic figure of merit for both material-dielectric interfaces and suspended films. Here, a material-vacuum interface is considered, since vacuum environment is crucial to realize NFRT as the dominant heat exchange mechanism. The polaritonic figure of merit \( Z \) is the product of \( Z_e \) and \( Z_p \), which respectively stand for the energy and propagation and confinement aspects, and defined as in Eq. (1) for material-vacuum interface (Ordonez-Miranda et al. 2017):

\[
Z_e = \frac{\chi^2}{2\varepsilon_i}, \ Z_p = \frac{\chi}{\varepsilon_i} (\varepsilon_i \eta_0 - \varepsilon_r \eta_+), \ Z = Z_e \cdot Z_p
\] (1)

In Eq. (1), \( \varepsilon_i \) and \( \varepsilon_r \) are respectively the imaginary and real parts of the relative permittivity of the film material. \( \chi \) and \( \eta_+ \) are defined as in Eq. (2) and (3) (Ordonez-Miranda et al. 2017):

\[
\chi = \sqrt{|\varepsilon| |\varepsilon + 1| + |\varepsilon|^2 + \varepsilon_r} \tag{2}
\]

\[
\eta_+ = \sqrt{1 + \varepsilon} \mp (1 + \varepsilon_r) \tag{3}
\]

The penetration depth (\( \delta \)) and the propagation length (\( \Lambda \)) of the surface phonon polaritons are calculated as in Eq. (4) and (5) (Ordonez-Miranda et al. 2017):

\[
\delta = \frac{1}{2Re(-k_0 \varepsilon/\sqrt{-1 - \varepsilon})} \tag{4}
\]

\[
\Lambda = \frac{1}{2Im[k_0 \sqrt{\frac{\varepsilon}{1 + \varepsilon}}]} \tag{5}
\]

In Eqs. (4) and (5), \( Re \) and \( Im \) denotes the real and imaginary parts of the corresponding parameter, while \( k_0 = \omega/c \) where \( c \) stands for the speed of light in vacuum. It is important to note that Eqs. (1-5) are here written specifically for dielectric material-vacuum interface, and their primary versions applicable to different configurations/cases are provided by Ordonez-Miranda et al. (2017).

The materials studied in this work (SiC, GaN, cBN) are known for their phononic behaviors. Polar materials support surface phonon polaritons in the IR and THz range, while exhibiting lower optical losses compared to plasmonic materials (i.e., metals) (Caldwell et al. 2015). Hence, realizing the behavior and performance of these common phononic materials is of importance for applications involving sub-diffraction confinement (Caldwell et al. 2015) and photon tunneling in the IR and THz spectral ranges.

2.2. Materials Studied in this Work

Plasmonic materials (e.g., metals) have had an important place in nano-photonics field. On the other hand, the optical losses associated with them limit their performances (Caldwell et al. 2015). Polar materials are known to support surface phonon polaritons (Narayanaswamy and Chen 2003) in the IR and THz range (Huber et al. 2008), and silicon carbide (SiC), cubic boron nitride (cBN), and gallium nitride (GaN) are among the leading phononic materials. Generation of surface phonon polaritons have found potential in microscopy and thermal emission applications, and phononic materials mostly exhibit high-temperature stability, and tunable polaritonic properties (Huber et al. 2008).

SiC is a wide bandgap semiconductor with high temperature stability (with a melting temperature >3000 K), high rigidity, high chemical and physical resistance (Huseynov 2020), and high breakdown field (Wondrak et al. 2001). GaN, similar to SiC, has high breakdown voltage and high electron velocity,
and is considered suitable in high power and high frequency applications (Mishra et al. 2008). GaN, as a nitride-semiconductor is also suitable for high temperature applications due to its wide energy bandgap and low carrier concentration in its undoped form (Pearson et al. 2002). For high-power electronics, the operating temperatures for 4H-SiC, 6H-SiC, and cBN were reported to be around 1230 K, 1200 K, and 1250 K, respectively (Lebedev and Chelnokov 1999). Apart from their individual usage, SiC and GaN can be paired for heteroepitaxial growth of GaN on SiC substrates (Talwar 2004). Cubic BN is the second hardest material and known to exhibit good thermal and chemical stability (Sharma et al. 2020). Bulk cBN crystals can also be doped to prepare p-type and n-type cBN (Litvinov et al. 1998), to fine-tune its properties. Due to high hardness of cBN, machining processes has been one of its prominent applications. These features combined with their phononic character makes SiC, GaN, and cBN highly worthy of investigation for nano-optics, nano-sensing, and nano-energy applications.

The frequency-dependent dielectric function of the materials is defined by the Drude-Lorentz Model, as in Eq. (6):

\[
\varepsilon_r(\omega) = \varepsilon_\infty \left( \frac{\omega^2 - \omega_{LO}^2 + i\Gamma \omega}{\omega^2 - \omega_{TO}^2 + i\Gamma \omega} \right)
\]

In Eq. (6), \(\omega\) stands for the angular frequency, while \(\varepsilon_\infty\), \(\omega_{LO}\), \(\omega_{TO}\), and \(\Gamma\) stand respectively for high frequency dielectric constant, longitudinal optical phonon resonance frequency, transverse optical phonon frequency, and damping factor. Values of these optical constants are provided in Table 1.

### Table 1. Optical constants of the studied materials.

| Material | \(\varepsilon_\infty\) | \(\omega_{LO}\) (rad/s) | \(\omega_{TO}\) (rad/s) | \(\Gamma\) (rad/s) |
|----------|------------------------|------------------------|------------------------|-------------------|
| SiC*     | 6.26                   | 1.849x10^{14}          | 1.500x10^{14}          | 0.56x10^{13}      |
| GaN**    | 5.3                    | 1.41x10^{14}           | 1.06x10^{14}           | 0.0152x10^{14}    |
| cBN***   | 4.46                   | 2.4551x10^{14}         | 1.9887x10^{14}         | 9.9512x10^{11}    |

Constants are taken from *Ordonez-Miranda et al. (2017), for 295 K, **Didari et al. (2018b), and ***Narayanaswamy and Chen (2003).

### 3. Results and Discussions

The overall polaritonic figure of merit (Z) calculated as in Eq. (1) for SiC, cBN, and GaN is plotted with respect to the angular frequency in Figure 2 (a). Figure 2 (b) is an alternate but more detailed version of the plot on the left (part (a)) by setting the y-axis in log-scale. It is seen that each of the materials show a single peak in their Z spectrum at a certain frequency. This peak is located at 1.501x10^{14} rad/s for SiC, 1.063x10^{14} rad/s for GaN, and 1.991x10^{14} rad/s for cBN. These peak locations correspond to locations very near to these materials’ transverse optical phonon frequencies, i.e., \(\omega_{TO}\). As discussed by (Ordonez-Miranda et al. 2017) for SiC, at around \(\omega_{TO}\) of the materials, surface phonon polaritons carry more energy. In addition, the dips in their Z spectrum (in Figure 2(b)) are located very near to the materials’ \(\omega_{LO}\), such that the minima of Z appeared at 1.829x10^{14} rad/s for SiC, 1.387x10^{14} rad/s for GaN, and 2.417x10^{14} rad/s for cBN.

When the overall polaritonic figure of merit peak magnitudes are compared, it is seen that SiC has the highest, followed by cBN and then GaN. In comparison, the peak magnitude of Z value of SiC is more than 4 times greater than that of cBN, and more than 17 times greater than that of GaN.
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Figure 2. (a) The overall polaritonic figure of merit ($Z$) of SiC, hBN, GaN, and cBN along the material-vacuum interface, (b) the plot in part (a) with y-axis in log-scale.

One of the components of the overall polaritonic figure of merit ($Z$) is the $Z_e$, which relates to the capacity of the structure to enhance the polariton thermal conductivity (Ordonez-Miranda et al. 2017). $Z_e$ is hence considered as the energy aspect of the overall polaritonic figure of merit. Figure 3 (a) shows the $Z_e$ trends for SiC, GaN, and cBN. The $Z_e$ values exhibit a minimum at 1.828x10$^{14}$ rad/s at SiC, 1.387x10$^{14}$ rad/s for GaN, and 2.412x10$^{14}$ rad/s for cBN. These angular frequency values point out to another material-specific quantity, resonant frequency ($\omega_{res}$), defined as in Eq. (7) for a single polar crystal-vacuum interface (Francoeur 2010):

$$\omega_{res} \approx \sqrt{\frac{\varepsilon_\infty \omega_{LO}^2 + \omega_{TO}^2}{\varepsilon_\infty + 1}}$$

With the light of the permittivity data given in Table 1 with Eq. (7), the $\omega_{res}$ values for SiC, GaN, and cBN are calculated respectively as 1.8049x10$^{14}$ rad/s, 1.3605 x10$^{14}$ rad/s, and 2.3765 x10$^{14}$ rad/s.

The results presented in Figure 3 (a) can be interpreted along with the polariton propagation lengths, which relate to the ability of the polaritons to travel long distances (Ordonez-Miranda et al. 2017).
As shown in Figure 3 (b) the polariton propagation lengths are at their minimum around the $\omega_{\text{res}}$ calculated for SiC, GaN, and cBN for single material-vacuum interface. This points out to the fact that when polaritons propagation length is small, they are highly confined at around these frequencies. Strong confinement of surface polaritons comes with high momentum (Dubrovkin et al. 2020), and they reflect their evanescent decay (as in Figure 1) and highly localized nature.

The other component of the overall polaritonic figure of merit is $Z_p$, which relates to the ability of the structure (here, the SiC-, GaN-, and cBN-vacuum interface) to support polaritons’ longitudinal propagation and transverse confinement. Figure 4 (a) depicts the $Z_p$ trends for the studied material-vacuum interfaces. When Figure 4(b), which is the detailed version of Figure 4(a) is examined, it is seen that the peaks of $Z_p$ are located very near the $\omega_{\text{TO}}$ of these materials, similar to the findings for overall polaritonic figure of merit, Z (in Figure 2), as also reported previously by (Ordonez-Miranda et al. 2017) for SiC. While this is the behavior with the $Z_p$ maxima, the dips of $Z_p$ are located at 1.829x10^{14} rad/s, 1.391x10^{14} rad/s and 2.415x10^{14} rad/s for SiC, GaN, and cBN, all of which are near the materials’ $\omega_{\text{res}}$ values.

These results can be better interpreted considering the penetration depth results in Figure 4 (b), which shows two dips in the penetration depth results in the studied spectral range, for SiC, GaN, and cBN. Examination of Figure 4 (b) reveals that the first and second dips as marked respectively with horizontal dashed and continuous lines are located at around $\omega_{\text{res}}$ and $\omega_{\text{TO}}$ values of the studied materials.

In magnitude, both the overall polaritonic figure of merit and the $Z_p$ values are the highest for SiC-vacuum interface, followed by cBN-vacuum interface, and GaN-vacuum interface, respectively. The spectral nature of the polaritonic behavior for these settings reveal the importance of the material selection by the peaks and dips in the polaritonic figure of merits’ spectral dependence on $\omega_{\text{res}}$, $\omega_{\text{TO}}$, and $\omega_{\text{LO}}$ values of the polar materials studied.
Figure 4. (a) The $Z_p$ component of overall polaritonic figure of merit for SiC, GaN, and cBN along the material-vacuum interface, (b) the plot in part (a) with $y$-axis in log-scale, (c) the phonon penetration depths at the SiC-, GaN-, and cBN-vacuum interface.

4. Conclusions

In this work, the polaritonic figures of merit derived by (Ordonez-Miranda et al. 2017) are solved for single material (SiC, cBN, or GaN)-vacuum interface systems. Solutions for SiC at two different temperatures (i.e., 295 K and 686 K) were previously presented by (Ordonez-Miranda et al. 2017). In this current work, the permittivity data of SiC given for 295 K by (Ordonez-Miranda et al. 2017) was utilized in recalculting single interface (SiC-vacuum) results for the figures of merit. The calculations are carried out for cBN and GaN this time for $Z_\alpha$, $Z_p$, $Z$, propagation length, and penetration depth. The findings for SiC are then compared to the results for cBN and GaN.

SiC, GaN and BN (both hexagonal and cubic BN) have been commonly investigated in NFRT systems, due to their support of surface phonon polaritons in the IR, leading to very high LDOS and spectral radiative fluxes by orders of magnitude greater than the blackbody limit at a given emitter temperature, at around the materials resonance frequency. This fact can help development of energy-efficient systems that are needed today and in the future. To the best of the authors knowledge, this work is the first to present polaritonic behavior of GaN and cBN, and compare them to those of SiC. Results showed that both $Z_p$ and $Z$ exhibited the highest peak magnitude for SiC, followed by cBN and then GaN, at around the resonance frequencies of these materials. The overall polaritonic figure of merit of SiC is greater.
than that of cBN by more than 4 times. The peak magnitudes of the overall figure of merit (Z) for SiC, cBN, and GaN are calculated respectively as $3.527 \times 10^7$, $7.368 \times 10^6$, and $2.061 \times 10^6$, centered around very near these materials’ $\omega_{TO}$. The dips in the Z spectrum are also material-specific, as they are located very near to the materials’ $\omega_{LO}$. When it comes to $Z_p$ component of the overall polaritonic figure of merit, Z, it is seen that the peak magnitude for SiC, cBN, and GaN are $4.086 \times 10^4$, $1.597 \times 10^4$, and $7.452 \times 10^3$, centered around very near these materials’ $\omega_{TO}$. The dips in the $Z_p$ spectrum are also material-specific, as they are located very near to the materials’ $\omega_{res}$.

This single interface (dielectric material-vacuum) system setting is fundamentally important as it involves one material and vacuum, fully reflecting the impact of the materials nature on the polaritonic performance. This is because polaritonic behavior of textured/corrugated surfaces and multi-layers (Didari et al. 2018a) implicitly include the effects of presence of such thin layers and nano-elements on the surfaces, their thicknesses, shapes, etc., in addition to the sole effect of the material. It is therefore important to understand the sole influence of the materials themselves in this fundamental configuration to develop new and more efficient systems. The results indicate that these merits are strongly influenced by the material type (and hence the materials dielectric properties) such that the performance merits’ maximum and minimum points are located very near the $\omega_{res}$, $\omega_{TO}$ and $\omega_{LO}$ values of the materials. Considering the fact that these semiconductors can be doped/mechanically modified to fine-tune their properties, and the technological nano-patterning opportunities, the material(s) can be selected with the desired polaritonic behavior, including optimum polariton propagation and confinement, for their potential use in opto-electronics, sensing, characterizing, and harvesting nano-devices. Such polaritonic performance evaluations of materials (both plasmonic and phononic) are expected to provide solid insight on development of systems including those harvest energy (e.g., NFRT), utilize surface polariton (plasmon, phonon, hybrid) resonance.

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