Influence of increasing SnO₂ content on the mechanical, optical, and gamma-ray shielding characteristics of a lithium zinc borate glass system

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A series of six samples were prepared based on the chemical composition of 65B₂O₃ + 20ZnO + (15-x)LiF + xSnO₂ (where x = 0, 0.25, 0.5, 0.75, 1, and 1.25 mol%) to study the role of SnO₂ on enhancing the optical and radiation attenuation capacity of the prepared glasses. The preparation of the glass series was performed using the melt quenching method at 1100 °C for 60 min. The density of the fabricated samples was measured using an MH-300A densimeter. The optical parameters of the fabricated glasses were calculated based on the spectrum recorded by a Cary 5000 UV–Vis–NIR double beam spectrophotometer in a wavelength range of 200 to 3000 nm. Furthermore, Monte Carlo simulation code and the XCOM online database were used to estimate the gamma-ray shielding capacity of the fabricated samples from 0.244 to 2.506 MeV. The results show enhanced gamma-ray shielding capacity due to the replacement of LiF by SnO₂. The linear attenuation coefficient at 0.244 MeV was enhanced from 0.352 to 0.389 cm⁻¹. The half-value thickness of the investigated glasses decreased from 1.967 to 1.784 cm when the increasing addition of SnO₂ from 0 to 1.25 mol%.

It is well understood that adding B₂O₃ to glass modifiers easily results in the formation of glass, and the resulting glass has several distinct physical and chemical properties, such as allowing visible light to pass through the glass and thus the possibility of using this type of glass in various optical applications. Borate-based glasses have notable characteristics, such as their high thermal stability, high chemical durability, low melting temperature, good mechanical stability and ease of fabrication. Furthermore, some heavy metal oxides are utilized to increase the chemical durability of these glasses and to lower their high phonon energy1,2. When ZnO is added to borate glasses, it plays dual roles, namely, a network former and network modifier; thus, ZnO enhances the physical characteristics of the host glass system, such as its thermal stability, and reduces the melting temperatures of the batches during fabrication.

Additionally, SnO₂ is a technologically important material because it has a wide range of applications, including in photocatalytic systems, gas sensors, and photovoltaic cells. Moreover, when SnO₂ is added to B₂O₃, a comparatively high density is observed, which promotes its use as an efficient shielding material. From the nineteenth century until the present, various radioactive isotopes emitting ionizing radiation have been widely employed in nuclear medicine, dentistry, agriculture, industry, sterilization, food, and other disciplines that are difficult to list3–5.

Borate-based glass systems are largely studied in the literature since they have a wide range of applications in different fields, such as medical treatment and research, industrial technology, and engineering. These applications, especially the medical applications, use ionizing radiation. Workers in these areas are exposed to different...
types of radiation at different doses and varying energies; thus, they may experience health problems, such as cancers.

As a result, there is an urgent need to develop protective radiation shields to safeguard nuclear power plant workers, medical professionals, and patients in radiotherapy facilities from the dangers of radiation.

Concrete and lead have historically been the most extensively utilized barriers for radiation protection since they have high photon attenuation ability. As a result, they are widely utilized as a protective barrier for gamma radiation and X-rays in a variety of forms, such as plates, tubes and bricks. However, concrete and lead are opaque to visible light; thus, they prevent light from passing through.

To overcome the opaqueness problem of these two materials, researchers have resorted to developing new types of radiation-protective materials using different types of glass because glass has the ability to transmit light. Recently, several research groups have developed multiple glass systems and studied the ability of these glasses to attenuate radiation and thus reduce its harmful effects on humans and the environment.

When studying radiation attenuation qualities, researchers identify fundamental quantities, such as the mass attenuation coefficient and some associated numbers. This amount may be calculated in a variety of ways, the most significant of which is a practical approach that involves exposing the sample to photons. A detector can be used to determine the number of photons that have penetrated the sample.

The Monte Carlo simulation approach is another option. This approach involves creating a virtual reality for the practical experiment taking place in the laboratory. This approach differs from the practical way in that it saves time and effort while also lowering the likelihood of individuals becoming exposed to photons. Its significance is heightened by the shortage of equipment and facilities, forcing more researchers to rely on simulation. In the last few years, several researchers have adopted Monte Carlo simulations to investigate the radiation attenuation performance of borate glasses with different additives. For example, Mahmoud et al. used MCNP5 to examine the influence of CdO on the attenuation shielding performance of alkali borate glasses. They found that the addition of more CdO enhanced the radiation attenuation performance of the selected glasses. Acikgoz et al. prepared new alumina borate glasses, and when using MCNP5, they determined the role of CeO2 and Er2O3 on the radiation shielding performance of the prepared glasses. The authors concluded that glasses that contained Er2O3 had a higher mass attenuation coefficient than glasses with CeO2.

Rammah et al. studied the impact of Bi2O3 and PbO on the radiation attenuation parameters of borate glasses. They used Monte Carlo simulation to investigate the attenuation factors at different energies ranging from 15 keV up to 15 MeV. The authors reported that the addition of PbO to borate glass enhanced the protective features of the samples to nuclear radiation. Stalin et al. prepared new lithium-borate glasses using the quenching method and simulated the radiation attenuation factors via the MCNP5 code. The authors examined the influence of replacing Bi2O3 with WO3 on the half-value layer and other related parameters. According to their results, the glass that contained 60 mol% Bi2O3 and 20 mol% Li2O and Bi2O3 had the smallest half-value layer. Asadi and Hosseini used the MCNP code to simulate the radiation attenuation coefficients of borate glasses doped with ZnO, Li2O and Bi2O3 in the energy range of 200 keV to 1.5 MeV. According to their results, glass with a composition of 20ZnO-50Bi2O3-15Li2O-15B2O3 had the shortest mean free path. Abou Hussein et al. simulated the radiation attenuation coefficients of borate glasses doped with one of the following oxides: V2O5, Cr2O3, Fe2O3 and TiO2. The authors concluded that a 3 cm-thick glass sample containing Cr2O3 was the best attenuator and could reduce the lifetime risk to cancer by five times.

In continuation of the previous efforts made by researchers to study the radioactive attenuation properties of borate glasses using Monte Carlo simulation, enrich the scientific community with more studies on the radiation attenuation properties of this kind of glass, and check the possibility of developing new shielding glasses based on borate glass, the Monte Carlo simulation code is utilized to estimate the gamma-ray shielding capacity of the novel lithium-zinc borate glass system. Moreover, to understand the optical properties of the developed glasses, the optical absorption band was measured in the wavelength range of 200 to 3000 nm. Additionally, the Matsuoka-Makimenzie model was applied to predict the elastic moduli and microhardness of the fabricated glasses.

### Experimental techniques

**Glass fabrication, characterization, and optical properties.** Tin-doped lithium zinc borate glass samples with nominal compositions of 65B2O3 + 20ZnO + (15-x)LiF + xSnO2 (where x = 0, 0.25, 0.5, 0.75, 1 and 1.25 mol%) were prepared. The glass composition and chemicals used to prepare samples along with glass code are shown in Table 1. Approximately 20 g of the batch composition of these chemicals was mixed well in an agate

| Sample | Chemical composition mol% | Density (gm/cm³) | Molecular weight (gm/mol) | Molar volume (cm³/mol) |
|--------|---------------------------|------------------|---------------------------|-----------------------|
| BZLSn0 | 65 20 15 0                | 3.025            | 65.426                    | 21.628                |
| BZLSn1 | 65 20 14.75 0.25          | 3.052            | 65.738                    | 21.539                |
| BZLSn2 | 65 20 14.5 0.5            | 3.083            | 66.049                    | 21.424                |
| BZLSn3 | 65 20 14.25 0.75          | 3.131            | 66.362                    | 21.195                |
| BZLSn4 | 65 20 14 1               | 3.199            | 66.674                    | 20.842                |
| BZLSn5 | 65 20 13.75 1.25         | 3.265            | 66.986                    | 20.516                |

**Table 1.** Chemical composition and density of the prepared BZLSn glasses.
mortar for 60 min and then preheated at 350 °C for 60 min. This homogenous mixture was then heated in a porcelain crucible at 1100 °C for 30 min before being cast onto a stainless steel mold to generate glass discs. After quenching, the samples were immediately transported to a muffle furnace set at 300 °C for annealing.

The densities of the glasses were measured using the MH-300A densimeter according to Archimedes’ rule, as shown in Eq. (1). Toluene was used at room temperature as an immersion liquid during the density measurements.

\[ \rho_{\text{glass}} = \frac{W_a}{W_a - W_b} \rho_{\text{tolune}} \]  

where \( W_a \) and \( W_b \) are the weights of the glass sample in air and in liquid toluene, respectively. The measured density, as well as the calculated molecular weight and molar volume, are listed in Table 1 for the fabricated BZLSn glasses.

The optical absorption data of the glass samples were obtained for the polished glass samples using a Cary 5000 UV–Vis–NIR double beam spectrophotometer.

**Gamma-ray shielding capacity evaluations.** Both Monte Carlo simulation and the XCOM theoretical program were used to predict and affirm the photon shielding capacity of current BZLSn glasses. The calculation using the XCOM program depended only on the chemical composition of the fabricated glass system. After that, the calculation for the attenuation coefficient depended on the mixture, as presented in Eq. (2).

\[ \mu_m \left( \frac{\text{cm}^2}{\text{g}} \right) = \sum_i \omega_i \times (\mu_m)_i \]  

where \( \omega_i \) is the fractional abundance of the \( i \)th element in the shielding material.

On the other hand, the Monte Carlo simulation code is based on many parameters, such as the sample composition, sample density, problem geometry, and ENDF/B-VIII library, from which the cross-section of interaction is extracted for various elements. The Monte Carlo simulation geometry consists of many cards that describe all components required for the simulation process, as illustrated in Fig. 1. The first important card is the cell card, which introduces the cell type, density, and boundaries. The mentioned figure shows many cells inside the prepared geometry, such as the collimators of lead with a density of 11.34 g/cm³, glass samples with their densities listed in Table 1, a radioactive source, and F4 detector tally to estimate the photons average track length per unit cell of the detector. The second is the surface card, which usually describes the shape and dimensions of the cells introduced in the first step. The dimensions of the various geometric components are illustrated in the figure. The third is the material card that introduces chemical composition of each cell in the arranged geometry. The source card also describes the source position, type, energy, decay probability, and emission distribution. The cutoff card is one of the physical cards used in the prepared geometry, and it is set to stop interaction after 106 historical cards.

**Results and discussion**

**Physical and optical properties.** The density and molar volume of the 65B₂O₃ + 20ZnO + (15-x) LiF + xSnO₂ glass system with various SnO₂ concentrations are shown in Fig. 2. The density of the sample improves as the SnO₂ content (x) is increased; on the other hand, the molar volume decreases. The density of the fabricated glasses slightly increase from 3.025 ± 0.151 to 3.265 ± 0.163 g/cm³, increasing the SnO₂ insertion ratio.
from 0 to 1.25 mol%. The substitution of Li\(^{2+}\) ions (\(\rho = 0.534 \text{ g/cm}^3\)) with Sn\(^{2+}\) ions (\(\rho = 7.31 \text{ g/cm}^3\)) causes a slight increase in the density of the fabricated glass. Additionally, the increase in the fabricated glass density may be related to a significant decrease in the interatomic distance due to the substitution of LiF by TeO\(_2\).

Figure 3 shows the UV–Vis absorption spectra of the produced SnO\(_2\)-doped lithium zinc borate glass samples, and the inset in Fig. 3 shows the absorption spectra from 350 to 800 nm. These measurements were carried out on samples with a thickness of 2.2–2.5 mm. The spectrum shows that the produced samples have significant UV absorption and that the strength of the absorption decreases as the concentration of SnO\(_2\) is increased\(^{27}\). As the SnO\(_2\) substitution ratio is increased, distinctive bands of SnO\(_2\) are observable in these glass samples: the peak at 445 nm (*) is attributable to a ligand in the metal charge assigning transition (Sn\(^{4+}\)).

Tauc’s rule, which Mott and Davis\(^{28,29}\) have modified, is shown in Eq. (3). This rule is used to estimate the optical energy band gaps for the investigated SnO\(_2\)-doped lithium zinc borate glass samples in this study:

\[
(\alpha(v)h\nu)^m = C(h\nu - E_g)
\]  

\(\alpha(v)\) is the absorbance, \(h\nu\) is the photon energy, \(E_g\) is the optical band gap energy, \(C\) is a constant, and \(m\) is a constant related to the type of transition. The value of \(m\) is 2 for direct allowed transitions and 0.5 for indirect allowed transitions.
where \( m \) represents the kind of electronic transition. In this case, \( m = 2 \) indicates a direct allowable transition and \( m = 0.5 \) indicates an indirect allowable transition.

Figures 4 and 5 show the dependence of \((\alpha h\nu)^2\) and \((\alpha h\nu)^{1/2}\) on \((h\nu)\) for the direct and indirect allowed transitions of the examined samples. Table 2 summarizes the \( E_g^{\text{optical}} \) values of both the direct and indirect transitions.

### Table 2. Optical parameters of the BZLSn glasses.

| Sample   | Direct | Indirect | \( E_g^{\text{optical}} \) | \( n \) | \( E_g \) (eV) |
|----------|--------|----------|-------------------|-------|-------------|
| BZLSn0   | 3.257  | 2.902    | 3.079             | 2.376 | 0.251       |
| BZLSn1   | 3.230  | 2.855    | 3.043             | 2.386 | 0.234       |
| BZLSn2   | 3.202  | 2.825    | 3.014             | 2.393 | 0.219       |
| BZLSn3   | 3.201  | 2.820    | 3.011             | 2.394 | 0.214       |
| BZLSn4   | 3.198  | 2.817    | 3.008             | 2.395 | 0.211       |
| BZLSn5   | 3.197  | 2.815    | 3.006             | 2.395 | 0.208       |

Figure 4. Dependence of \((\alpha h\nu)^2\) on \((h\nu)\) for the direct allowed transition of the examined glass samples.

Figure 5. Dependence of \((\alpha h\nu)^{1/2}\) on \((h\nu)\) for the indirect allowed transition of the examined BZLSn samples.
of undoped and doped SnO₂ glass samples. The $E_g$ optical values decrease during the direct and indirect transitions when the SnO₂ concentration in the produced glasses is increased.

The difference between the Eg of SnO₂ (3.6 eV) and the large band gap of LiF (14 eV) may be related to the decreasing optical energy values $E_g$ with increasing the content of SnO₂.

Equation (4) is used to compute the refractive index values ($n$) of the manufactured samples during the direct and indirect allowed transitions:

$$\left(\frac{n^2 - 1}{n^2 + 2}\right) = 1 - \frac{E_{optical}}{20}$$

Table 2 lists the values of ($n$) for all synthesized samples. Clearly, the optical energy band gaps are inversely related to the refractive index ($n$) of all the manufactured glasses. The refractive index increases from 2.376 to 2.395, indicating that the proposed glasses may be used as a promising material for optical filters and photonic applications.

The absorption coefficient ($\alpha$) can be estimated using Urbach's empirical formula:

$$\alpha = \alpha_0 \exp \left(\frac{h\nu}{E_u}\right)$$

where $\alpha_0$ is a constant and $E_u$ is the Urbach energy. Equation (5) can be written as:

$$\ln(\alpha) = \ln(\alpha_0) + \left(\frac{h\nu}{E_u}\right)$$

As a consequence, the slope of the straight line obtained by plotting $\ln(\alpha)$ vs. $(h\nu)$ may be used to calculate Urbach's energy ($E_u$). The change in $ln(\alpha)$ with photon energy $(h\nu)$ is exhibited in Fig. 6. As listed in Table 2, the Urbach energies decrease as the SnO₂ concentration is increased, indicating an increase in stability and decrease in disorder in the glass samples.

**Mechanical properties of the fabricated samples.** The effect of replacing LiF with SnO₂ on the dissociation energy ($G_t$) and packing factor ($V_t$) of the fabricated glasses was studied. Figure 7 depicts a linear decrease in $G_t$ as well as an increase in $V_t$ with an increasing replacement ratio of LiF by SnO₂. The $G_t$ values decrease slightly from 69.620 to 69.613 kJ/cm³, while the $V_t$ values increase from 0.732 to 0.777 when the SnO₂ insertion ratio is increased from 0.0 to 1.25 mol%, respectively. The mentioned decrease in the $G_t$ values is due to the replacement of the LiF compounds with $G_t = 38.801$ kJ/cm³ by SnO₂ with a comparable $G_t$ value where $(G_t)_{SnO2} = 38.264$ kJ/cm³. The increase observed in the $V_t$ values is due to the replacement of a low packing factor $V_t$ compound with a compound having a higher one; for instance, the $V_t$ of LiF is 4.894 cm³/mol while it is 13.862 cm³/mol for SnO₂.

Figure 8 shows the effect of replacing LiF with SnO₂ on the elastic moduli of the fabricated glasses. It is known that the elastic moduli are related and proven from the predicted values of both $G_t$ and $V_t$ presented in the previous section. According to the results presented in Fig. 8 A, the Young modulus increases from 101.949 GPa to 108.226 GPa when the SnO₂ insertion ratio is increased from 0.0 to 1.25 mol%, respectively. This increase in the Young modulus is directly related to the increase achieved for the $V_t$ of the fabricated glasses when LiF is replaced by SnO₂, where $Y = 2V_tG_t$. Additionally, Fig. 8B,C,D show that the bulk (B), shear (S), and longitudinal
(L) moduli follow the same trend as the Young modulus. The bulk modulus increases from 89.574 to 100.953 GPa, the shear modulus (S) increases from 38.903 to 40.953, and the longitudinal modulus exhibits the greatest values, increasing from 141.444 to 147.030 GPa when the SnO₂ ratio is increased from 0.0 to 1.25 mol%.

The Poisson ratio (σ) is a mechanical parameter that measures the ratio of glass deformation in a direction perpendicular to the loading direction. According to the theoretical calculations, the Poisson ratio is based on the predicted values of the Young (Y) and Shear moduli, where $\sigma = \frac{E}{2S} - 1$. In the present work, as illustrated in Fig. 9, the σ values increase from 0.310 to 0.321 when the SnO₂ ratio is increased from 0.0 and 1.25 mol%.

The microhardness (H, GPa) is a factor used to describe the hardness of the fabricated glass samples at the microscale, and it is proof of the Poisson ratio and Vₗ. The H calculations show that the H values slightly decrease from 4.920 to 4.878 GPa when the SnO₂ is increased from 0 to 1.25 mol% in the fabricated samples. This slight decrease is related the fabricated glasses’ Poisson ratio and Vₗ.

Figure 7. Dissociation energy (Gₚ, kJ/cm³) and packing density (Vₗ) versus the increasing Sn₂O ratio.

Figure 8. Effect of increasing SnO₂ on the elastic moduli of the fabricated glass system.
Gamma-ray shielding properties. The shielding capacity of SnO$_2$-based zinc lithium borate glasses was evaluated using Monte Carlo simulation$^{35}$ and the XCOM$^{36}$ theoretical program between 0.2234 and 2.506 MeV. First, the obtained average track length (ATL) of the gamma photons was evaluated based on the MC simulation. Then, the ATL was transferred to the linear attenuation coefficient (LAC, cm$^{-1}$) using Lambert–Beer’s law. The MC evaluated results were verified using the XCOM program at the same energies as the MC simulation.

Figure 10 illustrates that the LAC of the fabricated SnO$_2$-based borate glasses suffers a moderately exponential decrease as the energy is increased from 0.2234 to 2.506 MeV. This exponential decrease is due to Compton scattering (CS), representing the primary interaction in the selected energy region. The cross-section corresponding to CS decreases linearly with increasing photon energy, where $\sigma_{\text{CS}} \propto E^{-1}$.$^{37}$ When the photon energy is varied from 0.244 to 2.506 MeV, the LAC decreases in the following intervals: 0.352–0.116, 0.357–0.117, 0.362–0.118, 0.369–0.120, 0.379–0.122, and 0.388–0.125 cm$^{-1}$ for samples BZLSn0, BZLSn1, BZLSn2, BZLSn3, BZLSn4, and BZLSn5, respectively. The increase in the energy of the incident photon increases the frequency and penetration power of the incident photon. Thus, the photon cross-section of the interaction decreases. Therefore, the decreases in the interaction probability are associated with the significant decreases in the linear and mass attenuation values.

The effect of the added SnO$_2$ appears clearly in Fig. 11, where the LACs of the BZLSn glasses increase with increasing replacement of LiF by SnO$_2$. At a gamma energy of approximately 0.662 MeV, the LAC increases in the order of 0.226, 0.228, 0.230, 0.234, 0.239, and 0.244 cm$^{-1}$ when the amount of SnO$_2$ added into the glass
matrix is 0, 1, 2, 3, 4, and 5 mol%, respectively. This increase in the LACs is caused by the replacement of LiF (with a lower density and absorption cross-section ($\rho_{\text{LiF}} = 2.64 \text{ g/cm}^3$)) with SnO$_2$ (with a higher density and photon absorption cross-section($\rho_{\text{SnO}_2} = 6.95 \text{ g/cm}^3$)). As a result of the replacement of a light compound (LiF) with a dense compound, the probability of the interaction between the incident photons and the distributed atoms and electrons increases due to the increase in the interaction cross-section. This increase in the interaction cross-section is reflected in the LACs and MACs, where both the LAC and MAC increase with increasing SnO$_2$ ratio in the fabricated samples. The enhancement of the LACs and MACs is positively reflected in other shielding parameters, as illustrated in Fig. 11. The thickness required to decrease the source radioactivity to half (half-value layer, HVL) is enhanced by adding SnO$_2$ into the glass matrix. The HVL regularly decreases with increasing LiF substituted by SnO$_2$. Among the current study, thinner HVLs are achieved at 0.223 MeV. The HVL decreases from 1.967 to 1.784 cm when the SnO$_2$ concentration is increased from 0 and 5 mol%. On the other hand, the thicker layers achieved at 2.506 MeV decrease in the following order: 5.975, 5.923, 5.863, 5.773, 5.652, and 5.353 cm for the prepared BZLSn0, BZLSn1, BZLSn2, BZLSn3, BZLSn4, and BZLSn5 samples, respectively. Additionally, Fig. 11 shows that, at a gamma-ray energy of 0.662 MeV, the HVL decreases from 3.069 to 2.845 cm as the SnO$_2$ ratio is increased from 0 to 5 mol%. For all fabricated BZLS glasses, the HVL reduction is due to the LAC being enhanced with increasing SnO$_2$ addition, where HVL = 0.693/LAC.

Figure 11. Linear attenuation coefficient (LAC, cm$^{-1}$) and half-value layer (HVL, cm) with respect to the SnO$_2$ concentration at 0.662 MeV.
of photons inside the shielding layers. Therefore, the replacement of LiF with SnO₂ reduces the transmission of photons through the BZLSn glasses. The small variation in the TFs among the glass samples is related to the small amounts of LiF displaced by SnO₂, where the highest SnO₂ concentration is only 5 mol%.

The radiation protection efficiency (RPE, %) is used to describe the energy deposited inside the fabricated BZLSn glasses. Figures 14a and b depict the RPE variation with respect to the incident gamma photon energy and glass thickness. In the case of Fig. 14a, the RPE shows a linearly decreasing trend with increasing energy from 0.244 to 2.506 MeV. Among the studied energies, the best RPE is recorded at 0.244 MeV. Notably, the RPE varies, with values of 65.261, 65.743, 66.999, 67.936, 68.833% when the content of SnO₂ is 0, 0.25, 0.5, 0.75, 1.0, and 1.25 mol%, respectively (representing BZLSn₀, BZLSn₁, BZLSn₂, BZLSn₃, BZLSn₄, and BZLSn₅). These results indicate that at a gamma-ray energy of 0.244 MeV, the photons undergo many collisions with the glass atoms and lose a very large amount of their energy inside the glass layer. Thus, only a small number of photons can penetrate the glass and reach exposed workers. Then, with increasing gamma photon energy, the photon penetration power increases, the number of photon collisions decreases, and the energy transferred from photons to the glass layer decreases with increasing number of photons that can penetrate the glass layer. Hence, the number of photons reaching workers increases with a decrease in the glass thickness (i.e., RPE). Among the
studied energies, the lowest RPE is recorded at 2.506 MeV, and the values vary in the order of 29.390, 29.607, 29.858, 30.248, 30.780, and 31.317% for BZLSn0, BZLSn1, BZLSn2, BZLSn3, BZLSn4, and BZLSn5, respectively.

The glass thickness also greatly affects the protection capacity of the shielding material. As illustrated in Fig. 14b, the RPE increases linearly with increasing BZLSn glass thickness from 1.5 to 10 cm. Among the studied glass thicknesses, the lowest RPE is obtained at a thickness of 1.5 cm from the fabricated BZLSn glass samples. At 1.173 MeV, the recorded RPE for a thickness of 1.5 cm is 22.609, 22.777, 22.973, 23.281, 23.713, and 24.133 cm. In addition, the highest level of protection is obtained from the thicker samples with a thickness of 10 cm, where the RPE varies, with values of 81.889, 82.149, 82.449, 82.912, 83.544, and 84.139% for samples BZLSn0, BZLSn1, BZLSn2, BZLSn3, BZLSn4, and BZLSn5, respectively. It is clear that increasing the glass thickness causes an increase in the path length of the gamma photons inside the investigated samples, which forces the incident photons to undergo additional collisions with glass atoms and electrons. Thus, the energy deposited inside the glass layer increases, and the number of photons that can escape from the glass thickness decreases. Therefore, the RPE ratio increases.

To select the best fabricated glass that possesses suitable shielding and mechanical properties, as well as being inexpensive to fabricate, the variations of the LAC, HVL, microhardness, and fabrication cost versus the SnO2 concentrations were studied; the results are shown in Fig. 15. As mentioned in the last paragraphs, the insertion of SnO2 causes a significant increase in the µ values due to the significant increase in the packing factor Vt and bulk density associated with the replacement of LiF by SnO2. The significant increase in the packing factor causes an increase in the Poisson ratio associated with a decrease in the microhardness of the material. According to the mentioned figure, the best sample in the present study is has a SnO2 concentration of 0.75 mol% (i.e., BZLSn3). The µ value of the mentioned sample is 0.2338 cm⁻¹, and the HVL is 2.9638 cm at 0.662 MeV. Additionally, the mentioned sample contains a microhardness value of 4.9022 GPa. Furthermore, the fabrication cost for a sheet of BZLSn3 glass with dimensions of 100 cm × 100 cm, which can attenuate half of the incident photons when the energy is 0.662 MeV, is approximately $3151.46. This relatively high cost is due to LiF being expensive, which costs $0.384/g of fabricated sheet. Therefore, the replacement of LiF by SnO2 causes a significant decrease in the fabrication cost due to SnO2 having a relatively low price compared to LiF (SnO2 costs approximately $0.05/g).

**Conclusion**

In the present study, the effect of increasing SnO2 in a glass system based on zinc lithium borate was studied. The fabricated glass density increased from 3.025 and 3.265 g/cm³, while the molar volume decreased from 21.628 to 20.516 cm³/mol for the fabricated BZLSn glasses when the content of SnO2 was increased from 0 to 1.25 mol%.

According to the optical property calculations, the energy gap during direct and indirect transitions decreased from 3.257 to 3.197 eV and from 2.902 to 2.815 eV, respectively. The Monte Carlo simulation showed that BZLSn5 with a SnO2 composition of 1.25 mol% had the highest shielding capacity among the fabricated samples. The linear attenuation coefficient of the previously mentioned sample decreased from 0.389 to 0.125 cm⁻¹, and the half-value layer increased from 1.784 to 5.535 cm as the photon energy was increased from 0.244 to 2.506 MeV. The previously illustrated results showed that the replacement of LiF by 1.25 mol% SnO2 resulted in an enhanced shielding capacity, reaching 9.3% at 0.244 MeV.
Figure 15. Variation of the linear attenuation coefficient, half-value layer, manufacturing cost, and microhardness versus the SnO₂ insertion ratio.

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M.I.S.: write the main manuscript text, prepared figures A.H.A.: writing—review and editing, funding acquisition A.S.A.: write the main manuscript text, prepared the samples K.A.M.: write the main manuscript text, calculated the radiation shielding parameters using MCNP5 code.

Competing interests
The authors declare no competing interests.

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