Color management of semi-transparent nano-patterned surfaces

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Abstract. Many applications require colored surfaces while maximizing photons to pass through with no optical losses. As the perceived color depends on the spectral content and the number of reflected photons for each wavelength, it is necessary to consider not only the chromatic content but also the brightness. As a first application example, photovoltaic modules could see their acceptance in urban situations greatly improved if their color could be controlled while maximizing the number of photons used for energy conversion. In this case, it is important to limit light brightness to avoid glare when light coming from the sun is reflected directly. As another example, a colored and semi-transparent glass layer in the infrared could efficiently protect sensors while camouflaging them, such as lidar in automotive applications. We propose a theoretical study for multicolor applications, considering 2D partially etched grating waveguide structures that behave as a perfect broadband antireflection coating while the residual layer couples light in a waveguided mode in order to reflect only a particular photon band. This strategy allows to produce a large variety of colors and the use of non-absorbing materials avoids any optical loss. The width of the resonance peak is a crucial parameter for balancing brightness in multicolor applications. In the second part, we developed a method for a fast optimization of the geometrical parameters and we applied it to obtain the three primary colors with strong geometrical constraints, so that such structures could be manufactured in a single step by nano-imprint. Finally, we propose some improvements of these structures, in particular by reducing broadband reflection to obtain more saturated colors, or to improve the angular behavior of these structures. © 2021 Society of Photo-Optical Instrumentation Engineers (SPIE) [DOI: [10.1117/1.OE.60.5.055101](https://doi.org/10.1117/1.OE.60.5.055101)]

Keywords: CIE 1931 color space; photonic crystal; patterning; sRGB; brightness.

Paper 20201152 received Sep. 28, 2020; accepted for publication Apr. 6, 2021; published online May 13, 2021.

1 Introduction

Grating waveguide structures (GWSs) have been considered this last decade to obtain tiny resonant bandpass filters that can be used to control high chromaticity in the color printing field as shown in Refs. 1–3. In the indoor photovoltaic area, for example, merging elementary pixels based on GWS with different periods on top of the front glass would lead to the fabrication of colored panels without a strong degradation of the conversion efficiency. Performance items required for such filters are a broadband antireflection coating coupled to a sufficient narrow reflection resonance involving accurate perception of color, with limited color variations depending on the optical conditions as described in Ref. 4.

In this paper, we focus on a GWS that behaves both as a bandpass color filter and as a single layer anti-reflection coating (s-ARC). We will first assume that its reflected spectrum has only one peak which can be defined by its central wavelength, its width, and the level of background reflection. These three independent parameters impact the colors perceived by the human eye,

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such as brightness, saturation, and hue. In the first part, we carry out a theoretical study of these parameters in order to predict the overall shape of the reflected spectrum to produce a desired color. For multicolored applications, this part also aims to highlight the relationships to be considered between the different reflected spectra. In the second part, we focus on GWS structures and their simulation methodology and optimization to control the reflected spectrum and so the reflected color.

2 Range of Colors Accessible by Patterning the Front Glass Layer

To reduce the number of parameters to study an ARC on a glass substrate, we consider that the reflected spectrum has a Lorentzian shape centered on wavelength λ_c , with a width to mid height $\Delta\lambda_c$. In the case of a resonant mode, the maximum of the reflection is equal to 1, but a reflection on the entire visible spectrum R_0 will eventually be taken into account, leading to an amplitude of the reflected signal equal to $1 - R_0$,

$$R(\lambda) = R_0 + \frac{1 - R_0}{1 + \left[\frac{2 \cdot (\lambda - \lambda_c)}{\Delta\lambda_c} \right]^2}. \quad (1)$$

These assumptions restrict the study to only three parameters: λ_c , $\Delta\lambda_c$, and R_0 . Using this theoretical expression of $R(\lambda)$, it is possible to calculate the CIE (International Commission on Illumination) 1931 xyY components using the three color matching functions $\bar{x}$, $\bar{y}$, and $\bar{z}$ as explained in Ref. 4. However, this color space is not perceptually uniform. For this reason, we prefer Judd's Uniform Chromaticity Scale (UCS) diagram using Ref. 5,

$$\begin{cases} u = \frac{0.4661 \cdot x + 0.1593 \cdot y}{y - 0.15735 \cdot x + 0.2424} \\ v = \frac{0.6581 \cdot y}{y - 0.15735 \cdot x + 0.2424} \end{cases} \quad (2)$$

As the (u, v, Y) diagram is uniform, we can use the Euclidian distance between two colors $C_1(u_1, v_1, Y_1)$ and $C_2(u_2, v_2, Y_2)$ and calculate gamut coverage areas. Unlike other color spaces (CIE 1964 $U^*V^*W^*$, CIE 1976 $L^*a^*b^*$, or CIE 1976 $L^*u^*v^*$), it does not depend on a reference illuminant for its white point (WP): most often, it corresponds to daylight at temperature 6504K (D65) or 5003K (D50) and color coordinates are relative to the chosen WP. In addition, the chromaticity is contained in a 2D plane (as for the CIE 1931 xyY , CIE 1964 $U^*V^*W^*$, CIE 1976 $L^*u^*v^*$, and contrary to CIE 1976 $L^*a^*b^*$ and its derived color spaces), which makes color representation easier when Y stands for the corresponding brightness. Finally, we chose the model derived from Judd's work from Ref. 5 that has been historically less exploited than the simplified (for computational purposes) version of MacAdam proposed in 1937 in Ref. 6, which was adopted by CIE in 1960.

Finally, in this article, we consider that the reference spectrum I_0 is AM1.5D which represents the light emitted directly by the sun for an air mass of 1.5. As a consequence, the calculated colors correspond to an outdoor use and the resulting WP corresponds to $R_0 = 1$ at coordinates $(u = 0.3966, v = 0.4273)$.

2.1 Study of an Ideal Optical Response

In this part, we consider the ideal case $R_0 = 0$ in Eq. (1) in order to determine the optical properties for a very favorable case and for a single resonant mode. In that case, the calculated brightness is called Y_0 .

Figure 1 represents the chromaticity when the position λ_c of the Lorentzian varies for different values of $\Delta\lambda_c$. For each value of $\Delta\lambda_c$, the gamut starts toward the WP: when λ_c increases, the calculated color approaches the blue color and sweeps the rainbow colors in direction of the reddest point. Then, the curve goes back to the WP again.

Unsurprisingly, the gamut is the most extended for a narrow resonance ($\Delta\lambda_c = 1$ nm) and it gets closer to WP when the width of the peaks $\Delta\lambda_c$ increases. This behavior is a direct

Fig. 1 Different gamuts for an ideal case with no constant reflection in the visible range ($R_0 = 0$). Each gray contour line is obtained when varying the peak position λ_c for a given value of $\Delta\lambda_c$. The standard red–green–blue (sRGB) domain corresponds to the colors that can be produced by monitors.

Fig. 2 Color brightness Y_0 for the same Lorentzian functions as Fig. 1. On top of the figure, hue is given for illustration purpose.

consequence of a lower spectral selectivity of the reflected light, with a decrease of the saturation of the obtained color. A maximum coverage of 39% is obtained for the thinner peak. Last, it can be observed that the area above the purple line, containing shades of blue, magenta, and red colors, cannot be obtained even if it represents 61% of the gamut surface: for these colors, a photonic crystal (PC) with two simultaneously reflected peaks is necessary.

For each curve from Fig. 1, brightness Y_0 varies significantly with an envelope similar to that of the function $\bar{y}$ (as plotted in Ref. 4), as shown in Fig. 2. As the human eye is very sensitive to the green color, the width of the PC resonance must be increased when moving away from the green color. To obtain $Y_0 = 1$, every photon in the visible range must be reflected, leading to $\Delta\lambda_c = +\infty$ and corresponding to the WP in the gamut.

For an arbitrary target brightness $Y_0 = 0.4$, $\Delta\lambda_c$ in the 20- to 50-nm range is enough for the green color while it is necessary to consider wider than 200-nm peaks to balance with red and blue colors. According to Fig. 1, this brings the perceived color very close to the WP, with a poor saturation. On the other hand, we notice a relative symmetry around the wavelength corresponding to the green color (560 nm): to balance the brightness of a blue (473 nm) and a red surface (650 nm), we need $\Delta\lambda_c$ between 100 and 150 nm.

Fig. 3 Iso luminance curves for an ideal case with no constant reflection in the visible range ($R_0 = 0$). (a) Chromaticity diagram of with few iso luminance curves and (b) relation between the wanted color (λ_c) and the width of the reflected signal ($\Delta\lambda_c$). On top of the figure, hue is given for illustration purpose.

2.2 Maximum Brightness for a Single Mode PC

For multicolor applications, we may want to control the brightness of each color. For this purpose, we determined in Fig. 3(a) the gamut obtained for different constant values of the brightness Y_0 . This curve is very close to that obtained in Ref. 5 (Fig. 13, in the CIE 1931 xyY diagram) which gives curves of constant excitation purity, except in the bottom part of the gamut. The more we look for a significant brightness, the closer we get close to the WP by mainly reducing the extension of the achievable colors in the red and blue directions. However, if such a structure is directly illuminated by the sun, it is important not to dazzle humans. It is consequently advised not to exceed the ambient brightness, which corresponds to approximately $Y_0 = 10\%$ when we consider the diffuse level contained in the AM1.5G spectrum and AM1.5D as illuminant. The curve $Y_0 = 10\%$ shows that we can therefore theoretically reach and exceed the green and red colors of the sRGB domain, and come very close to the blue color.

Figure 3(b) represents the relationship between λ_c and $\Delta\lambda_c$ for the same Y constants as in Fig. 3(a). For wavelengths below 500 nm and greater than 650 nm, these two parameters vary almost linearly with each other. One may notice that when Y_0 increases, the blue and red colors become more and more difficult to obtain: for $Y_0 = 0.75$, shades of red, orange, yellow, and green can only be obtained for less than 200-nm-width peaks that cannot be currently fabricated with a PC since it exploits resonances. On the contrary, for very contrasting colors, a low brightness allows to get all the colors of the rainbow with a narrow single resonance. Last, this figure can easily be used as an abacus to obtain the stencils of the reflection spectra of each wanted color of a polychromatic drawing, each consisting of a PC for instance.

2.3 Impact of a Residual Reflection

In this part, we focus on the impact of $R_0 > 0$ since this constant represents a reflection on the whole visible spectrum. It can be the consequence of

- a patterned layer that does not behave as an efficient anti-reflective coating around the resonance;
- a parasitic reflection under the glass substrate (Fig. 4). For instance, considering the front glass layer constituting the encapsulation of a solar cell in its module, this constant allows to take into account the reflected photons on the solar cell which escape from the front surface of the module.

Since the brightness calculation is a scalar product (as detailed in Ref. 4), its linearity implies that the brightness Y_{R_0} can be expressed according to that (Y_0) when $R_0 = 0$ (previous paragraph),

Fig. 4 Schematic structure considering a backside reflection and a green dominating peak for the reflection.

$$Y_{R_0} = R_0 + (1 - R_0) \cdot Y_0 > R_0. \quad (3)$$

For small values of R_0 , this equation is approximated by

$$Y_{R_0} \approx R_0 + Y_0. \quad (4)$$

This approximation shows that the abacus in Fig. 3(b) can be used to calculate the brightness Y_{R_0} when $R_0 \neq 0$. For example, for $R_0 = 2\%$, just use the level curve $Y_0 = 3\%$ to get a brightness Y_{R_0} of 5% and adjust the wanted peaks properties (widths and respective positions).

In addition to impacting the color brightness, R_0 significantly modifies the obtained gamuts. As can be seen in Fig. 5 for $R_0 = 1\%$ and $R_0 = 2\%$, when Y_{R_0} is close to R_0 , the gamut is extremely small. Indeed, the photons responsible for the desired color are dominated by the number of photons emitted over the entire visible spectrum. For very low brightness ($Y_{R_0} \sim R_0$), the most saturated colors correspond to blue ones, since the $\bar{x}$ color matching function (plotted in Ref. 4) is narrower than the $\bar{y}$ and $\bar{z}$ ones. Last, when the total brightness Y_{R_0} increases and dominates R_0 , the gamut reaches a maximum area (around $Y_{R_0} = 10\%$ for $R_0 = 1\%$ and $Y_{R_0} = 20\%$ for $R_0 = 2\%$) before shrinking again. This behavior can be explained by the Center of Gravity Law of color mixture (see Sec. 2.4) since adding a constant reflection leads to a resulting color closer to the WP. Last, the dashed line between the “bluest” point, WP, and the reddest one delimits the border of the colors that can be achieved; this surface is smaller than with $R_0 = 0$ (Fig. 1) since it corresponds to 49.8% of the gamut (instead of 61%).

2.4 Taking the Most of Multimode PC

In the previous section, we have shown that for a given resonance, we obtain a color C in the form of a triplet (x, y, Y) or (u, v, Y) which determines the maximum brightness $Y = Y_{\max}$ accessible with a single resonant mode. To obtain a color C' of identical chromaticity (x, y) with an intermediate brightness Y in range $[0, Y_{\max}]$, we can mix n resonances giving colors from C_1 to C_n . Because of the Center of Gravity Law of color mixture detailed in Ref. 7, the resulting color C_r is a new triplet (x_r, y_r, Y_r) defined as follows:

$$x_r = \frac{\sum_{i=1}^n \frac{x_i Y_i}{y_i}}{\sum_{i=1}^n \frac{Y_i}{y_i}}, \quad y_r = \frac{\sum_{i=1}^n Y_i}{\sum_{i=1}^n \frac{Y_i}{y_i}}, \quad Y_r = \sum_{i=1}^n Y_i. \quad (5)$$

In the case of two resonant modes giving colors C_1 and C_2 , the resultant chromaticity (x_r, y_r) is on the segment connecting those of C_1 and C_2 both in the CIE 1931 xyY and the UCS color spaces. Thus, by taking two photonic resonances, centered at λ_{c_1} and λ_{c_2} of width $\Delta\lambda_{c_1}$ and $\Delta\lambda_{c_2}$, one can obtain a color in the central part of the gamut, with a controlled low brightness if each peak is very narrow (low values of $\Delta\lambda_{c_1}$ and $\Delta\lambda_{c_2}$). This property also makes it possible to obtain

Fig. 5 Iso luminance curves for a constant reflection in the visible range: (a) with $R_0 = 1\%$, (b) with $R_0 = 2\%$, and (c) with $R_0 = 5\%$.

colors in the blue–magenta–red area which is inaccessible with a single resonance. However, this means that we need PCs

- with a single resonance for the brightest or purest colors and
- with two (or more) resonances to access additional colors or to adjust the brightness to the wanted level.

In both cases, for a correct color, whatever the position of the observer, it is important that these modes are not (or very weakly) sensitive to the angle of incidence θ_0 .

3 Simulation Methodology to Size a s-ARC Using a 2D Photonic Crystal

3.1 Description of the Structure

In this part, we focus on the design of a s-ARC where light at normal incidence will be almost totally transmitted through the patterned layer, except for few wavelengths. Such a PC can be fabricated using a square array of cylindrical patterns, where the filling fractions (ff) represent the amount of surface that is not etched while patterning. For a single etched layer (or a stack of different dielectric layers), the main parameters are:

- the dispersion of each considered material. To avoid optical losses, none of the considered materials should be absorbing in the visible range;
- the thickness of the patterned layer t_{ARC} ;
- the ff of the PC, equal to $\pi \cdot (r/P)^2$ for pillars and $1 - \pi \cdot (r/P)^2$ for holes; and
- the period P of the PC.

Fig. 6 Example of a s-ARC reflecting the green color in case of a 2D array of pillars on glass made of a single dielectric layer with optical index n_{ARC} .

Figure 6 shows an illustration when a single dielectric layer (n_{ARC}) is patterned to obtain a 2D array allowing the green color to be reflected. This 2D array behaves as a bandpass filter for the reflected light if two conditions are satisfied.

1. Out of the resonant wavelength, the filter acts as a broadband ARC due to the Fabry–Perot cavity (vertical direction) leading to destructive interferences for the reflected light. To make a link with the previous part, a more realistic description of $R_0(\lambda)$ should be the optical response of this filter as illustrated in Ref. 8.
2. A sharp resonance of the reflection and transmission occurs due to the destructive interference between the specular transmission beam and the waveguided mode localized in the 2D periodic array. This effect, called a Fano resonance, produces a peak on the reflection spectrum at position λ_c . In the previous part, this peak has been modeled by a Lorentzian to control its shape with only two parameters instead of three.

From now, we propose to study the structure of Fig. 6 in two steps that can be calculated easily.

1. Choose the appropriate material for the s-ARC n_{ARC} and determine the thickness t_{ARC} of this layer and its filling factor ff .
2. Consider the wanted color and determine the right value of the period P .

3.2 Fabry–Perot Behavior in the Visible Range

In this part, we assume that the device is not encapsulated, so the ambient media is air, with an optical index $n_0 = 1$. As a wide range with a flat and low residual reflection is necessary, we focus on a single layer ARC. To easily size the structure, we define a linear equivalent optical index for the s-ARC layer (Fig. 7),

$$(n_{\text{ARC}}^{\text{eq}})^2 = n_{\text{ARC}}^2 \cdot ff + (1 - ff) \cdot n_0^2. \quad (6)$$

The ff can be adjusted so this optical index also satisfies the following equation to minimize the reflection over a wide range of wavelengths:

$$n_{\text{ARC}}^{\text{eq}} = \sqrt{n_g \cdot n_0}. \quad (7)$$

Fig. 7 Equivalent structure of Fig. 6 to size the optical behavior except at the PC resonance and get R_0 as low as possible.

As a consequence, to conciliate both Eqs. (6) and (7), we need

$$ff = n_0 \cdot \frac{n_g - n_0}{n_{\text{ARC}}^2 - n_0^2} \approx \text{constant}. \quad (8)$$

As ff may not be constant, we suggest to weight $ff(\lambda)$ by the color matching function $\bar{y}$ since this function measures the brightness and does not depend on the considered illuminating spectrum. Finally, a single quarter-wavelength coating leads to fix the thickness t_{ARC} since $t_{\text{ARC}} = \lambda_0 / (4 \cdot n_{\text{ARC}}^{\text{eq}})$, where λ_0 is chosen in the middle of the visible range. As the bandwidth of ARC is directly linked to the optical path, the study is limited to the smallest values of t_{ARC} providing the broaden ARC spectral range.

These values (t_{ARC} and ff) are used as initial guesses of a minimization routine based on a Levenberg–Marquardt algorithm (as detailed in Ref. 9) and the rigorous coupled-wave analysis (RCWA) method (as explained in Ref. 10). For this rigorous calculation, a sub-wavelength period of 100 nm is used in order to remove the reflection peak resonance since we can notice that for higher periods, the ARC behavior of the filter remains the same as the perturbation induced by the Fano resonance effect occurs only on a small range of wavelength in the vicinity of the reflection peak. Consequently, at this initial step, the resonant reflection peak is not present in the optical response so a wide spacing between the calculated wavelengths is enough (for instance, only 7 points between 300 and 900 nm), providing a fast optimization of these two parameters in only few minutes.

3.3 Selection of the Reflected Wavelength

As thickness and filling factor are now calculated, the different resonances must be tuned to the desired values that can be achieved by the geometry of the PC (pattern and period). To adjust the wanted reflected signal at wavelength λ_c , we need to consider the propagation in the plane of the s-ARC layer (Fig. 8). Then, light escapes on the front side with an angle θ_0 . On the glass substrate side, the emitted signal may interfere with the incident signal that has directly passed through the ARC layer.

Depending on the polarization of the propagative mode, the equivalent optical index of the central region is given at first approximation in Ref. 11,

$$\begin{cases} (n_{\text{ARC}}^{\text{eq}})^2 = n_{\text{ARC}}^2 \cdot ff + (1 - ff) \cdot n_0^2, & \text{TE polarization} \\ \frac{1}{(n_{\text{ARC}}^{\text{eq}})^2} = \frac{ff}{n_{\text{ARC}}^2} + \frac{1-ff}{n_0^2}, & \text{TM polarization} \end{cases} \quad (9)$$

We can notice that a more accurate formulation can be used involving the grating period P . In this study, we use the first formulation in order to illustrate that f mainly rules the effective optical indices while P controls the wanted reflected wavelength λ_c .

Then, the effective index method for an anti-symmetric three layers slab waveguide is used (as detailed in Ref. 12) to calculate $n_{\text{eff}(j)}^{\text{TE}}$ and $n_{\text{eff}(j)}^{\text{TM}}$. Depending on t_{ARC}/λ_c , different orders j may exist. Due to the grating equation and considering an arbitrary order m , the Bragg equation gives

$$n_{\text{eff}(j)}^{\text{TE/TM}} = n_0 \cdot \sin(\theta_0) \pm \frac{m \cdot \lambda_c}{P}, \quad (10)$$

where $\theta_0 = 0$ at normal incidence, and θ_0 corresponds to the angle at which the selected signal escapes from the PC. This equation leads to a very simple relationship between the selected wavelength λ_c and the period P of the PC,

Fig. 8 Illustration of the propagation of the wanted color (TE mode) in the equivalent ARC layer.

$$\lambda_c = \frac{n_{\text{eff}(j)}^{\text{TE/TM}} \cdot P}{k}. \quad (11)$$

For a given order of the propagating mode, $n_{\text{eff}(j)}^{\text{TE}} > n_{\text{eff}(j)}^{\text{TM}}$ as demonstrated in Ref. 13. Consequently, once the period is chosen, if the system is sized for the TE mode, another color for the TM mode may create another peak at a shorter wavelength. From the analysis in the previous part, this property makes it possible to access the blue–purple–red side of the gamut depending on the chosen material and chosen geometry (t_{ARC} , ff , P). In addition, by sacrificing range of the low reflectivity with a thicker ARC layer, it is possible to increase the number of modes j in the visible range, and thus adjust the brightness of the colors obtained.

The chosen geometry is then used as a starting point for more rigorous calculation with home-made RCWA.

3.4 Leakage Control of the Selected Mode

Whatever the polarization, Eq. (9) leads to $n_{\text{ARC}}^{\text{eq}}$ in between n_0 and n_g . Consequently, the selected colors correspond to leaky modes and part of the incident light is therefore transmitted through the substrate, reducing the reflected signal and so the color saturation. To avoid this unwanted effect, different solutions can be considered:

- partially etch (depth t_s) the glass substrate [Fig. 9(a)], so light is confined in the media with optical index n_{ARC} . Here, t_g must be high enough to avoid leakage, leading to few hundred nanometers. The partial etch of the glass layer may induce an effective index lower than the substrate itself, leading to an increase of the background reflection R_0 ;
- partially etch the ARC layer [Fig. 9(b)] so that a very thin ($t_{\text{slab}} < 20$ nm) layer remains with optical index $n_{\text{ARC}} > n_g$.

In the first case, the issue may be the ARC behavior, since for all other wavelengths, the device looks like a double-layer ARC, that may reduce the transmission plateau. In the second case, the behavior of the ARC is slightly affected, and thus the study exposed in the whole of this article remains valid.

To illustrate the impact of the slab, Fig. 10 compares the reflection for different t_{ARC} in the visible range with (b) and without (a) a 20-nm remaining TiO_2 layer. Without the slab (a), the reflection coefficient does not exceed 1.8% for $t_{\text{ARC}} = 95$ nm. As the equivalent optical index is lower than the one of the glass substrate, the device does not support waveguided modes without losses giving Fano resonances with poor contrast. On the contrary, partially etching the TiO_2 layer (b) results in a residual TiO_2 layer of higher refractive index than the one of glass substrate. This layer reduces the losses of the waveguided mode involved in the Fano resonance. The thickness of this layer must be as low as possible to preserve the ARC properties of the structure. Calculation of the effective index of the mode with a standard optical planar waveguide mode solver as detailed in Ref. 14 shows that a residual thickness t_{slab} of 20 nm is enough to obtain an effective index greater than the refractive index of the glass substrate ($n_g = 1.46$). The evolution of the reflection coefficient versus the wavelength and t_{ARC} with this residual layer is given in

Fig. 9 Different ways to reduce leakage for the wanted colors. (a) Overetching of the glass substrate and (b) partial etching of the ARC layer.

Fig. 10 Evolution of R_0 as a function of the wavelength and the thickness of the ARC for $n_{ARC}^{eq} = 1.21$. (a) $t_{slab} = 0$ and (b) $t_{slab} = 20$ nm.

Fig. 10(b). We can notice that the reflection coefficient remains below 2.9% in the visible range for an optimized t_{ARC} value of 75 nm.

3.5 Design and Optic Optimization by RCWA

Reflection coefficient of 2D cylindrical periodic structures is calculated with home-made RCWA code as detailed in Ref. 10. Optimization of the s-ARC follows the following procedure.

1. We calculate the optical response of the structure for sub-wavelength grating ($P = 100$ nm).
2. Due to a Levenberg–Marquardt algorithm, we determine the thickness t_{ARC} of the remaining TiO_2 layer and its filling factor ff corresponding to the lowest background reflection coefficient R_0 .
3. The period P of the filter is set to match the Fano resonance with the wanted peak that requires the highest period. This particular choice leads to optimize the s-ARC with the biggest leakage coefficient.
4. A thin layer t_{slab} on the bottom of the grating is added in a such way that the leakage of the waveguided mode matches with the coupling coefficient of the GWS (as shown in Ref. 15) giving a maximum of the reflection resonance close to 1. At this step of the optimization process, adding t_{slab} involves an increase of the background reflection coefficient R_0 : different successive iterations between steps 2 and 4 provide a trade-off between the contrast of the resonant peak and the ARC behavior of the filter. The brightness and saturation of the wanted colors are controlled at this step.
5. The period P of the filter is obtained in a last step to match the Fano resonance with all the wanted wavelengths.

All the simulations are performed with refractive indexes of Tables 1 and 2.

Table 1 Experimental Sellmeier coefficients for TiO_2 deposited by sputtering with a maximum error of 1.5 % in the visible range.

i	A_i	λ_i^2 (μm^2)
1	3.150	6.922×10^{-2}
2	1.312	4.829×10^{-4}
3	0.607	155.3

$$n(\lambda) = \sqrt{1 + \sum_i \frac{A_i \lambda_i^2}{\lambda^2 - \lambda_i^2}}$$

Table 2 Sellmeier coefficients for the amorphous SiO₂ from Palik (data available in Ref. 18).

i	A_i	λ_i^2 (μm^2)
1	1.103	8.300×10^{-3}
2	0.364	43.69

As an example, Fig. 11 summarizes the value of the different parameters to obtain the primary colors. In this optimization process, we assume that thickness's t_{slab} and t_{ARC} are the same for all the wanted colors: this condition would correspond to different colored surfaces fabricated in a single step, for example by a single nano-imprint process and a single mold with different patterns. This condition strongly impacts the brightness and saturation, since it impacts the width of the different peaks and the background reflection coefficient as illustrated in Fig. 12. To independently tune the brightness and saturation, one may consider more complex PCs or use a multistep process to set t_{slab} and t_{ARC} at the best values for each wanted color.

The lowest reflection is obtained without the slab ($t_{\text{slab}} = 0$) since in the vertical direction the optical stack behaves like a perfect Bragg mirror. With the 8-nm-thick slab, reflection is higher

Parameter	Value		
ff (%)	$\leftarrow 38 \rightarrow$		
t_{ARC} (nm)	$\leftarrow 73 \rightarrow$		
t_{slab} (nm)	$\leftarrow 8 \rightarrow$		
λ_c (nm)	650	530	470
P (nm)	445	360	310
Expected color*	UCS Judd coordinates		
u	0.4860	0.3383	0.3389
v	0.4356	0.4883	0.3284
Y	0.0802	0.0592	0.2704
	sRGB approximation		
R	138	96	74
G	82	161	84
B	74	33	142
Calculated color			

*: using AM1.5D as illuminant.

Fig. 11 Parameters of the designed samples with a TiO₂ s-ARC layer. **Fig. 12** Intermediate and optimized spectrum for the three primary colors.

($R_0 \sim 5.6\%$). Then, increasing the period P locates the position of λ_c for each wanted color. The reflection due to the slab limits the saturation of each color (Fig. 11), and each is positioned in the sRGB triangle of the UCS Judd gamut.

This method provides the advantage of decorrelating the geometrical parameters of the grating. Furthermore, as the reflection background spectrum out of the resonance is a slowly varying curve, minimization is performed with a reduced number of points and a wavelength resolution of 50 to 100 nm is enough to ensure convergence of the Levenberg–Marquardt algorithm. Then, the computation time of this minimization procedure is comparable to that of the calculation of one of the curves including the Fano resonance with a resolution of 0.5 nm.

4 Discussion and Conclusion

GWS make it possible to obtain structures with a very selective optical behavior. By using non-absorbing materials, it is therefore possible to design nano-patterned layers that are capable of reflecting only certain photons at a wanted wavelength and transmitting all the others in direction of the bottom surface, which is particularly suitable to produce photovoltaic modules with a colored front glass.

The impact of the resonance properties of these GWS has been studied theoretically in the first part. If the width of the resonant peak affects both the brightness and saturation, the parasitic reflection very quickly limits the saturation of the obtained color, at the exception of blue. Despite a discontinuous solar spectrum, the width of the red and blue peaks can be almost linearly related to the chosen wavelength as soon as brightness is chosen: this property makes it possible to very quickly optimize the shape and the position of the optimum resonance for a selected brightness. Due to the high brightness sensibility of the human eye around the green range [450 to 650 nm], an extremely narrow resonance must be used to balance blue or red colors with a same brightness.

In a second part, we have explained how to design a GWS to obtain a reflection spectrum with a single peak and a transmission close to 1 over the whole visible spectrum. Instead of directly using a simulation software and of scanning the different parameters to obtain saturated colors, we have explained how to quickly obtain the geometry and optimize such a structure by decorrelating the different geometric parameters. The first optimization step only requires a 1D optical mode solver and makes possible to size the filling factor of the PC and the thickness of every layer in order to maximize the amount of light transmitted across all the visible range. The PC is partially etched to obtain a residual layer as thin as possible to provide a guided mode with a moderate increase of the background reflection. Optimization with RCWA is particularly fast since the previously calculated thicknesses are accurate initial guesses, and computation for only few wavelengths is required. Last, increasing the period of the PC helps in positioning the Fano resonance peak at the right wavelength for color management: more generally, lateral size of the pattern controls the reflected spectrum properties.

Producing multicolor semi-transparent surfaces with a single nano-patterning step implies that the same thicknesses and partial etching depth must be considered for all the chosen colors. This condition greatly limits the maximum color saturation since the obtained solution corresponds to $R_0 > 7\%$. For outdoor photovoltaic applications, this broadband reflection also reduces the optical transmission and so conversion efficiency. However, one could theoretically reduce this parasitic reflection using a double-layer ARC and materials with suitable dispersion curve such as different silicon oxynitrides. The main drawback of this solution is the narrowing of the optical filter bandpass that may also impact the conversion efficiency of a solar cell because of a high reflection in the infrared domain.

Last, to reduce the impact of the sun position and of the human being toward the patterned surface, one may consider more complex PCs as demonstrated in Ref. 16 or satinated front surfaces as proposed in Ref. 17.

Acknowledgments

The authors would like to thank the NanoLyon technological platform at INL for TiO₂ layers processing and their optical measurement.

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