

Towards high-efficiency industrial p-type mono-like Si PERC solar cells

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ABSTRACT

The cast-grown monocrystalline-like silicon (mono-like Si) technology has been reactivated recently for the manufacture of high-efficiency solar cells at low cost. In this paper, we have provided a progressive research, both experimentally and theoretically, to improve the efficiency of mono-like Si passivated emitter and rear cells (PERCs) through production lines. By employing rear surface passivation of AlO_x /triple- $\text{SiN}_x\text{:H}$ layers, the average efficiency has increased to 21.87% with a champion efficiency of 22.32% from the standard mono-like Si PERCs of 21.72% efficiency with AlO_x capped with double- $\text{SiN}_x\text{:H}$ layers. We have further optimized the rear electrode contact patterns and introduced the selective emitter (SE) technology, yielding an absolute efficiency improvement of 0.19% and 0.26%, respectively. Integrating all approaches of the triple- $\text{SiN}_x\text{:H}$, rear contact pattern optimization and SE technology, we have demonstrated that the mono-like Si PERCs can reach the average efficiency of 22.2% with a maximum efficiency of 22.5%, similar to the current Czochralski monocrystalline Si counterparts.

1. Introduction

For the recent five years, there is a growing tendency of crystalline silicon solar cells that the conventional aluminum back surface field (Al-BSF) design is in a transition to passivated emitter and rear cells (PERCs). It is expected that the production capacity of PERCs will be over 100GWp in China by the end of 2019, and nowadays PERCs have become the mainstream of the highly competitive photovoltaic (PV) market. The concept of PERCs was first proposed by Blakers et al. [1], in 1989. In the past 30 years, a significant subsequent motivation for the shift from conception to mass production is associated with the rapid development of manufacturing technologies [2,3], e.g., rear dielectric coating [4,5], wet chemical polishing etch [6] and laser-based selective emitter (SE) [7] and contact ablation [8]. An aluminium oxide/silicon nitride ($\text{Al}_2\text{O}_3/\text{SiN}_x\text{:H}$) structure was proposed for the rear passivation in 2009 [9,10], which has long been adopted in the industry until now. In our previous work, a feasible silicon oxide (SiO_2) passivation technique for industry-related PERCs was also investigated [11] as a way of improving the long-wavelength response and reducing rear surface recombination. In 2011, a 19.5% cell efficiency of multicrystalline silicon (mc-Si) PERCs was achieved by Q-Cells [12]. Next year, Suntech

Power successfully manufactured advanced PERCs with a laser-based SE [13], instead of other complicated approaches such as ion implantation [14] and printable dopant [15]. While LONGi Solar has declared the realization of world-record PERC efficiency of 24.06% in 2019 [16], the average efficiency for Czochralski monocrystalline silicon (Cz mono-Si) PERCs in current mass production is in the range of $22.2 \pm 0.2\%$.

It should be noted that the PV industry as a whole is dominated by Cz mono-Si PERCs due to in principle the reduced cell efficiency of mc-Si PERCs as well as their limited space to improve the efficiency [17]. Nevertheless, another challenge is indispensable, originating from the fact that the cost of Cz mono-Si wafers is still 30–40% higher than that of mc-Si ones at the present stage [18], causing a significant influence on the price of high-efficiency PERCs. This is unfavorable for the leveled cost of electricity (LCOE) and resolutely enhancing PERC efficiency without increasing the cost should be a decisive strategy. In this situation, the ever popular cast-grown monocrystalline-like silicon (mono-like Si) technology in 2010 has been reactivated [19], which was first proposed by Cizek et al. [20], in 1979. The rapid development of mono-like Si wafers provides the possibility to fabricate high-efficiency PERCs at a relatively low cost. Since 2017, a great variety of works has been carried out to investigate the growth and material features of

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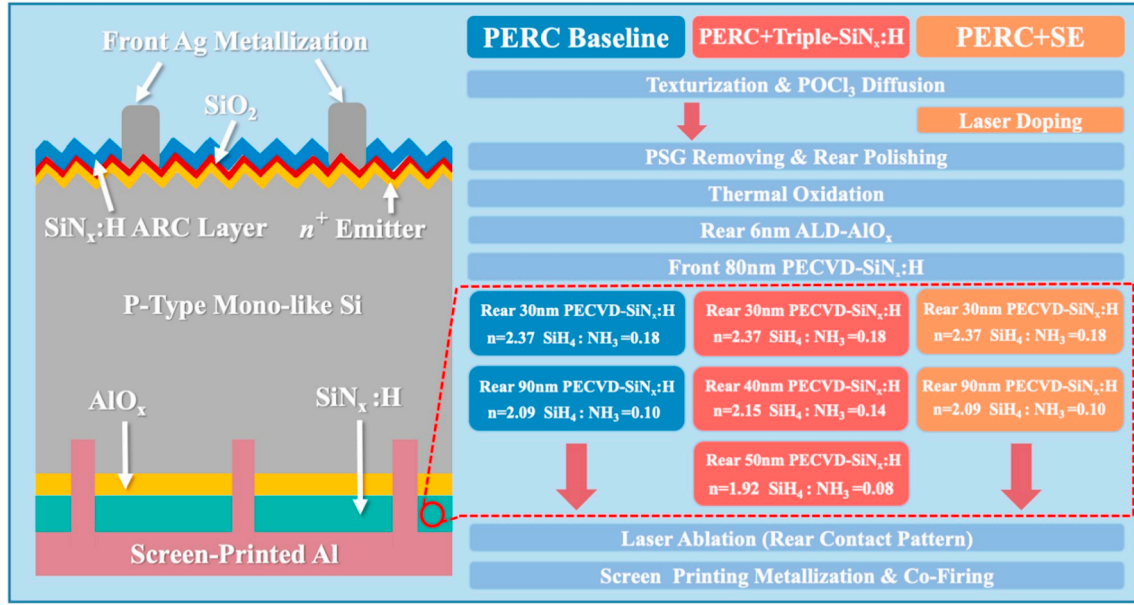


Fig. 1. Schematic of the standard PERC structure (left) and the process flow for the PERC baseline, PERC + Triple-Si_xN_y:H (passivated with triple layers of Si_xN_y:H at the rear side), PERC + SE (combined with SE in the emitter) fabricated in our study together with details of rear AlO_x and Si_xN_y:H layers as well as the front Si_xN_y:H ARC layer (right).

mono-like Si, e.g., impact of impurities [21], defect-related recombination [22] and dislocations [23]. Liu et al. [24], proposed a phosphorous diffusion process to improve the electronic quality of highly defective mono-like Si. Lan et al. [25], introduced a simple approach to control the ingot quality and cell appearance of mono-like Si. Particularly, GCL Poly has played a significant role in the research and development of mono-like Si material, which is strongly supported by several solar cell manufacturing companies. In 2019, Canadian Solar has reported the world-record efficiency of 22.28% for the mono-like Si PERCs [26]. As a promising alternative to the conventional Cz mono-Si PERCs, the production capacity of mono-like Si PERCs is expected to reach 6GWp in 2019 with an average efficiency of $21.8 \pm 0.2\%$ in mass production currently [27].

In contrast to the quick pushing in the PV industry, there is almost no relative report on the mono-like Si PERCs in academia, except three proceedings papers by Xu et al. [28], by Dullweber et al. [29], in 2013 and by Chang et al., in 2014 [30], respectively. We notice that the record cells achieved by either LONGi Solar or Canadian Solar were only certified by the third-party laboratory without any technical analysis. This paper is aimed at providing a progressive research, both experimentally and theoretically, to extend the overall efficiency of mono-like Si PERCs through the production lines. Based on the standard PERCs with AlO_x capped by double-Si_xN_y:H layers at the rear surface with the average efficiency of 21.72%, we have integrated the optimized rear AlO_x/triple-Si_xN_y:H passivation and contact patterns with front laser-based SE technology. With our choice of these feasible technologies, the average efficiency of mono-like Si PERCs has reached 22.2%, with a champion efficiency of 22.5%, which is almost the same as the Cz mono-Si counterparts.

2. Experimental and simulation

We fabricated three types of PERC structures, i.e., standard PERC (baseline with conventional double-Si_xN_y:H passivation layers), PERC with triple-Si_xN_y:H passivation layers (shortened as triple-Si_xN_y:H) and PERC with integration of the SE technology (shortened as SE). The wafers used for this work were p-type mono-like Si wafers with a <100> grain orientation, a size of $157.75 \times 157.75 \text{ mm}^2$, a thickness of $180 \pm 20 \mu\text{m}$ and a resistivity of about $0.8 \Omega \text{ cm}$. Fig. 1 illustrates the

PERC structures and the corresponding fabrication processes, together with the details of rear AlO_x and Si_xN_y:H layers as well as the front Si_xN_y:H antireflection coating (ARC) layer.

2.1. Standard baseline PERCs

The wafers underwent the following production line treatments: 1) standard damage etching, wet-chemical alkaline texturing like Cz wafers to achieve the random pyramid texture with a lower reflectance compared to acidic texturing, which is considered to be more advisable for mc-Si wafers with random grain orientations [31], and RCA clean (InOxSide, RENA); 2) diffusion using POCl₃ as precursor gas to form n⁺ emitter with the sheet resistance (R_{sh}) of approximately $70 \Omega/\square$ (M5111-4WL/UM, CETC 48th Research Institute); 3) edge isolation, rear side polishing and removal of phosphorous silicate glass (PSG) in HF/HNO₃ solution (InOxSide, RENA); 4) thermal oxidation (AS-300E) at the temperature of 680°C for 20min; 5) the 6 nm-thick AlO_x film by thermal atomic layer deposition (ALD) on the rear side (ALD, Ideal Energy); 6) double-Si_xN_y:H layers deposition with thickness of 80 nm on the n⁺ emitter surface by plasma enhanced chemical vapor deposition (PECVD) system (M82200-6/UM, CETC 48th Research Institute); 7) capped with double-Si_xN_y:H passivation layers through the same PECVD system with thickness of 120 nm on the rear side after ALD-AlO_x deposition; 8) laser ablation with a line pattern to form a local contact opening line width of $40 \pm 2 \mu\text{m}$ (DR-LA-Y40, DR Laser), screen printing (PV 1200, DEK) and co-firing technique (CFSeries, Despatch) to ensure front and rear side metallization.

2.2. Triple-Si_xN_y:H PERCs

The fabrication of triple-Si_xN_y:H PERCs was similar to that of the baseline ones, except for the step 7 in Section 2.1. Note that the initial overlapping 120 nm Si_xN_y:H films on the rear surface of baseline PERCs can be divided into the inner Si_xN_y:H layer with a thickness of 30 nm and a refractive index of 2.37 (at $\lambda = 632 \text{ nm}$) while the outer Si_xN_y:H layer with a thickness of 90 nm and a refractive index of 2.09. The different procedures during rear Si_xN_y:H deposition of triple-Si_xN_y:H PERCs are as follows: 1) keep the inner Si_xN_y:H layer fixed; 2) reduce the thickness of initial outer Si_xN_y:H layer from 90 nm to 40 nm, as a new middle layer

Table 1

The main parameter values for the Quokka 2 simulations.

Quokka2 Parameters	Rear electrode contact ratio	23.3%	34.4%	43.4%	53.8%	86.6%
Main	Cell thickness (μm)	170				
	Front width (μm)	1500				
	Rear width (μm)	600				
	Cell depth (μm)	1054	1037	1054	1051	1082
Front (Emitter)	Contact shape	Line				
	Half width (μm)	28.75				
Emitter 1	Sheet resistance (Ω/\square)	70				
	Junction depth (μm)	0.32				
Contact	J_{0e} (A/cm^2)	1.2E-12				
	J_{02} (A/cm^2)	0				
	Contact resistivity ($\Omega\bullet\text{cm}^2$)	1.33E-3				
Non-contact	J_{0e} (A/cm^2)	4.7E-14				
	J_{02} (A/cm^2)	0				
Bulk	Doping type	P-type				
	Resistivity ($\Omega\bullet\text{cm}$)	0.83				
	Background lifetime (μs)	150				
Rear (Base)	Contact shape	Rectangle				
	Number of contacts	1&3				
	Half width X (μm)	34				
	Half width Y (μm)	246	337	457	566	938
BSF 1	Sheet resistance (Ω/\square)	15				
	Junction depth (μm)	0.8				
Contact	J_{0BSF} (A/cm^2)	7.24E-13				
	J_{02} (A/cm^2)	0				
	Contact resistivity ($\Omega\bullet\text{cm}^2$)	2.20E-3				
Non-contact	$J_{0\text{-AlOx/SiNx}}$ (A/cm^2)	7.15E-15				

with a refractive index of 2.15; 3) add another $\text{SiN}_x\text{:H}$ layer as a new outer layer with a thickness of 50 nm and a refractive index of 1.92. The thickness and refractive index of the overlapping $\text{SiN}_x\text{:H}$ films were optimized by Wafer Ray Tracer software (PV Lighthouse) [32].

2.3. SE PERCs

The manufacturing processes of SE PERCs were identical to those of the baseline ones, except for the emitter formation procedure, which involved an additional step: laser doping (DR-LA-Y40, DR Laser) performed on wafers after POCl_3 diffusion and followed by the removal of PSG. Laser doping was included to achieve so-called highly doped emitters (n^{++} emitters) and lightly doped emitters (n^+ emitters), leading to the formation of emitters selectively. Among them, PSG produced during POCl_3 diffusion served as a doping source. The R_{sh} of the n^{++} emitter was $60 \pm 5 \Omega/\square$, while the R_{sh} of the n^+ emitter was $120 \pm 10 \Omega/\square$.

2.4. Symmetrical samples for passivation characteristics

Both saturation current density (J_{0e}) and effective minority carrier lifetime (τ_{eff}) effectively investigate the passivation quality of solar cells in the form of symmetrical structures. For baseline PERCs, the J_{0e} experiment was divided into two parts: the passivation zone and the metal zone. Regarding the passivation zone, $\text{SiN}_x\text{:H}$ passivation layers were deposited on both the textured and phosphorus-diffused surfaces of wafers, with R_{sh} of $70 \Omega/\square$, which forms the symmetrical $\text{SiN}_x\text{:H}/\text{Si}/\text{SiN}_x\text{:H}$ structure. Prior to $\text{SiN}_x\text{:H}$ deposition, textured and diffused wafers underwent the conventional HF cleaning and thermal oxidation processes. The wafers were further treated with sintering process without metallization. Regarding the metal zone, prior to sintering process, the textured and diffused wafers underwent the similar processing sequence as mentioned above, together with Ag printing on both full-area surfaces verse sintering process, with the symmetrical $\text{Ag}/\text{SiN}_x\text{:H}/\text{Si}/\text{SiN}_x\text{:H}/\text{Ag}$ structure. After that, the wafers were soaked in aqua regia solution ($\text{HCl}:\text{HNO}_3 = 3:1$) to remove the Ag paste. As for the τ_{eff} measurement, the sequence is the same as the J_{0e} experiment for the passivation zone with a symmetrical $\text{SiN}_x\text{:H}/\text{Si}/\text{SiN}_x\text{:H}$ structure.

For SE PERCs, the J_{0e} experiment was divided into three parts: the n^+

emitter zone, the n^{++} emitter zone and the metal zone. Regarding the n^+ emitter zone, the fabrication of symmetrically structured samples was similar to that of the samples tested for the passivation zone of baseline PERCs, except for the R_{sh} of about $125 \Omega/\square$. Regarding the n^{++} emitter zone, prior to $\text{SiN}_x\text{:H}$ deposition, full-area laser doping was added to the fabrication process on both surfaces of textured and phosphorus-diffused wafers as mentioned above, with R_{sh} of $65 \Omega/\square$, the other steps were exactly the same. Regarding the metal zone, the samples just used to test for the n^{++} emitter zone underwent the similar treatments as the metal zone of baseline. For the measurement of τ_{eff} , $\text{SiN}_x\text{:H}$ films were deposited on both surfaces of the textured and phosphorus-diffused wafers, with R_{sh} of $125 \Omega/\square$, similar to the process as mentioned above.

2.5. Characterization

The thickness and refractive index of AlO_x and $\text{SiN}_x\text{:H}$ films were measured by spectroscopic ellipsometry (SE400adv-PV). The microgram of different rear contact patterns were obtained by a 3D microscope (Zeta-20, ZETA Instruments). Dopant concentration profiles of the emitters were determined by electrochemical capacitance voltage (ECV) profiling (CVP21, WEP). Sheet resistance R_{sh} and resistivity of the wafers were investigated by four-point probes (280I Series, Four Dimensions Inc.). The interstitial oxygen (O_i) of the mono-like Si wafers was investigated by FTIR (Bruker IFS-113v). τ_{eff} , J_{0e} and implied open circuit voltage (V_{oc}) of the emitters were measured by a quasi-steady-state photoconductance (QSSPC) tester (WCT-120, Sinton Instruments). The reflectance and quantum efficiency of solar cells were measured by QEX10 (PV Measurements) system. Finally, the efficiency of the solar cells was tested under AM1.5 spectrum at 25°C by Halm 3600 in the production line.

2.6. Simulation of optical absorption and reflection

To understand the effect of rear triple- $\text{SiN}_x\text{:H}$ layers on the optical absorption and reflection, the online Wafer Ray Tracer software was employed to optimize thickness and refractive index distribution. Both the baseline PERCs and triple- $\text{SiN}_x\text{:H}$ PERCs were set coated with double- $\text{SiN}_x\text{:H}$ layers on the front surfaces, with thickness of 55 nm and refractive index of 2.03 for outer layer as well as thickness of 25 nm and

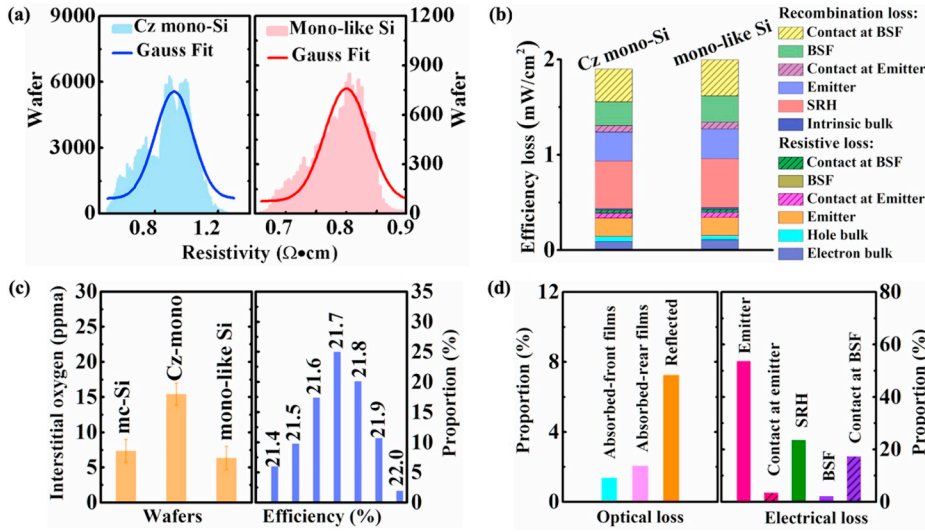


Fig. 2. (a) Resistivity distribution curve of Cz mono-Si and mono-like Si wafers. (b) FELA simulation results of Cz mono-Si and mono-like Si PERCs with the same structure, the measured bulk lifetime was used for the simulation, with a real value of 150 μ s for mono-like Si wafers and 200 μ s for Cz mono-Si wafers. (c) Comparison of interstitial oxygen of the mc-Si, Cz mono-Si and mono-like Si wafers (left) and efficiency distribution of our baseline PERCs fabricated in mono-like Si wafers in the range of 21.4–22.1% (right). (d) Simulated optical loss in baseline PERCs (left) and electrical loss in baseline PERCs (right).

refractive index of 2.37 for inner layer, respectively. With regard to the rear surfaces, fix the total $\text{SiN}_x\text{:H}$ layer thickness of 120 nm and the inner film thickness of 30 nm, change the thickness distribution of the other two layers. The refractive index of outer $\text{SiN}_x\text{:H}$ layer varied from 1.91 to 2.03 while the inner and middle layers remained 2.37 and 2.15, respectively.

2.7. Simulation of electrical performance and free energy loss analysis

Quokka 2 was used to study and verify the performance of three types of PERCs with different real ratios of rear electrode contact patterns. Table 1 lists the simulation parameters of the solar cells.

3. Results and discussion

3.1. Baseline mono-like Si PERCs

Before presenting the results of baseline mono-like Si PERCs, the intrinsic properties of Cz mono-Si and mono-like Si wafers were compared, both experimentally and simulatively. We show in Fig. 2(a) the resistivity distribution of these two types of wafers. There is an evident fluctuation found in the resistivity of Cz mono-Si wafers ranging from 0.5 Ω cm to 1.3 Ω cm, while the concentrated resistivity of mono-like Si wafers with a stable value of $0.80 \pm 0.1 \Omega$ cm is more preferable for the PERC production process. As illustrated in Fig. 2(b), we make a comparison between Cz mono-Si PERCs and mono-like Si PERCs with the same cell structure in efficiency and free energy loss analysis (FELA) by Quokka 2 simulation. It is worthy to mention that the measured bulk lifetime was used for the simulation, with a real value of 150 μ s for mono-like Si wafers and 200 μ s for Cz mono-Si wafers. According to the simulation results, it is found that the energy loss of Cz mono-Si PERCs is lower than that of mono-like Si PERCs, leading to a cell efficiency improvement of 0.03%. The main reason lies in the higher bulk recombination of mono-like Si PERCs. However, mono-like Si wafers exhibit much lower O_i with the value of 6.34ppma, while the O_i of Cz mono-Si wafers is 15.43ppma as listed in the left graph of Fig. 2(c). The lower O_i in mono-like Si wafers is instrumental in reducing the light-induced degradation (LID) [29] of PERC modules. The right graph of Fig. 2(c) depicts the mass production efficiency distribution of the baseline mono-like Si PERCs with the manufacturing process flow listed in Fig. 1. It is found that the cell efficiency (η) is mainly concentrated at 21.80% with a share of 25%, the efficiency of 21.9% second, yielding an average η of 21.72%. There is only 12.6% share with the η of more than 22.0%, indicating the large improvement room for the mono-like Si

PERCs as compared with the Cz mono-Si counterparts.

It is well known that the loss of solar cells could be categorized as optical and electrical losses. Fig. 2(d) shows the related losses from a series of simulations to straightforwardly clarify the direction of improvement. On the one hand, the optical losses by Wafer Ray Tracer software indicate that the reflection loss is the dominating factor to limit the carrier collection of solar cells, and the improvement of spectral response is restricted unless the rear films are optimized with an alternative routine for a long optical transmission length. The absorption loss from rear cell is also higher than that of front cell, causing a negative influence on the long-wavelength response. On the other hand, electrical losses with the help of Quokka 2 simulation reveal the serious front emitter loss, which could be effectively reduced by the SE technology. Apart from the Shockley-Read-Hall (SRH) loss, the loss from contact at BSF also accounts for a considerable proportion, which can be reduced by optimal rear laser cut groove to achieve a low resistive and recombination loss. The above simulation results guide us to investigate the influence of emitter and rear structure on cell performance below.

3.2. Effect of rear triple- $\text{SiN}_x\text{:H}$ passivation films

It is well known that the optical loss in the long wavelength is mainly caused by the reflection and absorption losses from the rear metal layer [33], leading to the reduction of short circuit current (I_{sc}). To further improve the long-wavelength response of cells, it is necessary for rear passivation films to obtain a higher reflection when light arrives at the rear, leading more long-wavelength photons to be reflected into the Si wafers to be fully utilized again. According to the reflection theory, triple- $\text{SiN}_x\text{:H}$ instead of conventional double- $\text{SiN}_x\text{:H}$ in the industry (i.e., the baseline) can effectively increase the reflection because of their multiple surfaces, so that the light fraction has another chance to enter the wafers [34]. Prior to the experiments, it is essential to figure out the optimal thickness and refractive index for triple- $\text{SiN}_x\text{:H}$ PERCs by means of simulation. Here, the thickness distribution of each $\text{SiN}_x\text{:H}$ layer from the inner to the outer was categorized as two types: 30nm/30nm/60 nm and 30nm/40nm/50 nm, while the thickness distribution of baseline is 30nm/90 nm. We used the Wafer Ray Tracer software to calculate the optically generated current density (J_G) with different refractive index of triple- $\text{SiN}_x\text{:H}$ PERCs (see section 2.6 for details). It's worth to mention that the passivation of $\text{SiN}_x\text{:H}$ films enhances with an increase in refractive index [35]. This is due to the higher density of Si:H bond determined by the ratio of SiH_4 to NH_3 gas flow in the PECVD system [36,37]. The larger the ratio of SiH_4 to NH_3 gas flow, the more the Si:H bond density and thereby the higher the refractive index of $\text{SiN}_x\text{:H}$.

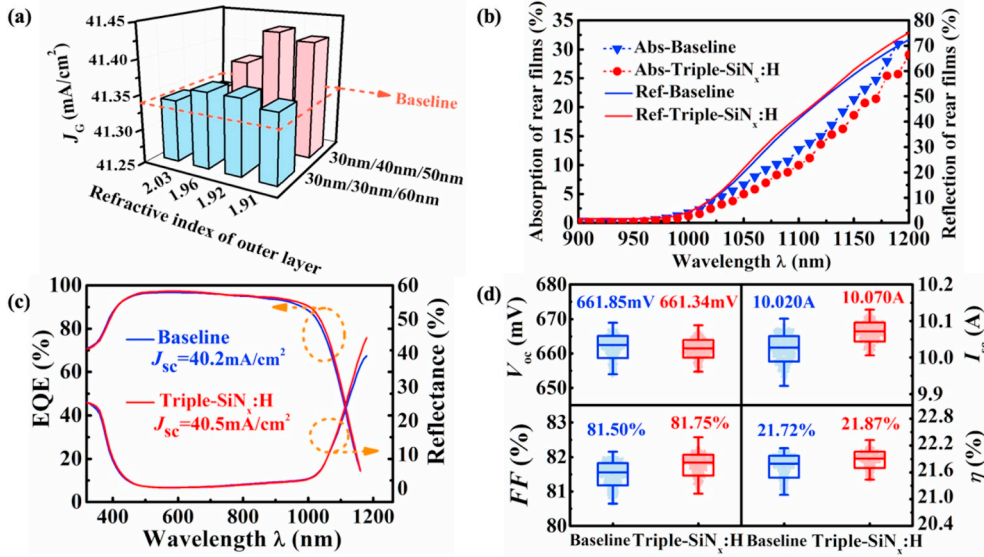


Fig. 3. (a) Simulated values of J_G for triple-SiN_x:H samples. 30nm/30nm/60 nm and 30nm/40nm/50 nm types represent the thickness of each layer from inner to outer. The refractive index of outer layer for each type varied from 1.91 to 2.03. The dotted line represents the J_G value of baseline. (b) Simulated optical absorption and reflection from rear films. (c) EQE and reflectance spectra measured with a mono-chromatic approach. (d) Measured I - V parameters (J_{sc} , V_{oc} , FF and η) with 400 pieces of solar cells.

However, the lower refractive index of SiN_x:H results in improving the reflection, thereby increasing the long-wavelength response of PERCs. This explains our choice of refractive index for triple-SiN_x:H PERCs.

Fig. 3(a) illustrates the calculated J_G characteristics, where there is an obvious change after triple-SiN_x:H replacing the double-SiN_x:H baseline (the J_G value for the baseline is 41.34 mA/cm²). The J_G s for the 30nm/30nm/60 nm type are 41.34, 41.36, 41.36 and 41.35 mA/cm², while for the 30nm/40nm/50 nm type, the corresponding J_G s are 41.33, 41.38, 41.43 and 41.42 mA/cm² with refractive index of 2.03, 1.96, 1.92 and 1.91 for outer SiN_x:H layer, respectively. This result clearly indicates that the triple-SiN_x:H PERCs exhibit higher J_G s overall compared with those of baseline. The maximum J_G of 41.43 mA/cm² is achieved at the refractive index of 1.92 in terms of the 30nm/40nm/50 nm type, thereby considered to be the best case for triple-SiN_x:H designs. In other words, the following discussion with respect to the triple-SiN_x:H PERCs is based on the 30nm/40nm/50 nm type with refractive index for outer SiN_x:H layer of 1.92. Although big difference of the refractive indices (2.37 and 1.92) in double-SiN_x layers (30 nm and 90 nm) can yield the similar optical effect, the introduction of intermediate refractive index of 2.15 in the triple-SiN_x scheme can save the industrial cost of rear SiN_x

fabrication in the PECVD system.

Obvious differences could be observed in Fig. 3(b) between triple-SiN_x:H and baseline PERCs from the absorption and reflection in the long wavelength from 900 nm to 1200 nm by Wafer Ray Tracer software. The surface of light propagation has increased for triple-SiN_x:H PERCs, leading to a slight increase in reflection and thereby a lower absorption. Another explanation for the improvement might be that there is an increase in extinction coefficient with increasing ratio of SiH₄ to NH₃ gas flow [38], resulting in a reduced absorption in the long wavelength of triple-SiN_x:H PERCs with a lower refractive index of 1.92 for outer SiN_x:H layer, in contrast to the baseline with a refractive index of 2.09 for outer SiN_x:H layer. The above argument reveals that the reflection and absorption of long-wavelength photons are significantly modified when the rear surfaces are passivated by triple-SiN_x:H films.

Fig. 3(c) depicts the experimental EQE and reflectance spectra of triple-SiN_x:H PERCs against the baseline ones. It is almost the same in the short wavelength of both EQE and reflectance for the reason that short-wavelength spectral response mainly shows the recombination and loss from the front cell, which is identical between triple-SiN_x:H and baseline PERCs. However, in the long wavelength, triple-SiN_x:H PERCs

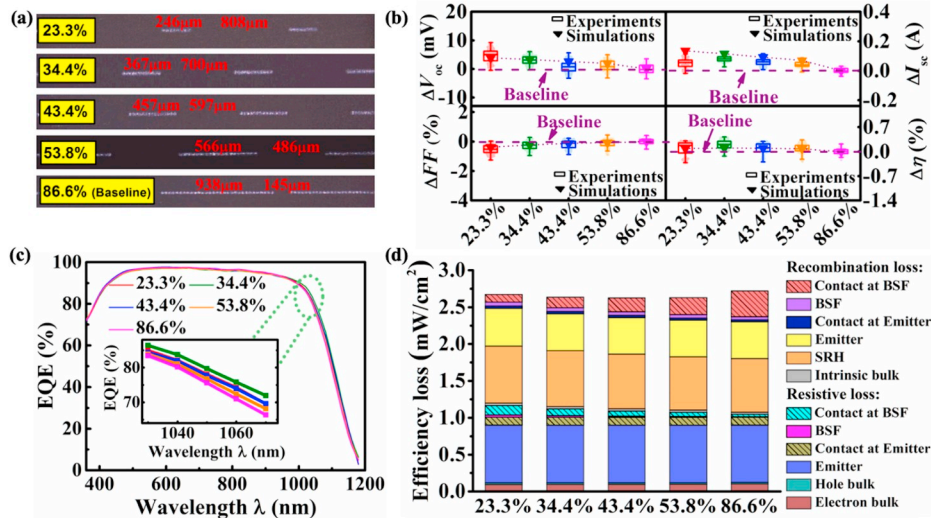


Fig. 4. Different rear contact patterns ($L^R/L = 23.3\%$, 34.4% , 43.4% , 53.8% and 86.6%): (a) 3D microgram, (b) measured and simulated I - V parameters (relative value of I_{sc} , V_{oc} , FF and η compared with baseline), (c) EQE spectra and (d) simulated FELA results.

achieves a better performance on EQE and reflectance than the baseline ones, demonstrating superior light utilization and carrier collection efficiency at the rear side. As a result, the yielded short circuit current density (J_{sc}) from EQE spectra rises from 40.2 mA/cm² to 40.5 mA/cm².

Fig. 3(d) shows the comparison of electrical parameters between the two types of PERCs. Obviously, the average I_{sc} and fill factor (FF) of triple-SiN_x:H PERCs exceed the baseline ones, leading to higher η . The average η reaches 21.87% for triple-SiN_x:H PERCs, accompanied by a corresponding increase of 0.15% over baseline ones with an average η of 21.72%. The increase in the I_{sc} of 50 mA is consistent with the improvement of EQE and reflectance in Fig. 3(c). It is also worth to mention that the value of series resistance (R_s) is 0.86 mΩ and 0.92 mΩ for triple-SiN_x:H and baseline cases, respectively. We assume that the higher FF in triple-SiN_x:H PERCs is due to a nitrogen-rich film at the Al/SiN_x:H interface because of a lower refractive index [39], resulting in attenuating the corrosion penetration of Al paste and thereby a lower R_s . However, no obvious change occurs in V_{oc} as a result of the equivalent total thickness of SiN_x:H films, meaning that the passivation effect is hardly influenced by the triple-layer films. Therefore, it could be concluded that the increased η is practically attributed to the improvement of I_{sc} and FF , which is mainly affected by the lower refractive index at outer layer, enhanced reflection and reduced absorption in the long wavelength. In summary, the rear triple-SiN_x:H film structure is confirmed to be a promising solution to enhance the long-wavelength spectral response for PERCs.

3.3. Influence of rear electrode contact patterns

Since the AlO_x/SiN_x:H stack is considered as an insulating layer, laser ablation is necessarily required for electrode formation at the rear side, that is, aluminum-silicon contact can be realized in the ablation portion [40,41]. There is no doubt that the electrode contact patterns should be well designed. Note that the contact patterns in production lines for baseline PERCs are dotted lines with length of 938 μm and lateral spacing of 145 μm, respectively, as shown in Fig. 4(a). Here, we define a new quantity, real ratio (L^R/L , where $L = L^R + L^I$, L^R is the length, and L^I is the lateral spacing), to simplify the line patterns description, fixing the values of longitudinal spacing 600 μm and width 40 μm. To investigate the influence of rear contact patterns, a series of PERCs with different real ratios L^R/L were fabricated in mono-like Si wafers. Fig. 4(a) shows the image of rear contact patterns with L^R/L ranging from 23.3% to 53.8%, and the pattern of baseline ($L^R/L = 86.6\%$) is also displayed for comparison. The corresponding ratio of electrode contact area to the total area of rear surface is about 1.1%, 1.8%, 2.3%, 2.9% and 4.7%, respectively. We aim at obtaining the optimal L^R/L , which can provide the highest output performance compared with baseline PERCs.

The relative values of experimental parameters (ΔV_{oc} , ΔI_{sc} , ΔFF and $\Delta \eta$) between four set samples ($L^R/L = 23.3\%$, 34.4%, 43.4% and 53.8%) and baseline are illustrated in Fig. 4(b). It is easy to find that with the increase of L^R/L , the V_{oc} diminished progressively. By contrast, obvious improvement in FF is observed with increasing L^R/L . The R_s of every set sample (23.3%, 34.4%, 43.4%, 53.8% and 86.6%) is 0.92, 0.78, 0.69, 0.66 and 0.60 mΩ, respectively. The lower the R_s , the larger the FF , thus the decline of FF is a result of the increasing R_s , which is mainly affected by contact resistance and lateral transport resistance [42,43]. The above trends can be adequately explained by the dependence of lateral transport resistance on L^R/L , which follows the adverse tendency as FF . However, an increase in the contact area means a reduction in the passivation area [44], causing the V_{oc} to drop clearly. Hence, it is not preferable to improve the performance by increasing the laser cut groove area. One can see that the I_{sc} of the four set samples exceed the baseline significantly. There is a nearly 47, 81, 59 and 41 mA improvement with L^R/L of 23.3%, 34.4%, 43.4% and 53.8%, respectively, and this could be attributed to differences between current transmission channels and current collection at the rear surface. Also, a homologous trend can be observed in η that the conversion efficiency of

four set samples is greatly better than the baseline absolutely increased by 0.09%, 0.19%, 0.08% and 0.08%, respectively. It indicates that the higher V_{oc} and greater I_{sc} would compensate the FF loss accounting for the fact that the sample with L^R/L of 34.4% exhibits the best electrical performance in comparison with other set samples.

Fig. 4(b) also shows the simulated electrical parameters by Quokka 2 software for verification. In order to ensure an effective and feasible simulation, the experimental data was adopted in the calculation, such as sheet resistance, junction depth, contact resistance, J_{0e} and so on. Note that trends of ΔV_{oc} , ΔI_{sc} , ΔFF and $\Delta \eta$ depicted by Quokka 2 with L^R/L varied from 23.3% to 53.8% are in good agreement with those through experiments. Simulation results reveal that I_{sc} of the 23.3% set samples is better than that of the 34.4% set ones. This may be due to the identical contact resistivity of rear side selected for each set samples during the simulations. Indeed, the smaller the L^R/L , the lower the contact resistivity [45]. Overall, compared with the simulated baseline counterparts, the optimal results of the electrical performance are obtained by Quokka 2 at $L^R/L = 34.4\%$ with an absolute increase in efficiency η of 0.23%, while the maximum improvement η of 0.19% has been achieved experimentally for the PERCs with real ratio L^R/L of 34.4%.

In order to evaluate the improvements of I_{sc} more convincingly, Fig. 4(c) depicts EQE spectra of all these set samples. One can clearly observe that the set samples with $L^R/L = 34.4\%$ exhibit a stronger spectral response in the long wavelength but nearly no difference in the short and middle wavelength compared with other set samples. This could be associated with the sufficiently high quality of rear surface passivation and good ability of rear current collection, rather than the influence of the emitter and the base. What's more, the EQE tendency of these five set samples is in accordance with the corresponding experimental I_{sc} listed in Fig. 4(b).

To straightforwardly analyze the influence of real ratio L^R/L on electrical losses, the FELA was also carried out by Quokka 2, as illustrated in Fig. 4(d). According to the FELA, these five set samples are obviously different in two items, the recombination loss and the resistivity loss of contact in BSF. This is because the higher L^R/L of rear contact area indicates that the wafer is subject to higher surface recombination velocity as well as greater laser damage, leading to more serious recombination losses of contact in BSF. However, the smaller L^R/L of rear contact area possesses a higher transverse resistance, resulting in much more losses. It is required to make a trade-off between the reduction of resistivity loss and combination loss [42], and $L^R/L = 34.4\%$ is confirmed to satisfy above requirements commendably. In conclusion, optimal rear electrode contact pattern should not only have high quality passivation, but also low transverse resistance. Too large and too small L^R/L are not advisable to improve the performance of the PERCs with the following two reasons: (i) the decrease of V_{oc} and I_{sc} is inevitable owing to the loss of passivation regions and enhanced absorption of photons by the electrode contact when L^R/L is too large; (ii) the decline of FF resulting from the excessive R_s appears as well as the deterioration of aluminum-silicon contact performance when L^R/L is too small.

From the FELA in Fig. 4(d), one can also note that the top-three energy losses for PERCs are the emitter resistance loss, the SRH bulk recombination loss and the emitter recombination loss. Therefore, the SE technology is a necessary and cost-effective strategy to further optimize the performance of PERC cells, which will be discussed in Section 3.4.

3.4. Analysis of SE technology

Great efforts have been made on optimizing the manufacturing process through production lines with regard to rear sides of the PERCs in our previous sections. Here we will investigate the gains from the front sides combined with SE technology. SE is a structure in which an n^{++} emitter is formed at the front metal finger, together with an n^{+}

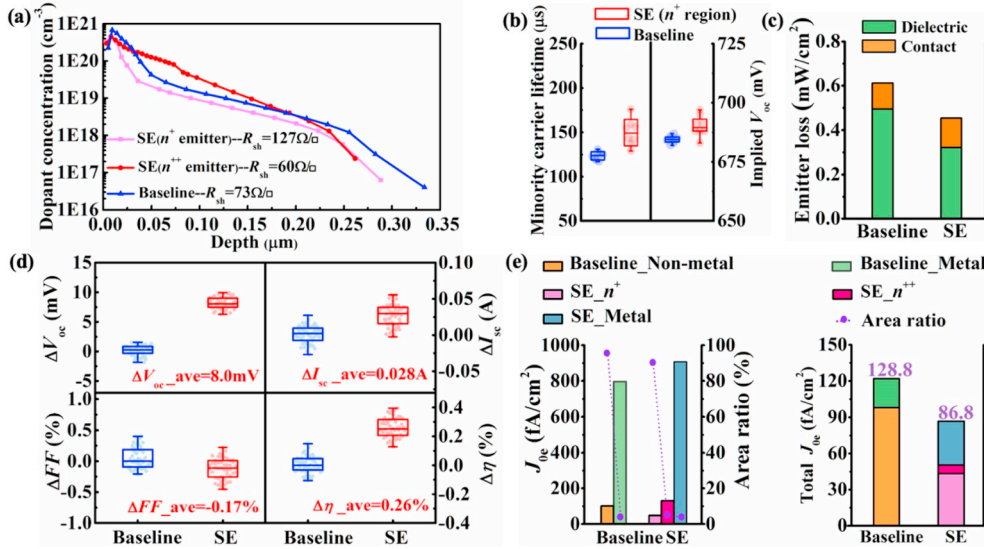


Fig. 5. (a) Phosphorous dopant concentration profiles of the emitters for baseline and SE samples. (b) Measured τ_{eff} and implied V_{oc} for baseline and the n^+ emitter of SE at an injection level of 10^{15}cm^{-3} with a symmetric structure. (c) Simulated efficiency loss in the emitter for baseline and SE. (d) Measured I - V parameters (relative value of I_{sc} , V_{oc} , FF and η compared with baseline). (e) Measured J_{0e} of different regions (metal and non-metal for baseline and n^+ emitter, n^{++} emitter and metal for SE) together with the area ratio of the corresponding regions (left), the calculated total J_{0e} as a function of J_{0e} of each region and the corresponding area ratio (right).

emitter to remain high quality passivation. On the one hand, the n^+ emitter has a completely lower dopant concentration and junction depth, which contributes to less serious recombination in the non-metallized areas, i.e., the Auger recombination and SRH recombination [46]. On the other hand, the higher dopant concentration of the n^{++} emitter form high-performance Ohmic contact among the emitter and metal electrodes to reduce the contact resistance [47].

Fig. 5(a) shows the phosphorous dopant concentration profiles of the emitter with two types of solar cells, baseline and SE technology, respectively. For the baseline, the value of R_{sh} is $73\Omega/\square$, with a phosphorous surface concentration of about $6.8 \times 10^{20}\text{cm}^{-3}$ versus a junction depth of $0.28\mu\text{m}$. For the case of SE technology, the R_{sh} of the n^+ emitter is $127\Omega/\square$, with a phosphorous surface concentration of about $4.25 \times 10^{20}\text{cm}^{-3}$ and a junction depth of $0.24\mu\text{m}$. A strong thermal effect due to the laser doping not only causes the phosphorous atoms to quickly penetrate into silicon wafers, but also melts the PSG so that the phosphorous atoms in the PSG precursor layer are almost completely utilized and penetrate into the deeper of the wafers [48]. Therefore, the n^{++} emitter was obtained with R_{sh} of $60\Omega/\square$, based on the n^+ emitter with R_{sh} of $127\Omega/\square$. It is expected that the SE technology possesses advantageous quality for emitters.

Fig. 5(b) displays the τ_{eff} and implied V_{oc} with and without SE. For a convenient and rapid comparison, the τ_{eff} was extracted at an injection level of 10^{15}cm^{-3} as well as the implied V_{oc} . Here, we just consider the passivation regions of SE for comparison, namely the n^+ emitter. It can be observed that the average value of τ_{eff} increases from approximately $127\mu\text{s}$ – $154\mu\text{s}$ and the implied V_{oc} raises from 683mV to 690mV after SE process, respectively. Reasons for the improvement can be evaluated as the follows: a notable reduction of surface Auger recombination and defect density from the associated low level of dopant concentration of the n^+ emitter. This supports the importance of SE technology as a result of good chemical passivation. To estimate whether or not the passivation quality of emitters could be further enhanced in combination with the SE structure, Quokka 2 software was implemented to achieve a deep insight to the recombination loss of the cell front, as schematically shown in Fig. 5(c). The loss from the cell front was mainly attributed to two components: dielectric and contact areas. It's clear that recombination from the contact between baseline and SE technology is almost the same, which means that recombination from the dielectric areas might be the primary determinant of the emitter quality.

As illustrated in Fig. 5(d), the experimental results of SE technology show an increase in the V_{oc} of 7.98mV , and an increase in the I_{sc} of 28mA , accompanied by a corresponding drop in the FF of 0.17%

compared to the homogeneous emitter baseline cells. There is now a general consensus that the improvement in V_{oc} is significantly correlated with the rewarding chemical passivation [49]. The higher I_{sc} is due to the reduced emitter recombination to the profitable impact of the short-wavelength spectral response. The maximum η improvement of 0.26% has been obtained on the PERCs combined with SE technology.

In order to further quantitatively explore front emitter quality of both cells, we constituted a valid comparison of recombination current density J_{0e} to evaluate with different regions, i.e., the metal zone, the n^{++} emitter zone, the n^+ emitter zone for the SE case, and the metal zone, the non-metal zone for the baseline one, respectively. It is worth to mention that J_{0e} of different regions were measured with the symmetrical structures (see section 2.4 for details). Here, the proportion of each region to the total area of cells is also displayed for the purpose of calculating the total recombination current density of cells, as given in Fig. 5(e) (the left graph). Note that the J_{0e} of the SE case in three regions (including n^+ emitter, n^{++} emitter and metal) is 48 , 130 and $907\text{fA}/\text{cm}^2$, whose area accounts for 90.6% , 5.4% and 4.0% , respectively, while the J_{0e} of the baseline case in two regions (including non-metal and metal) is $101\text{fA}/\text{cm}^2$ and $796\text{fA}/\text{cm}^2$, accounting for 96% and 4% , respectively. The presence of SE technology results in a higher level of J_{0e} in the metal regions. This may be understood by the fact that metal contact of the SE case is performed in the n^{++} emitter region, that is to say the dopant concentration of SE metal regions is higher in comparison with that of the baseline. Another explanation might be that the SE front surface would suffer unavoidable damage caused by the laser doping, thereby resulting in the increase of the Auger recombination in the emitter.

The total J_{0e} of both cells was calculated as a function of J_{0e} of different regions as well as the corresponding area ratios. In other words, the total J_{0e} of the SE case is the sum of the J_{0e} of three areas (n^+ emitter, n^{++} emitter and metal) multiplied by the respective area ratio, while the calculation of the baseline case was similar to that of the SE. Fig. 5(e) (the right graph) summarizes the values of the total J_{0e} for the baseline case of $128.8\text{fA}/\text{cm}^2$ and for the SE one of $86.8\text{fA}/\text{cm}^2$, suggesting a good enough front surface passivation. In conclusion, through the integration of SE technology, we have achieved good quality emitter, which is an essential prerequisite for high performance PERCs.

To verify the feasibility and superiority of our roadmap, we have further combined the efficiency enhancements of triple-SiN_x-H, rear contact pattern of $L^{\text{R}}/L = 34.4\%$ and SE technology for mono-like Si PERCs. As displayed in Fig. 6, the average results of the electrical parameters ($V_{\text{oc}} \sim 668.5\text{mV}$, $I_{\text{sc}} \sim 10.2\text{A}$, $FF \sim 81.6\%$ and $\eta \sim 22.2\%$) are

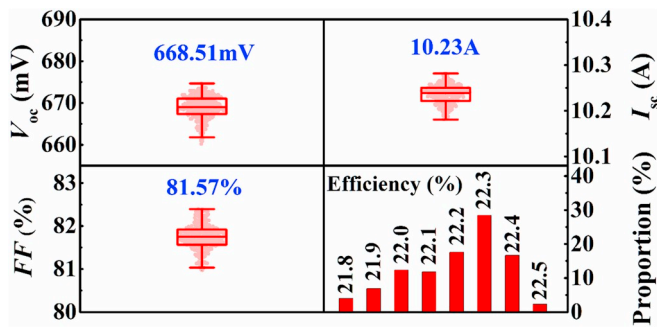


Fig. 6. Measured I - V parameters (absolute values of V_{oc} , I_{sc} , FF and η of mono-like PERCs combined with triple-SiN_x:H, rear contact pattern of $L^R/L = 34.4\%$ and SE technology).

obtained with a maximum efficiency of 22.5%. It is obvious that the η is mostly concentrated at 22.2%, 22.3% and 22.4%, which is almost the same as the Cz mono-Si counterparts.

4. Conclusions

In summary, we have investigated both experimentally and theoretically the application of mono-like Si wafers in the mass production of mainstream PERCs. By employing triple-SiN_x:H in combination with AlO_x at the rear surface to replace the standard AlO_x/double-SiN_x:H stacks, more long-wavelength photons can be reflected into the Si wafers to be utilized again, resulting in an increase in the I_{sc} of 50 mA. The rear triple-SiN_x:H also performs a high FF attributed to a nitrogen-rich film at the Al/SiN_x:H interface. Consequently, the average η of our mono-like Si PERCs has been increased to 21.87% from the existing baseline η of 21.72%. We have further demonstrated the effect and role of rear contact patterns, highlighting the necessity of keeping the balance between the passivation and series resistance. The optimal contact pattern of $L^R/L = 34.4\%$ exhibits a higher performance in both V_{oc} and I_{sc} , and the average η increases by more than 0.19% absolutely, compared to the standard PERCs with $L^R/L = 86.6\%$. Finally, the emitter saturation current density J_{0e} has been dramatically reduced from 128.8 fA/cm² to 86.8 fA/cm² by the implementation of SE technology, yielding an absolute average η improvement of 0.26%. Finally, we have combined these feasible technologies in the process of baseline mono-like Si PERCs. An average η of 22.2% accompanied by a maximum η of 22.5% was obtained, which is almost the same as their Cz mono-Si counterparts.

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