

Devices and Circuits Using Novel 2-D Materials: A Perspective for Future VLSI Systems

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Abstract—Here, we review the most recent developments in the field of 2-D electronics. We focus first on the synthesis of 2-D materials, discussing the different growth techniques currently available and assessing their strengths and weaknesses. Moreover, we describe a possible roadmap to enable CMOS compatible integration of 2-D materials. We then shift our attention to 2-D devices and circuits and review the state of the art. Among the plethora of device concepts, we look closely at 2-D tunnel FETs (TFETs) and negative-capacitance FETs (NC-FETs) for low-power applications. We also put a particular emphasis on doping-free polarity-controllable systems that use electrostatic doping to eliminate the need for physical or chemical doping. We conclude with an analysis of simulations of scaled devices and discuss the possibilities enabled at circuit level by 2-D electronics.

Index Terms—2-D materials, beyond CMOS, doping-free, growth, integration, low power, polarity control, scaling, transfer, transition metal dichalcogenides (TMDCs).

I. INTRODUCTION

RESEARCH on 2-D materials has experienced remarkable growth in the last decade. Major semiconductor companies are looking with interest at this novel class of materials, in the hope of addressing the shortcomings that are making scaling of silicon-based electronic devices increasingly difficult.

Despite the diversity in conduction properties and atomic composition, all 2-D materials are composed by covalently bonded in-plane layers that are held together by weak van der Waals interactions to form a 3-D crystal. Each layer has a uniform thickness ranging from 0.3 to 0.7 nm, depending on its atomic structure [e.g., graphene is composed by a single

layer of carbon atoms, while, in general, a transition metal dichalcogenide (TMDC) layer with a general formula of MX_2 is composed by a transition metal (M), sandwiched between two chalcogen atoms (X) [1], [2]. The weak van der Waals interaction is present in the out-of-plane direction thanks to the pristine surface of each layer, i.e., the absence of dangling bonds. This peculiar layered structure, together with certain optical and electrical properties, had already been discovered in 1947 for graphite [1] and in 1969 for TMDCs [2]. However, it was not until the pioneering work of Novoselov *et al.* [3] in 2004 that a graphite monolayer (now commonly known as graphene) was isolated and studied, effectively marking the beginning of 2-D electronics. However, the semimetallic nature of graphene prevents it to be used as the primary vehicle for novel digital transistors and circuits. Among the TMDC materials, the most studied are the ones formed by group IV [zirconium (Zr) and hafnium (Hf)] and group VI metals [molybdenum (Mo) and tungsten (W)]; however, other materials formed by different metals, such as platinum (Pt), rhenium (Re), and tin (Sn), are currently gaining interest. These materials show a semiconducting behavior and have exhibited excellent electrical properties [4]–[6], which will be further analyzed in Section IV. The presence of a sizeable bandgap (1–2 eV) makes the TMDC materials appealing for electronics applications, as it allows us to realize devices with low leakage level and high ON/OFF current ratios [7]–[10]. Among the other remarkable features of TMDCs, their layered structure provides 2-D films of controllable uniform thickness with dangling-bond-free interfaces. Moreover, their extreme thinness and low in-plane dielectric constant alleviate short-channel effects (SCEs) and drain-induced-barrier-lowering (DIBL) [11], [12], which are detrimental to device performance. The high effective mass of charge carriers (especially with respect to III–V materials) helps reducing direct source-to-drain tunneling at ultrascaled dimensions [13], [14], providing a better control of the device’s OFF state by the gate terminals. The 2-D materials are appealing for future VLSI systems, and there are several areas in which they could bring novel functionality in electronic systems. However, several challenges still remain for high-quality growth and successful integration, and experimental demonstrations are still far from simulated devices performances.

This review is organized as follows. In Section II, we focus on the synthesis of 2-D materials. We highlight the different

Manuscript received February 6, 2019; revised March 18, 2019; accepted March 26, 2019. Date of publication June 4, 2019; date of current version June 26, 2019. This work was supported in part by the IMEC beyond CMOS Program and in part by the National Science Foundation (NSF) under Career Award 1751064. (Corresponding author: Giovanni V. Resta.)

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Digital Object Identifier 10.1109/TVLSI.2019.2914609

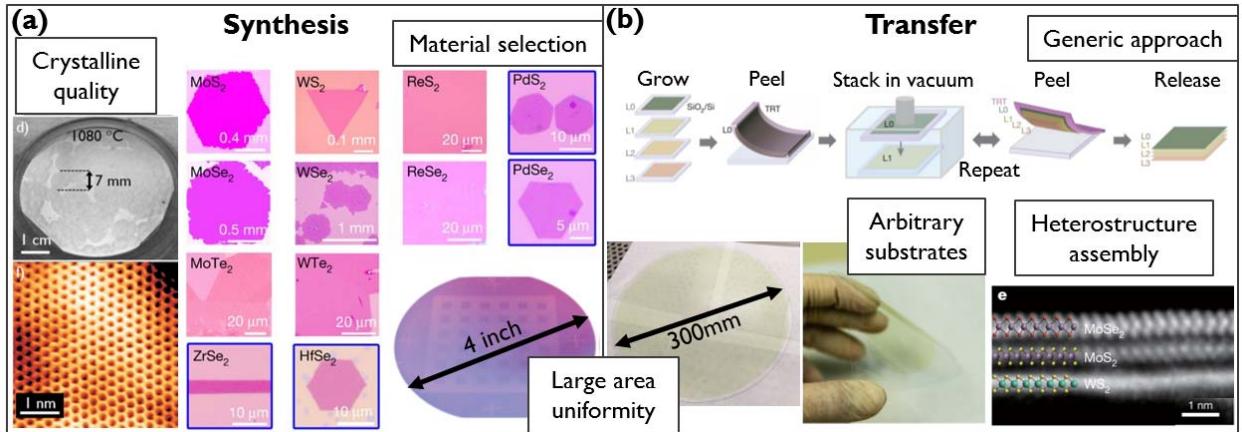


Fig. 1. Relevant metrics for synthesis and transfer of 2-D materials, with notable examples. (a) Crystalline quality and wide material selection are important to allow HP devices. (b) Transfer is a generic approach to allow integration in arbitrary substrates or assembly of heterostructures. Large area uniformity is essential to both synthesis and transfer, allowing low variability in material properties over wafer scales. Adapted with permission from [19], [46], [53], and [61].

approaches that are considered for the growth of 2-D materials on existing CMOS logic, back-end-of-the-line (BEOL) integration and discuss the need for the development of an efficient transfer technique. We also focus on the possibility of synthesizing lateral and vertical heterostructures, which are appealing for TFETs applications. In Section III, we describe a possible roadmap to enable CMOS compatible integration of 2-D materials, highlighting the main challenges that still have to be overcome. In Section IV, we shift our attention to 2-D devices and circuits and review the state of the art. Among the plethora of device concepts, in Section V, we look more closely at 2-D tunnel FETs (TFETs) and negative-capacitance FETs (NC-FETs) for low-power applications. Section VI is dedicated to doping-free polarity-controllable systems that use electrostatic doping to eliminate the need for physical or chemical doping on 2-D materials. Finally, in Section VII, we look at scaling opportunities for different types of 2-D devices and evaluate their projected performances.

II. SYNTHESIS AND TRANSFER OF 2-D MATERIALS

The first works with monolayer (1L) and multilayer (ML) 2-D materials were all done on exfoliated flakes, from the pioneering work on graphene [3], which led to a Nobel Prize in physics, to several demonstrations with TMDCs [7], [21]–[23]. Exfoliation results in high crystalline quality flakes that are used to understand the fundamental properties of 2-D materials and pave the way for future applications. The nondeterministic nature of mechanical exfoliation, however, leads to fluctuations in flake thickness, lateral size, and shape, all of which make it nonscalable to industrial integration. Logic circuits have been demonstrated on such exfoliated flakes [8], [24]–[27], but most reports involve a small number of devices and low level of integration. If the 2-D materials are to be considered for CMOS integration, large-area, uniform, and high-quality synthesis becomes a critical aspect to realize their full potential for future electronic circuits. Fig. 1(a) presents relevant metrics for the synthesis of 2-D materials, namely the crystalline quality (e.g., grain size) of the grown layers, the possibility of growing a number of different materials, and the large area

uniformity of the synthetic layer. Two main synthesis approaches exist: growth directly on the device wafer, as often done for TMDCs [15], [17], [28] or growth on an optimized substrate, with a subsequent transfer step [18], [29] for integration. The optimal growth substrate changes with respect to the 2-D material grown, with metal films being used for graphene [16], [30] and crystalline hexagonal-boron nitride (h-BN) [31], while the best results for TMDCs have been obtained with sapphire [32], [33] or silicon dioxide (SiO_2) [34]. Direct growth is limited by the maximum temperature allowed at different process steps, especially when 2-D materials are cointegrated with a standard silicon CMOS technology. This poses a challenge since heterogeneous cointegration is an attractive route to enhance the speed or functionality of silicon chips by employing 2-D materials alongside the current architectures. Examples include using graphene boosting for interconnects [35], where the presence of a graphene layer leads to a resistivity reduction in the metal line, and 2-D TMDC transistors at the BEOL [36], which can provide further functionalities in the framework of 3-D scaling. However, if the growth is limited to 450 °C, for BEOL integration, the resulting material, particularly 2-D TMDCs, is highly defective [37], [38]. Besides prohibitive temperatures, the growth environment (corrosive or metal-containing gases) can also result in lower reliability of the dielectric where the 2-D material is grown. When using a dedicated growth substrate, the temperature limitations are lifted, which allows this approach to achieve high crystalline quality [30], [32]. The growth substrate can be carefully selected and tailored to achieve the best possible synthesis, where the layer-by-layer growth has been demonstrated with precise atomic control and reduction of strain and doping effects [39]. This temperature flexibility comes at the expense of an additional transfer step, which decouples the growth parameters from the device integration.

A transfer allows high-quality 2-D materials to be integrated into arbitrary substrates, which ranges from the standard CMOS wafers [either in the front-end-of-the-line (FEOL) or BEOL] to flexible substrates [18]. It is, in principle, a generic

approach, which allows the manipulation of different 2-D materials with the same method [20]. In the context of logic devices, uniform transfer of large-area 2-D materials is a key challenge. Transfer relies on interface interactions between the original substrate, the 2-D material, and the target wafer and is normally done using water- or acid-assisted intercalation [18], [29], [40], [41], which can introduce additional contamination or material modifications [42]–[44]. Understanding those interactions and being able to develop an all-dry process is critical to achieve a uniform and repeatable transfer, which has minimal impact on the 2-D material. By alloying 300-mm growth capabilities with standard wafer bonding tools, a uniform 300-mm dry transfer was recently reported [19]. Fig. 1(b) shows the transfer process, with notable examples of each of these metrics. Besides transferring the 2-D materials onto the arbitrary substrates, understanding the interactions present after transfer and the impact of different substrates is also necessary. During this integration step, the 2-D material is transferred onto a wafer surface, which may have a series of nonidealities, such as roughness, topography, surface chemistry, and contamination, among others. Due to the atomically thin nature of the 2-D material, it is highly sensitive to the surrounding environment, and these nonidealities can cause unintentional doping, strain, and increased defectivity. This surface must be carefully engineered in order to result in stable and reliable transistor operation. Current understanding is that other 2-D materials are the ideal substrates for both graphene and TMDC devices [45]–[47].

In the context of transistor fabrication, heterostructures are inherently required, either between TMDCs and conventional metals and oxides or more exotic TMDCs/graphene, TMDC/hBN, or TMDC/TMDC heterostructures [22], [48]. These can be realized by the above-discussed two techniques, either *in situ* growth or assembly *ex situ* through transfer. The growth of vertical (WSe₂/SnS₂, MoS₂/WSe₂, WS₂/MoS₂, graphene/h-BN, and so on) [49]–[52] and lateral (WS₂/WSe₂, WSe₂/MoS₂, WS₂/MoS₂, and so on) [50], [53] heterostructures presents encouraging results. The main figure of merit in terms of heterostructures is atomically abrupt interfaces, being them vertical or lateral, since this is an important requirement for applications such as TFETs (as will be discussed in Section V-A). This is why, the *in situ* growth of the heterostructure—as opposed to assembly *ex situ*—is an appealing technique, where the presence of contamination at the interfaces is drastically reduced. In vertical heterostructures, atomically sharp junctions are easier to be obtained since the layered nature of 2-D materials can be exploited, but achievements in terms of growth techniques also allow atomic stitching of lateral heterostructures [53], [54]. Nonetheless, high-quality, large-area heterostructure assembly has also been demonstrated either by sequential transfers of different 2-D materials one on top of the other [55] or by using a 2-D material to pick up a different one [20], [56]. Beyond their use in established transistor architectures, heterostructures also present interesting physical phenomena, such as tightly bound excitons [57], [58] and plasmonic effects [59], [60], which can lead to exciting new device concepts in the future. These considerations in terms of growth and transfer of 2-D

materials lead us to the discussion of what are the required milestones to be achieved for their integration with current CMOS technology.

III. ROADMAP FOR CMOS INTEGRATION OF 2-D MATERIALS

When evaluating the integration of 2-D materials on future consumer and industrial applications, several different aspects should be taken into account. Their atomically thin nature and high mechanical strength make them perfect candidates for high-performance (HP) flexible electronics [25], [62]. Their electronic structure, featuring an indirect-to-direct bandgap transition, makes them promising for optoelectronics [63]. Their large surface-to-volume ratio results in great potential for both energy storage [64] and sensing [65] applications. While other roadmaps include these aspects [66], [67], we will primarily focus on the potential of 2-D materials as speed and functionality boosters alongside the standard CMOS technology. The 2-D materials are excellent candidates for this application due to their ultrathin and self-passivated nature and for featuring high mobilities even at the atomic thickness limit.

For that to turn into reality, the 2017 IEEE IRDS roadmap [68] mentions several requirements that have to be fulfilled, which are listed in the following. The list combines requirements for graphene (●), TMDCs (○), and shared requirements (◐). We would like to add that some of the aspects IRDS mention only for graphene, are also challenges for TMDCs, and are added as shared requirements in the list in the following:

- bandgap formation with high mobility and ON/OFF ratio greater than 10⁴;
- armchair/zigzag nanoribbon formation;
- improvement of contact resistance;
- ◐ large-area synthesis with low defect density;
- ◐ growth or transfer at controlled locations;
- ◐ growth or transfer of high-quality material at low temperature;
- ◐ thickness control;
- ◐ defect-free contacts;
- ◐ techniques for doping;
- ◐ characterization and imaging of nanoscale structures and composition.

Fig. 2 outlines the core steps of technology development for 2-D material integration in the CMOS process flow. With hundreds of 2-D materials currently being isolated or synthesized, selecting a few materials from this growing library becomes an important integration decision. Material selection can be done by benchmarking against current silicon technology, as proposed by Agarwal *et al.* [69], or can be done based on application-specific requirements, with TFETs being the most attractive example [70]–[72]. Another integration option is to select based on the stability of the material in ambient and process availability of high-quality large-area materials. Irrespective of how to select them, a few materials have already attracted special attention, such as graphene, MoS₂, WS₂, WSe₂, hBN, and black phosphorus (BP).

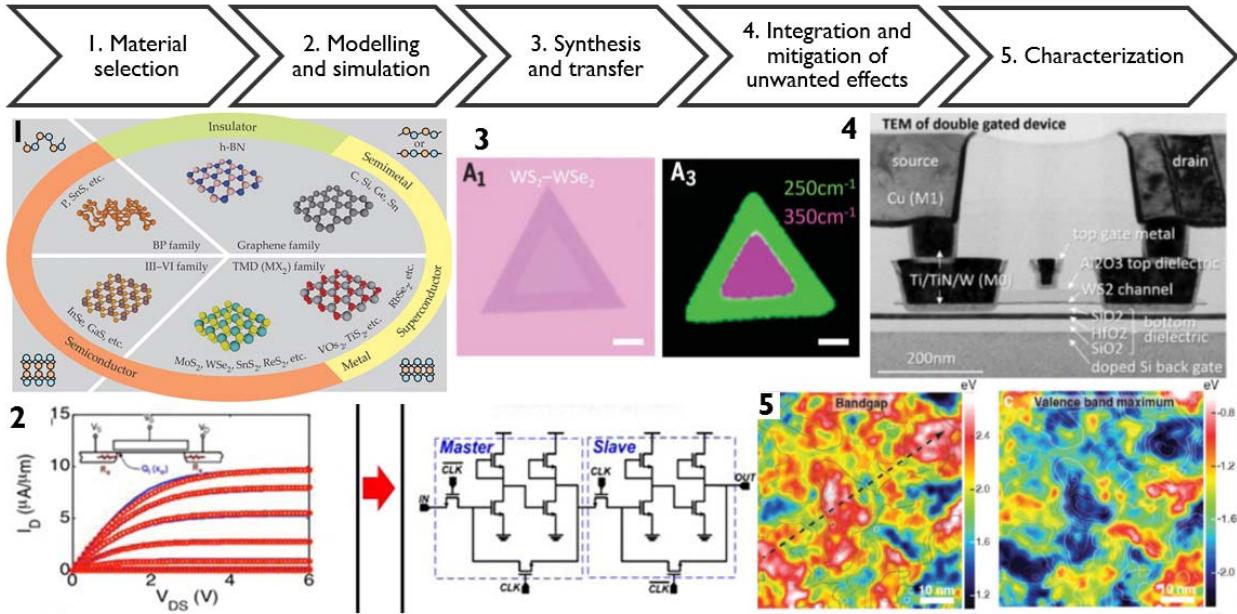


Fig. 2. Roadmap for integration. Required steps for 2-D material to reach product stage. Adapted with permission from [19], [46], [53], and [61].

Modeling and simulation are the key elements for device and circuit design and fabrication and will be the focus of Section VII. Modeling is essential to understand device functionality, especially the impact of contact barriers, non-idealities, and possible defects on the transistor functionality [61], [73]–[77]. Important advances are being made in terms of fundamental material and device understanding [78]–[80]. Besides the important design guidelines that it generates, simulation tools are also essential ingredients if widespread circuit design with 2-D materials is the objective.

Synthesis and transfer, as discussed in detail in Section II, are challenging steps, which aim to achieve materials with a high crystalline quality, controlled thickness at the atomic scale, and integration with arbitrary substrates. Encouraging results from the growth of large single crystals and precise heterostructures [53] were demonstrated. When 300-mm integration is the goal, two main routes for transfer emerge: wafer-to-wafer transfer on complete 300-mm blankets [19] or die-to-wafer positioning of small crystals [81].

The integration core step includes not only upscaling successful lab-scale processes to industry-compatible 300-mm platforms but also mitigating of unwanted effects arising from the substrates, encapsulating materials, or device processing. Some works have explored the mitigation of unintentional doping and strain for both graphene and TMDCs [44]–[46]. Linked to integration is also the precise modulation of the electrical properties of the devices, such as threshold voltage (V_{TH}) and polarity. The precise control of V_{TH} of different devices is critical for correct circuit operation, and while great efforts have been made in the direction of doping, a standard is yet to emerge [82]–[85]. Furthermore, while some materials present either n-type or p-type conduction, others combine both, opening interesting routes for doping-free devices and optimized circuit design, as will be discussed in depth in Section VI [27], [86], [87].

For most of the processing steps, characterization techniques are essential to assure yield and reliability. While the current 2-D material characterization allows probing several properties, wafer-scale characterization techniques are still not matured. The core techniques used to probe 2-D material quality are Raman spectroscopy, photoluminescence, and atomic force microscopy [88]–[90]. Techniques to understand interface properties are not fully developed and are essential for achieving reliable and stable transistors [46], [91], [92]. In-line metrology will surely need additional techniques in order to fully probe the reduced dimensions and specific properties of 2-D materials. Importantly, the development of a standardized characterization toolkit capable of measuring from material quality to electrical characteristics, including interface properties, is still pending.

In conclusion, in order to realize the full potential of 2-D materials as boosters in VLSI systems, several technological developments are still necessarily spanning from a better quality of the grown material to the development of an efficient transfer technique to the upscaling of processes to 300-mm platforms. However, the recent progresses shown in the understanding of the properties of 2-D materials and the demonstration of novel devices keep the research field vibrant and appealing to industries and universities. We now review the state of the art for 2-D electronics based on TMDCs and graphene providing insights into the most recent developments in the field.

IV. STATE OF THE ART FOR EXPERIMENTAL DEVICES AND CIRCUITS

Since the first demonstration of a monolayer (1L) MoS_2 transistor with high mobility and high ON/OFF current ratios in 2011 by Radisavljevic *et al.* [7], the field of 2-D electronics has gained worldwide traction with a constant increase in the number of publications and a growing interest of

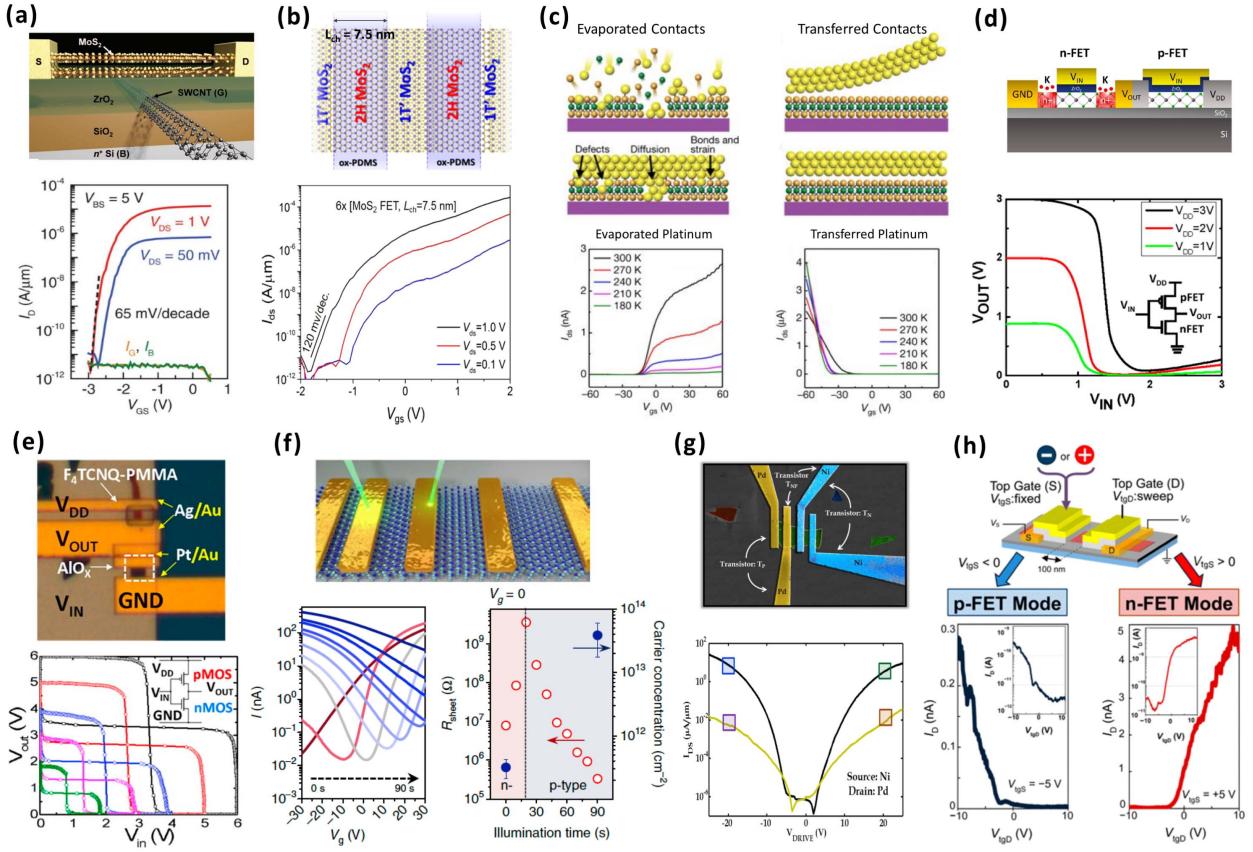


Fig. 3. 2-D-TMDCs electronics. (a) Schematic and transfer characteristics of a MoS₂ device gated by a single carbon nanotube. Adapted with permission from [93]. (b) Ultra-scaled MoS₂ devices fabricated with lithium-induced change to the metallic phase. Adapted with permission from [94]. (c) Comparison between the transfer characteristics of a MoS₂ device with evaporated and transferred platinum contacts. It is shown how p-type conduction can be achieved on MoS₂. Adapted with permission from [95]. (d) Demonstration of a high-gain complementary inverter realized with chemically doped WSe₂. Adapted with permission from [10]. (e) Chemical doping and different metal contacts, providing better carrier injection for electrons and holes, are used to demonstrate logic gates on WSe₂. Adapted with permission from [9]. (f) Light-induced defects on MoTe₂ cause p-type doping of the 2-D semiconductor, allowing for the development of complementary logic. Adapted with permission from [96]. (g) Ambipolar characteristics of WSe₂ device. Adapted with permission from [97]. (h) Ambipolarity and polarity control demonstrated on exfoliated MoTe₂. Adapted with permission from [98].

the research community. Among the several semiconducting materials belonging to the TMDCs family, and introduced in Section II, the most appealing for electronic application have been the molybdenum (MoS₂, MoTe₂, and MoSe₂) and tungsten compounds (WSe₂ and WS₂). These semiconducting materials are stable in the air; even in their monolayer form, they are easily exfoliated from commercially available bulk crystal and are also grown with high quality in a large area, as shown in Section II. Several major results presented in this review have been achieved using mechanical exfoliation, which despite being a nonscalable approach, provides high-quality flakes that are used to gain insights into the properties of these materials and assess their potential for electronic applications. Fig. 3 summarizes a few of the most noticeable experimental results of the last years.

The most known TMDC material, MoS₂, has proven to be a viable solution for the realization of nMOS transistors [7], [8], [24], [99], [100] and ultrascaled devices that have been recently demonstrated [93], [94], [101] [see Fig. 3(a) and (b)]. In Fig. 3(a), a single carbon nanotube was used to gate

a transistor with MoS₂ semiconducting channel, proving the superior electrostatic control achievable thanks to the ultrathin MoS₂ [93]. In Fig. 3(b), the devices with gate length below 10 nm were realized by inducing a change to the metallic phase of the MoS₂, with a lithium solution. This process also allows for a reduced contact resistance thanks to the metallic MoS₂ forming a seamless contact scheme with the 3-D metal [102]. To date, the largest circuit reported on 2-D MoS₂ is a 1-bit microprocessor composed of 115 transistors [103] and other small circuits that have also been previously demonstrated [8], [24], [104]. However, due to considerable difficulties in achieving the p-type behavior in MoS₂ [105], these circuits all adopt a noncomplementary nMOS logic that is not power-efficient. It has recently been demonstrated how p-type conduction can be achieved on MoS₂ by transferring metal contacts on top of the 2-D material, rather than evaporating it [95] [see Fig. 3(c)]. Using this innovative technique, it has been possible to show an almost linear relationship between the work function of the contact metal and the height of the Schottky barrier created, and

p-type conduction has been achieved with high work function metals (such as gold and platinum). Being able to develop both n- and p-type devices on the same semiconducting material is extremely important in order to achieve complementary operation of logic gates and circuits. Conventional CMOS uses ion implantation to physically dope silicon creating low-resistance ohmic contacts and irreversibly setting the polarity of the fabricated device (n- or p-type) according to the dopant atoms used [arsenic (As) for electron and phosphorus (P) for hole doping]. Since a reliable physical doping technique for 2-D materials is still lacking, chemical doping techniques have been explored to both reduce contact resistance (thus achieving higher ON-current) and to achieve complementary behavior [9], [10], [106]–[109]. These techniques are often tailored to the specific material used, in order to dope it against its natural carrier concentration (i.e., n-type doping for WSe₂ and p-type doping for MoS₂ and WS₂). The growth conditions also determine the natural carrier concentration of the materials, and it is important to develop techniques that allow the growth of high-quality, intrinsic materials in order to precisely tune the specific doping processes. Using WSe₂ as a semiconducting material, complementary operations of inverters and other small logic gates have recently been shown [see Fig. 3(d) and (e)] [9], [10]. However, chemical doping is often nonscalable, nonstable, and noncompatible with conventional CMOS fabrication. An innovative doping strategy has recently been proposed where a strong light source is used to locally create defects in MoTe₂ semiconducting channels, which upon oxidation, induce p-type doping [96] [see Fig. 3(f)]. This technique has allowed the realization of photovoltaic cells and bipolar-junction-transistor (BJT) arrays [96]. The possibility of using two different 2-D semiconductors to separately develop n- and p-type transistors has been explored using MoS₂ to fabricate nMOS device and monolithically integrate WSe₂ p-type devices on top, demonstrating complementary logic gates [26]. This paper shows the potential of 2-D materials to be used in combination with CMOS for 3-D monolithic integration. An alternative to the use of either chemical or physical doping is the exploitation of the ambipolar behavior, which refers to the capability of a semiconductor to conduct both charge carriers, and has been shown on several 2-D semiconductors, such as WSe₂, MoTe₂, WS₂, and MoSe₂ [97], [110]–[113] [see Fig. 3(f) and (g)]. A more comprehensive discussion on the opportunity to use the ambipolar behavior to realize doping-free and polarity-controllable devices is presented in Section VI. Other 2-D materials that are being explored include the members of the TMDCs family, such as HfSe₂ and ZrSe₂, which are appealing since their native oxides HfO₂ and ZrO₂ are CMOS compatible [114] and also BP. BP, in particular, has recently drawn considerable interest thanks to its high carrier mobility, ambipolar nature, and lower semiconducting bandgap compared with TMDCs [25], [86], [115]–[117]. However, BP is nonstable in ambient condition and only a few minutes of air exposure results in a complete oxidation of the semiconducting layer, making it challenging to integrate into a conventional VLSI fabrication flow [118]. While TMDCs and BP have been extensively studied for the

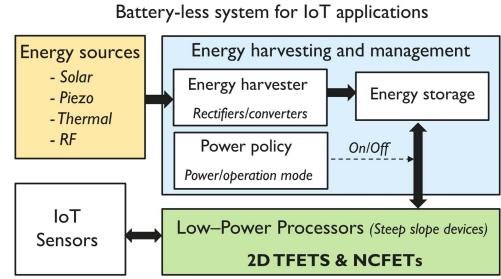


Fig. 4. Overview of self-sustainable energy harvesting system using low-power device alternatives. Adapted with permission from [128].

fabrication of novel electronic devices, graphene, due to its semimetallic nature and high carrier mobility [119], is mainly considered to be integrated into VLSI systems for interconnect scaling [120], [121]. Monolayer graphene is explored as a diffusion barrier and capping layer to extend scaling of conventional copper (Cu) interconnects [120], [122], while multilayer graphene, in the form of nanoribbons, is looked at for entirely replacing Cu interconnects [120], [123]. One of the major challenges is the development of a technique that would allow to grow high-quality graphene at BEOL-compatible temperatures (i.e., 450 °C). Recently, a low-temperature process to grow graphene has been demonstrated, and upon doping with FeCl₃ graphene, nanoribbons of 20-nm width have shown lower resistivity than copper. It is expected that by improving the fabrication process and the doping technique, multilayer graphene interconnects could provide at least 4× smaller circuits delay compared with conventional Cu interconnects [123]. Graphene-based interconnects are also considered for future fully 2-D systems [124] and flexible electronics [125]–[127].

Following this general overview, we focus on the experimental results of two classes of devices: TFETs and NC-FETs for low-power applications and polarity-controllable FETs for doping-free circuits.

V. DEVICES FOR LOW-POWER APPLICATIONS

In order to scale the supply voltage in the future technology nodes and provide more energy-efficient systems, several alternatives based on low-power devices are currently being explored. These devices are targeted to reduce the supply voltage to less than 0.5 V by reducing the subthreshold swing below 60 mV/decade. Besides the benefit in power consumption for VLSI systems, low-power devices could enable the realization of power-efficient processors powered by energy harvesters for the Internet-of-Things (IoT) applications [129]–[131] as well as low-voltage static random access memories (SRAMs) [130]. Fig. 4 shows a schematic of a self-sustainable (batteryless) energy harvesting system, which uses an ambient energy source, such as solar, piezo, or thermal to power a processor based on 2-D TFETs or ferroelectric FETs, also known as NCFETs [128]. NC-FETs are also being explored for energy-efficient nonvolatile computing, which is attractive for IoT applications [131]–[133]. This section

focuses on TFETs and NCFETs, as several promising demonstrations of these classes of devices has been shown using 2-D materials.

A. Tunnel FETs

TFETs are one of the candidates to replace conventional CMOS for future technology nodes due to their low supply voltage (V_{DD}) requirement and steep subthreshold slope (SS) at room temperature. The capability to achieve an SS lower than the thermionic limit of 60 mV/decade is enabled by the different conduction mechanism used in these devices. In fact, current flows in a TFET thanks to quantum mechanical tunneling of carriers from the valence band maximum of one material to the conduction band minimum of the other material, while in conventional MOSFETs, conduction is dominated by drift-diffusion mechanisms [140]. In recent years, much research has focused on realizing TFETs with semiconductors from group IV and group III-V. It has been shown that materials with a direct bandgap, such as III-V compounds, can provide higher ON-currents with lower supply voltage compared with indirect bandgap materials. This is because indirect bandgap materials require additional energy to compensate the mismatch in momentum when tunneling from the conduction band to the valence band [141]. Moreover, the III-V materials have low bandgaps and high electron mobilities, which are ideal for TFET performances. A good figure of merit to measure the TFET current performance is the I_{60} current, which is the maximum current value up to which the SS is less than 60 mV/decade [142]. In order for TFETs to compete with CMOS technology, the minimum required I_{60} current is $10 \mu\text{A}/\mu\text{m}$ and an ON-current of at least $100 \mu\text{A}/\mu\text{m}$ [141]. However, experimental demonstrations are still far away in reaching these goals. The main reason is that the III-V materials show a high density of trap states both at the dielectric and at the heterojunction interface, due to lattice mismatch, which degrades the band-to-band tunneling (BTBT) mechanism [143]–[145]. Therefore, the monolayer 2-D materials appear to be an ideal candidate for the realization of vertical heterostructures TFET applications thanks to their atomically thin body, direct bandgap, and self-passivating surfaces, which could result in no dangling bonds at the interfaces. The advantage of using monolayer 2-D materials is evident from simulation studies [146], [147], and however, it has not yet been confirmed by the experiments, which often use few-layer semiconductors [134]–[139].

Most experimental demonstrations with 2-D materials aim at achieving BTBT by forming a vertical heterostructure between a 2-D n-type material, such as MoS₂, WS₂, ReS₂, and SnSe₂, and a 2-D p-type material, such as WSe₂, MoTe₂, and BP, to create a vertical tunneling heterojunction. Heterostructure stacks, such as MoS₂–WSe₂ [136], [138], [148], [149], MoS₂–MoTe₂ [137], SnSe₂–WSe₂ [149], SnSe₂–MoTe₂, MoS₂–BP [84], [139], [150], have all shown to exhibit BTBT. Fig. 5 displays a summary of the state of the art for 2-D TFETs. Fig. 5(a) shows a MoS₂–BP heterostructure TFET fabricated using an electrolyte gating with a steep subthreshold of 55 mV/decade as shown in the I_D – V_G characteristics [134]. Fig. 5(b) shows an example of

a 3-D–2-D-based tunneling approach where germanium (Ge) is stacked with MoS₂ to form a TFET. The I_D – V_G characteristics of Ge–MoS₂ [135] are shown with an average SS of ~ 31 mV/decade, using electrolyte gating. This SS is the lowest reported on TFETs to date. The advantage of using a bulk material is that they can be easily doped (not yet possible with 2-D materials, as discussed in Section IV) to provide a lower band offset. This, combined with a 2-D material, can lower scattering and provide improved electrostatics, thus leading to more efficient tunneling [151]. However, as previously mentioned, great care needs to be taken when dealing with the 3-D–2-D interface in order to minimize the trap states at the surface. In Fig. 5(c), a MoS₂–WSe₂ TFET is fabricated with two separate gates, enabling the individual modulation of the n- (MoS₂) and p-type (WSe₂) materials [136]. The device also shows SS less than 60 mV/decade. Fig. 5(d) displays a MoS₂–MoTe₂ TFET that also follows the previous approach of having an isolated bottom gate that modulates MoTe₂ and not MoS₂. The I_D – V_G characteristics shows a temperature-dependent BTBT current, which occurs due to the recombination/generation currents present at higher temperatures [137]. The devices in Fig. 5(a) and (b) follow a complete top/bottom gate configuration approach, where both the n- and p-type materials are modulated with the same gate potential. However, a major drawback is that this device architecture can hinder the possibility of creating a broken band alignment for tunneling if the materials are not doped enough. Moreover, the Schottky contacts are also modulated simultaneously with the heterostructure, which can degrade the steep switching in the device [137]. The isolated gate approach shown in the devices in Fig. 5(c) and (d) has the advantage of modulating the carrier concentration of the individual n- and p-type materials when neither of the materials is highly doped. However, this configuration is more complicated to fabricate, as it requires two gates per transistor, which is not ideal in an integration point of view. Fig. 5(e) shows a MoS₂–WSe₂ heterostructure FET that shows the phenomena of negative differential transconductance (NDT) in the I_D – V_G characteristics [152]. A similar behavior is also shown by a MoS₂–BP heterostructure FET [see Fig. 5(f)]. The NDT phenomena can be exploited to create devices that support multivalued logic (MVL), and MVL inverters have been demonstrated with 2-D TFETs [138], [139], [153]. Thanks to the higher number of logic states and higher data storage density, MVL logic is considered to be a promising alternative to traditional binary logic.

Despite the large amount of experimental work on 2-D TFETs, not many devices have shown a steep SS of less than 60 mV/decade. This is possibly due to several factors, such as the presence of Schottky contacts at the metal–semiconductor heterojunction [137], [154], the high density of interface traps at the semiconductor–oxide interface, and the presence of defects and grain boundaries in the 2-D materials. Further experimental work is needed to improve the performances of 2-D TFETs and enable their integration in VLSI systems.

B. Negative Capacitance FET

NCFETs have recently shown promising performances for steep switching application. Unlike TFETs that obtain the

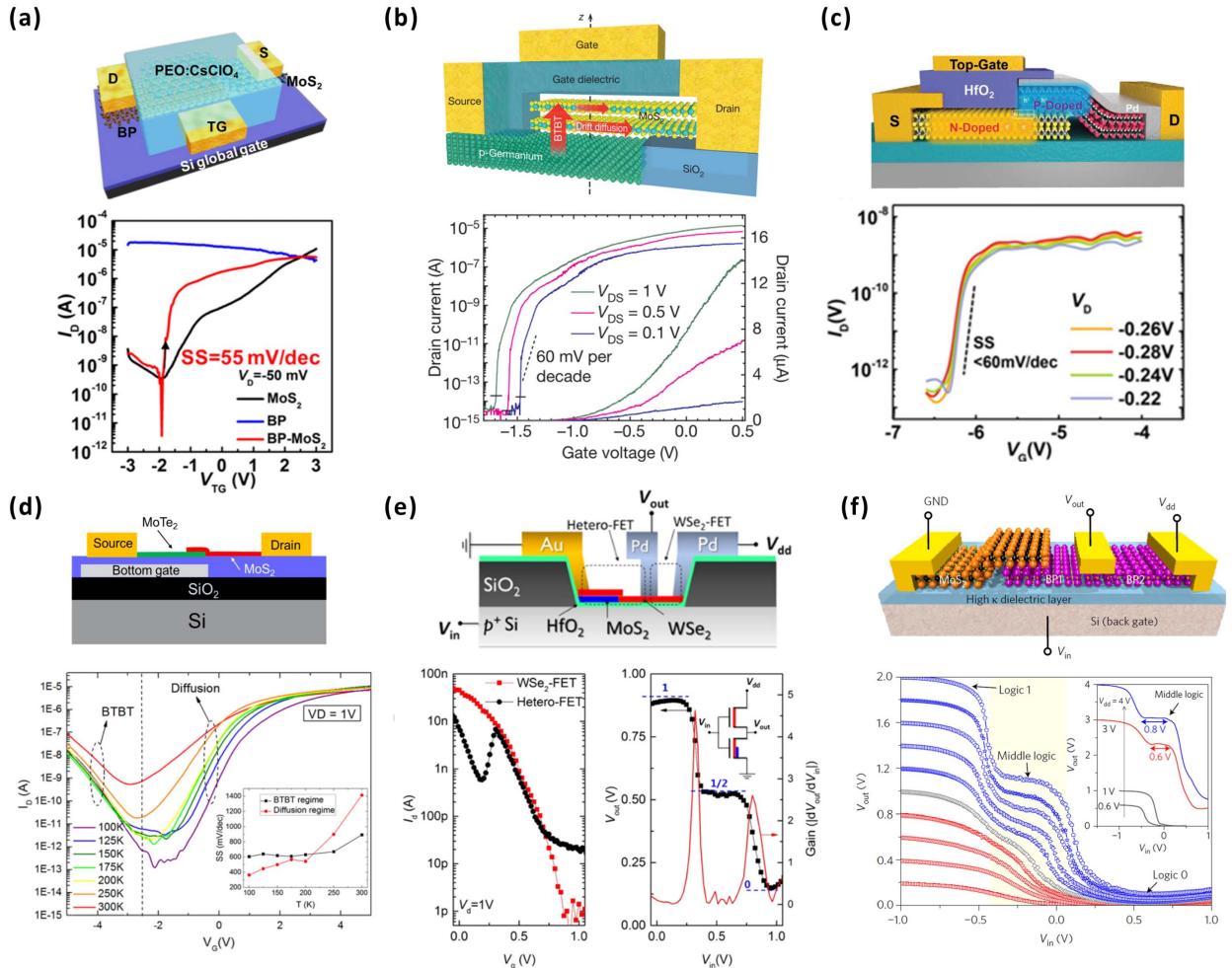


Fig. 5. Experimental results of 2-D TFETs. (a) Device schematic and I_D - V_G curve of a MoS_2 -BP TFET with electrolyte gating. Adapted with permission from [134]. (b) Device schematic and transfer characteristics of a Ge - MoS_2 TFET with electrolyte gating. Adapted with permission from [135]. (c) Schematic cross section and I_D - V_G of a MoS_2 - WSe_2 TFET with an isolated gate device. Adapted with permission from [136]. (d) Schematic cross section and transfer characteristics of MoS_2 - MoTe_2 with an isolated bottom gate. Adapted with permission from [137]. (e) Device cross section and I_D - V_G curve of a MoS_2 - WSe_2 heterostructure FET showing NDT and MVL operation. Adapted with permission from [138]. (f) Device schematic and transfer characteristics of a MoS_2 -BP heterostructure FET showing the NDT behavior for MVL applications. Adapted with permission from [139].

steep switching through BTBT, NCFETs still use a conventional drift-diffusion conduction mechanism, paired with a ferroelectric gate dielectric. The steep switching is enabled by the phenomena of negative capacitance that arises from the transient response of the ferroelectric dielectric while switching its polarization. A ferroelectric material has two polarized states, which can be altered from one state to another with an external applied bias [159]. When the polarization state is switched, the charge-to-voltage variation becomes negative (i.e., $dQ/dV < 0$). This negative value of dQ/dV corresponds to a negative differential capacitance and, when inputted in the equation of SS, results in a subthreshold swing of less than 60 mV/decade [160], [161]. The phenomena of negative capacitance could act as a performance booster for any CMOS platform, even conventional silicon FinFETs, by improving both the SS and the overdrive voltage [162]. However, 2-D materials are attractive for NCFETs because,

thanks to their atomically thin body and high mobility, they can respond to a change in polarization more rapidly than a bulk semiconductor.

Several works have focused on the integration of ferroelectric materials in Si MOSFETs and FinFETs to achieve steep slope performance, and CMOS compatible hafnium-oxide-based ferroelectric materials have been demonstrated by doping hafnium oxide (HfO_2) with materials, such as Al, Zr, and Si. The 2-D-based NCFETs using Al- and Zr-doped HfO_2 and other ferroelectrics have been shown to have steep switching characteristics, as discussed in the following [165]–[167]. Moreover, there has been considerable interest in the recent demonstration of an NCFET using a 2-D semiconductor (MoS_2), as well as a 2-D ferroelectric oxide CuInP_2S_6 (CIPS) [168]. A considerable and yet unsolved issue with NCFET is the large hysteresis that arises from the ferroelectric material while shifting the polarization state [159].

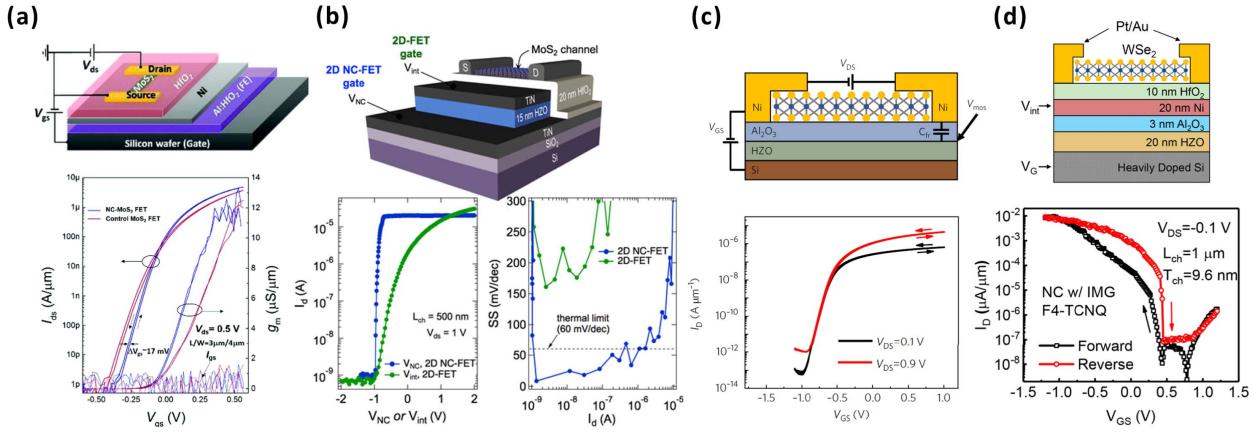


Fig. 6. Experimental studies result in 2-D NCFET based on different 2-D materials and ferroelectric gate-stacks, respectively. (a) Structure schematic and I_D - V_G curve of the MoS₂-NCFET with ferroelectric Al:HfO₂ layer. Adapted with permission from [155]. (b) Structure schematic and I_D - V_G and SS curves of the MoS₂-NCFET with ferroelectric HZO layer. Adapted with permission from [156]. (c) Schematic cross section and I_D - V_G characteristics of the MoS₂-NCFET with ferroelectric HZO without an IMG layer. Adapted with permission from [157]. (d) Schematic cross section and transfer characteristics of the WSe₂-NCFET with ferroelectric HZO layer. Adapted with permission from [158].

TABLE I
STATISTICAL TABLE OF EXPERIMENTAL RESULTS ON 2-D NCFETs

No.	2D Material	FE Material	Type	L_{ch}	V_d	I_{D0}	I_{ON}	lowest SS	ref
2	MoS ₂	HZO	MFMIS	500 nm	1 V	5 μ A	5 μ A	6.07 mV/dec	[157]
3	MoS ₂	Al:HfO ₂	MFMIS	3 μ m	0.5 V	500 pA	13 μ A	57 mV/dec	[156]
4	MoS ₂	P(VDF-TrFE)	MFIS	80 nm	0.5 V	10 nA	100 μ A	24.2 mV/dec	[164]
5	MoS ₂	HZO	MFMIS	1 μ m	0.5 V	100 pA	10 μ A	37.6 mV/dec	[165]
6	MoS ₂	HZO	MFIS	2 μ m	0.5 V	N/A	510 μ A	52.3 mV/dec	[158]
7	WSe ₂	HZO	MFMIS	1 μ m	0.1 V	200 pA	10 nA	14.4 mV/dec	[159]

Hysteresis could be reduced by capacitance matching between the ferroelectric layer and the dielectric layer. This complicated process aims at achieving a single energy minimum in the energy landscape of the resulting capacitor, thus reducing or eliminating hysteresis, which originates from the presence of two energy minimum [169], [170]. Design guidelines for NCFETs should thus include specifications for the tolerance range of hysteresis, as addressed in [171]. However, in order to achieve a steeper SS, most of the NCFETs reported in the literature are fabricated with an internal metal gate (IMG), placed in between the ferroelectric and dielectric layers, which makes capacitance matching impossible to obtain and enhances the hysteretic behavior [172].

Fig. 6 displays several 2-D NCFETs fabricated with and without an IMG architecture. Fig. 6(a) show a schematic cross section of a MoS₂ NCFET using Al:HfO₂ dielectric with a Ni IMG layer [155]. The I_D - V_G curve shows a low hysteresis with a slight improvement in SS of 57 mV/decade compared with when only gating it with regular HfO₂. Fig. 6(b) shows a schematic cross section of a MoS₂ NCFET using hafnium-zirconium oxide (HZO) with a TiN IMG layer [156]. The I_D - V_G curve displays a very steep SS with a minimum of 6.07 mV/decade when compared with the regular MoS₂ FET. However, a very large hysteresis is also observed for this device. Fig. 6(c) shows a MoS₂ NCFET with HZO as

ferroelectric without an IMG layer [157]. The I_D - V_G curve has very small hysteresis, however, with the tradeoff of a larger SS of 52.3 mV/decade. For this device, the drain ON-current level, I_{ON} , of 510 μ A/ μ m is among the highest values for the state-of-the-art steep-slope FETs reported so far [162]. A p-type NCFET using WSe₂ was also realized, as shown in Fig. 6(d), using HZO ferroelectric with an IMG layer [158]. The device dimensions and their performances in terms of SS and ON-current are summarized in Table I, which also includes other notable examples. The devices are separated into two types: metal–ferroelectric–metal–insulator–semiconductor (MFMIS), the devices that contain an IMG layer, and metal–ferroelectric–insulator–semiconductor (MFIS), devices without an IMG layer. The devices have shown to obtain a steep SS below 60 mV/decade and have shown more promise compared with its TFET counterpart. All the reports used exfoliated flakes with a majority of the work done using MoS₂ and different ferroelectric materials.

Apart from improving performances of CMOS devices, NCFETs have other potential applications such as to realize nonvolatile flip-flops for IoT applications, provided that the polarization of the ferroelectric layer is maintained in the absence of gate voltage [132]. As mentioned earlier, the hysteresis created by the introduction of the ferroelectric material is detrimental to purely logic applications, but if properly

controlled could improve the noise margin on SRAMs [131]–[133]. Moreover, the possibility of dynamically tuning the hysteresis to the zero-gate-voltage region will create novel memory-logic interchangeable circuits for energy harvesting processor and in-memory computing [128]. As a final remark, it is worth mentioning that NCFETs offer extensive dynamic voltage–frequency scaling (DVFS) with a wide range of power-performance options [173], [174]. This feature can be applied to low-voltage near-threshold computing, which is an active area of research with attractive energy efficiency [174], [175].

VI. POLARITY-CONTROLLABLE DOPING-FREE DEVICES AND LOGIC GATES ON WSe₂

Physical doping, through ion-implantation, of the 2-D material has not proven to be successful due to the extreme thinness of the 2-D semiconductors, with ions just implanting in the substrate. The possibility of introducing dopant atoms [such as rhenium (Re) or niobium (Nb) for MoS₂] during the growth of the 2-D material has been reported but lacks selectivity and does not allow for any control of the doping profile [176]–[180]. Several chemical and molecular doping techniques have been developed and have been addressed in Section IV, but are often nonstable and non-CMOS compatible, thus not allowing the integration of chemically doped 2-D devices in the BEOL fabrication of VLSI circuits [9], [10], [106]. A device concept that would not rely on any physical doping and use undoped materials would be of great interest in this regard. When no physical or chemical doping is introduced (i.e., an undoped material is used), contact to a 2-D semiconductor usually results in the creation of a Schottky barrier at the source and drain contacts. The height of the Schottky barrier determines the conduction properties of the device, as undoped materials are, for the most part, able to conduct both charge carriers (i.e., electrons and holes). However, it is challenging to control the OFF state of a purely ambipolar device when using a standard gate configuration, as we cannot selectively suppress the conduction of a specific type of charge carriers [97], [181]. In order to control the polarity of the transistor and achieve either n- or p-type behavior, a separate gating of different channel regions is introduced, which we refer to as a double-independent-gate (DIG) structure [181]. A second gate is added, named polarity gate (PG), which induces electrostatic doping at the contact interfaces, effectively modulating the height of the Schottky barrier and blocking the injection either of electrons or holes. It is then possible to dynamically reconfigure the device polarity at run time, by reversing the bias applied to the PG. A conventional gate, named control gate (CG), acts in the central region of the channel and controls the ON/OFF state of the device. The possibility to dynamically flip the polarity at a run time enhances the switching property of the devices, enabling a doping-free, highly flexible design for logic circuits [182], [183]. Differently, from the TFETs and NCFETs, DIG-FETs do not inherently provide a way to reduce the SS below the limit of 60 mV/decade, although steep switching has been demonstrated on Si FinFETs with a DIG structure by increasing

the V_{DS} bias and achieving weak-impact ionization of charge carriers [184]. Moreover, the application of ferroelectric oxide to a DIG-FET structure could be foreseen in future research.

As mentioned in Section IV, ambipolar behavior has been demonstrated on several 2-D materials (such as WSe₂, MoTe₂, MoSe₂, and BP), with WSe₂ showing high mobility and high current densities for both charge carriers [97], [185]–[188]. Recent progress in the demonstration of 2-D polarity-controllable doping-free devices and circuits has been made using WSe₂ and is summarized in Fig. 7 [27]. A schematic cross section of a single device presented in Fig. 7(a), highlighting the position of the PG and CG as well as the semiconducting channel. The devices are fabricated experimentally on exfoliated WSe₂, and the demonstration of polarity control is shown in Fig. 7(b). The inset shows a schematic of the band profile at the source contact for two opposite biases of the PG. When the PG is biased at 0 V or below, holes are favorably injected in the channel, and the device behaves as a p-type transistor. Conversely, the transistor is programmed to function as an n-type device when the PG is biased at 3 V and above. In both cases, the CG is able to turn on and off the transistor when swept between 0 and 2 V. A standard-cell library has been demonstrated with doping-free polarity-controllable WSe₂ FETs, including inverters, NAND, NOR, XOR, and MAJ gates [see Fig. 7(c) and (d)]. The DIG structure results in a more expressive switching function of the fabricated devices, which act as comparison driven switches (i.e., the device is conducting when both PG and CG are controlled by the same logic value). Thanks to the higher expressivity of the single device, it is possible to realize compact two-input XOR gates using only four transistors (not counting the inverters needed to generate the negate signals), while in CMOS, eight transistors would be necessary. The peculiar switching properties of the devices also enable the fabrication of three-input XOR and three-input MAJ gates using only four transistors and replacing the V_{DD} and GND terminals with logic inputs [27], [183]. The possibility of realizing highly compact XOR and MAJ gates has been explored in the synthesis of arithmetic circuits, as will be discussed in Section VII. The challenges that need to be addressed to enable the use of polarity-controllable devices in VLSI systems include achieving a symmetric behavior of the n- and p-type branches by tuning the Schottky barrier at contacts, reach the p-type behavior without the need to apply negative gate voltages (to ensure cascadability), and operate both gate terminals with the same voltage bias.

The experimental results presented in Sections IV–VI show the potential of 2-D semiconductors for several applications, but simulations are essential to better evaluate the strength and weakness of different devices. In Section VII, we will focus on the most recent simulation results for 2-D devices and circuits.

VII. PERFORMANCES PROJECTION AT SCALED DEVICE DIMENSIONS

The need for modeling and simulation to assess the performances of devices and circuits at scaled gate dimensions has already been highlighted in Section III as one of the critical steps in the roadmap. It is crucial to rely on accurate simulations to understand the scaling possibilities for different

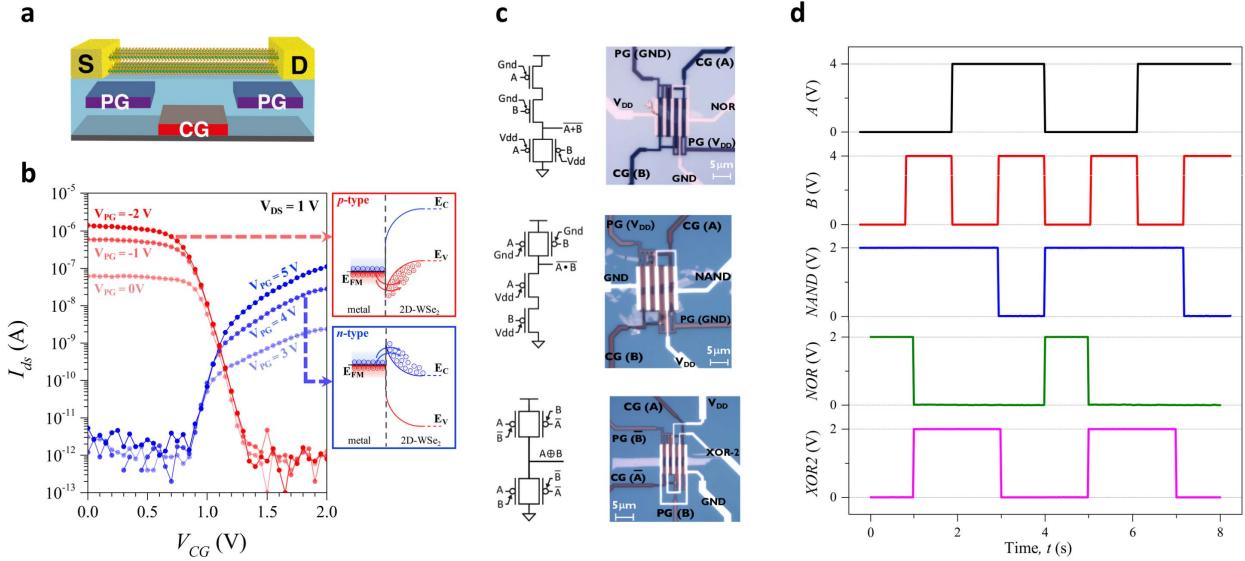


Fig. 7. Polarity-controllable devices and circuits. (a) Schematic cross section of a single polarity-controllable device, highlighting the presence of the PG acting at the contact interface and the CG placed in the central region of the channel. (b) Transfer characteristics demonstrating the polarity-controllable behavior. Insets: schematic of the band-bending at the contacts for n- and p-type operation. (c) Optical micrograph and circuit schematic of the fabricated NAND, NOR, and 2-input XOR gates. (d) Quasi-static measurements of the doping-free logic gates, demonstrating proper operation. Adapted with permission from [27].

classes of devices, compare their performances, and assess the circuit-level metrics. Simulations can also provide guidance and insights into improving the fabrication process (e.g., which metal provides the lowest contact resistance or which oxide lowers the phonon scattering). Scaled device characteristics for 2-D materials have mainly been studied using *ab initio* methods, where the band structure extracted from density functional theory (DFT) calculations [192] is used within a nonequilibrium Green's function (NEGF) formalism to derive the current–voltage characteristics of a device [193]. These simulations require a high computational cost, which increases tremendously when several conduction and valence bands are considered in the Hamiltonian and when scattering is introduced in the simulations [194]. They are also only suitable for single scaled-devices simulations. In order to reduce the computational cost and extend the simulations to circuit level, several compact models have been proposed that are able to reproduce the results of NEGF simulations or experimental results for both doped-FETs and TFETs [61], [73]–[77], [195]. Using these compact models, circuit simulations have shown that 2-D materials could theoretically outperform conventional Si-FinFETs and meet ITRS performance and power-delay requirements [61], [196]–[199]. Fig. 8 summarizes some of the main simulation results both at the device and circuit level for the three main classes of devices presented in this paper (i.e., conventional doped 2-D FETs, 2-D TFETs, and doping-free 2-D FETs). We excluded considerations on NC-FETs, as simulation results on scaled devices and circuit perspectives are still lacking for 2-D materials.

A. Doped 2-D FETs

For 2-D-doped FETs, it was found that a double-gated structure, like the one presented in Fig. 8(a), provides better

electrostatic control on the channel, particularly when using a few-layer 2-D material. Work from Cao *et al.* [11] and Szabó *et al.* [194] have focused on understanding the impact of charge and phonon scattering on the device performances [see Fig. 8(b)] and have shown that even at ultrascaled dimensions ($L_G = 8$ nm), scattering can reduce the ON-current considerably. As mentioned in Section III, the material choice can be guided by several factors, such as intrinsic device performances, fabrication simplicity, or stability of the 2-D materials. In this context, several simulations have been performed to assess which 2-D material can provide the best device performances [11], [12], [14], [69], [189], [198] [see Fig. 8(c)]. Thanks to the possibility of low-temperature integration (see Section II), the performance projection of monolithically integrated 2-D-FETs has been studied [see Fig. 8(d)] [69]. Simulations including parasitic capacitances and wire loads have shown how devices based on WS₂ could be able to provide lower energy-delay-product (EDP) than conventional Fin-FETs in future technology nodes [69]. Moreover, a 32-bit commercial processor core has been implemented with 5-nm design rules to assess the benefits of 2-D transistors on a large-scale design [196]. It was found that 2-D-FETs can provide $\sim 2 \times$ better EDP with respect to conventional Si Fin-FETs, provided that a contact resistivity $< 6 \times 10^{-8} \Omega \cdot \mu\text{m}^2$ can be achieved [196].

B. Two-Dimensional TFETs

The schematic of a scaled TFET exploiting band-to-band vertical tunneling between the p- and n-type semiconductors is shown in Fig. 8(e), with a schematic of the band levels and effective masses [147]. As shown in Fig. 8(f), the transfer characteristics of the scaled devices appear promising reaching

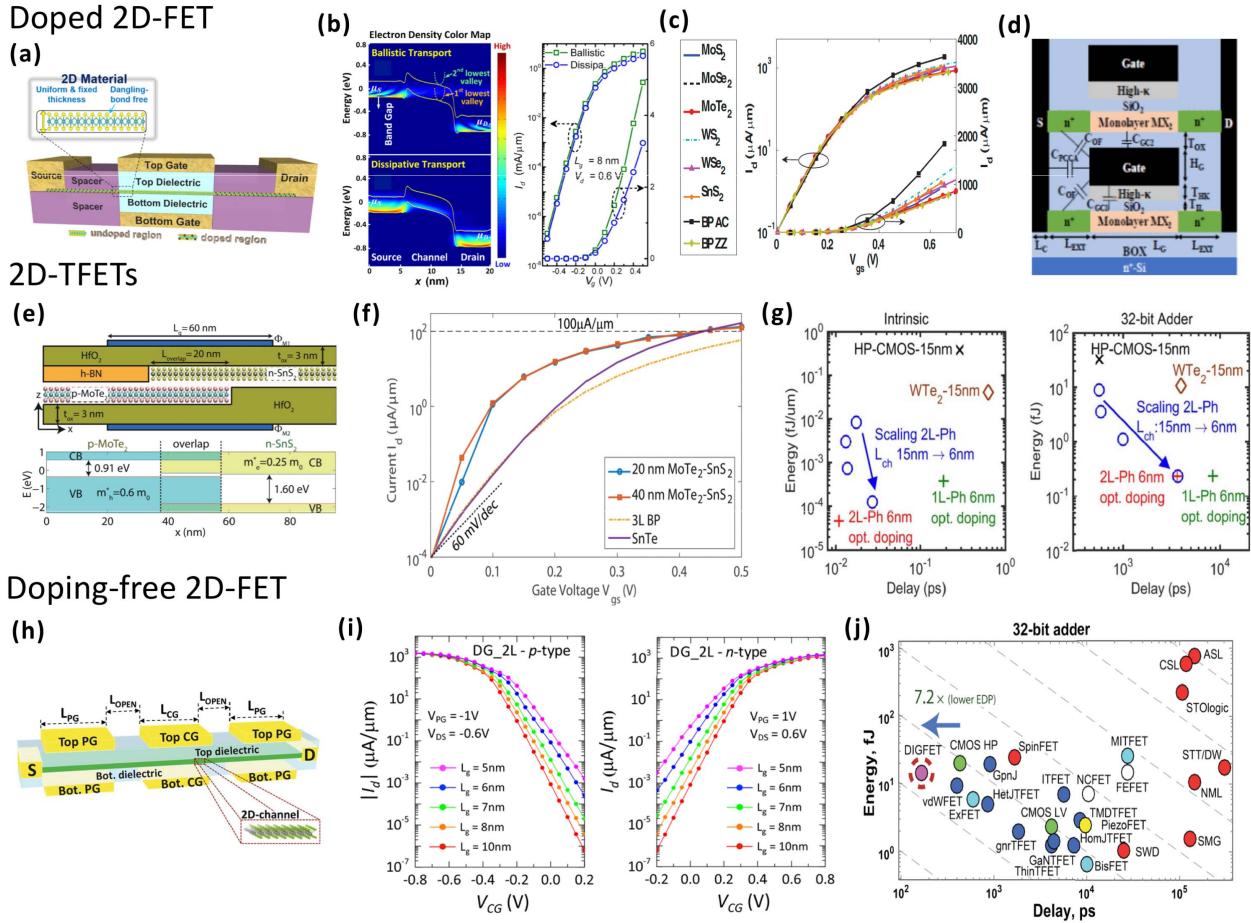


Fig. 8. Device and circuits performances at scaled gate lengths. (a) Schematic of a double-gate FET with doped contacts. Adapted with permission from [11]. (b) Electron density color map and transfer characteristics of the scaled device in (a), showing the effect of scattering on the performances of the device. Adapted with permission from [11]. (c) Comparison of transfer characteristics for several 2-D TMDs and BP. Adapted with permission from [189]. (d) Schematic of a monolithically integrated device with MX₂ materials, which is benchmarked against conventional Si-FinFETs. Adapted with permission from [69]. (e) Schematic cross section of a scaled TFET device, showing the band alignment between the 2-D semiconductors. Adapted with permission from [147]. (f) Transfer characteristics of the device in (e), showing steep-switching and relatively high ON-current. Adapted with permission from [147]. (g) Study of EDP improvements both at the device and circuit level for ultrascaled BP-TFETs. Adapted with permission from [70]. (h) Three-dimensional schematic of the simulated polarity-controllable FET. (i) Transfer characteristics for the device in (h), showing the feasibility of polarity-control down to 5-nm gate lengths. Scaling of the channel length results in an increased SS, as also demonstrated for conventional doped 2-D FETs in [11]. Adapted with permission from [190]. (j) EDP projection for a doping-free 32-bit adder realized with polarity-controllable 2-D FETs. The use of polarity-controllable 2-D FETs enables a $\sim 7 \times$ reduction in EDP compared with scaled high-power CMOS. Adapted with permission from [191].

ON-currents greater than $100 \mu\text{A}/\mu\text{m}$, having I_{60} of around $1 \mu\text{A}/\mu\text{m}$ and operating at $V_{DD} = 0.5 \text{ V}$. These simulations, although promising, are purely ballistic and not include any effect arising from scattering, traps, and contact resistance. Other works have also investigated different geometries and materials [200]–[203]. As mentioned in Section V-A, BP is a material that has attracted considerable interest for TFETs applications and its performances at scaled gate lengths have been studied [70] and compared with HP CMOS as in the beyond-CMOS benchmark [204]. In this case, no heterostructure is used, and the BTBT takes place between the heavily p-doped source and the heavily n-doped drain through an intrinsic region (i.e., p-i-n structure). Fig. 8(g) shows the switching energy versus delay, a metric usually referred to as EDP, for the intrinsic-scaled BP device and for a 32-bit

adder realized with 2-D-BP TFETs. It is demonstrated that thanks to the lower bandgap, two-layer BP with asymmetric optimized doping provides the most improvement in EDP. However, it can be seen in Fig. 8(g) that the improvements in EDP for the 32-bit adder are not as significant as the ones promised by the intrinsic EDP. This is caused by the presence of interconnects and parasitic capacitances that in the sub-10-nm regime, play a critical role.

C. Doping-Free 2-D FETs

Using a scaled double-gated structure [see Fig. 8(h)], the performances of doping-free polarity-controllable devices have been assessed by Resta *et al.* [190], using ballistic NEGF simulations. It is shown how a reduction in the semiconducting

bandgap of the MX_2 material can lead to higher ON-currents (due to improved tunneling at the contacts) while maintaining sufficient electrostatic control over the channel to switch OFF the device [see Fig. 8(i)]. For a bandgap energy of 0.8 eV, I_{ON} can reach $1.5 \mu\text{A}/\mu\text{m}$ while keeping I_{OFF} well below $10^{-2} \mu\text{A}/\mu\text{m}$ down to $L_G = 5 \text{ nm}$. It should be noted that due to the presence of the additional PG, the device is inherently longer than conventional CMOS, and in the simulations, L_G refers to the length of each gated segments, either PG or CG. The simulated transfer characteristics for the 10-nm devices were used to simulate the behavior of a 32-bit adder, which is compared with other technologies of beyond-CMOS benchmark [191], [204]. Thanks to the design possibilities enabled by polarity-controllable devices at the circuit level, e.g., compact realization of XOR and MAJ gates, it was shown a $\sim 7 \times$ reduction in EDP when compared with HP-scaled CMOS [see Fig. 8(j)] [191].

VIII. CONCLUSION AND PERSPECTIVE

This paper provided an overview of the recent advances in the field of 2-D electronics. We focused on three different aspects that are equally relevant for the evaluation of novel materials and novel devices considering material synthesis, experimental demonstrations, and projections of device performances. We looked closely at novel device concepts that have been boosted by the properties of 2-D materials, such as TFETs and NCFETs, and also provided an in-depth discussion on the possibility of realizing doping-free 2-D circuits thanks to polarity-controllable devices. For each topic, we highlighted the main achievements and evaluated the challenges that remain to be addressed in order to integrate the 2-D materials into VLSI systems. This integration can be fruitful by using 2-D materials alongside the conventional silicon CMOS, where we could envision 2-D-TMDCs circuits being integrated into the BEOL and graphene nanoribbons used for interconnects. A truly 3-D nanosystem could be enabled by the low-temperature fabrication process required by the 2-D materials leading to dense 3-D monolithic VLSI systems, integrating memory, sensors, and electronics on the same chip [205]. Moreover, the foldable and flexible electronic system could also greatly benefit from the thinness and large area availability of 2-D materials, enabling more compact and portable wearable systems.

ACKNOWLEDGMENT

The authors would like to thank C. Huyghebaert and F. Catthoor for their valuable input and stimulating discussions.

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