

11-Surfaces, Interfaces and Thin Films

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PS-11.01.28 AN INVESTIGATION ON (100) MONOEPI-
TAXIAL DIAMOND MONOCRYSTALLINE FILMS. By Z.M.
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For diamond's excellent microelectrical prop-
erties and the rapid development of CVD tech-
niques, the diamond high temperature semicon-
ductor materials have been received increasing
attention. In our work, we have obtained quite
smooth homoepitaxial monocrystalline films de-
posited on (100) oriented HT-HP synthesized
diamond by microwave PCVD method. The microwa-
ve frequency was 2.45 GHz, internal dimension
of the quartz horizontal reaction tube was 35mm,
the reaction gases were acetone and hydrogen,
their flow rate was 100 SCCM and pressure 4-6.5
KPa. Microwave power was about 300W. For the
characterization, SEM, Raman Spectroscopy, Dou-
ble Crystal Diffraction and HEED were used. The
results show that the acetone concentration di-
rectly affected the morphology and structure of
the films, as well as the growth process. For
lower concentration (0.9 Vol%), the growth pat-
tern was essentially in line with the spiral
dislocation growth mechanism, when raised to
1.7 Vol%, the epitaxial surface became more
smooth, but the accumulation of defects in grown
film led to the formation and growth of parasitic
nuclei, and with higher concentration (2.6
Vol%), the cellular structure was locally formed.
In addition, it is the advantage to enhance two-
dimension growth and depress parasitic nucleation
if argon was added, in this case, quite smooth
monocrystal epitaxial layers formed.

PS-11.01.29 APPLICATION OF THE X-RAY STANDING
WAVE TECHNIQUE TO ANALYSIS OF
HETEROINTERFACES WITH MULTI-SITE OCCUPANCY

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The X-ray Standing Wave (XSW) technique
is a powerful structural tool to study
surfaces and interfaces. The intensity
modulation of events excited by the
electromagnetic field when Bragg diffraction
takes place, allows the knowledge of both
the modulus (named F) and the phase (named
P) of an XSW structure factor relative to
the excited atoms whose radiative decay is
measured as a function of the Bragg angle.
If the atoms under study occupy one well
defined position relative to the diffracting
planes, the measured P and F represent
respectively the occupied position and the
fraction of the total atoms which occupy
that position, (1-P) being randomly
distributed. However, often several sites
are occupied at the same time and the
resulting P and F are functions of the
individual positions p_i 's and the relative
occupancies f_i 's.

In this contribution we will present some
examples of X-ray Standing Wave studies of
metal/semiconductor
and semiconductor/semiconductor
heterointerfaces where multi-site occupancy
took place. It will be shown that under
appropriate assumptions the XSW can
accurately describe the structure even when
several sites are involved. This is the
case, for example, of the alkali metal/Si
interface. In the (111) orientation seven
different sites have been considered. With
two reflections and the assumption of a
covalent bond the distribution of the alkali
metal atoms into the seven sites have been
determined.
Other examples will concern with III-V
compounds and Si/Ge heterointerfaces.

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X-RAY STANDING WAVE STUDY OF THE
FE/SI(111) INTERFACE. By B.Capelle*, A.Taccoen, J.-C.
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At room temperature, α -iron grows on the (111) surfaces of
silicon with the following epitaxial relations and with an
abrupt interface:

$$\text{Fe}(111)\parallel\text{Si}(111)$$

$$\text{Fe}[1\bar{1}0]\parallel\text{Si}[1\bar{1}0]$$

$$\text{Fe}[11\bar{2}]\parallel\text{Si}[11\bar{2}]$$

The growth of the first three monolayers ($1\text{ML}=0,78 \cdot 10^{14}$
 at/cm^2) of iron on silicon has been studied by the X-Ray
Standing Wave (XSW) method using the synchrotron
radiation at the Laboratoire pour l'Utilisation du
Rayonnement Electromagnétique (Orsay, France). Samples
were elaborated in ultra high vacuum at room
temperature. As XSW experiments occurred at air, iron
layers were covered by amorphous silicon to avoid any
oxidation. For this study, new monolithic monochromators
with four reflections (111, 220 and 004) were developed and
inclined reflections (11 $\bar{1}$, 220 and 004) have been used
(Boulliard J.C, Capelle B., D. Ferret, A. Lifchitz, C.
Malgrange, J.-F. Pétroff, A.Taccoen and Y.-L. Zheng,
J.Phys.I France, 2, 1992, 1215-1232).

The following growth model fits well with the XSW
experiments: iron atoms of the first layer adsorbate onto the
 T_4 sites with a contraction of the surface of the silicon (5 to
6%). Distance between iron atoms of the first layer and the
nearest silicon atoms is between 2,30 and 2,40 Å. The three
first layers grow with islands whose height is limited to
three layers and which are completed before a fourth layer
appears.