## MATRL 100A: Structure and Properties I, assignment 4

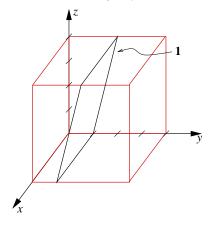
This is a practice midterm exam. In the exam, you would be allowed a simple periodic table (not with the configurations) and a calculator.

Each question carries ten points. Please read the questions carefully.

- 1. Iridium (Ir) is one of the heaviest elements, with a density  $\rho = 22.55$  g cm<sup>-3</sup>. The cubic unit cell parameter of Ir is a = 3.84 Å.
  - (a) Use the data to determine the number of Ir atoms in the unit cell. [7]
  - (b) Is the cell simple (*P*) cubic, body-centered (*I*) cubic, or face-centered (*F*) cubic? [3]

(You need the atomic weight of Ir, which is 192.22 and the Avogadro number is  $N_{\rm A} = 6.022 \times 10^{23}\,{\rm mol}^{-1}$ .)

2. In the following object:



- (a) Identify the Miller plane 1. [4]
- (b) Using the same coordinate system, sketch the  $(1 \bar{1} 0)$  and (2 2 2) Miller planes. [6]
- 3. The *I*-cubic (bcc) structure:
  - (a) Sketch this structure [3]

[3]

- (b) How many neighbors does each atom have, and what is the distance between an atom and its nearest neighbor in terms of the cell parameter a?
- (c) How many next-near neighbors does an atom in this structure have? What is the distance from an atom to its next-nearest neighbor in terms of a? [4]

## 4. Bonding:

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| (a) | What kind of bond would you find in the hydrogen molecule $(H_2)$ ? How is the noble gas configuration achieved and which is the noble gas? | [2] |
|-----|---|-----|
| (b) | What kind of bond is found in LiH (lithium hydride). How is the noble gas configuration achieved? Which is the noble gas or gases?          | [2] |
| (c) | What kind of bonding would you find between the atoms in (i) solid Ar (ii) HCl (iii) NaCl   | [3] |
| (d) | Write the electronic configuration of the following species: (i) nitrogen in $Na_3N$ (ii) phosphorous in $PF_3$ (iii) $S$ in $Na_2S$        | [3] |

## Table of electronegativities:

| Н   | Li  | N   | F   | Na  | P   | S   | Cl  | Ar |
|-----|-----|-----|-----|-----|-----|-----|-----|----|
| 2.1 | 1.0 | 3.0 | 4.0 | 0.9 | 2.1 | 2.5 | 3.0 | _  |

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