

VEIN BATCH 2027

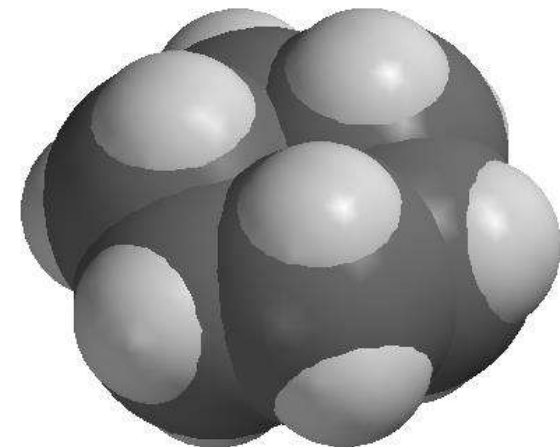
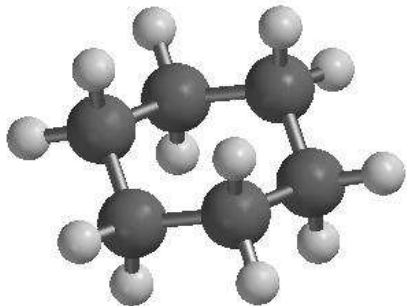


Sub: Organic المادة:

Lecture: 2 المحاضرة:

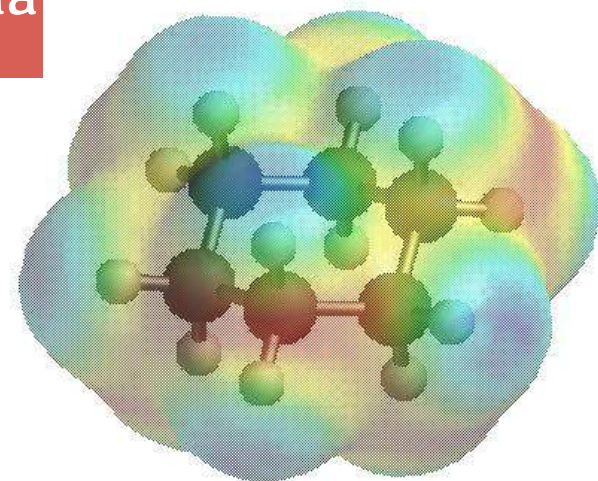
By: Johainah Taha إعداد:

Edited: تعديل:



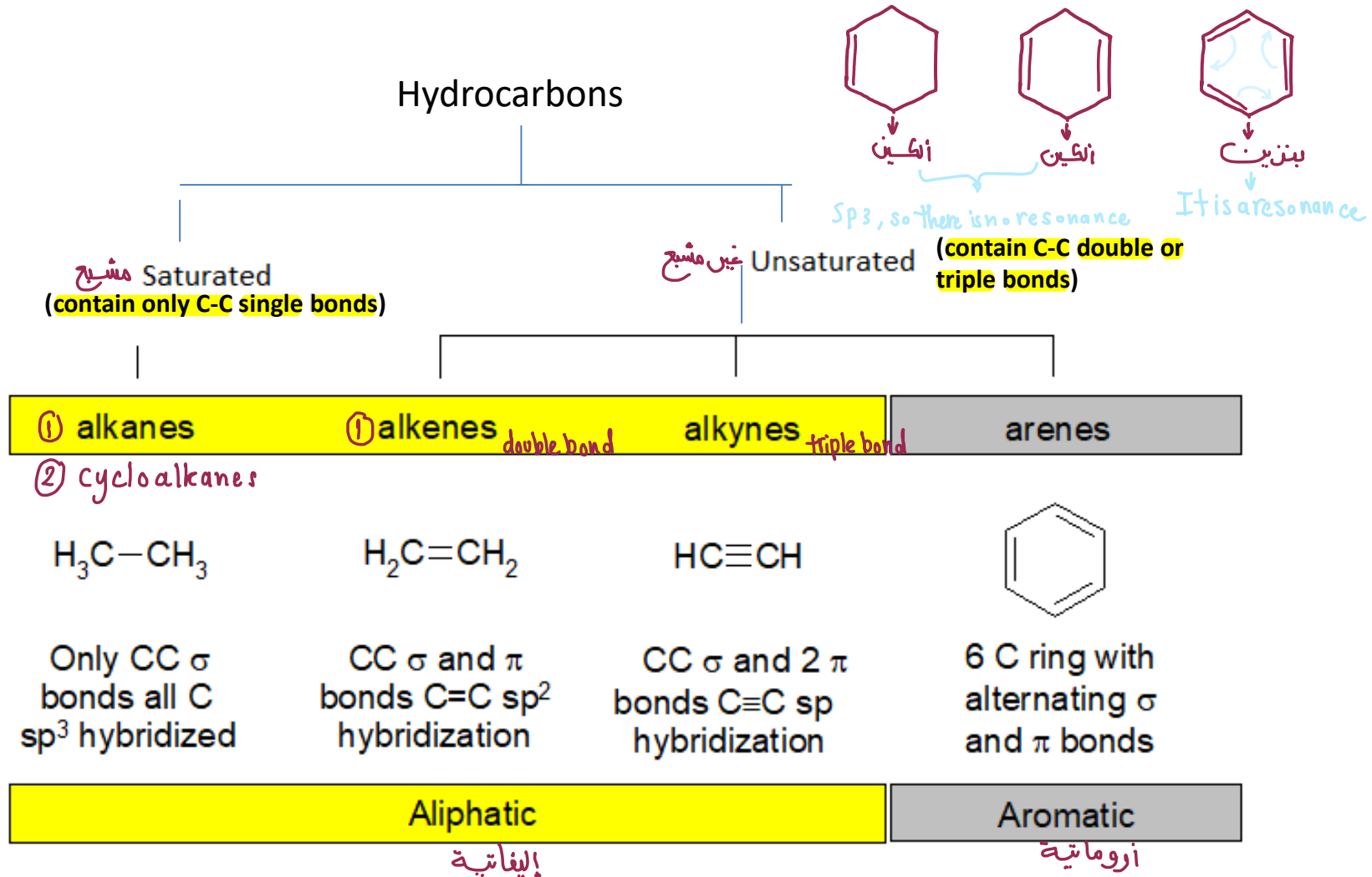
Chapter 2: Alkanes and Cycloalkanes: Conformational and Geometric Isomers

Done by : Johainah Taha



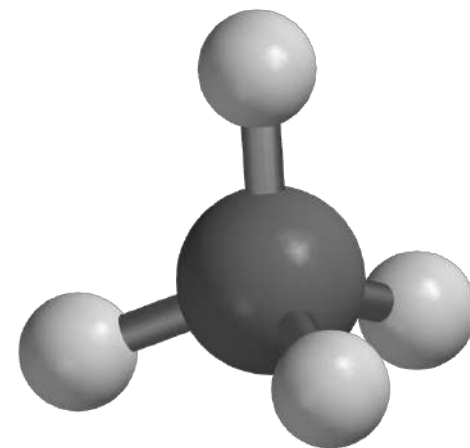
Types of Hydrocarbons

Hydrocarbons are compounds that only contain C and H atoms. سؤال؟ شو الفرق بين البنزين والالكين؟

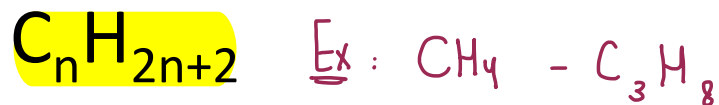


Structure of Alkanes

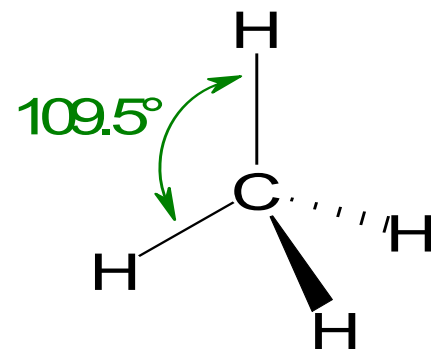
Alkanes are **saturated** hydrocarbons, that is they contain the **maximum** number of H atoms possible for the **number of C atoms present**.



The generic formula for an alkane is:



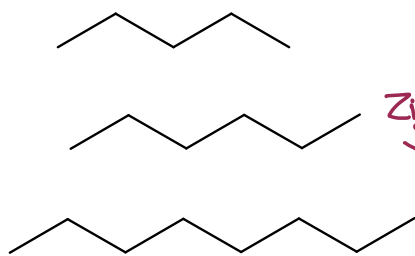
This means every C atom is **sp³** hybridized with bond angles of **~109.5°**



*Tetrahedral molecular geometry.

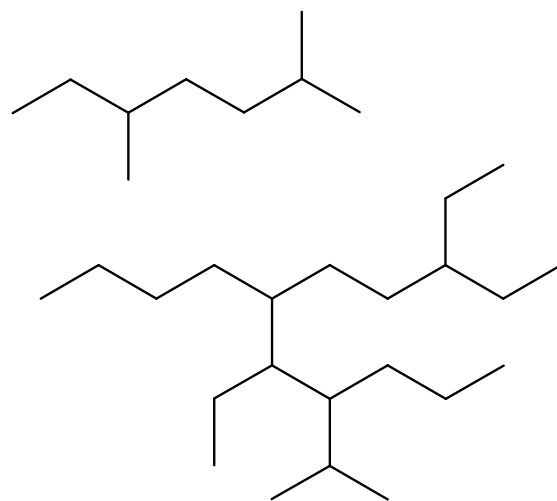
Structure of Alkanes (cont'd)

Alkanes can come in two forms, normal (or linear) and branched. Branched alkanes can have branched branches. Therefore the number of isomers possible grows quickly



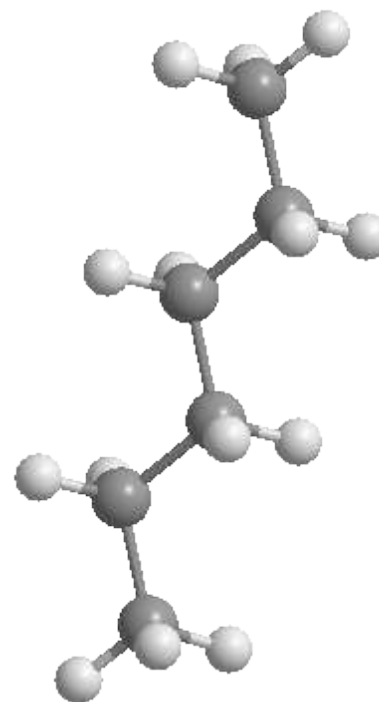
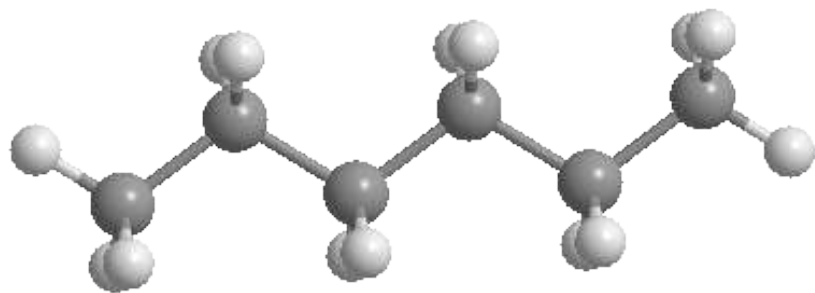
• الرسمة بتكون Zig-Zag
وهي فعليا تمثل الشكل
الحقيقي لرضه المرادفات

Normal, unbranched (or linear)



branched

Structure of Alkanes (cont'd)



التسمية Nomenclature of Organic Compounds

A. **Common names** In the early days of organic chemistry, each new compound was given a name that was usually based on its source or use.

B. IUPAC

The IUPAC name of any compound contains **3 parts** :

IUPAC name = ^③ Prefix + ^① parent + ^② suffix

Prefix : What and where substituents ^{التفرعات} → third step.

Parent (Root) : longest chain → first step

Suffix : functional group , whether it is alcohol, or alkyl etc. → Second name

Table 2.1 ■ Names and Formulas of the First Ten Unbranched Alkanes

Name	Number of carbons	Molecular formula	Structural formula	Number of structural isomers
¹ <u>methane</u> ^{alkan}	1	CH ₄	CH ₄	1
² <u>ethane</u> ^{Alkan}	2	C ₂ H ₆	CH ₃ CH ₃	1
³ <u>propane</u> ^{Alkan}	3	C ₃ H ₈	CH ₃ CH ₂ CH ₃	1
butane	4	C ₄ H ₁₀	CH ₃ CH ₂ CH ₂ CH ₃	2
pentane	5	C ₅ H ₁₂	CH ₃ (CH ₂) ₃ CH ₃	3
hexane	6	C ₆ H ₁₄	CH ₃ (CH ₂) ₄ CH ₃	5
heptane	7	C ₇ H ₁₆	CH ₃ (CH ₂) ₅ CH ₃	9
octane	8	C ₈ H ₁₈	CH ₃ (CH ₂) ₆ CH ₃	18
nonane	9	C ₉ H ₂₀	CH ₃ (CH ₂) ₇ CH ₃	35
decane	10	C ₁₀ H ₂₂	CH ₃ (CH ₂) ₈ CH ₃	75

حفظ
الاسماء

Note: The repeating group is (-CH₂-) is Methylene group

طريقة سهلة لحفظهم كنت أتبعها بالمدرسة فه

ميتان و إيثان بنات العم ، تزوجوا بروبان وسكنهم

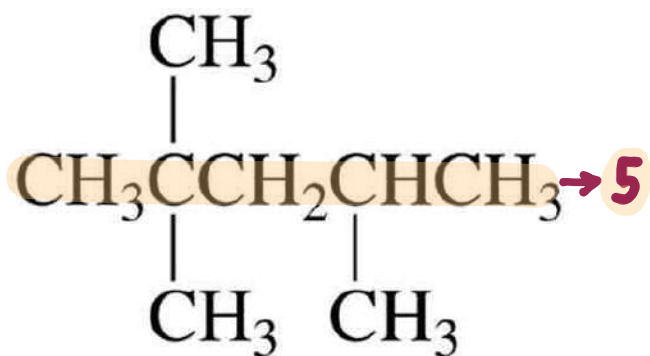
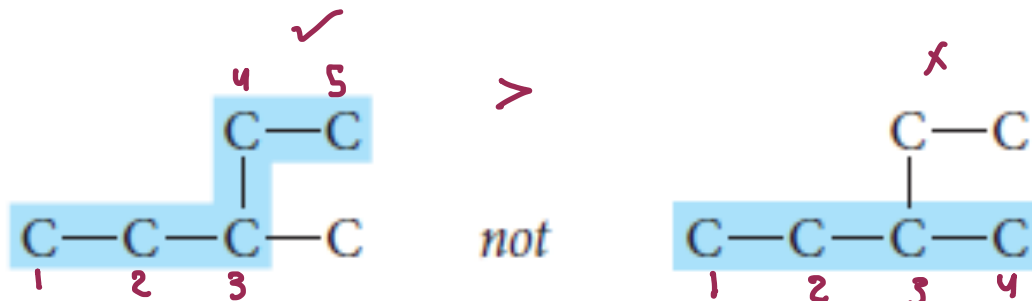
في بيوتان ، وأنجبتا بنتان ، الأولى اسمها هكسان والثانية

هبتان ، ولما كبروا تزوجتا أوكتان و نونان ، ومهرهم

كان ديكان

IUPAC Rules for Naming Alkanes

1. First identify the longest continuous chain (parent name)

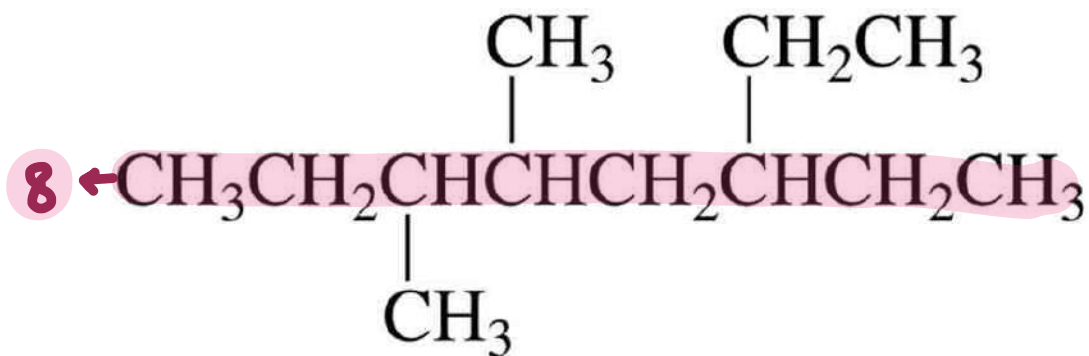


2,2,4-trimethylpentane

not

2,4,4-trimethylpentane

because $2 < 4$



6-ethyl-3,4-dimethyloctane

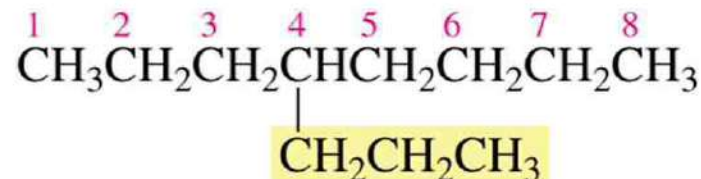
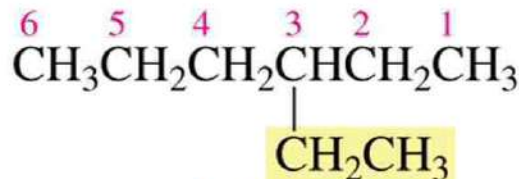
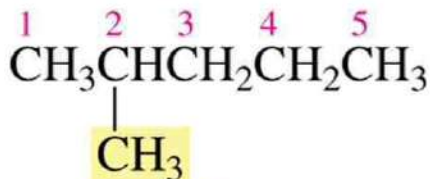
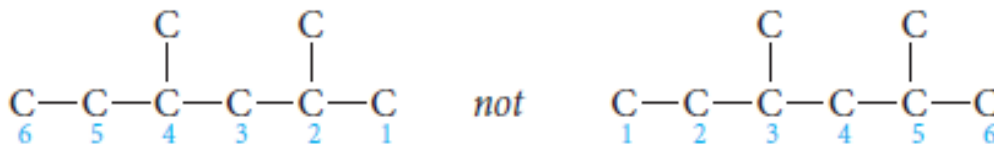
not

3-ethyl-5,6-dimethyloctane

because $4 < 5$

2. Number the chain in the direction that gives the substituent as low a number as possible

نختار الأقرب الى
التفرع

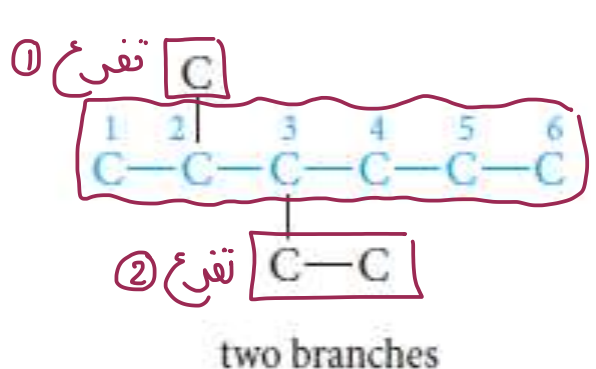


②-methylpentane
not
④-methylpentane

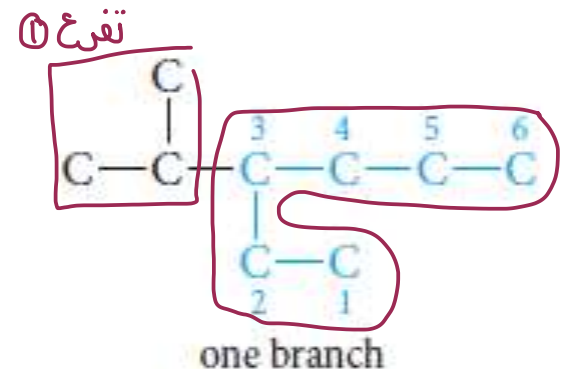
③-ethylhexane
not
④-ethylhexane

④-propyloctane
not
⑤-propyloctane

Note: If there are two equally long continuous chains, select the one with the most branches. For example:



not



* $\text{CH}_4 \rightarrow$ methan

$\text{CH}_3 \rightarrow$ methyl

* $\text{C}_2\text{H}_6 \rightarrow$ Ethan

$\text{C}_2\text{H}_5 \rightarrow$ Ethyl

* $\text{C}_3\text{H}_8 \rightarrow$ prop an

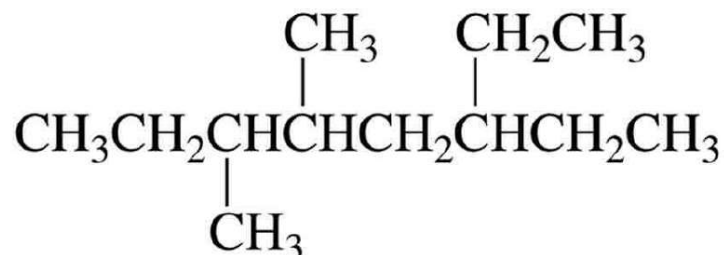
$\text{C}_3\text{H}_7 \rightarrow$ propyl

etc ...

If there is a branch equidistant from each end of the longest chain, begin numbering nearest to a third branch



- **How to name a substituent?**
- **It can be a branch in the chain**
- **A branch (alkyl substituent): is (Alkane - H)**



Replace "ane" of alkane with "yl."

CH_3-
methyl group

CH_3CH_2-
ethyl group

$\text{CH}_3\text{CH}_2\text{CH}_2-$
propyl group

$\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2-$
butyl group

$\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2-$
pentyl group

$\text{R}-$
any alkyl group

-There are two propyl Groups

a secondary carbon

a primary carbon



remove a hydrogen



normal ← n-propyl group

→ -Pr رمزها

↓
linear صلت

حفظ



isopropyl group

→ i-Pr- رمزها

↓

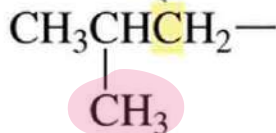
-There are Four Butyl Groups

primary carbon



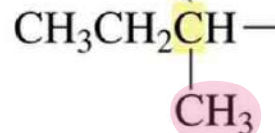
butyl group

primary carbon



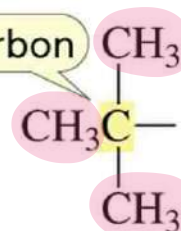
isobutyl group

secondary carbon



sec-butyl group

tertiary carbon

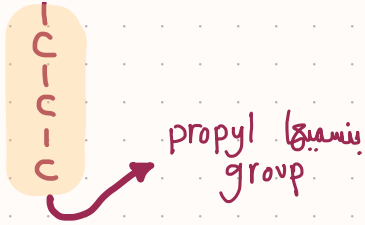


tert-butyl group

Note: These names for the alkyl groups with up to four carbon atoms are very commonly used, so you should memorize them.

شرح مني ت : مثل ما حكينا البروبيل هو C_3H_7 يعني ٣ كربونات ، والبروبيل انه 2 Shapes :-

مكان ارتباطه بالتركيب هما جان سواء اكان هكسان ، اوكتان ... الخ



مثال :-



الشكل الاول

اسمه :- propyl group

رمزها :- -pr



مثال :-



الشكل الثاني

اسمه :- isopropyl group

رمزها :- -i-pr

الملاحضه

أشكال البروبيل

iso propyl



ارتباط من هون

normal propyl



ارتباط من هون

أما بالنسبة لـ البيوتال فهو C_4H_9 ، وفيه 4 Shapes

الشكل الأول



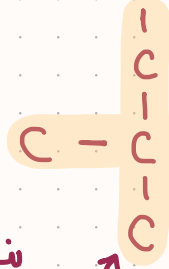
butyl group

الشكل الثالث



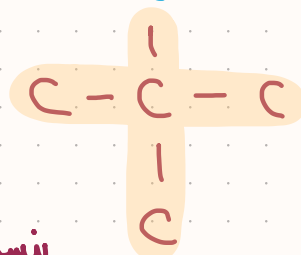
sec-butyl group

الشكل الثاني



isobutyl group

الشكل الرابع

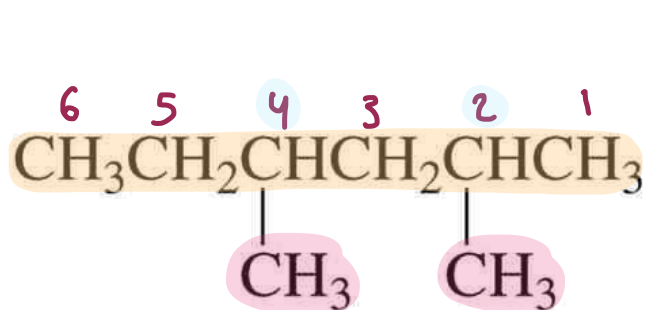


tert-butyl group

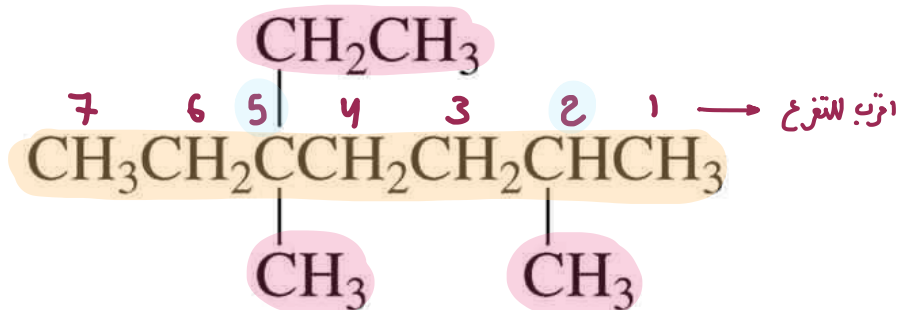
Examples:

Note 1: di-, tri-, and tetra- are used for the same **Multiple Substituents**

Note 2: Substituents are listed in alphabetical order.
(di, tri, tetra, sec, tert are not alphabetized)

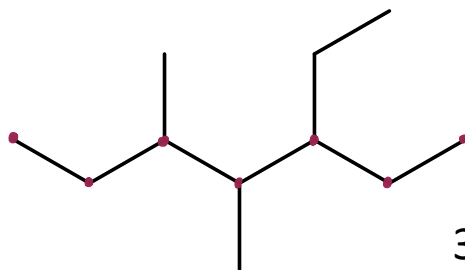


2,4-dimethylhexane

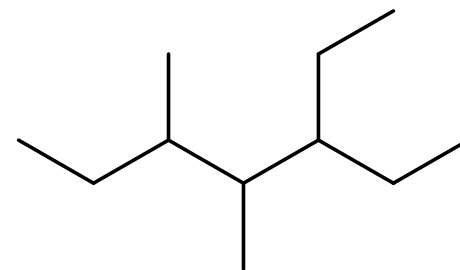


5-ethyl-2,5-dimethylheptane

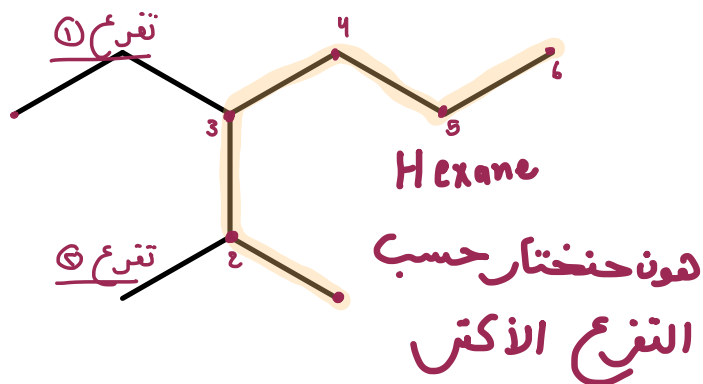
رتبناهم حسب الأعراف (e قبل m)
numbers are separated by a comma;
a number and a word are separated by a hyphen (-)



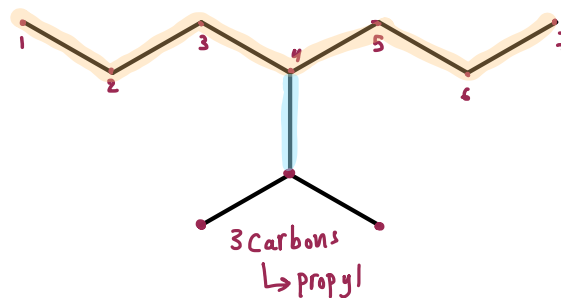
3-ethyl-4,5-dimethylheptane



Examples



3-ethyl-2-methylhexane



4-isopropylheptane

اذا كان عندك أكث من خيار ببعطوا ال
longest chain ، فلانم نختار الخيار الاكثر

تفرع

اذاً الاولويات
1) الاطول
2) الاكثر تفرع

2.4: Alkyl and Halogen Substituents (R-X)

هالوجينات
↓
الكيل

Alkyl Halide يكونوا

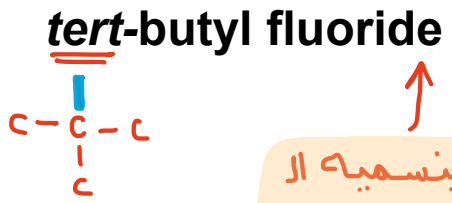
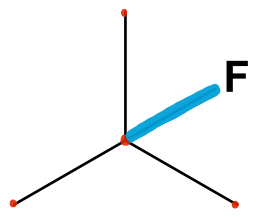
in simple compounds

note
الفرق بين
Common - IUPAC

Common name : Alkyl halide

eg. $\text{CH}_3\text{CH}_2\text{Br}$

Ethyl bromide
Alkyl halide



هالوجينات
Common name

IUPAC:

Halogen substituents are named by changing the *-ine* ending of the element to *-o*.

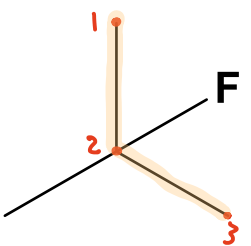
F: fluoro

Cl: chloro

Br: bromo

I: iodo-

على وزن هالو



longest chain
2-fluoro-2-methylpropane

هالوجينات
IUPAC name

افزادها قیادہ
الدكتور طارق
عيلبر

Sources of Hydrocarbons

Hydrocarbons are found in crude oil and natural gas. Both are mixtures of different hydrocarbons and they are separated by “fractional distillation” a process where the different compounds are separated in a long distilling column based on their boiling points. Heavy fractions can be “cracked” into small lighter, hydrocarbons using heat and/or catalysts.

انتقل الخام

خليط

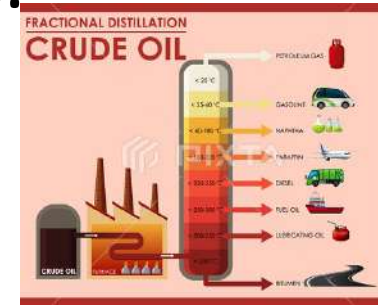
التقطير الجزئي

قطارة

غليان

متشققا ومكسور

مُحفز



2.7: Physical Properties of Alkanes and Nonbonding Intermolecular Interactions

A physical property is any property of matter or energy that can be measured.

When it changes, the chemical composition of the object does not change مثلاً لو وضعت الماء بالفرن حتى يجف Solid ولكن كيميائية chemical composition ما تغيرت

غير قابل للذوبان

Alkanes are insoluble in water. This is because water molecules are polar, whereas alkanes are nonpolar (Like dissolves like: Polar/ionic solvents dissolve polar/ionic solutes and non-polar solvents dissolve non-polar solutes)

درجة الحرارة التي يتساوى فيها External pressure ج
verbal pressure (درجة الغليان) ← من الروابط بين molecules

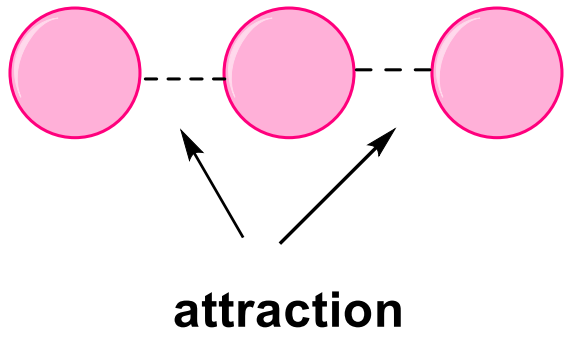
Alkanes have lower boiling points for a given molecular weight than most other organic compounds. This is because they are nonpolar molecules.

على ماذا تعتمد؟؟

The physical properties of molecules are in part dependent on the type's of intermolecular forces (IMF) present.

between

- Alkan properties:
- 1- insoluble in water
- 2- non polar
- 3- lower boiling point

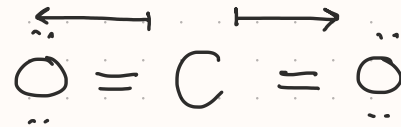


NOTE للفائدة ، ميب قَطْبِي ومبينك ٩٩

خطي ، لهم نفس الكهروسلبية ← غير قطبي



كهروسلبية مختلفة ، الجزيء كامل غير قطبي



لأنه قوة الجذب يتلغى بعضه ولكن

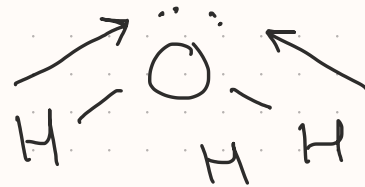
بين C و O ذرتين لوحدهم يوجد

فرق بالكهروسلبية ولهذا بينهم قطبي

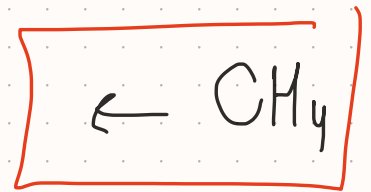
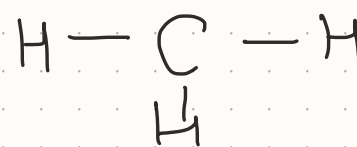
المركب قطبي



المركب قطبي



غير قطبي ، القوى يتلغى بعضه



Intermolecular Forces

The physical properties of molecules are in part dependent on the type's of intermolecular forces (IMF) present.

Boiling points (BP) are also dependent on the **mass of the molecule**.

الذوبان

Solubility, the ability to dissolve into a solvent is dependent on IMFs.

The strength of the interaction between molecules is also dependent on the **overall shape** of the molecule.

Intermolecular Forces

There are 3 types of IMFs, by decreasing strength they are:

1) Hydrogen bonding الأتوري

2) Dipole-dipole

3) Van der Waals or London Dispersion الأضعف
 ← المركبات ال nonpolar الأضعف

Dipole-dipole

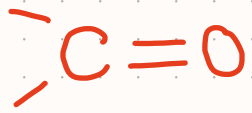
Dipole-dipole forces arise from the attraction of oppositely charged atoms (other than H) in molecules. These molecules may have a **permanent** dipole moment. Generally in organic molecules they result from the presence of C-X bonds where X is more electronegative than C.

١. الشحنة
٢. polar
٣. الترتيب
قائمة

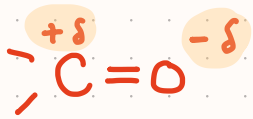
These are generally weaker than H-bonding, ranging from about 5-10 kJ/mol.

→ الشرح

شرح dipole - dipole



- what is the type of the bond?
covalent bond.
- Is it polar or non polar?
polar.
- Where is the direction of polarization?
toward oxygen.
- charges?



*these are dipole (ثنائين) *
لهم قيمة ثابتة ومصدرة
فعلقتهم مع other molecule
بتكون في نفس الترتيب



interaction

تأثر من dipole-dipole
ويعتبر strong

إذاً هي رابطة بتعبر بين المركبات القطبية حيث
كل مركب فيه عدد من الذرات كل ذرة لها
شحنة ثابتة وخصائص ثابتة من الغير ممكن إن
تتغير.

نوع الرابطة : تساهمية
قوة التجاذب dipole-dipole

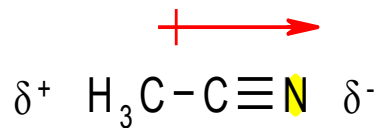
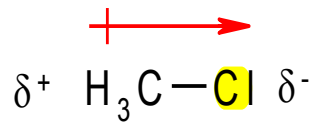
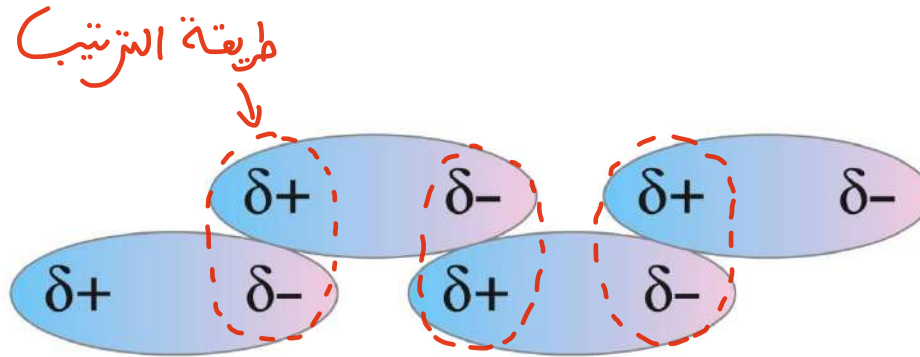


مثل ما اتفق ملاحظين المركب قطبي لأن Cl أعلى كهروسالبية
ولرندا :-

Cl ← شحنة جزئية سالبة H ← شحنة جزئية موجبة

فينشأ قوة تجاذب بين كل ذرتين شحنتهم متعاكسة

Dipole-dipole

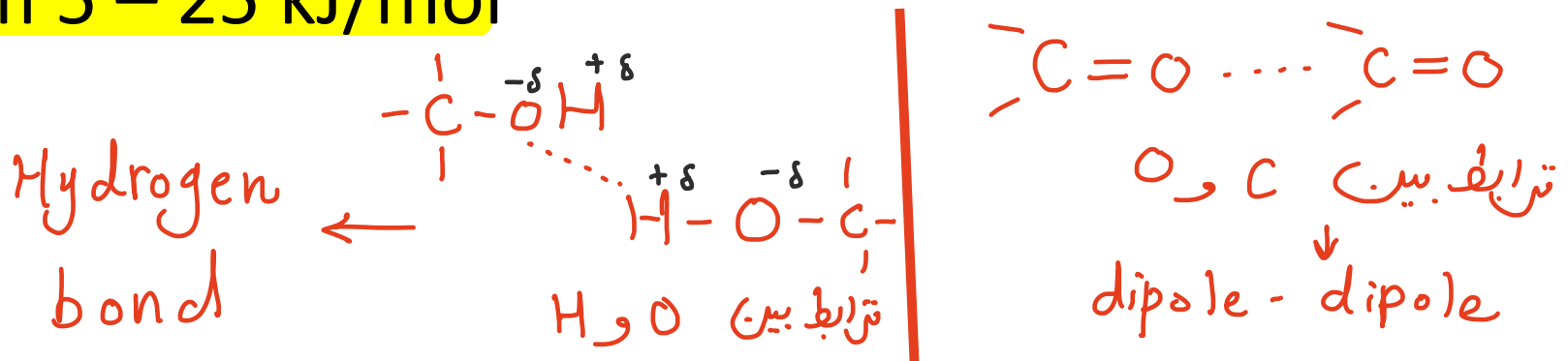


Hydrogen Bonding

N, O or F with H

Hydrogen bonding is a complex interaction that includes dipole-dipole, as well as orbital interactions and the transfer of electron density ^{انتقال} between molecules.

These are the strongest of the IMFs and range from 5 – 25 kJ/mol



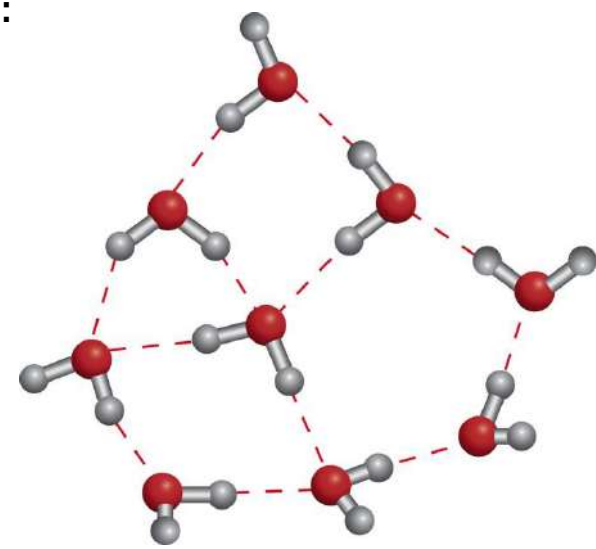
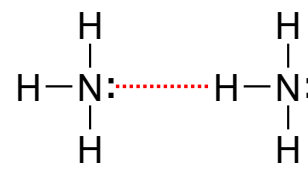
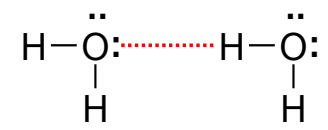
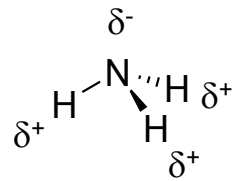
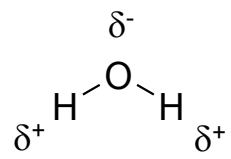
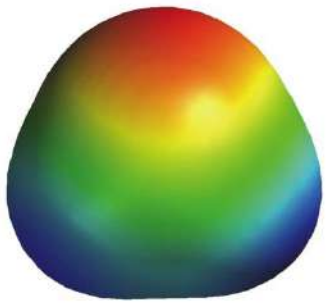
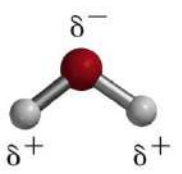
Hydrogen Bonding

علاقة طردية بين EN و Strong

نوف
١ ٢ ٣

Occur primarily between OH, NH and FH. The more EN the atom the stronger the interaction. (The atom H is attached to usually has a lone pair of e⁻)

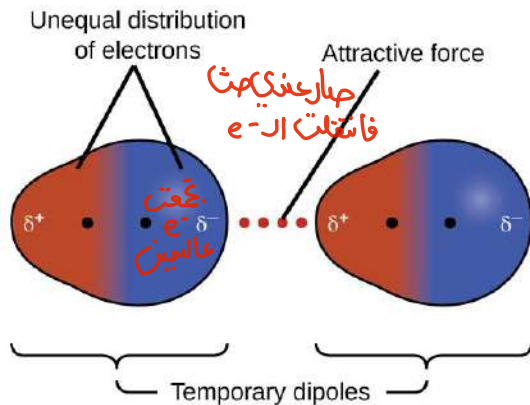
وهاد هوسيب انو الهيدروجينا اقوى وحدة



* موهودة في جميع المركبات الحنوية → Van der Waals (dispersion)

Van der Waals or (London) **dispersion** forces arise from the **movement of electrons** within a molecule. This natural motion can produce an uneven distribution of the electrons (polarization of the distribution) resulting in a **temporary dipole moment** in the molecule. This will **induce** the movement of electrons in adjacent molecules producing a dipole moment in them. These “induced” dipole moments are very brief as they disappear when the electrons move to new locations within the molecule, so they forces are very **brief and weak, only 2-5 kJ/mol.**

* نلاحظ ان لندن والهيدروجينية ليس الهم انتقال في e^- .



Very Weak attraction

London dispersion, < dipole-dipole, < Hydrogen bonding

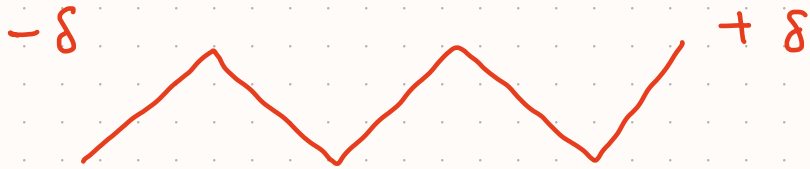
London force

* تتمثل أهميتها في non-polar مثل Alkane dipole ما عليه \uparrow

الذات إلكترونات هنا ليست حالة ثابت
انما حالة حركة \uparrow



لوحدها عنا ازاحة رح تولد dipole



وحتكون قيمته صغيرة



temporary dipole



other molecule \rightarrow induce \rightarrow يجعل \rightarrow فبيترتبوا بشكل انه اد \rightarrow opposite مع بعض

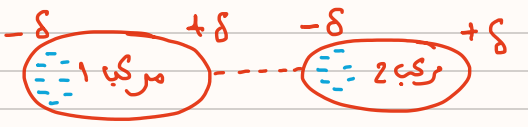
* هاد شرح الدكتور، كما ان شوي
شرحته لكم بطريقة مفهومة
أكثر

اقرأوه كيش
بفنيكم

مقارنة سريعة بين intermolecular forces

LONDON

قوة تجاذب وتزاييل بين الجزيئات غير القطبية ، ويكون فيط فرق الكهروسلبية يساري صفر .
وتتكون تساهمية .

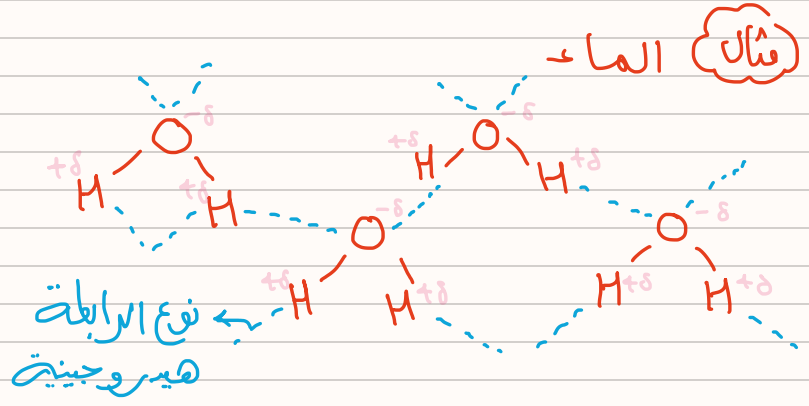


فبمير بيزم حد ، فيعاد تنظيم الشحنات السالب ينقل لجهة والموجب ينقل لجهة .

الانتقال حسب حركة الانتقال e^-

Hydrogen bond

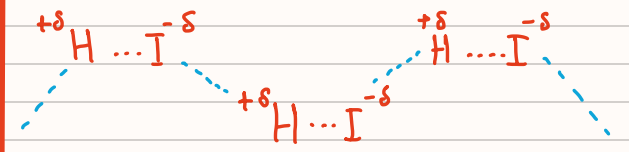
بتمير بركبات قطبية بين عنصر الهيدروجين و نوب (N, O, F) ، اي الكهروسلبيتهم عالية



وسبب الكهروسلبية العالية ، الاكترونات تنسحب باتجاه الاكسجين ، و تتكون بينهم رابطة قوية بسبب الكهروسلبية العالية و كما ان بسبب انتقال الاكترونات في ذرة حتكسي $+δ$ و ذرة $-δ$

Dipole-Dipole ثنائية القطب

تحت في الجزيئات القطبية .



اي بيحدث كالتالي ، الموجب والسالب يجذبوا بعضا برابطة بنسجورا ثنائية القطب .
فبمير تزايل بين الجزيئات .

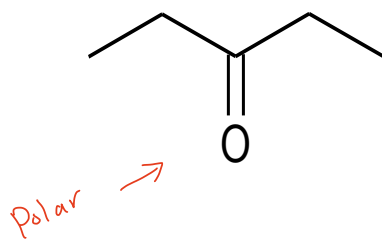
* ملاحظة *

الوقت بينك وبين لندن انه هاي قطبية ولندن غير قطبية وايضا بار dipole $+δ$ و $-δ$ ثابتين ما بتغير مكانهم مثل لندن

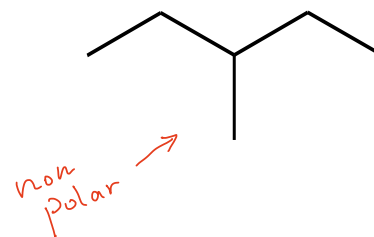
بالمركبات العضوية بنلا تير بين C مع X و لا يتكون اكثر كهروسلبية لانظر من المجموعة السابعة

Effects on Physical Properties (cont'd)

Dipole forces: much strong IMFs so the BP is higher when dipole-dipole interactions are present, i.e.



mass = 86 amu
BP = 101.7 °C



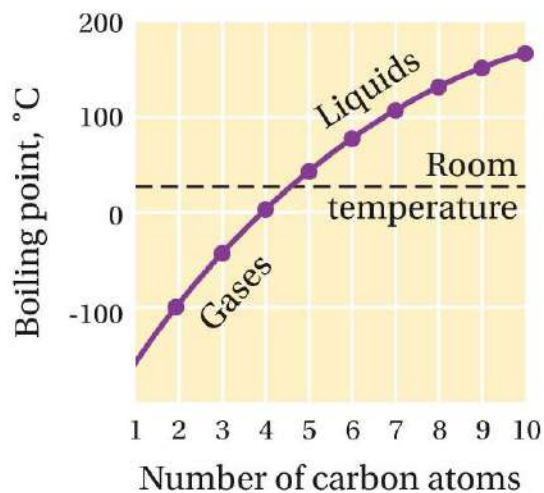
mass = 86 amu
BP = 63.3 °C

سبب الاختلاف انه الي الشمال dipole dipole واي عليمين london .

كلما كانت الرابطة اقوى كلما زادت درجة الغليان
و dipole dipole اقوى من London

Alkanes: no H-bonding or dipole moment (C & H have nearly the same EN ∴ not polar).

Main force
↓
London



Name	Formula	Boiling point, °C
pentane	$\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_3$	36
2-methylbutane (isopentane)	$\begin{array}{c} \text{CH}_3\text{CHCH}_2\text{CH}_3 \\ \\ \text{CH}_3 \end{array}$	28
2,2-dimethylpropane (neopentane)	$\begin{array}{c} \text{CH}_3 \\ \\ \text{CH}_3 - \text{C} - \text{CH}_3 \\ \\ \text{CH}_3 \end{array}$	10

no branch

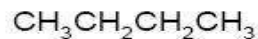
1 branch

4 branch

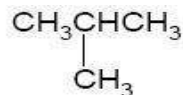
- Boiling point for alkanes increases with increasing size (mass)
- Boiling point for alkanes decreases with increasing branching

↑ mass
↓
↑ attraction
↓
↑ Boiling point

Example:



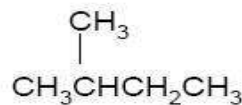
n-butane, 0°C



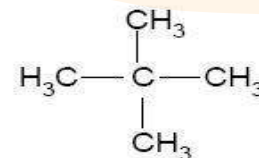
Isobutane, -12°C



n-pentane, 36°C



Isopentane, 28°C

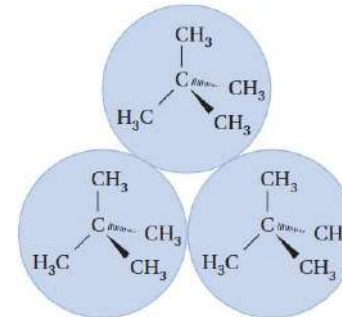
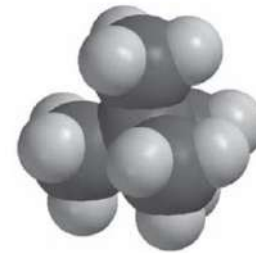


Neopentane, 9.5°C

↑ branching → ↓ surface area → ↓ interaction ↓ boiling

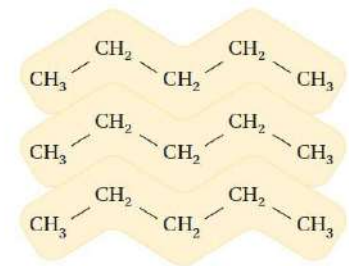
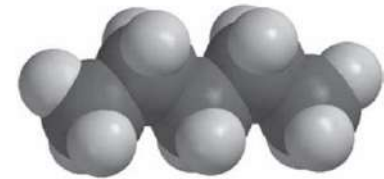
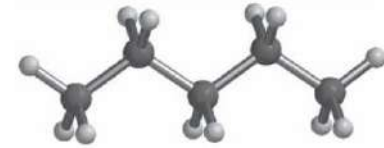
Structural Effects on IMFs

The strength of the IMFs depend on the amount of contact between the molecules, especially for dispersion forces. Hence the shape of the molecule can affect the surface area of contact, long thin molecules have more surface in contact than spherical molecules.



2,2-dimethylpropane
bp 10°C

(a)

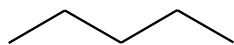


pentane
bp 36°C

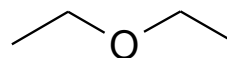
(b)

Effects on Physical Properties (cont'd)

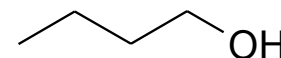
Hydrogen bonding:
strongest
intermolecular
forces so BP are
very high for
equivalent MW
compounds, i.e.



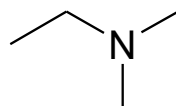
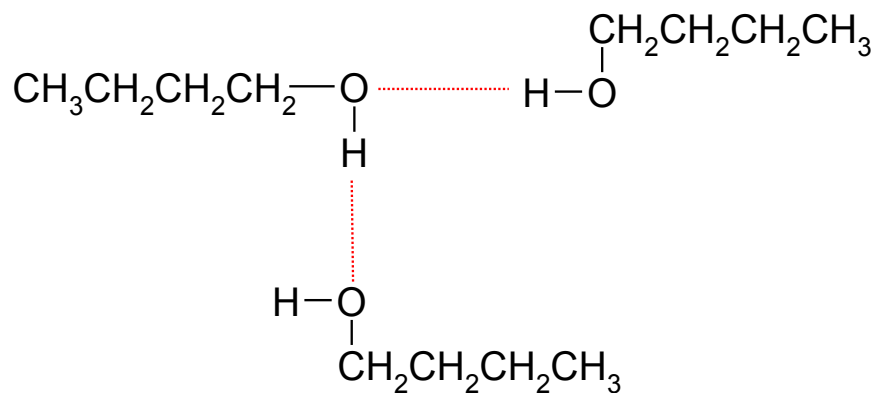
mass = 72 amu
BP = 36.1 °C



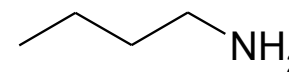
mass = 74 amu
BP = 35 °C



mass = 74 amu
BP = 117 °C



mass = 73 amu
BP = 36 °C



mass = 73 amu
BP = 78 °C

New terms

- saturated: مشبع
- unsaturated: غير مشبع
- generic formula: الصيغة العامة
- substituent: التفرعات
- crude oil: النفط الخام
- fractional distillation: التقطير الجزئي
- insoluble: غير قابل للذوبان
- solvent: مذيب
- dissolve: يذيب
- permanent: دائم و ثابت
- temporary: مؤقت
- Induce: الحث

Remember...

It's ok to have a bad day

It's ok to make mistakes

Set back is not failure

It's ok to take a break

Nothing is perfect

You are stronger than you think you are

Asking for help is a strength

Small steps are also progress

Good Luck