This paper describes objective technical results and analysis. Any subjective views or opinions that might be expressed in the paper do not necessarily represent the views of the U.S. Department of Energy or the United States Government.

Topological Effects on Separation of Alkane Isomers in Metal-Organic Frameworks

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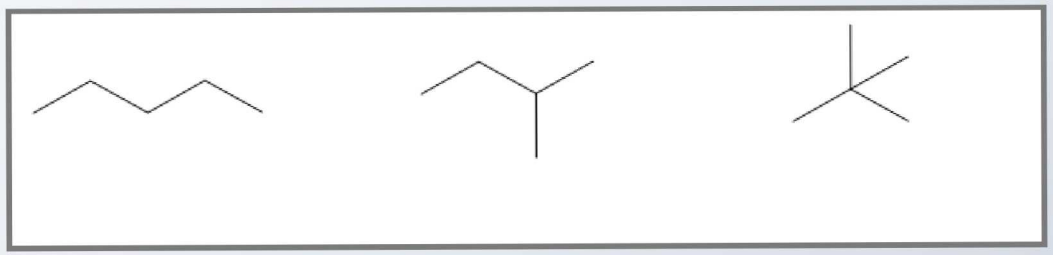
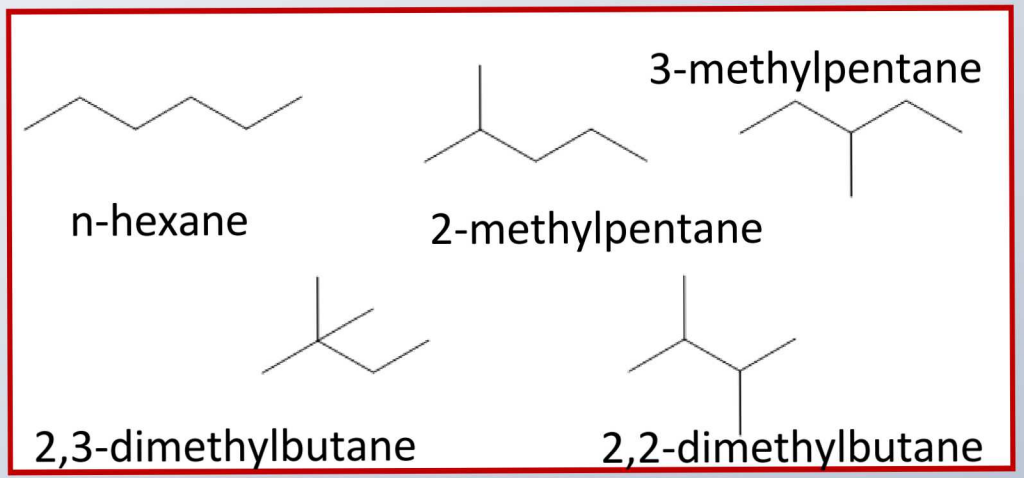
Alkane Isomers

n-butane isobutane

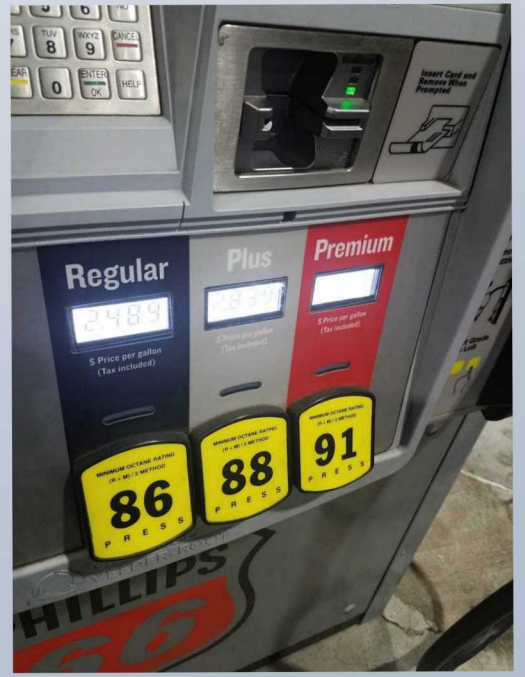


n-pentane 2-methylbutane 2,2-dimetylpropane (neopentane)

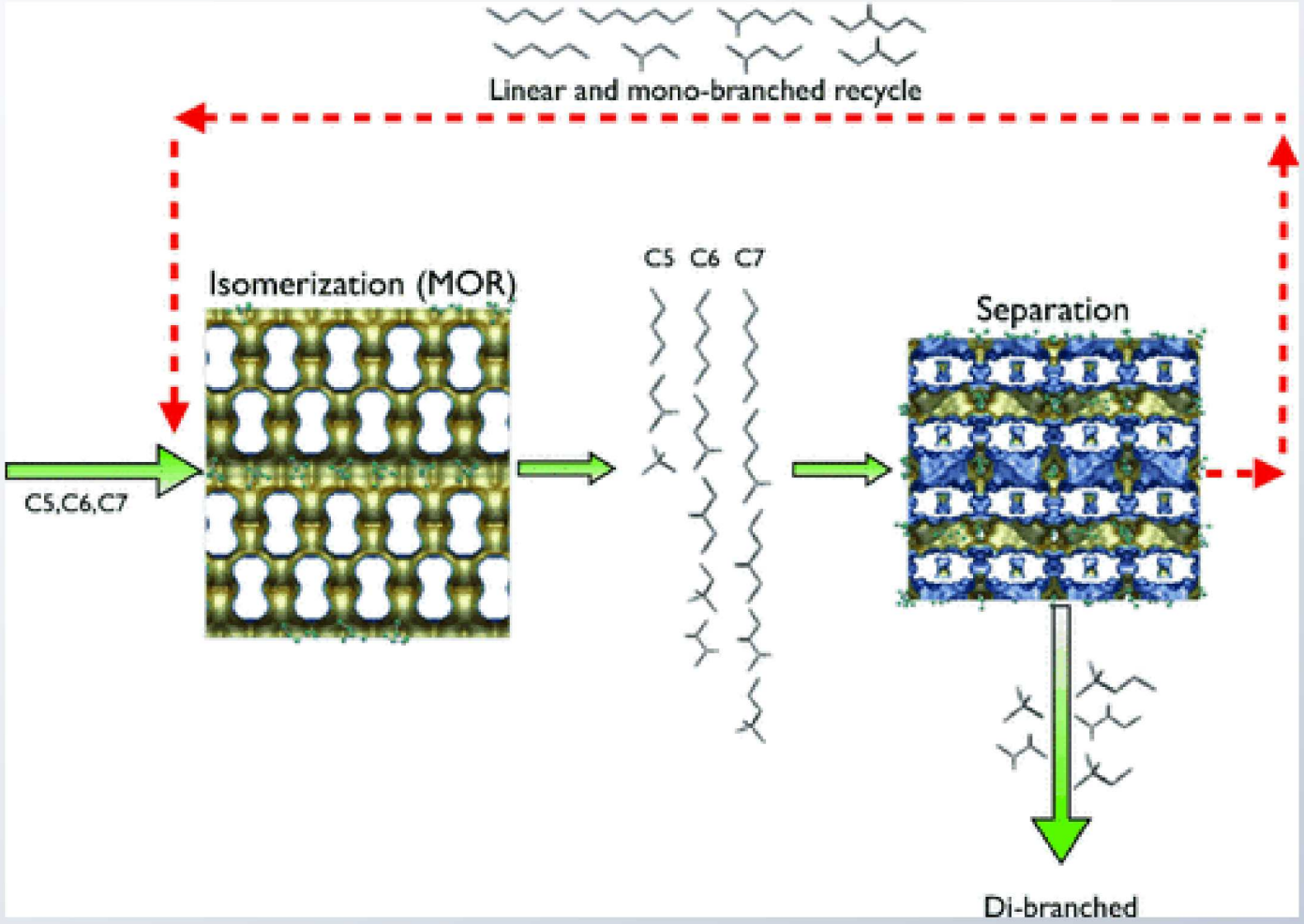


Isomer	Research Octane #
n-hexane	25
2-methylpentane	73
3-methylpentane	75
22-dimethylbutane	92
23-dimethylbutane	104

J. Combustion, ID 8406754, 2018



Branched isomers of hexane are key component of high-octane gasoline.



Dubbeldam, Angew. Chem. Int. Ed., 51, 11867-11871, 2012

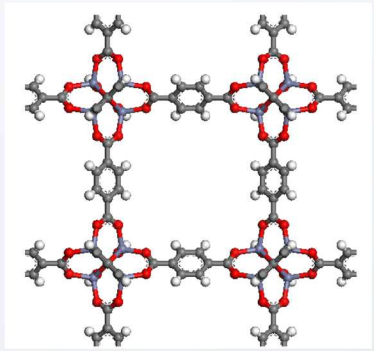
Gasoline is enriched by separating linear and branched alkanes and feeding back into an isomerization process.

It is advantageous to find a solid sorbent (eg MOF, zeolites) that can achieve this separation with a high degree of selectivity and capacity.

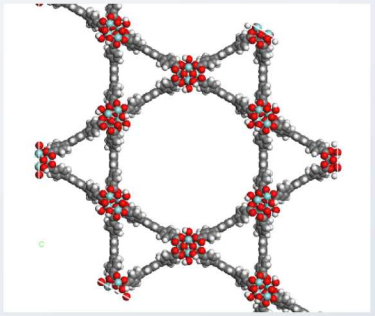
Isomer	B.P (˚C)
n-hexane	68.7
2-methylpentane	63.3
3-methylpentane	60.3
22-dimethylbutane	58.0
23-dimethylbutane	49.7

Distillation is difficult and expensive.
10-15% of world energy consumption.

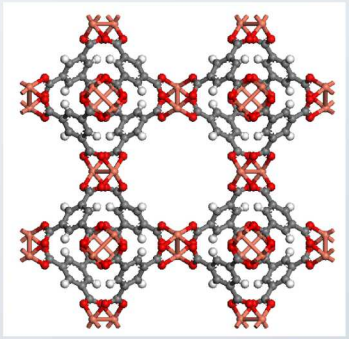
MOFs offer diverse pore shapes and sizes. We can choose a MOF to separate linear and branched isomers selectively based on shape.



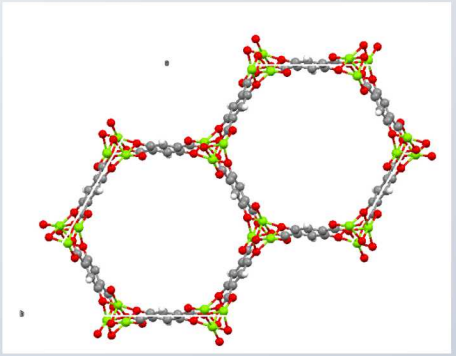
MOF-5



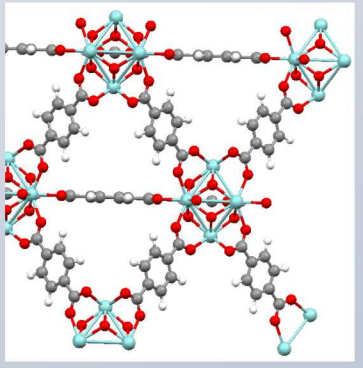
NU-1000



HKUST-1



MOF-74

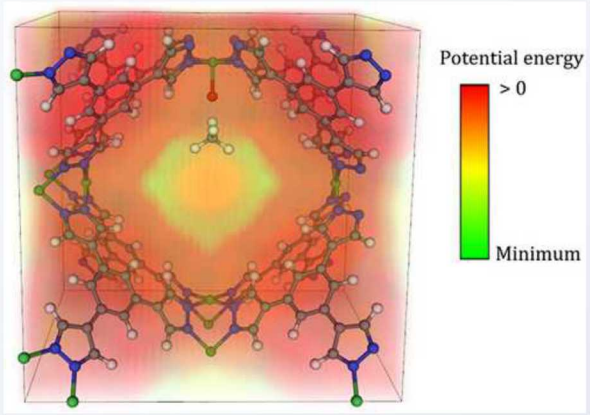


UiO-66

Shape-selective pores

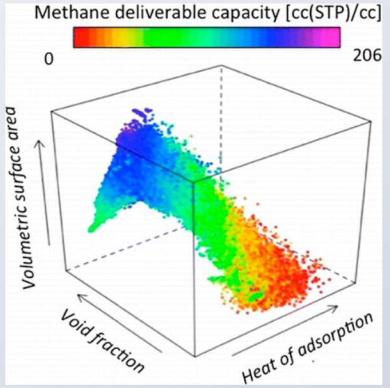
Mason, future chemical engineer

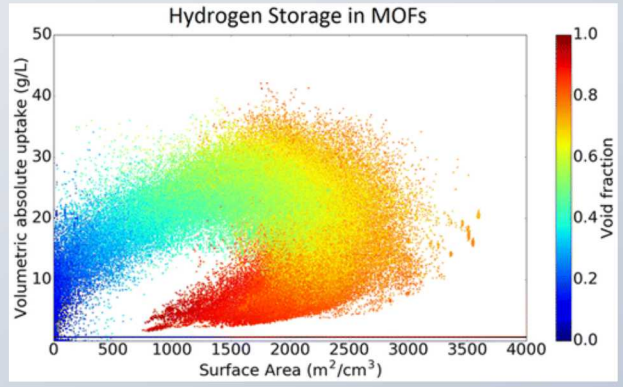
Catalysis



Rosen, J. Comp. Chem., 40, 1305-1318, 2019

Gas Storage





Gomez-Gualdron, J. Phys. Chem. C., 118, 6941-6951, 2014
Bobbitt, J. Phys. Chem. C, 120, 27328-27341, 2016

Nodes/Metals

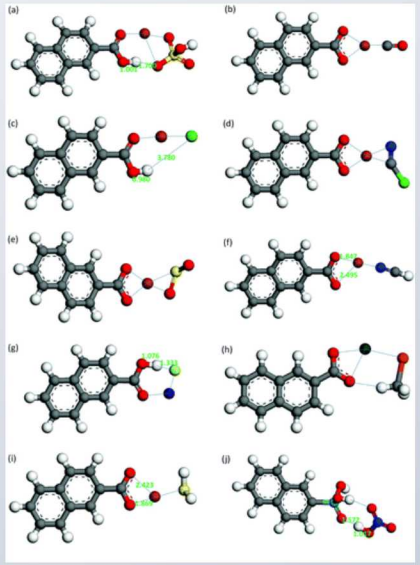
Linkers

TUNABLE PROPERTIES

Functional Groups

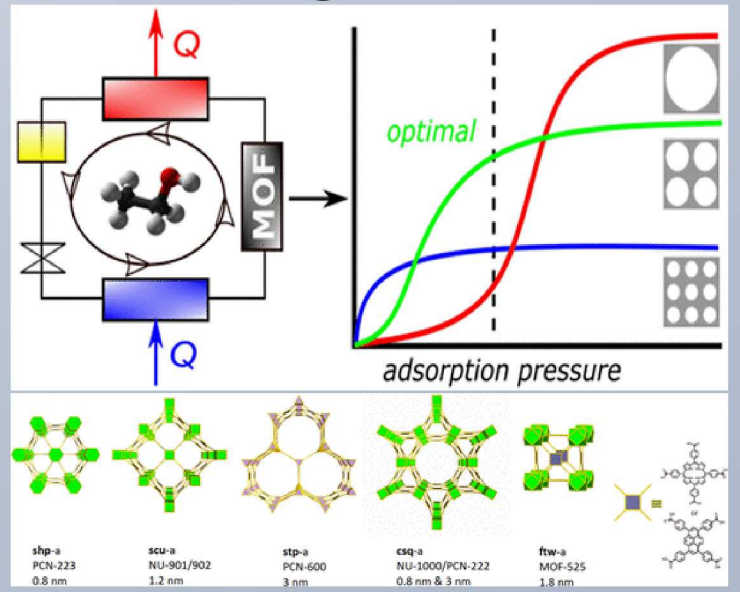
Topology

Toxic Gas Capture



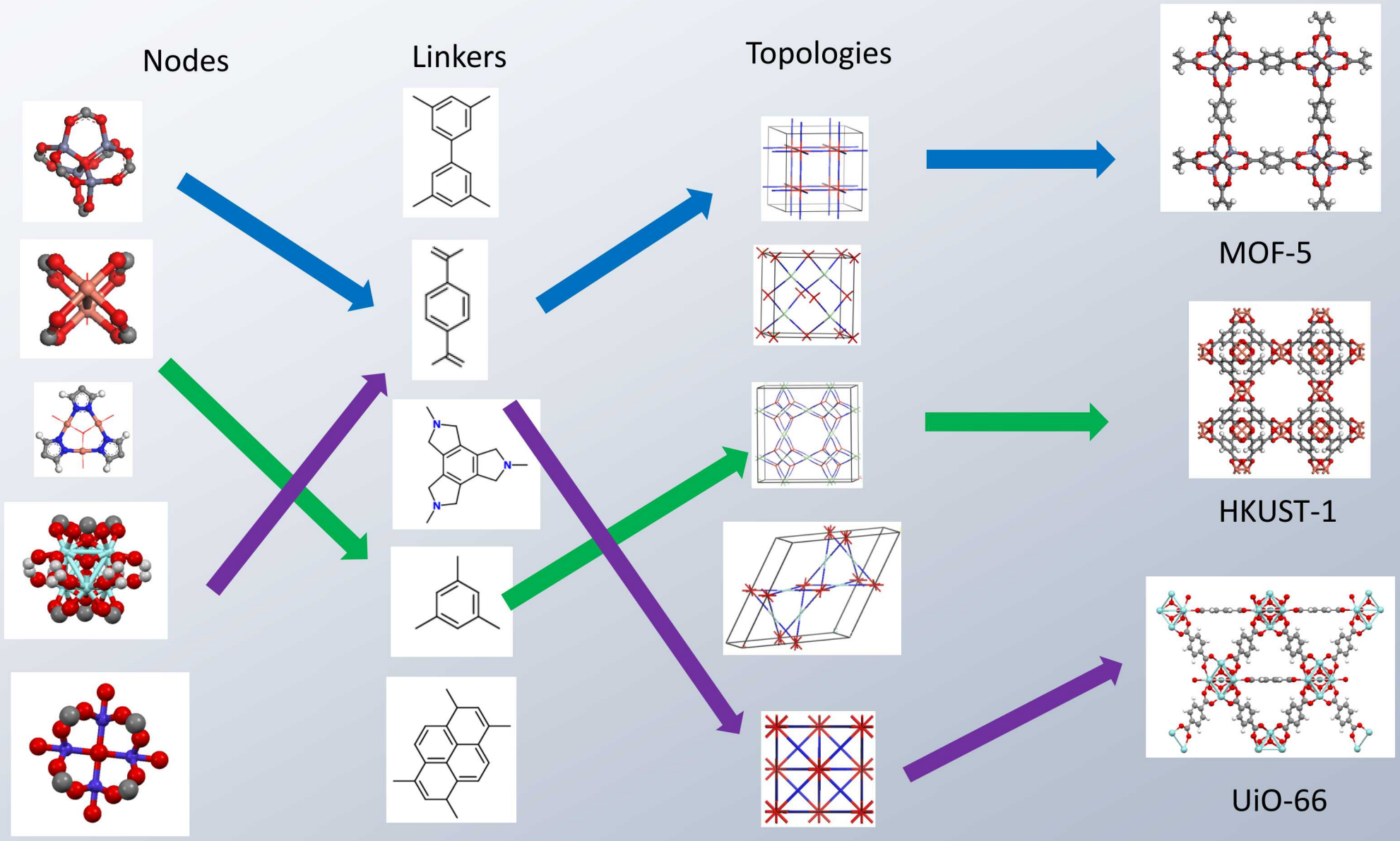
Kim, PCCP, 19, 31766-31772, 2017

Refrigeration



Chen, Chem. Mater., 31, 2702-2706, 2019

Many different combinations of nodes, linkers, and topologies can be used to form unique MOFs.

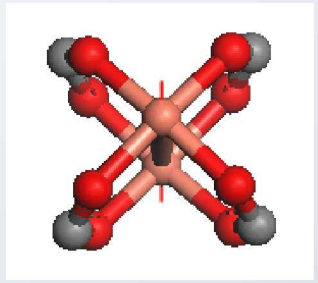


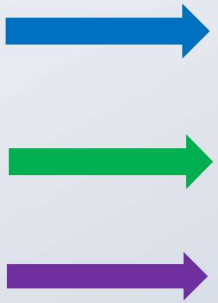
The same nodes and linkers can be combined in different topological networks to form unique MOFs with the same chemical formula.

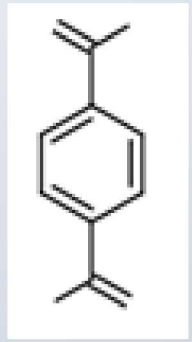
Nodes

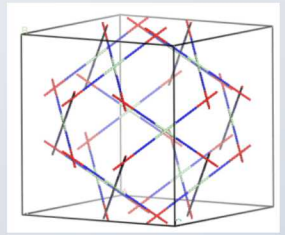
Linkers

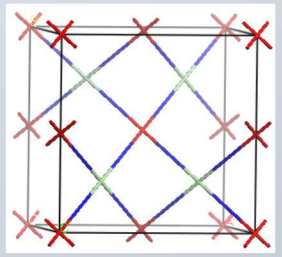
Topologies

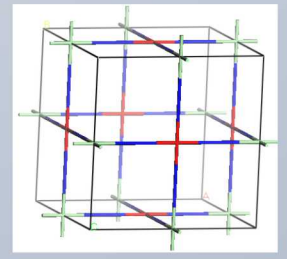


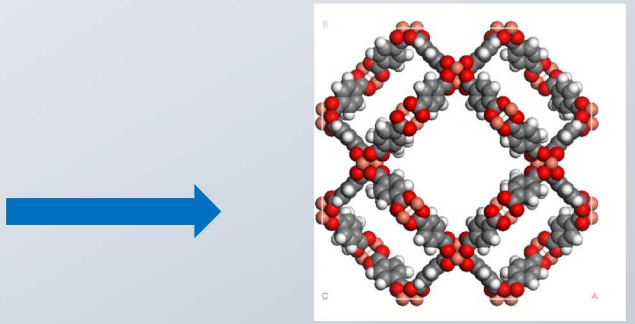


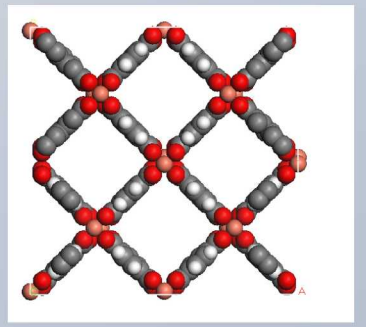


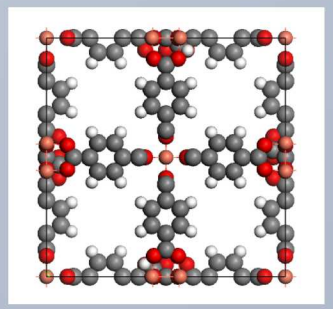












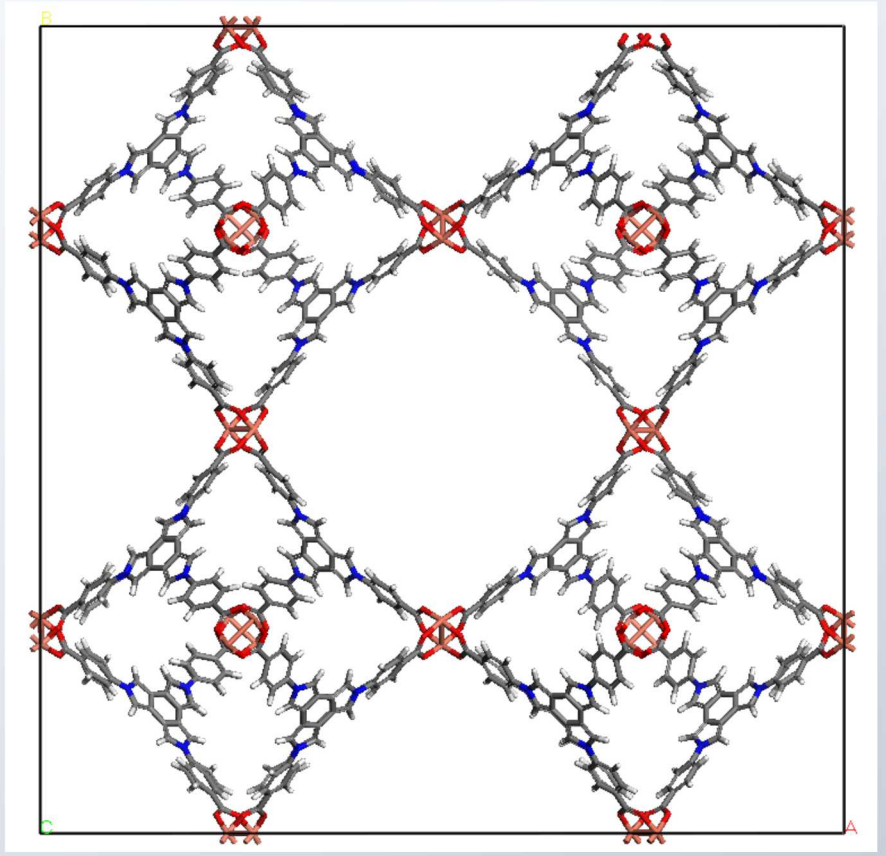
$CuC_8H_4O_4$

$CuC_8H_4O_4$

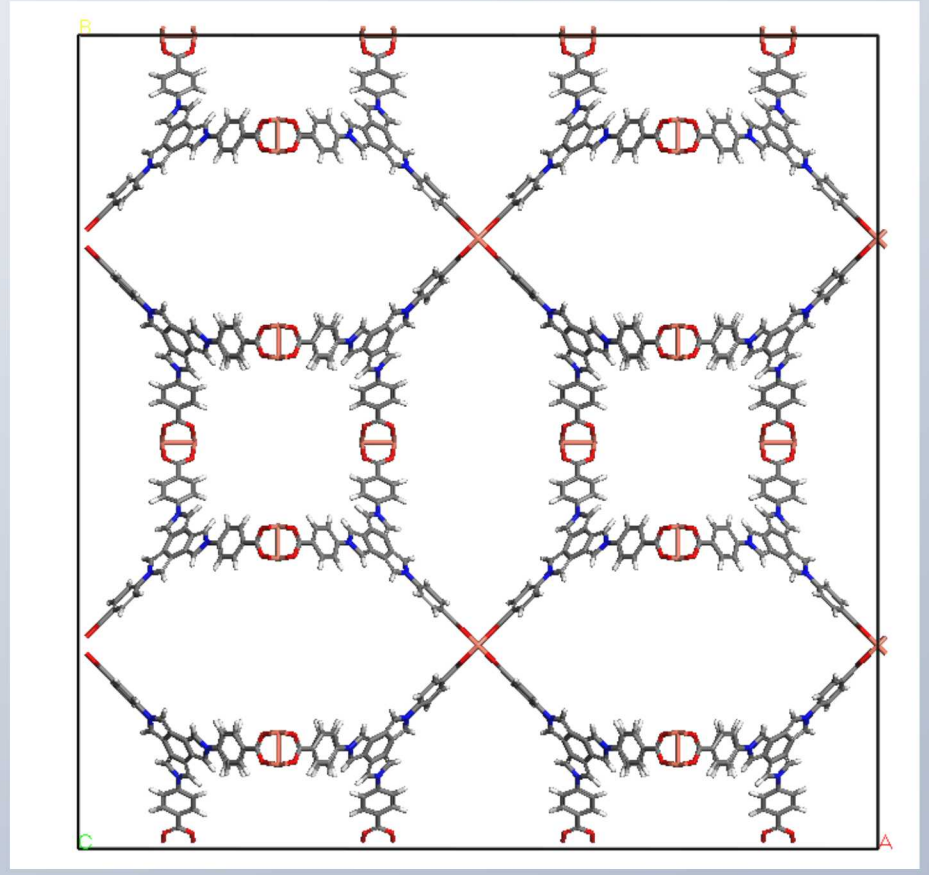
$CuC_8H_4O_4$

Polymorphs

Polymorphs are families of MOFs that contain the same nodes and linkers but arranged in different topologies.

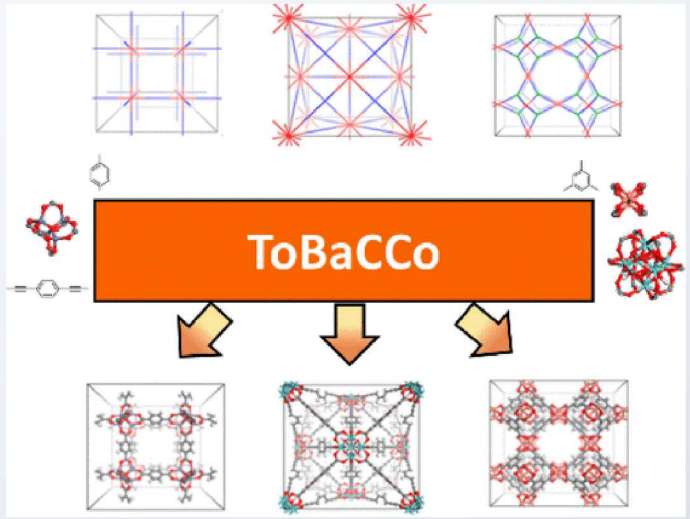


tbo



pto

I have selected four families of polymorphs that consist of 38 MOFs from the Tobacco database.



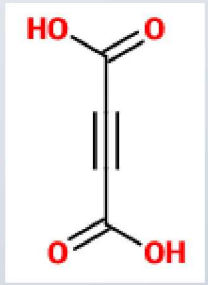
Colon, Gomez-Gualdron, Snurr, Cryst. Growth Des., 17, 5801-5810, 2017

Polymorph families:

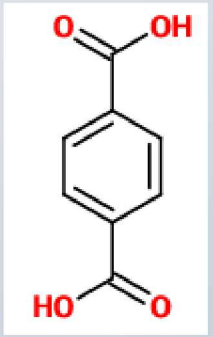
dia/lcs/sod

lvt/nbo/rhr

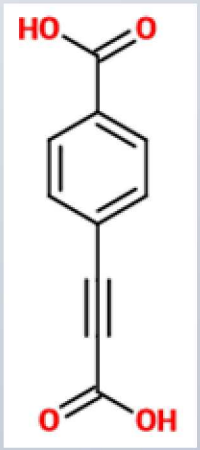
tbo/pto



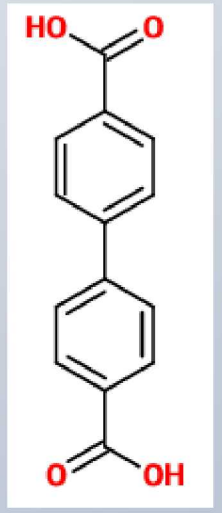
L3



L12

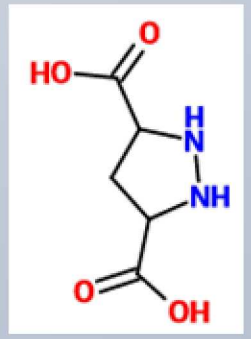


L20

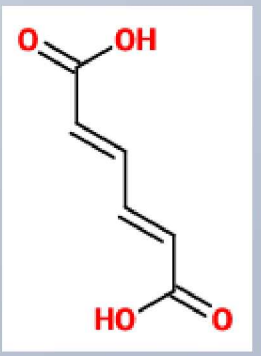


L24

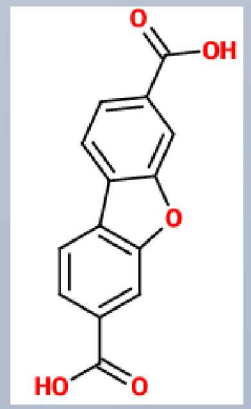
ssa/ssb



L4



L15



L21

Computational details

1. Monte Carlo simulations were performed in RASPA

2. We used the TraPPE model for alkanes and LJ parameters from UFF for MOFs.

3. MOF structures are kept rigid.

4. 20,000 cycles for Henry calculations and 100,000 total cycles (50k initialization-50k production) for GCMC using configurational biasing

For further details see Chung et al., Chem. Mater. 29, 6315-6328, 2017

Henry constants
mol/kg/Pa

nbo_L12 shows great affinity for n-hexane over branched isomers

MOF	C4	isobutane	C5	neopentane	2-methylbutane	C6	2-methylpentane	3-methylpentane	22-dimethylbutane	23-dimethylbutane
dia_L_12	7.6E-05	6.3E-05	1.8E-04	5.9E-05	1.3E-04	4.0E-04	3.1E-04	2.8E-04	1.8E-04	2.3E-04
dia_L_20	5.4E-05	4.5E-05	1.1E-04	4.2E-05	8.6E-05	2.3E-04	1.8E-04	1.7E-04	1.1E-04	1.4E-04
dia_L_24	7.0E-05	5.8E-05	1.5E-04	5.3E-05	1.2E-04	3.3E-04	2.6E-04	2.3E-04	1.5E-04	2.0E-04
dia_L_3	4.9E-05	4.2E-05	1.0E-04	4.0E-05	8.1E-05	2.1E-04	1.7E-04	1.6E-04	1.1E-04	1.4E-04
lcs_L_12	8.6E-05	7.1E-05	2.1E-04	6.6E-05	1.6E-04	5.1E-04	3.9E-04	3.6E-04	2.2E-04	3.0E-04
lcs_L_20	5.5E-05	4.6E-05	1.2E-04	4.3E-05	9.0E-05	2.4E-04	1.9E-04	1.7E-04	1.2E-04	1.5E-04
lcs_L_24	7.0E-05	5.8E-05	1.6E-04	5.4E-05	1.2E-04	3.5E-04	2.7E-04	2.5E-04	1.5E-04	2.1E-04
lcs_L_3	5.4E-05	4.6E-05	1.2E-04	4.5E-05	9.3E-05	2.5E-04	2.1E-04	1.9E-04	1.3E-04	1.6E-04
lvt_L_12	3.5E-03	3.4E-03	2.5E-02	9.1E-03	3.0E-02	1.6E-01	1.9E-01	2.4E-01	2.0E-01	3.1E-01
lvt_L_20	2.4E-04	2.1E-04	8.7E-04	2.5E-04	6.8E-04	3.1E-03	2.5E-03	2.3E-03	1.5E-03	1.9E-03
lvt_L_24	1.5E-04	1.4E-04	4.3E-04	1.6E-04	3.8E-04	1.2E-03	1.1E-03	1.0E-03	7.4E-04	9.3E-04
lvt_L_3	6.3E-04	6.3E-04	3.1E-03	9.6E-04	2.9E-03	1.4E-02	1.4E-02	1.3E-02	9.5E-03	1.2E-02
nbo_L_12	2.8E-02	1.9E-04	4.2E-01	2.1E-04	7.9E-04	**6.3E+00**	4.5E-03	3.2E-03	1.4E-03	2.3E-03
nbo_L_20	4.5E-05	3.9E-05	9.7E-05	3.7E-05	8.1E-05	2.0E-04	1.8E-04	1.6E-04	1.2E-04	1.4E-04
nbo_L_24	9.0E-05	8.1E-05	2.2E-04	8.0E-05	1.9E-04	5.1E-04	4.4E-04	4.1E-04	3.0E-04	3.6E-04
nbo_L_3	4.5E-05	4.0E-05	1.2E-04	5.2E-05	1.0E-04	2.9E-04	2.8E-04	2.8E-04	2.1E-04	2.6E-04
pto_L_12	2.1E-04	1.9E-04	6.2E-04	2.4E-04	6.1E-04	1.7E-03	1.7E-03	1.8E-03	1.4E-03	1.8E-03
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rhr_L_12	1.3E-03	2.3E-03	4.8E-03	6.2E-03	1.2E-02	1.8E-02	3.2E-02	4.3E-02	9.2E-02	9.4E-02
rhr_L_24	1.3E-04	1.1E-04	3.8E-04	1.2E-04	3.1E-04	1.2E-03	9.4E-04	8.3E-04	5.5E-04	6.9E-04
rhr_L_3	5.6E-05	2.4E-05	1.5E-04	1.8E-05	5.4E-05	4.0E-04	1.6E-04	1.3E-04	5.3E-05	7.9E-05
sod_L_12	7.5E-05	6.2E-05	1.8E-04	5.9E-05	1.3E-04	4.0E-04	3.2E-04	2.9E-04	1.8E-04	2.4E-04
sod_L_20	5.7E-05	4.8E-05	1.2E-04	4.5E-05	9.2E-05	2.5E-04	2.0E-04	1.8E-04	1.2E-04	1.5E-04
sod_L_3	5.6E-05	4.8E-05	1.3E-04	4.8E-05	1.0E-04	2.9E-04	2.3E-04	2.2E-04	1.4E-04	1.9E-04
ssa_L_15	6.8E-04	4.7E-04	3.7E-03	5.6E-04	2.4E-03	1.9E-02	1.4E-02	1.3E-02	6.3E-03	9.4E-03
ssa_L_21	5.5E-04	4.5E-04	2.9E-03	4.8E-04	2.4E-03	1.2E-02	1.3E-02	1.3E-02	6.6E-03	1.2E-02
ssa_L_4	2.0E-03	1.5E-03	1.6E-02	1.6E-03	1.1E-02	1.1E-01	1.0E-01	9.0E-02	3.4E-02	6.6E-02
ssb_L_15	3.1E-04	1.8E-04	1.2E-03	1.3E-04	5.9E-04	4.4E-03	2.5E-03	1.9E-03	8.0E-04	1.1E-03
ssb_L_21	4.3E-04	3.7E-04	1.8E-03	4.7E-04	1.8E-03	7.9E-03	7.2E-03	7.8E-03	5.0E-03	8.3E-03
ssb_L_4	3.1E-04	2.7E-04	1.2E-03	3.4E-04	1.0E-03	4.5E-03	4.0E-03	3.8E-03	2.5E-03	3.4E-03
tbo_L_12	1.9E-04	1.5E-04	6.5E-04	1.8E-04	5.1E-04	2.3E-03	1.9E-03	1.8E-03	1.0E-03	1.5E-03
tbo_L_20	6.0E-05	5.2E-05	1.3E-04	5.0E-05	1.1E-04	2.9E-04	2.5E-04	2.3E-04	1.6E-04	2.0E-04
tbo_L_24	9.9E-05	8.9E-05	2.4E-04	8.8E-05	2.1E-04	5.8E-04	5.0E-04	4.7E-04	3.4E-04	4.2E-04
tbo_L_3	1.2E-04	1.1E-04	4.2E-04	1.7E-04	3.9E-04	1.4E-03	1.5E-03	1.5E-03	1.1E-03	1.4E-03

GCMC results at 300 K, 0.0001 bar
Heat of adsorption in kJ/mol

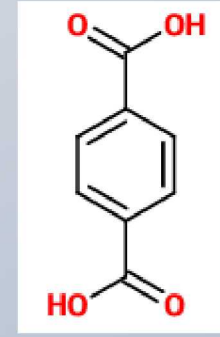
				DIFF [C5 - X]	
MOF	C5	2-mb	neopent	2-mb	neopent
dia_L_12	-10.1	-8.7	-6.3	-1.5	-3.8
dia_L_20	-8.7	-7.1	-3.0	-1.6	-5.6
dia_L_24	-9.3	-8.3	-5.4	-1.0	-3.9
dia_L_3	-5.9	-6.5	-5.8	0.6	-0.1
lcs_L_12	-11.1	-10.5	-6.9	-0.6	-4.2
lcs_L_3	-7.7	-7.8	-4.2	0.1	-3.6
lvt_L_12	-27.6	-28.5	-24.5	0.9	-3.0
lvt_L_20	-16.6	-16.1	-12.9	-0.5	-3.7
lvt_L_24	-12.8	-13.3	-10.6	0.5	-2.2
lvt_L_3	-23.0	-23.4	-20.6	0.5	-2.4
nbo_L_12	-52.1	-18.1	-13.2	**-34.0**	**-38.9**
nbo_L_20	-9.0	-9.7	-4.6	0.7	-4.3
nbo_L_24	-13.2	-13.2	-9.9	0.0	-3.3
nbo_L_3	-8.9	-9.2	-6.9	0.3	-2.0
pto_L_12	-14.5	-16.3	-12.5	1.7	-2.1
pto_L_20	-9.3	-9.9	-7.1	0.6	-2.1
pto_L_24	-11.0	-10.9	-7.2	0.0	-3.8
pto_L_3	-10.0	-9.9	-7.9	0.0	-2.1
rhr_L_3	-19.4	-11.9	-2.2	-7.5	-16.1
rhr_L_12	-27.2	-31.6	-30.0	**4.4**	**2.7**
rhr_L_24	-16.2	-15.6	-12.0	-0.6	-4.2
sod_L_12	-10.2	-9.7	-5.4	-0.5	-4.8
sod_L_3	-8.9	-7.2	-4.2	-1.7	-4.7
ssa_L_15	-27.0	-25.3	-19.7	-1.6	-7.3
ssa_L_21	-27.7	-28.0	-22.4	0.3	-5.3
ssa_L_4	-35.7	-34.7	-26.8	-1.0	-8.9
ssb_L_15	-24.8	-21.9	-12.2	-2.9	-12.6
ssb_L_21	-23.8	-24.1	-19.6	0.3	-4.2
ssb_L_4	-19.7	-19.6	-16.0	-0.1	-3.7
tbo_L_3	-13.5	-14.4	-12.0	0.8	-1.5
tbo_L_12	-16.2	-15.7	-	-0.5	-
tbo_L_20	-9.6	-9.0	-6.3	-0.6	-3.3

BLUE = favors branched isomer
RED = favors linear isomer

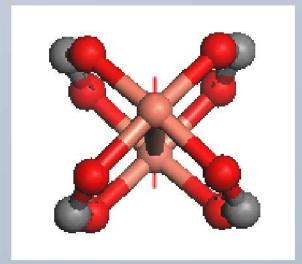
GCMC simulations show that nbo_L12 highly favors linear n-pentane over branched isomers.

Polymorph family:

lvt/nbo/rhr

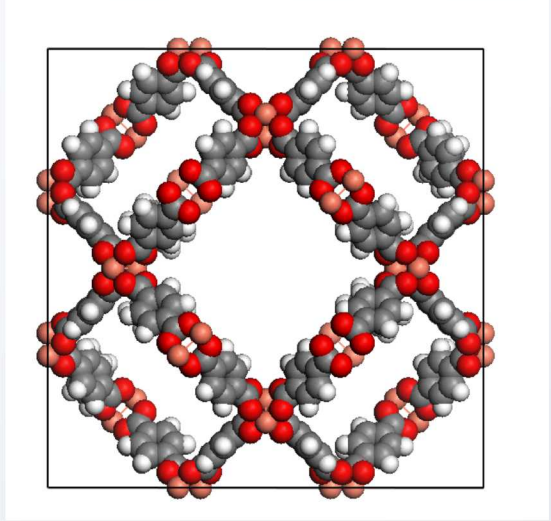


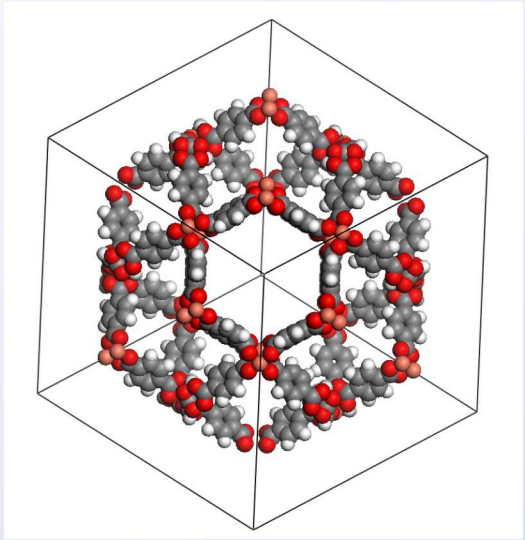
L12



Cu paddlewheel node

rhr_L12 polymorph shows slight selectivity for the branched isomers

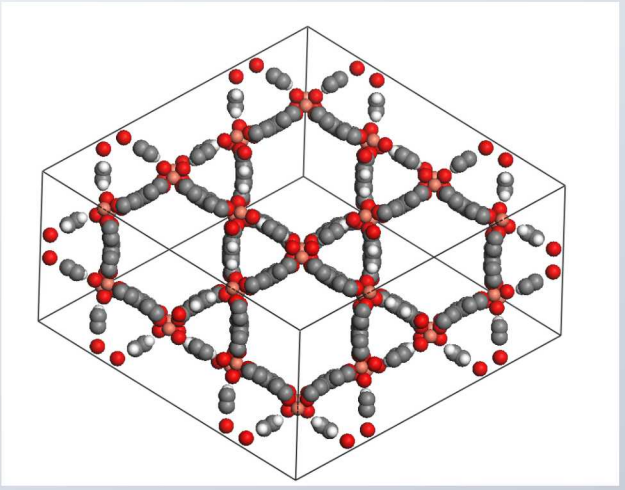




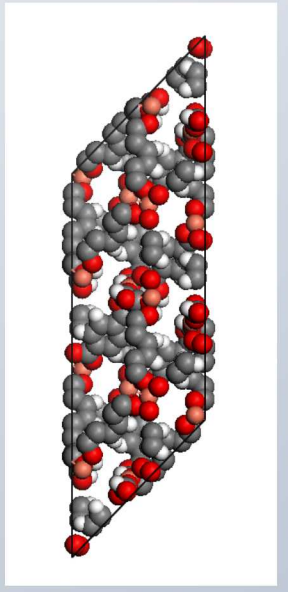
rhr_L12

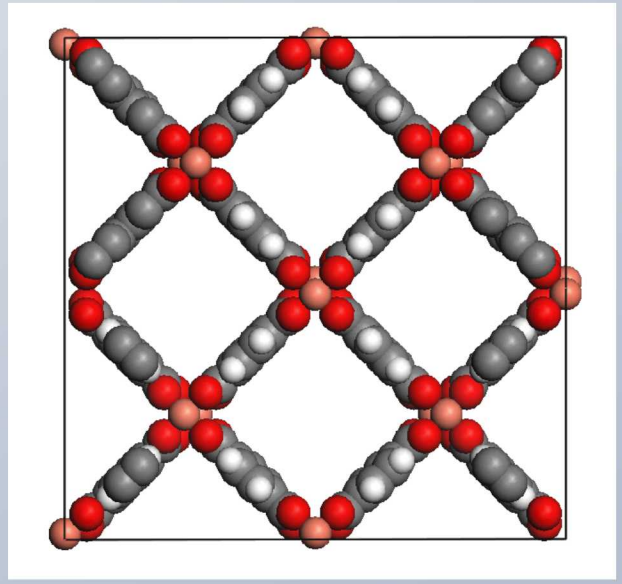
We will focus on these three polymorphs.

MOF	Pore size (Å)
rhr_L12	28.2
nbo_L12	4.6, 13.6
lvt_L12	7.4, 8.6

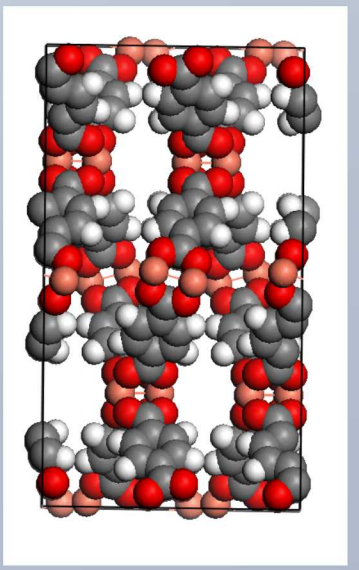


nbo_L12



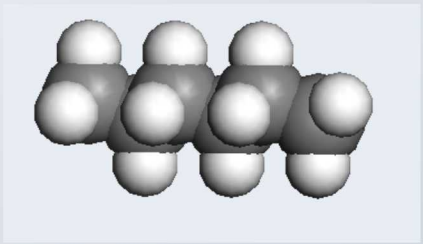


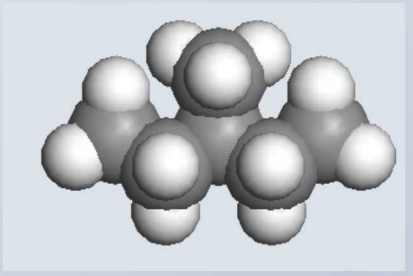
lvt_L12



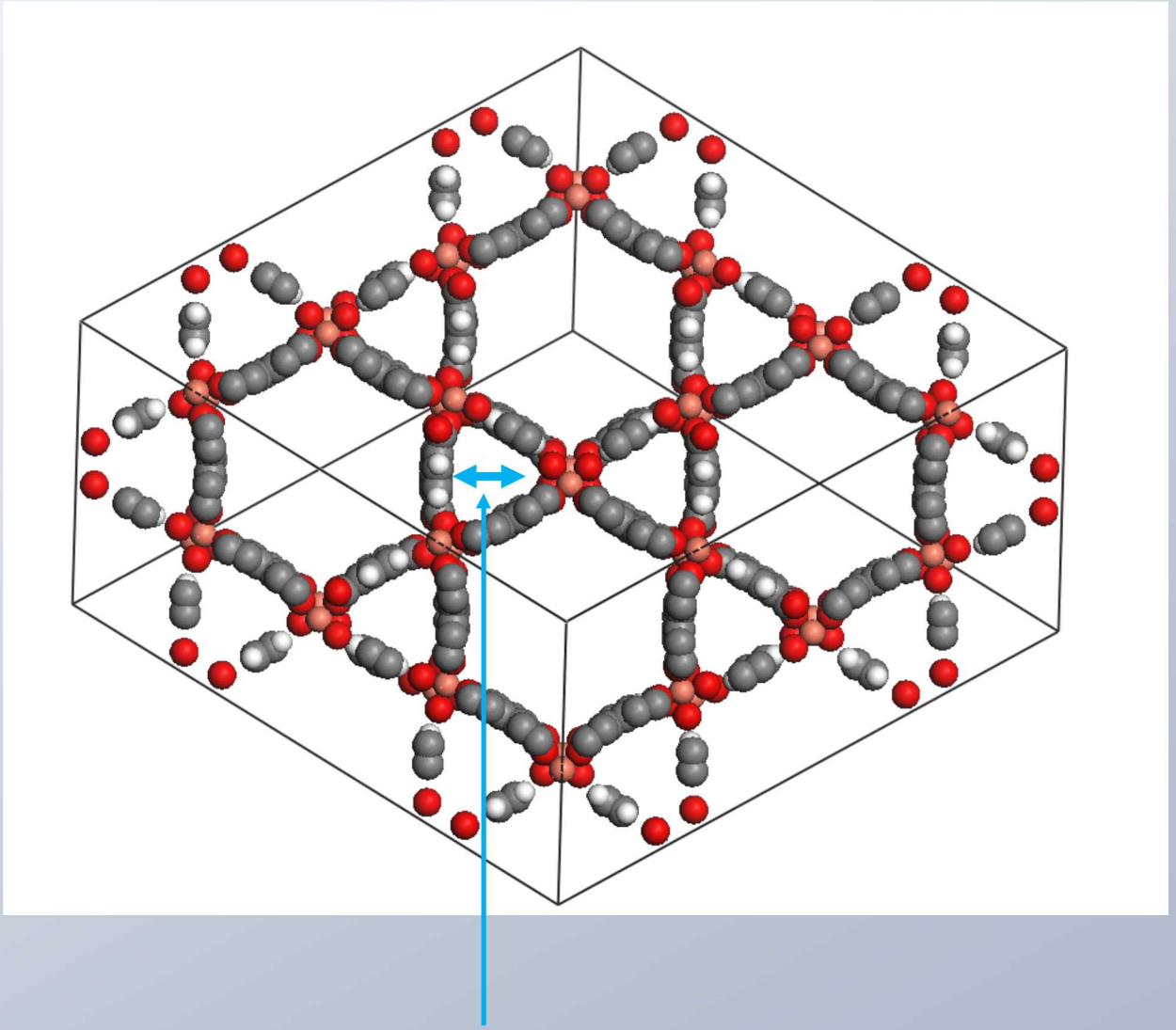
The small pores in nbo_L12 will accommodate n-hexane but exclude bulkier branched isomers.

Alkane	X	Y	Z	Kinetic Dia. (Å)
n-hexane	9.7	4.5	4.0	4.3
2-methylpentane	9.2	6.2	5.2	5.0
3-methylpentane	9.3	6.2	5.2	5.0
2,2-dimethylbutane	8.0	6.7	5.9	6.2
2,3-dimethylbutane	7.8	6.7	5.3	5.6





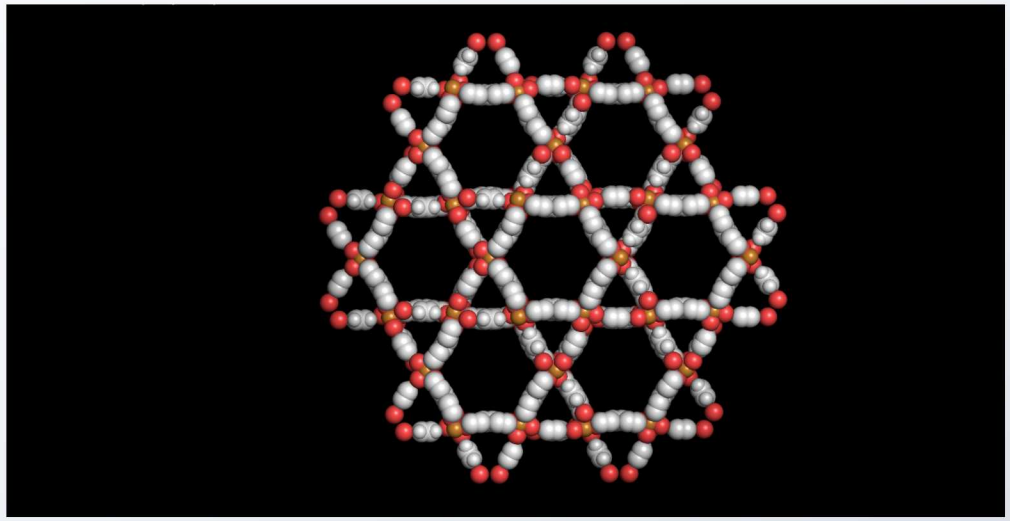
nbo_L12



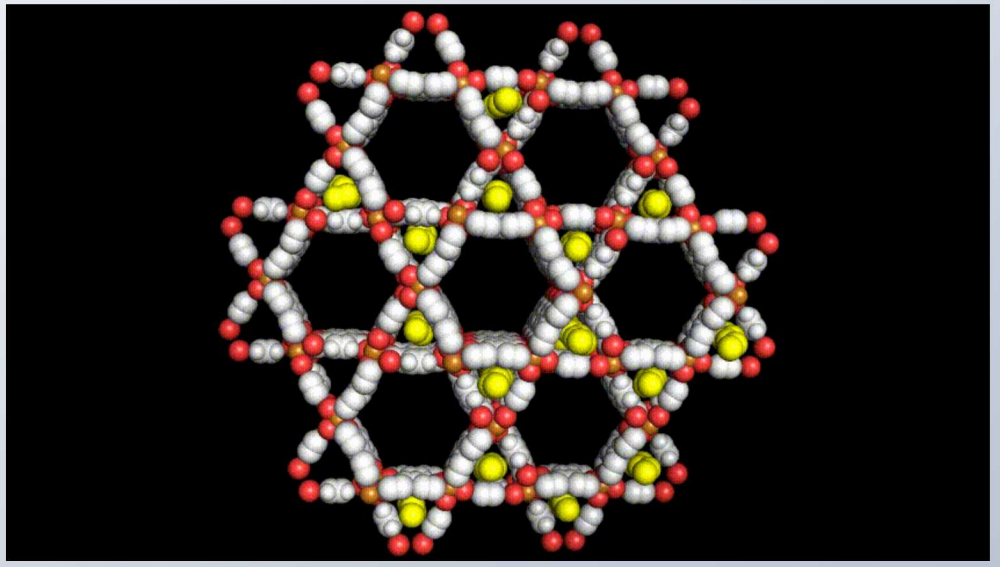
Alkane diameters: Peralta, Ind. Eng. Chem. Res., 2012,51, 4692-4702

n-hexane
300 K

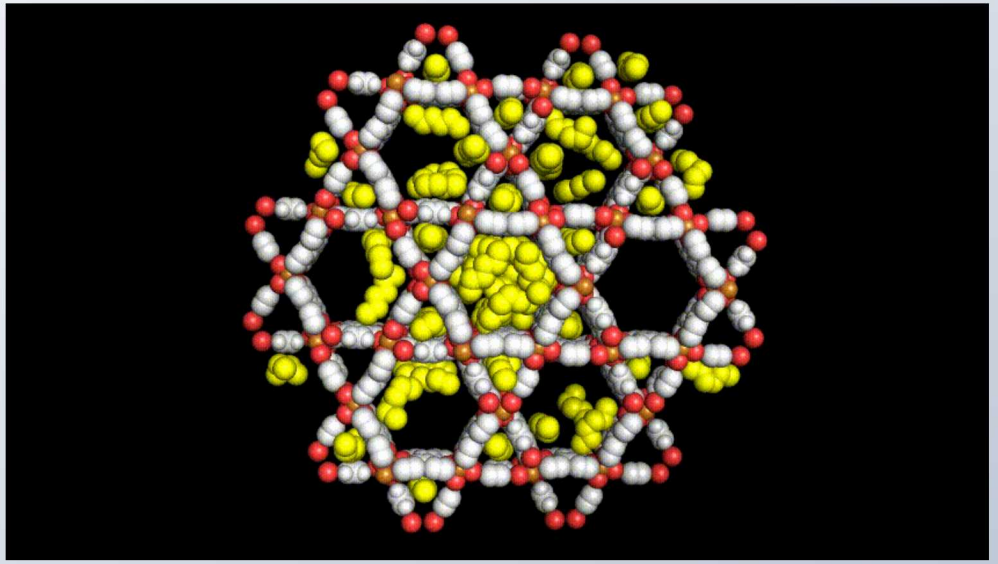
Linear molecules prefer the small pores at low pressure.



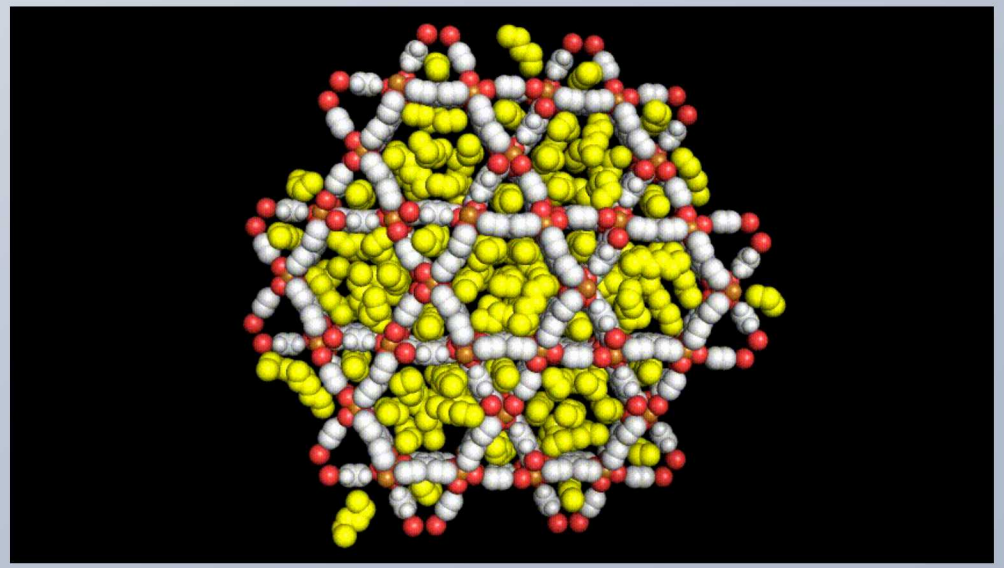
No adsorbate



10 Pa

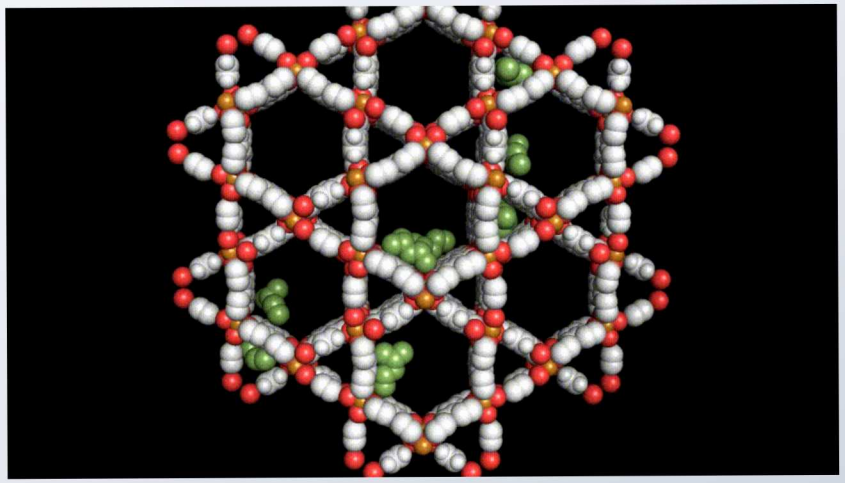


100 Pa



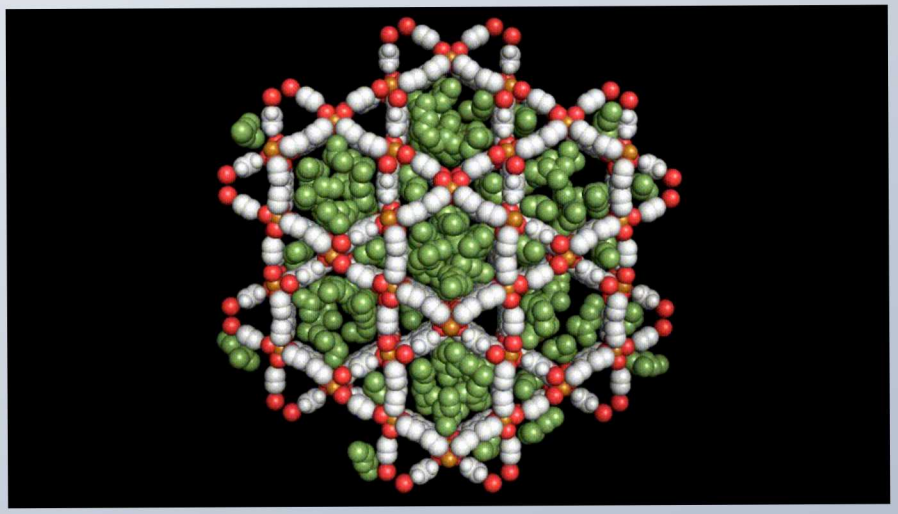
1000 Pa

At low pressure, the monobranched isomer prefers the large pores



100 Pa

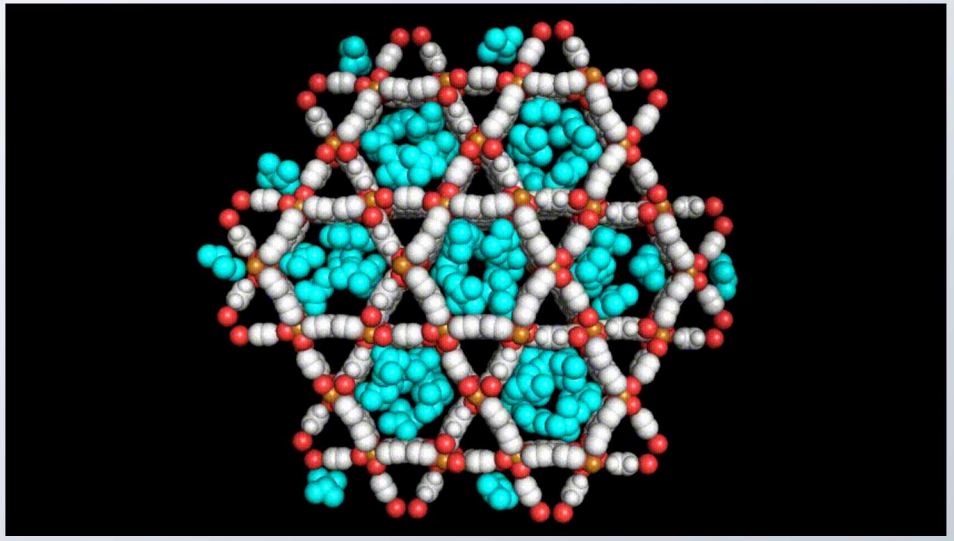
At saturation, 2-mp can be forced into the small triangle pore



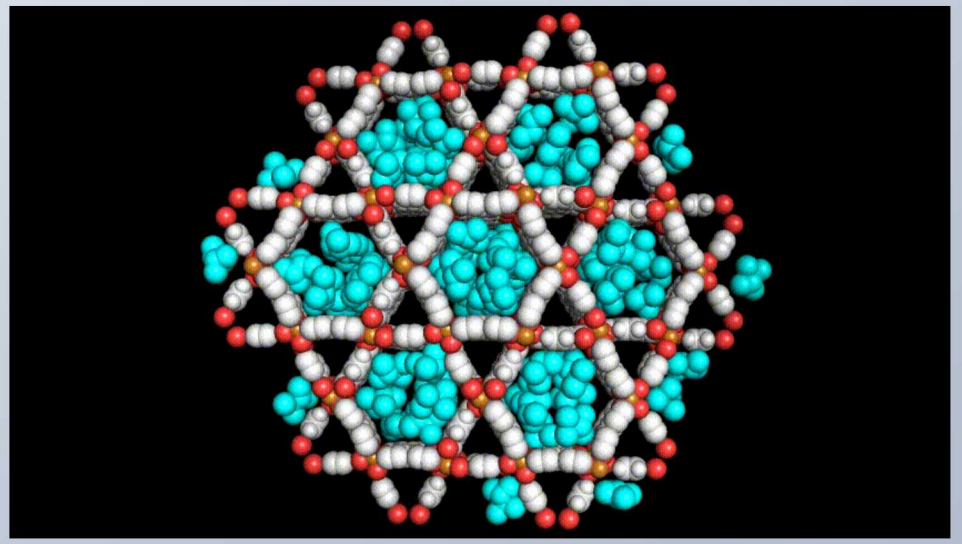
1000 Pa

2-methylpentane
300 K

23-dmb cannot go into the small pore even at high pressure



1000 Pa

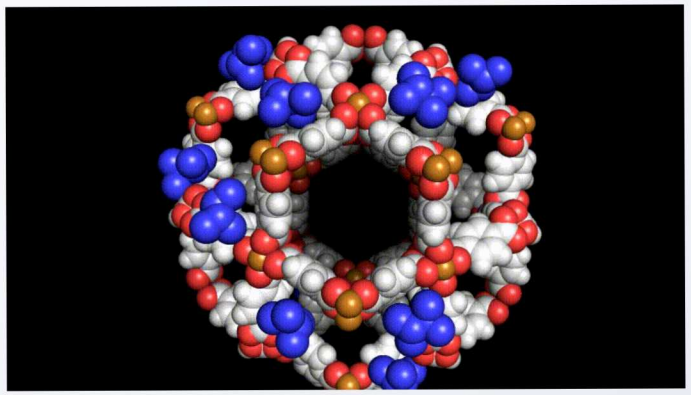


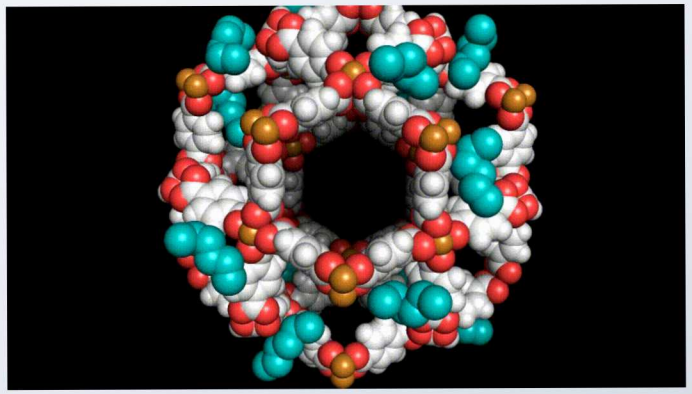
10000 Pa

2,3-dimethylbutane
300 K

2,3-dimethylbutane

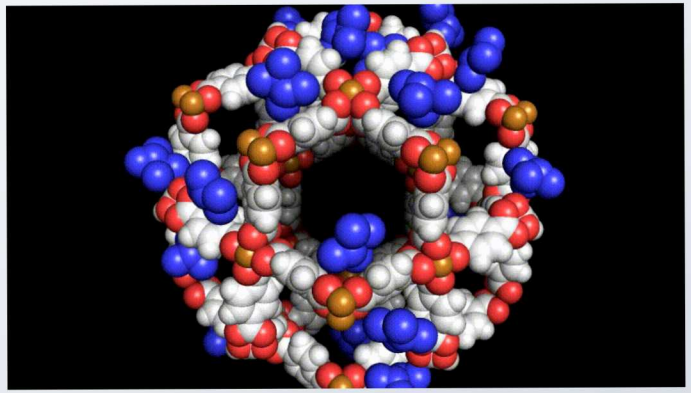
n-hexane

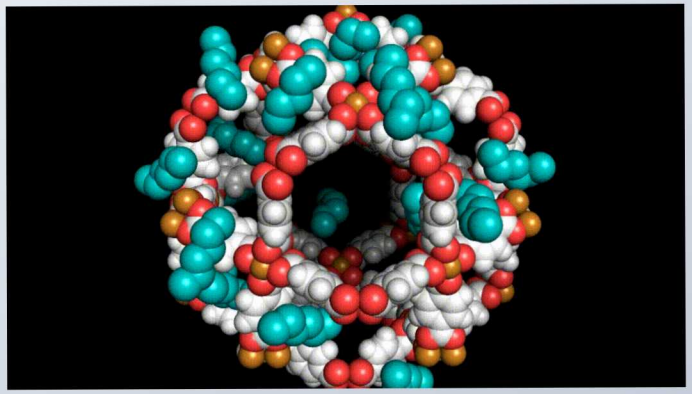




100 Pa

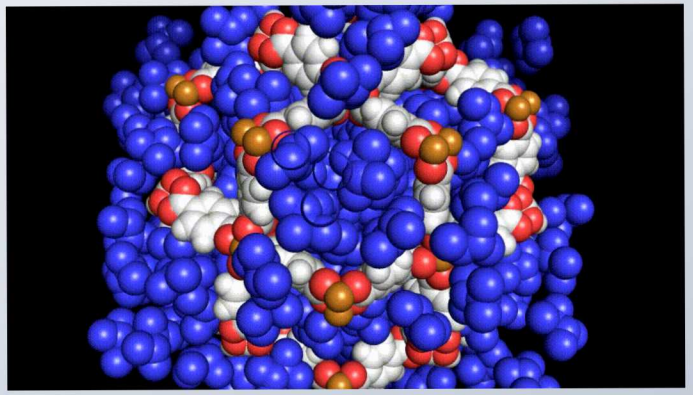
In rhr_L12, the molecules first go to the windows surrounding the large spherical pore

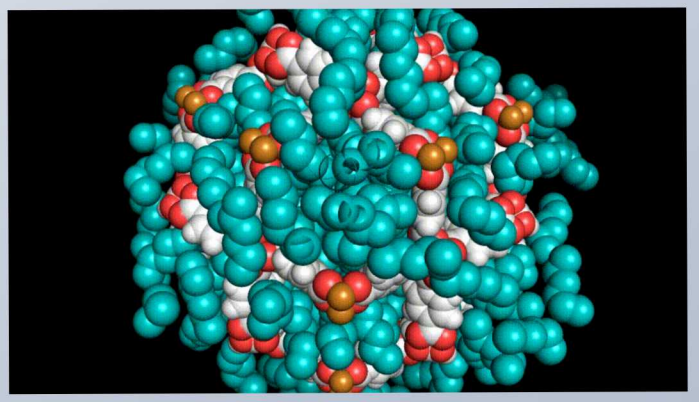




1000 Pa

At higher pressure, the pore is filled.

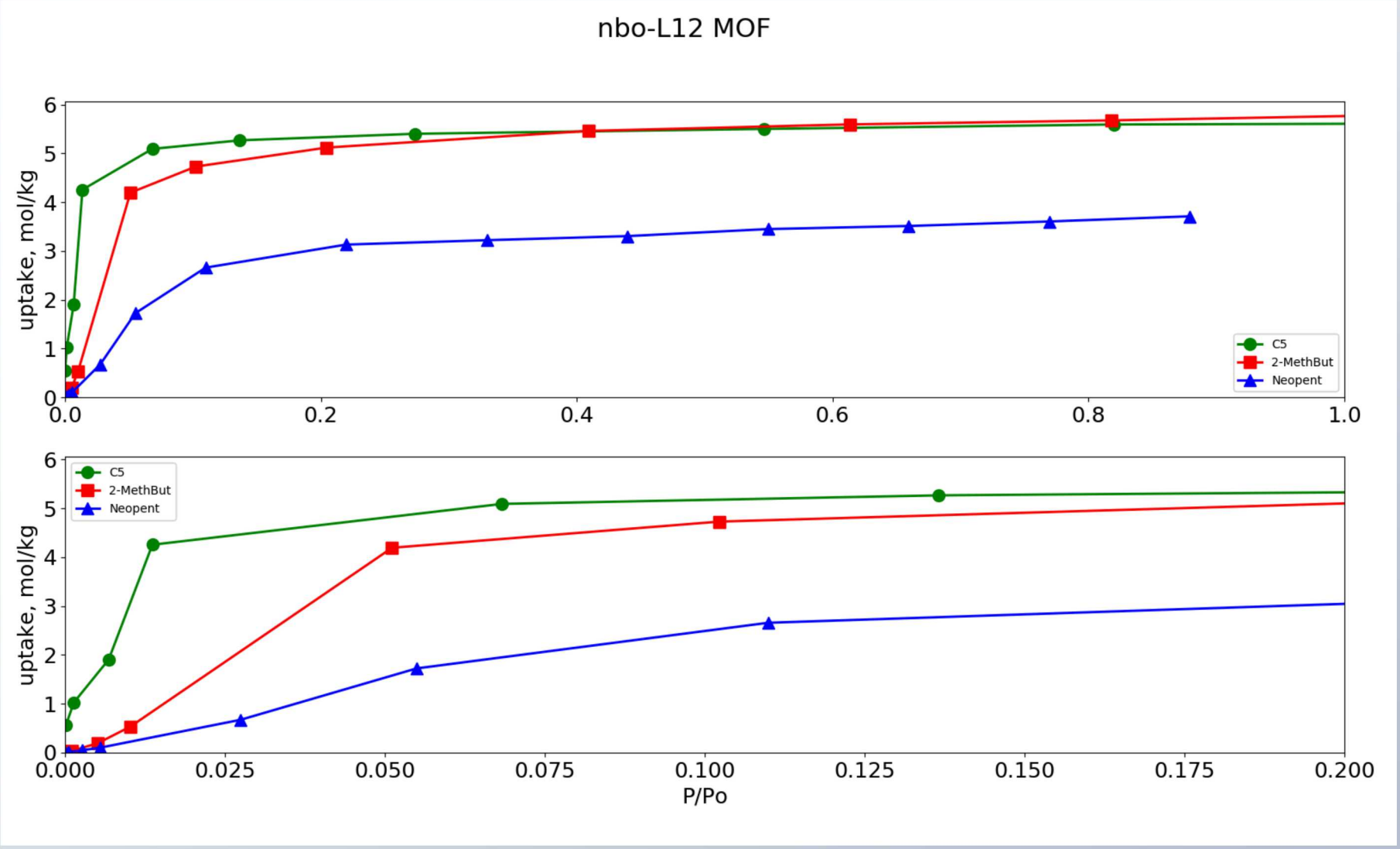




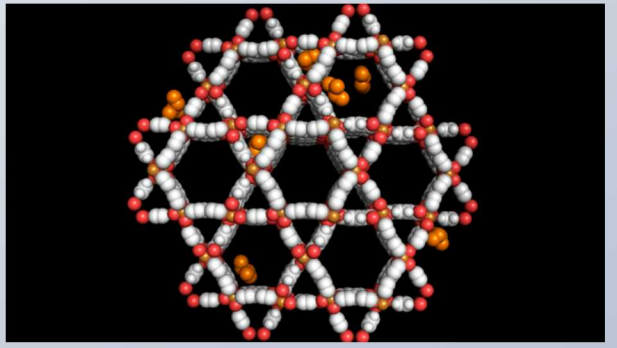
10,000 Pa

MOF nbo-L12 shows selectivity for n-pentane at low pressure. 2-methylbutane is adsorbed into the small pores.

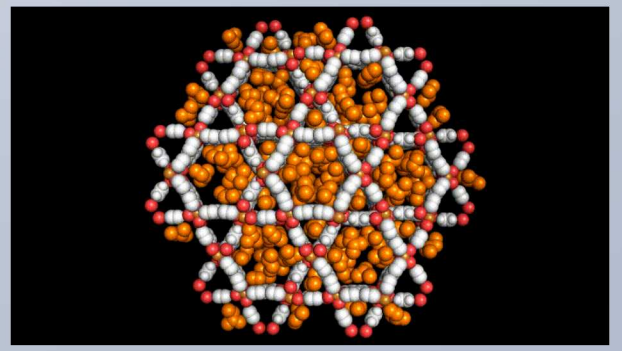
At saturation, there is less neopentane because it is excluded from the small pores.



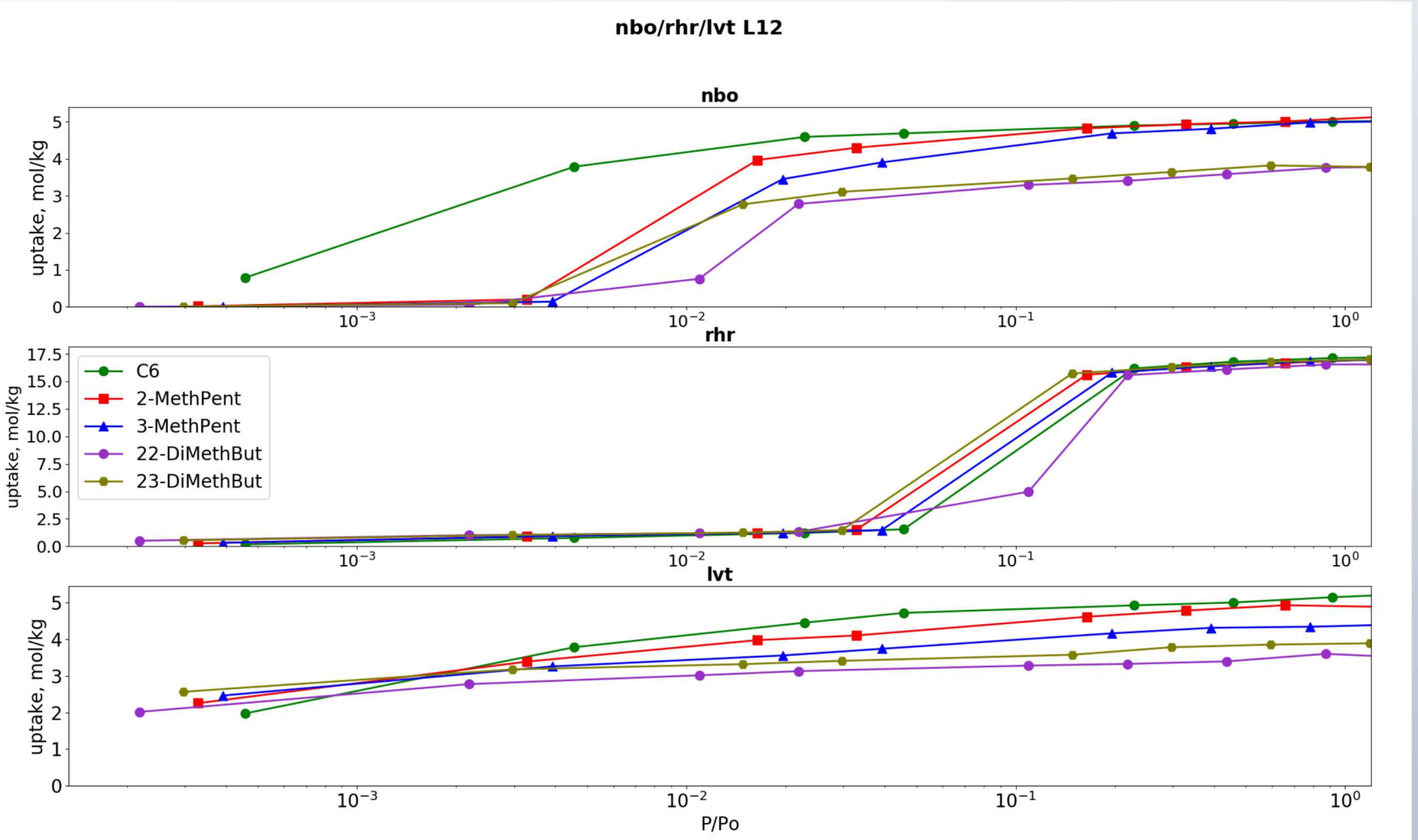
0.1 bar



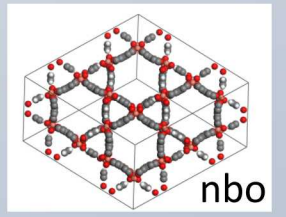
1.0 bar



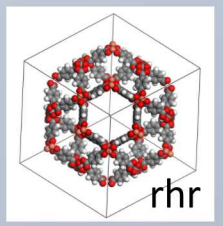
2-methylbutane
300 K
Single component isotherms



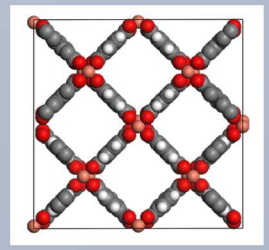
In nbo MOF, we see a separation of linear and 1-methyl isomers from dimethyl isomers.



No separation in the large pore rhr MOF.



lvt MOF shows reduction in bulkier molecules due to packing in the square pores



C6 isomers
300 K

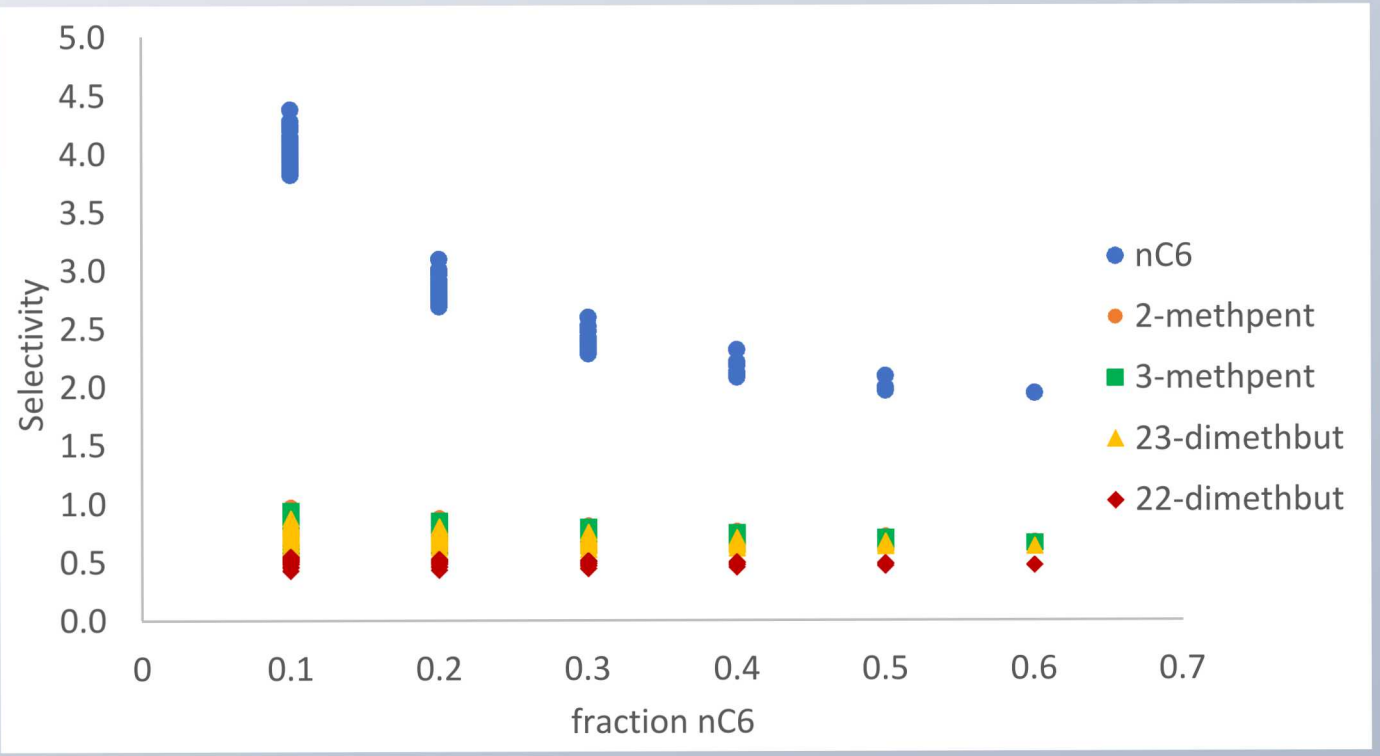
Multi-component Adsorption

GCMC simulations for a 5-component mixture containing all C6 isomers.
433 K, 1 bar 100k/100k cycles

Adsorbate	Selectivity
nC6	2.8
2-methpent	0.79
3-methpent	0.76
23-dimethbut	0.74
22-dimethbut	0.51
L+MP/DMB	1.86

Equimolar mixture

Linear + monobranched selectivity: 1.86

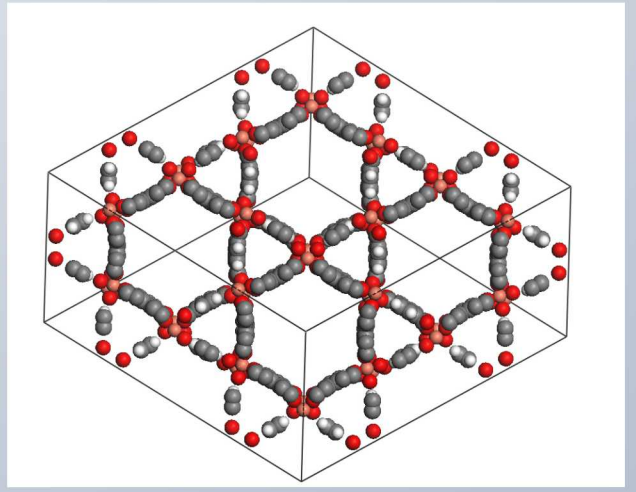


nC6 selectivity goes down when it forms a higher fraction of mixture

Thermodynamics

	H_ads				ΔH_ads [linear-X]		
	300 K	kJ/mol			300 K	kJ/mol	
	nbo_L12	rhr_L12	lvt_L12		nbo_L12	rhr_L12	lvt_L12
C5	-65.3	-36.9	-36.1		0.0	0.0	0.0
2-methylbutane	-30.1	-40.7	-36.7		35.2	-3.8	-0.6
neopentane	-23.7	-38.7	-33.0		41.6	-1.8	3.1
C6	-76.7	-41.6	-42.5		0.0	0.0	0.0
2-methylpentane	-40.4	-44.6	-42.6		36.4	-2.9	-0.1
3-methylpentane	-36.7	-45.7	-42.9		40.0	-4.1	-0.4
22-dimethylbutane	-29.9	-47.8	-42.1		46.8	-6.1	0.5
23-dimethylbutane	-31.8	-48.3	-43.5		44.9	-6.7	-1.0

RED: more favored than linear alkane
BLUE: less favored than linear alkane



Changing the topology can flip the trends in selectivity

Enthalpy of adsorption shows that nbo MOF is more selective for n-hexane by ~40 kJ/mol

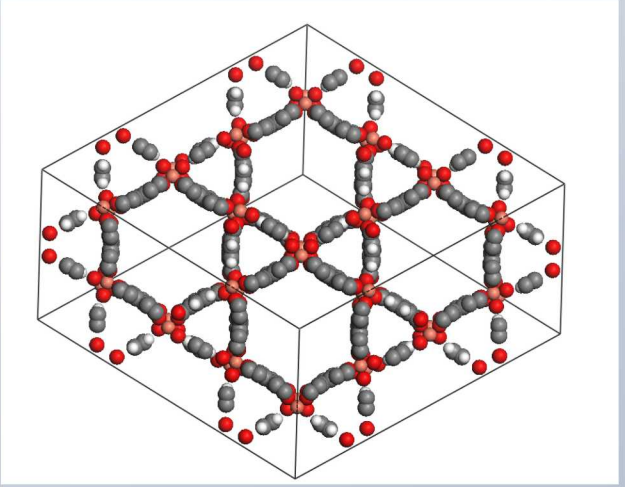
rhr MOF shows some slight selectivity for the branched isomers over the linear isomer

There is very little selectivity for the lvt MOF

Thermodynamics

Calculated from Henry constant data.

	nbo_L12				ΔH_ads [linear-X]		
	300 K	kJ/mol			300 K	kJ/mol	
	H_ads	-TΔS	A_ads		H_ads	-TΔS	A_ads
C5	-65.3	33.8	-34.0		0.0	0.0	0.0
2-methylbutane	-30.1	14.8	-17.8		35.2	-19.0	16.3
neopentane	-23.7	11.0	-15.2		41.6	-22.7	18.8
C6	-76.7	38.2	-41.1		0.0	0.0	0.0
2-methylpentane	-40.4	20.1	-22.7		36.4	-18.0	18.3
3-methylpentane	-36.7	17.1	-22.1		40.0	-21.1	19.0
22-dimethylbutane	-29.9	12.4	-20.0		46.8	-25.7	21.1
23-dimethylbutane	-31.8	13.2	-21.1		44.9	-25.0	19.9



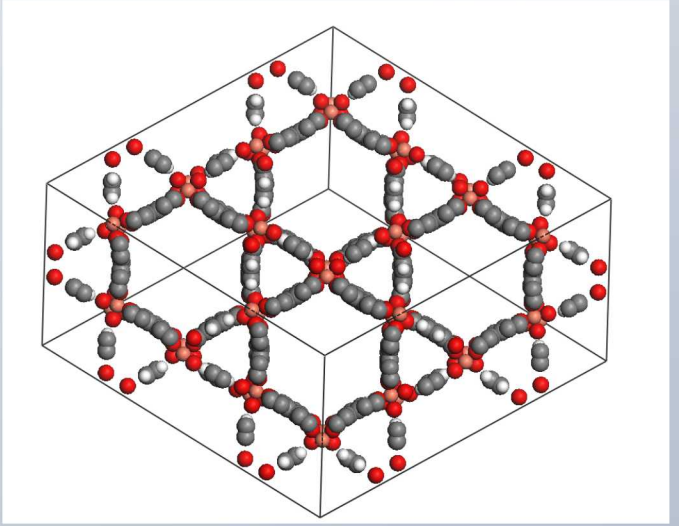
There is a large entropic penalty for the n-pentane and n-hexane molecules to go into the small pore in the nbo MOF.

Also the free energy for the mono and dibranched isomers is about the same due to the larger cost for the monobranched isomers.

Thermodynamics

Calculated from Henry constant data.

	TdelS				A_ads		
	kJ/mol, 300 K				kJ/mol, 300 K		
	nbo_L12	rhr_L12	lvt_L12		nbo_L12	rhr_L12	lvt_L12
C5	33.8	18.4	11.6		-34.0	-21.0	-27.0
2-methylbutane	11.0	19.6	11.0		-15.2	-21.6	-24.5
neopentane	14.8	20.7	11.6		-17.8	-22.5	-27.6
Spread [max - min]	22.7	2.3	0.7		18.8	1.5	3.0
C6	38.2	19.8	13.3		-41.1	-24.3	-31.7
2-methylpentane	20.1	21.5	13.0		-22.7	-25.5	-32.1
3-methylpentane	17.1	21.7	12.7		-22.1	-26.5	-32.7
22-dimethylbutane	12.4	22.0	12.4		-20.0	-28.3	-32.2
23-dimethylbutane	13.2	22.4	12.7		-21.1	-28.4	-33.3
Spread [max - min]	25.7	2.6	1.0		21.1	4.1	1.6

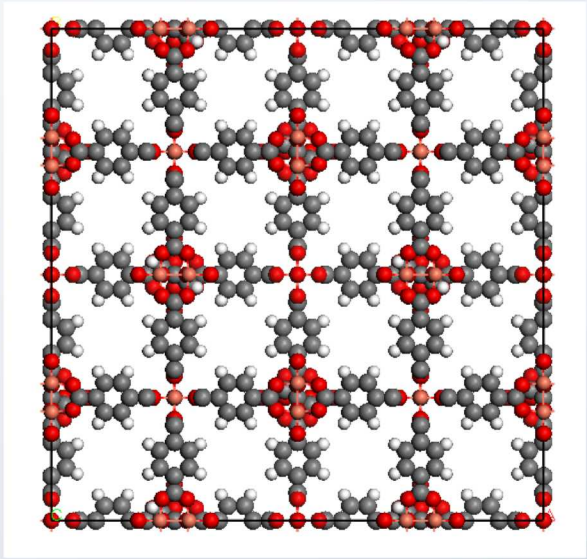


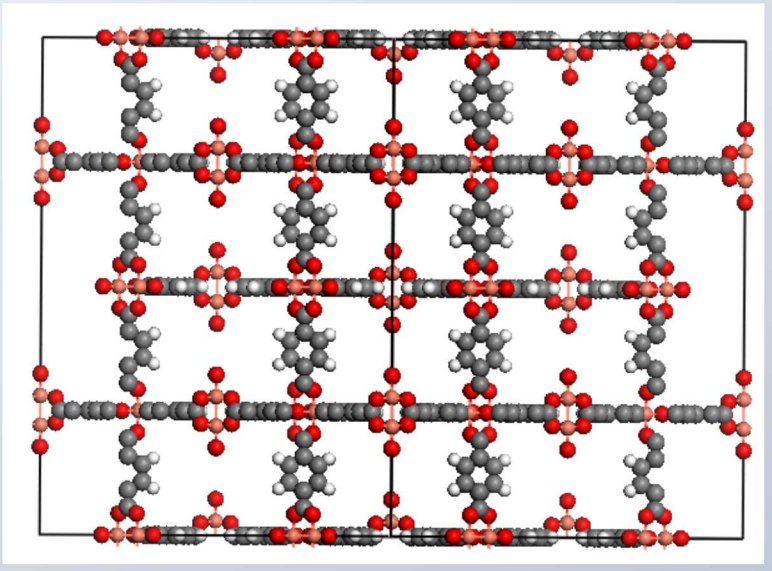
The rhr and lvt MOFs show little differentiation based on entropy.

The slight selectivity for branched isomers in the rhr MOF is driven by enthalpy.

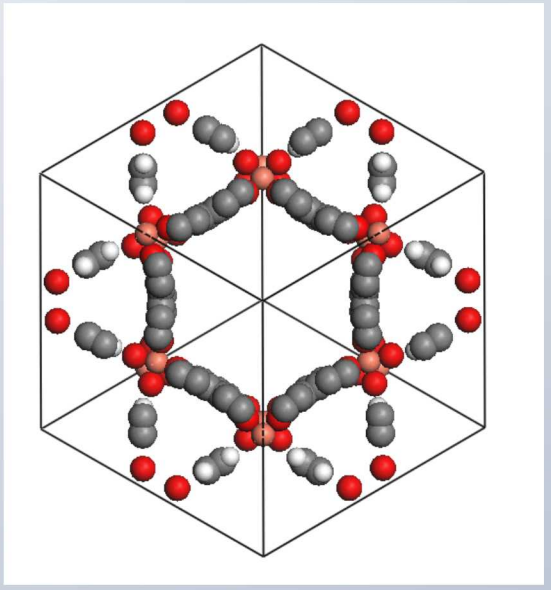
The nbo topology has two distinct conformations that are the same topology.

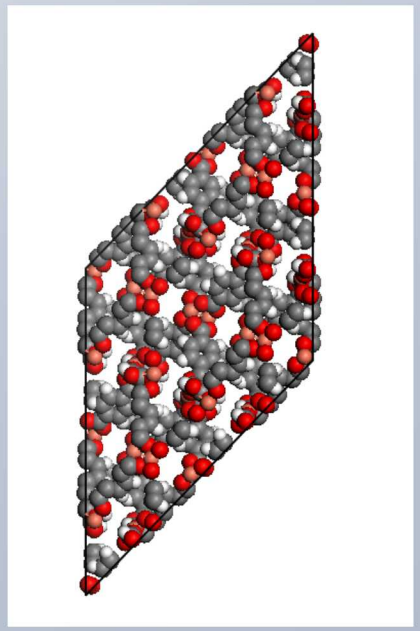
nbo cube



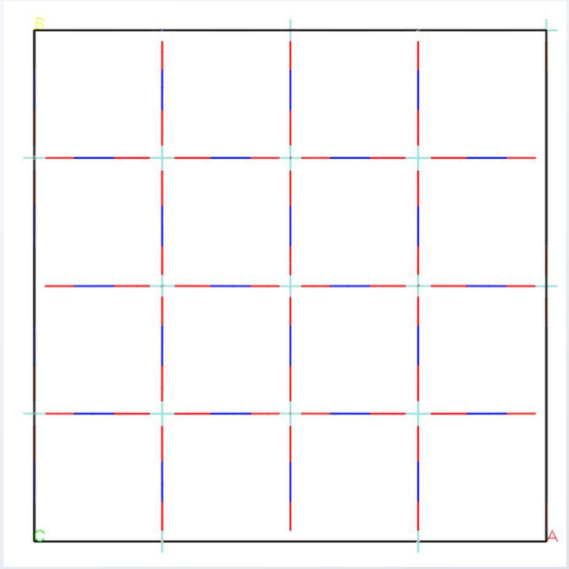


nbo star



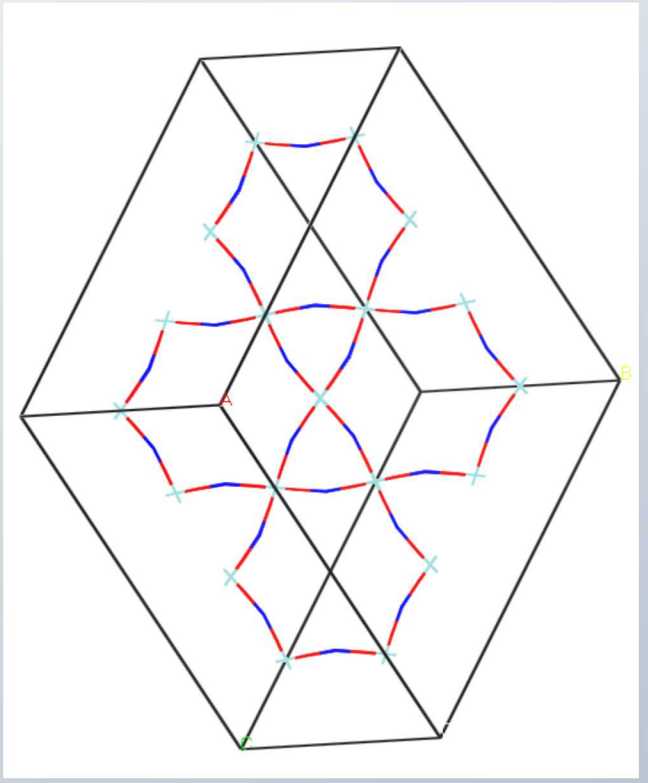


simplified nbo topology

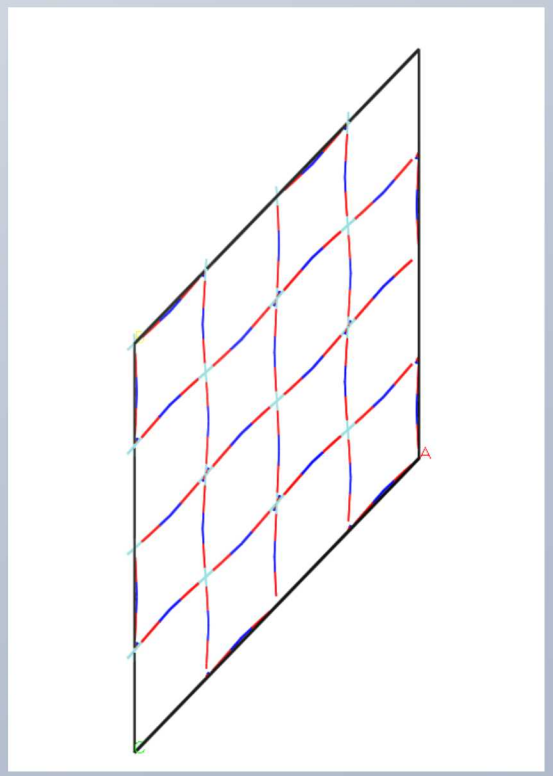


The connectivity is the same in each conformation. The simplified maps are equivalent.

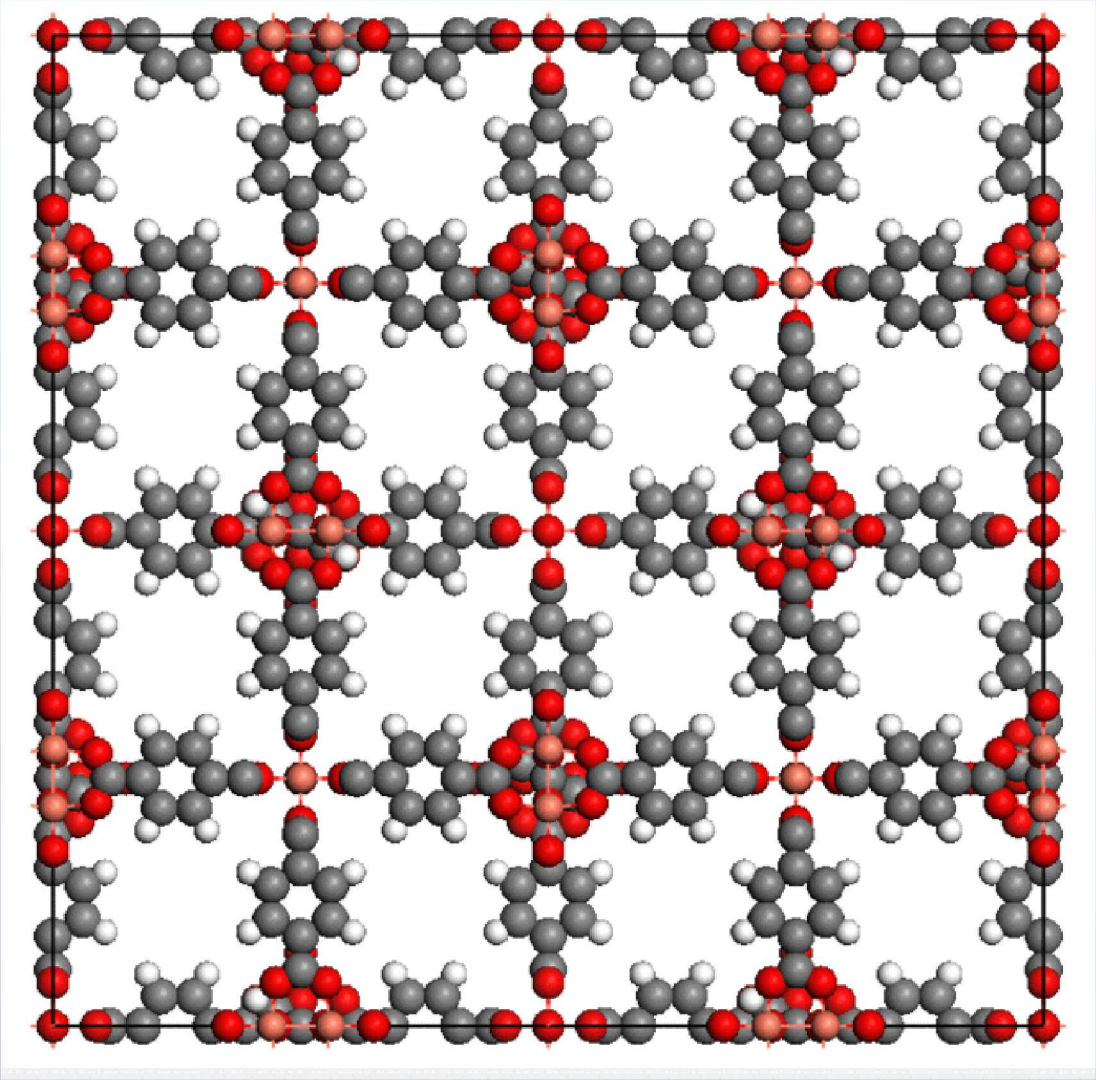
nbo cube



nbo star

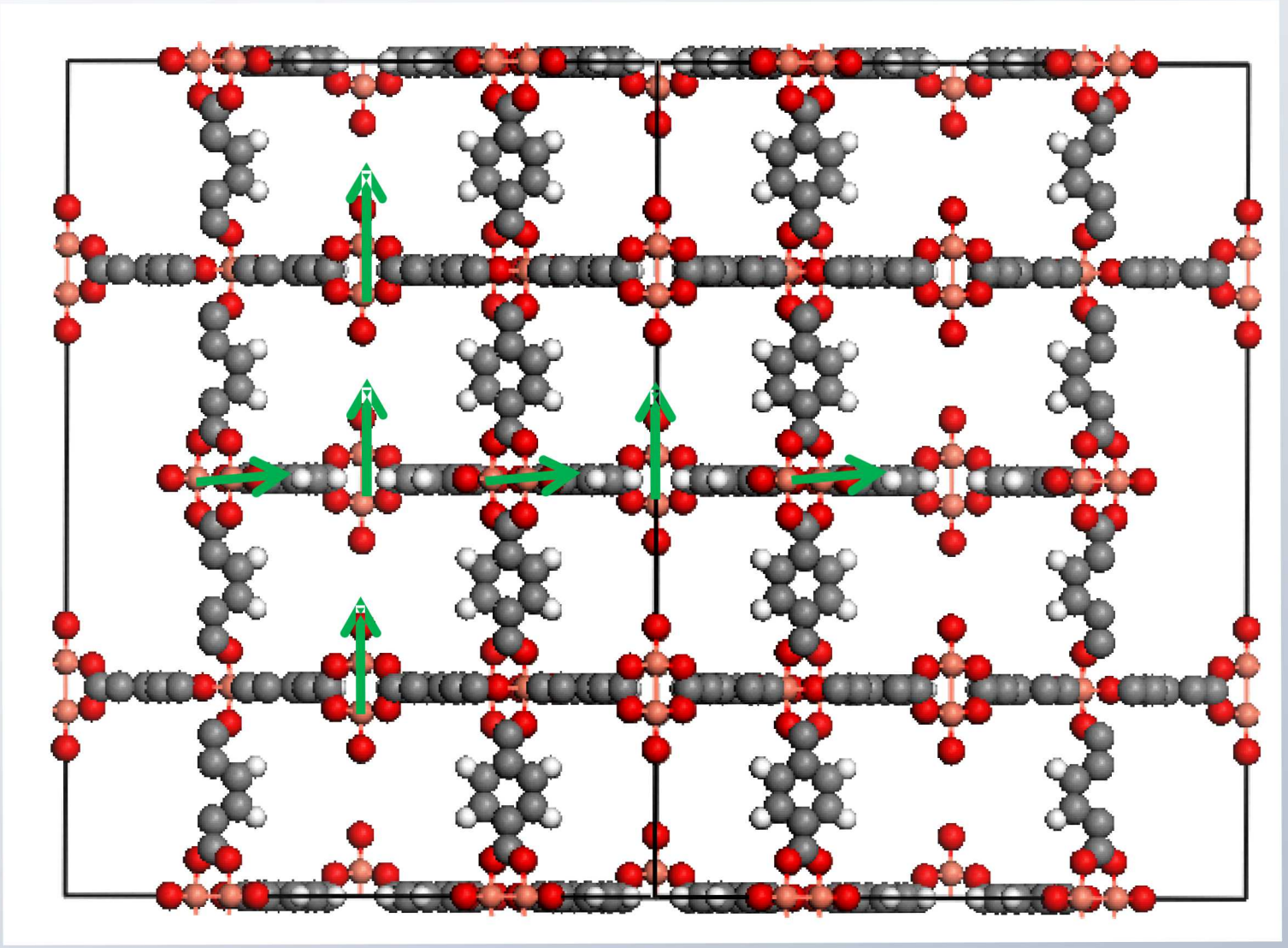


You can see how the star confirmation forms by looking down the [111] view of the cube confirmation.

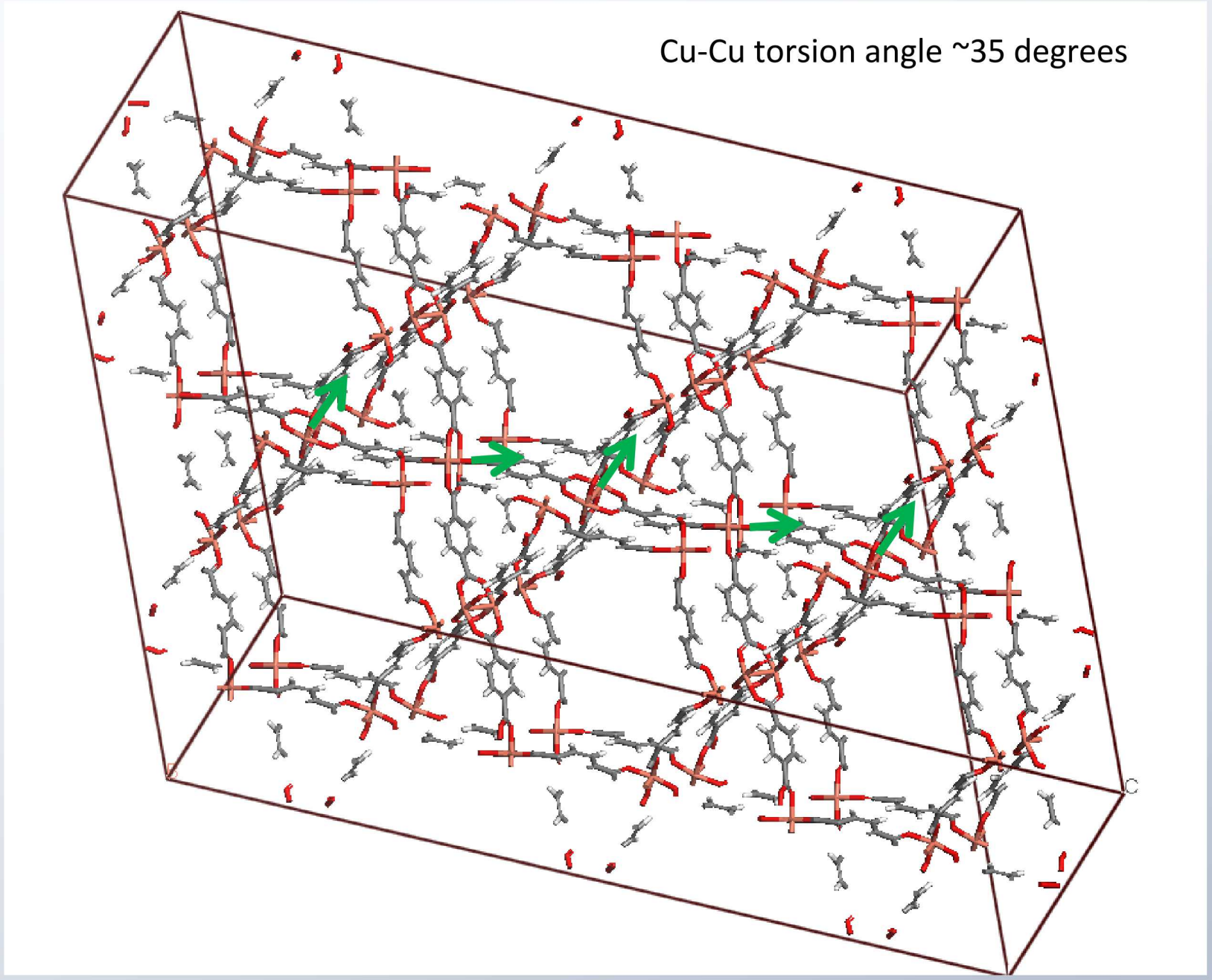


The star confirmation seen in MOF nbo_L12 is formed by compressing the MOF along the [111] axis so the

Cu-Cu vectors in alternating directions

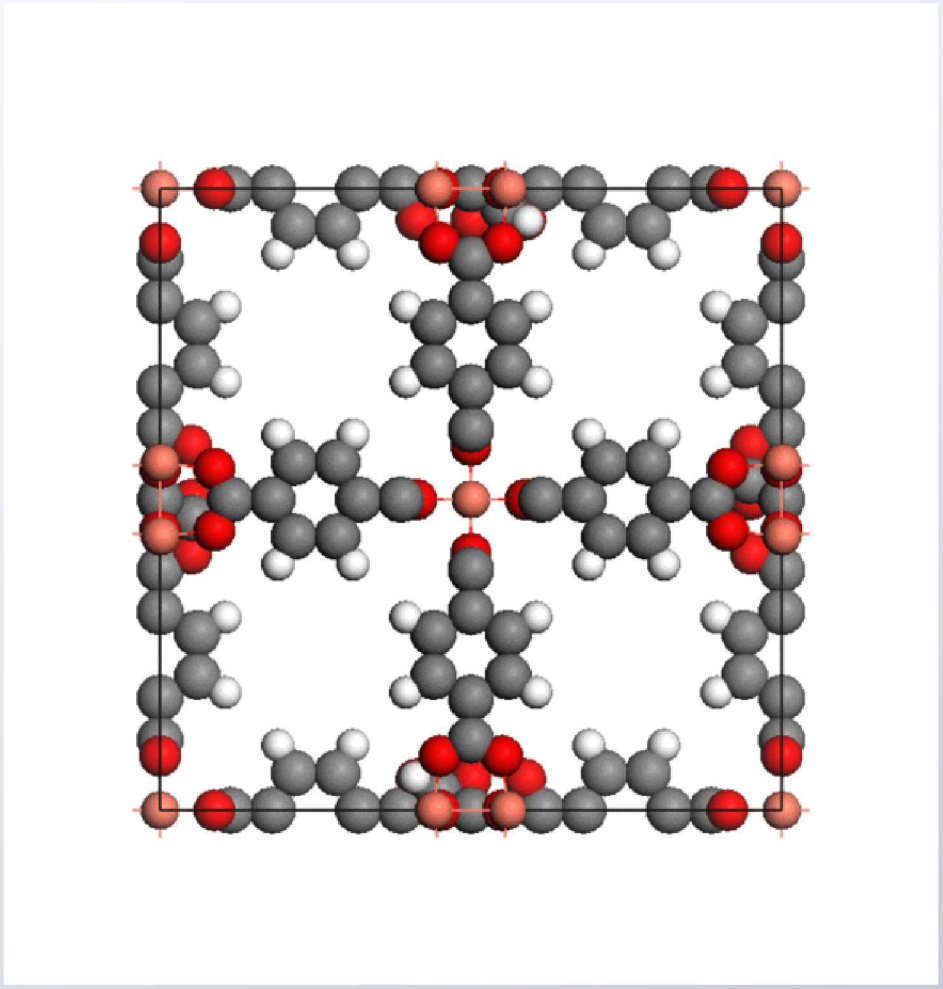


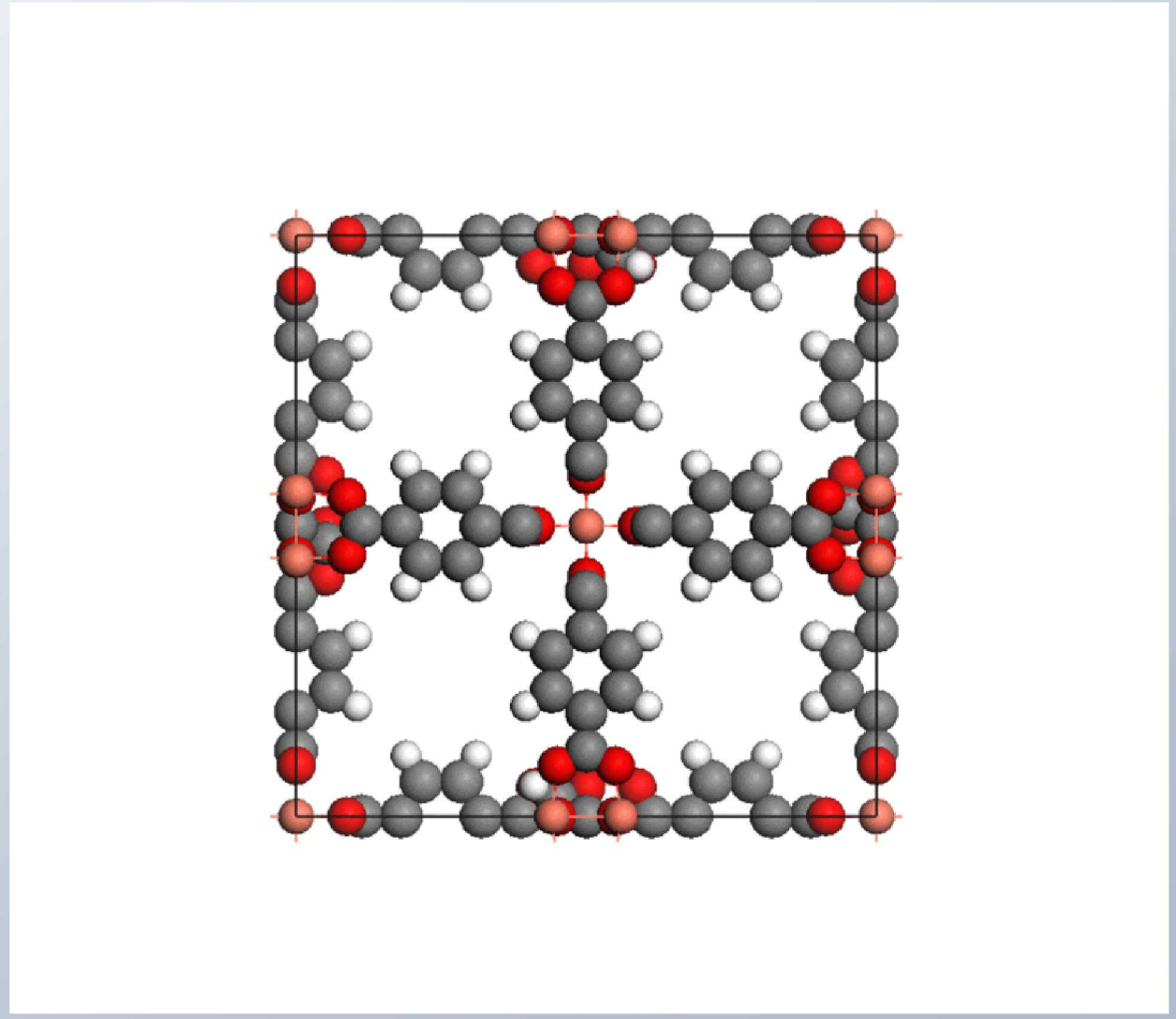
In square nbo confirmation the Cu paddlewheel nodes are aligned normal to their nearest neighbor node.



In the nbo star confirmation the Cu paddlewheel nodes are twisted to a 35 degree angle from their neighbors.

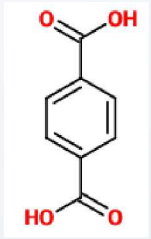
Geometrical minimization of MOF nbo_L12

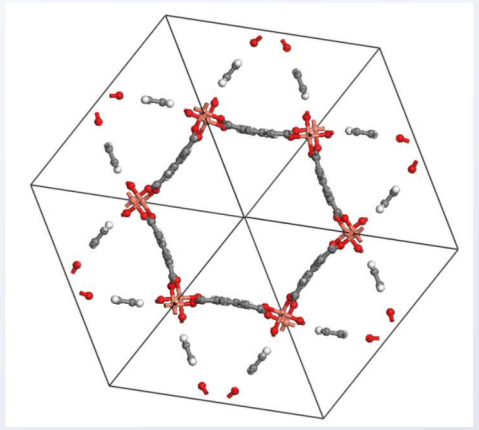




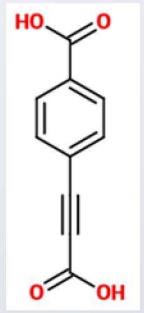
The conformation is determined by steric hindrance between the linker and node.

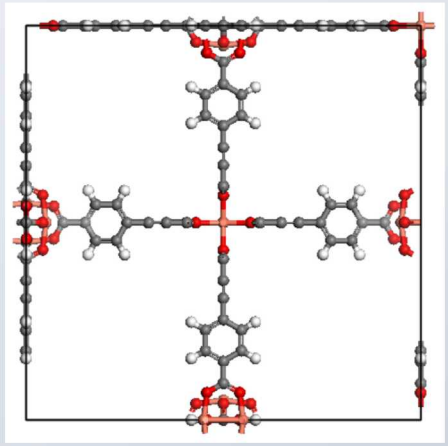
L12





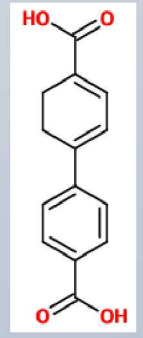
L20



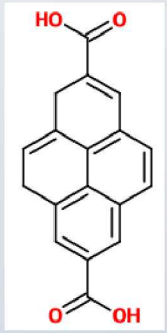


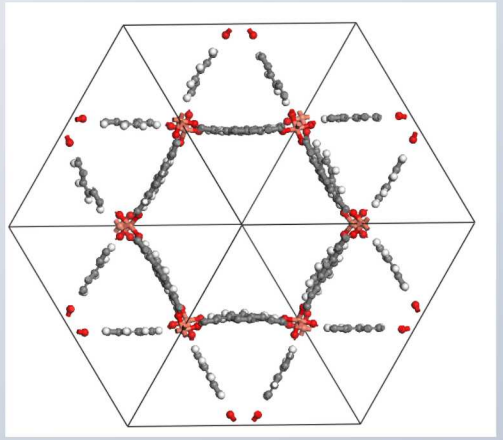
When the nodes are weakly coupled by hindrance the MOF takes on a slightly skewed confirmation.

L24

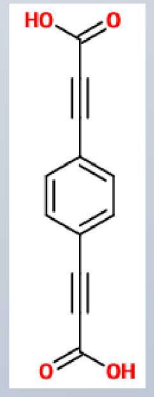


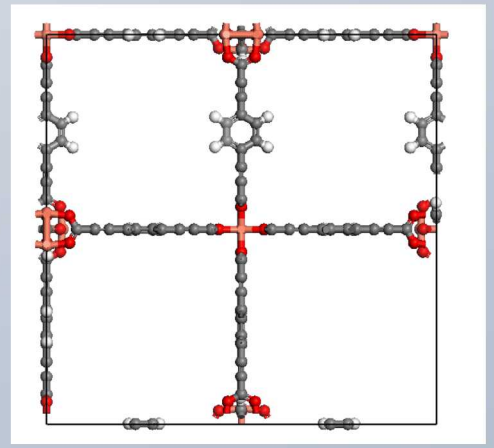
L22

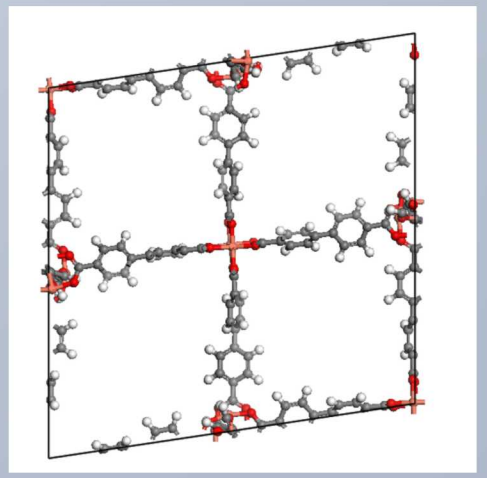




L34

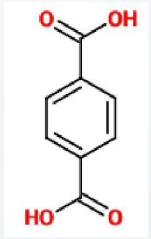


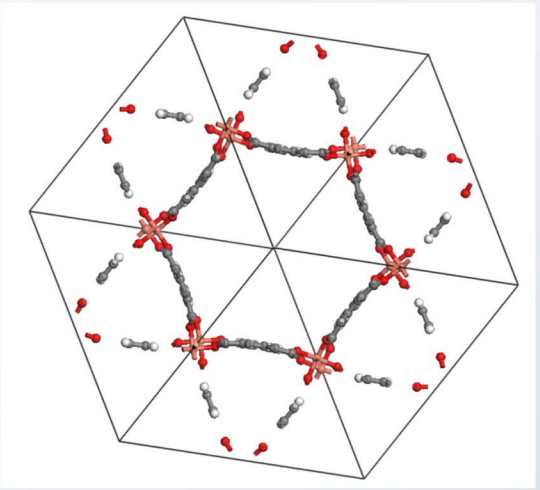




DFT calculations show that the star confirmation is more favorable for certain linkers.

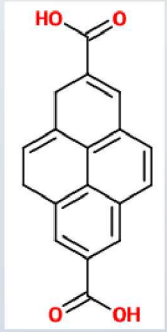
L12

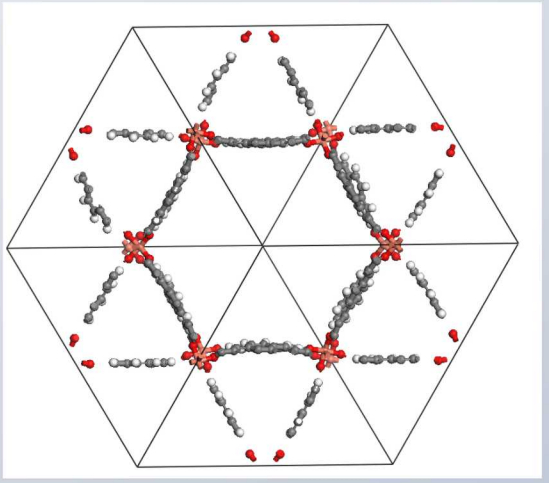




189 kJ/mol (per unit cell) more favorable than square

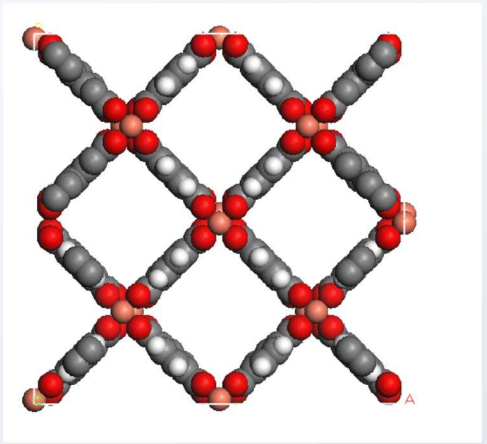
L22



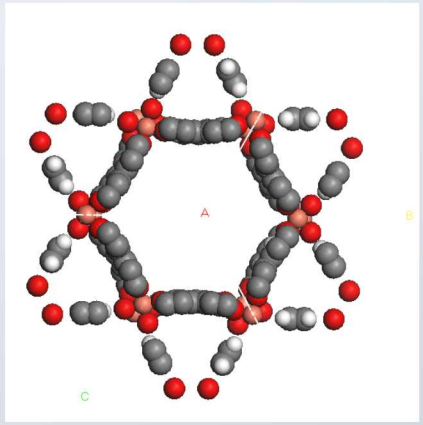


83 kJ/mol (per unit cell) more favorable than square

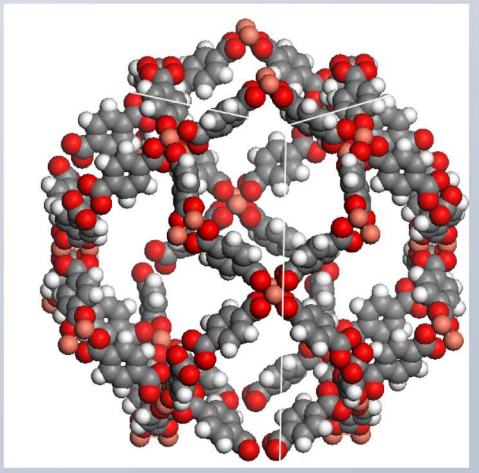
lvt topology is most favorable, similar to nbo
rhr is the least favorable.



lvt



nbo



rhr

	kJ/mol diff
Topology	**$CuC_8H_4O_4$**
lvt	0
nbo	1.2
rhr	17.7

*based on UFF optimization in Materials Studio

Conclusions

1. Topology can greatly affect MOF pore size and shape, which has implications for separations

2. nbo/lvt/rhr polymorph family has different linear/branched selectivity depending on topology

3. nbo MOFs with phenyl ring linker show high selectivity for linear alkanes due to small pores

4. nbo topology can form different conformations (cube vs star) depending on steric effects of the linker

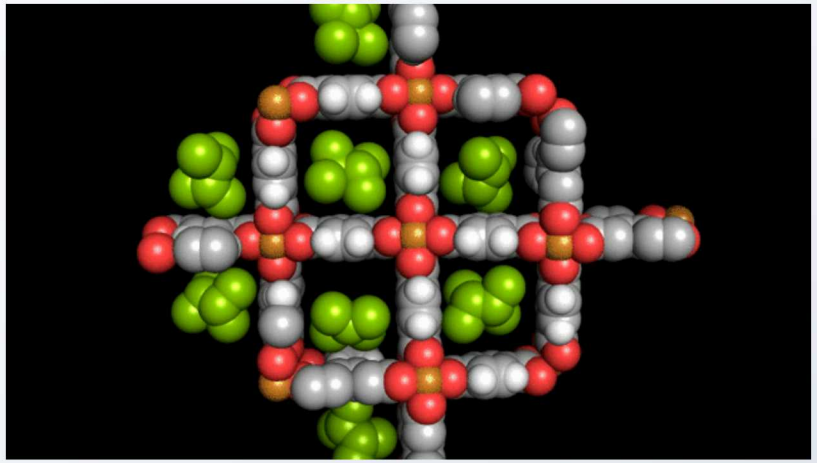
Acknowledgements:
This work was supported by the US Department of Energy under Award DE-FG02-08ER15967.

Vapor pressure at 300 K

Taken from Antoine equation

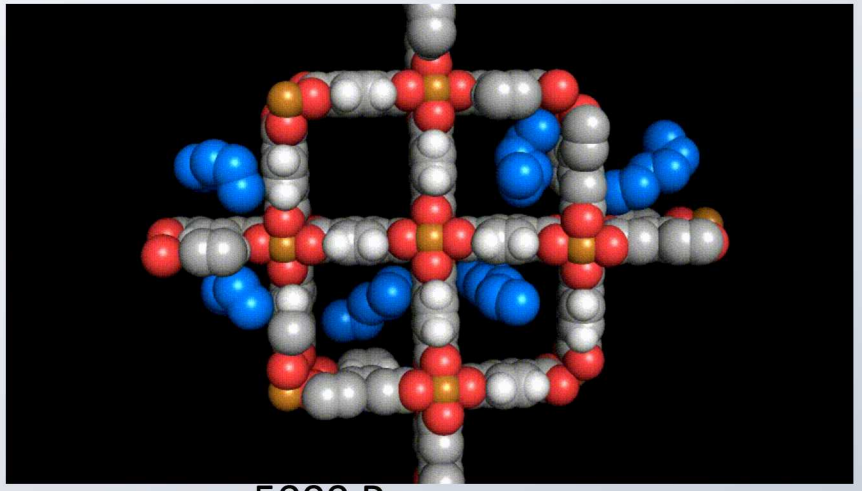
	kPa	bar
[C1]	-	-
[C2]	4398	43.98
[C3]	1023	10.23
[C4]	261	2.61
[C5]	73	0.73
[C6]	22	0.22
[isobutane]	377	3.77
[neopentane]	182	1.82
[2-methylbutane]	98	0.98
[2-methylpentane]	30	0.30
[3-methylpentane]	25	0.25
[22-dimethylbutane]	46	0.46
[23-dimethylbutane]	34	0.34

dimethyl-butane



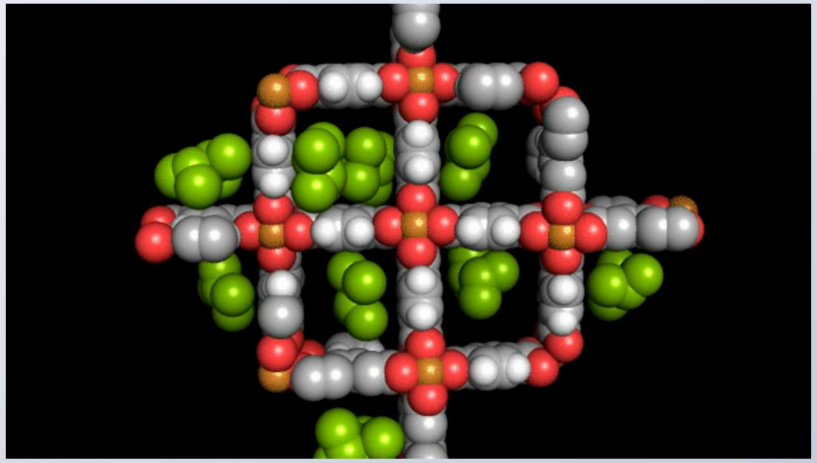
1000 Pa

n-hexane

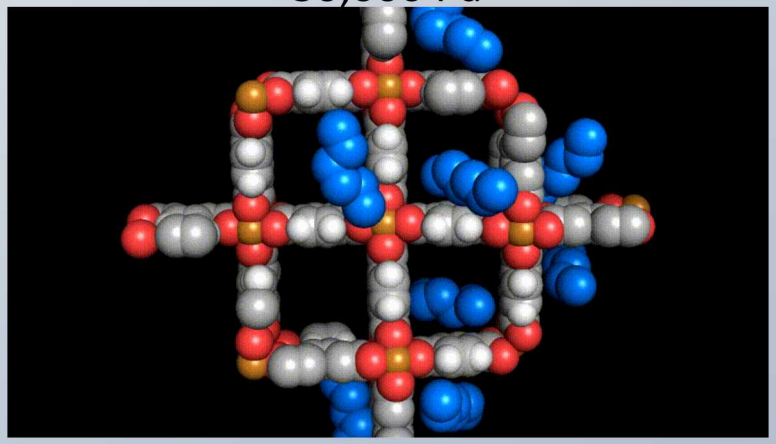


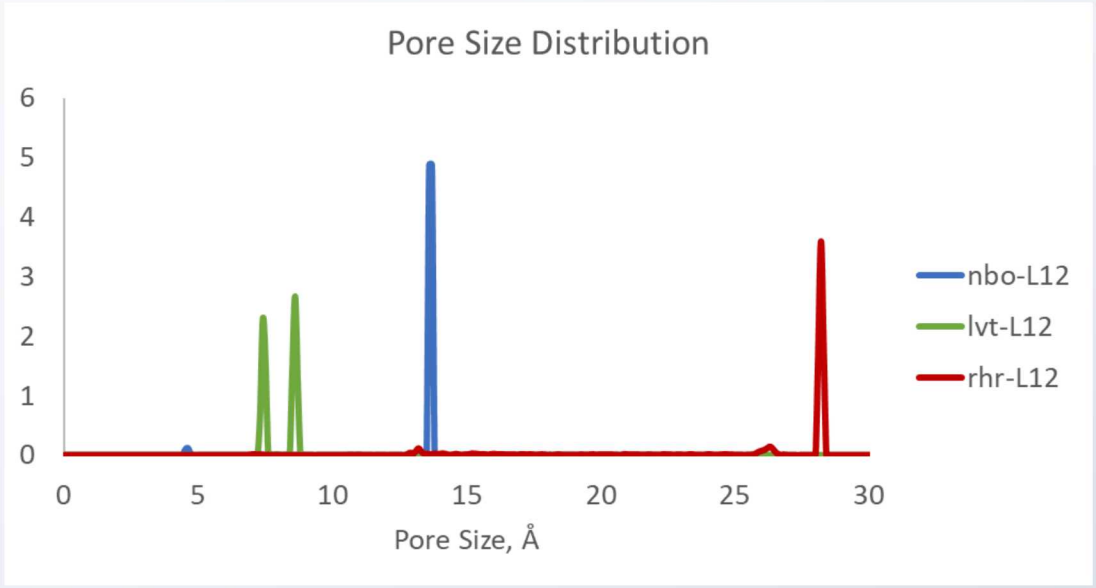
5000 Pa

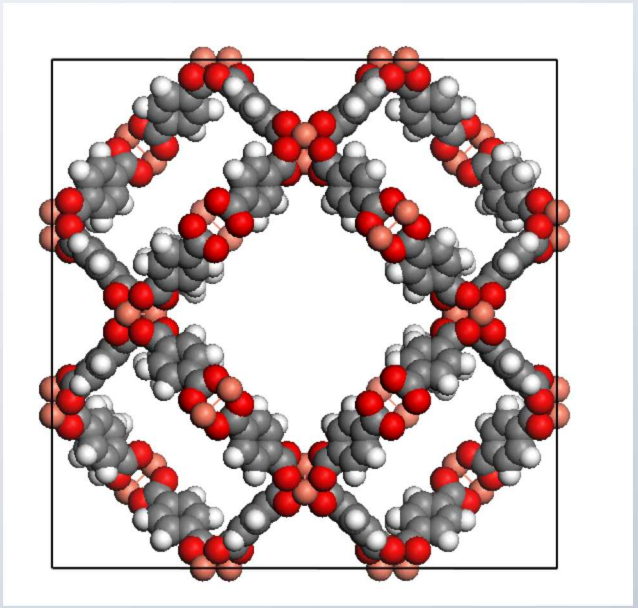
30,000Pa



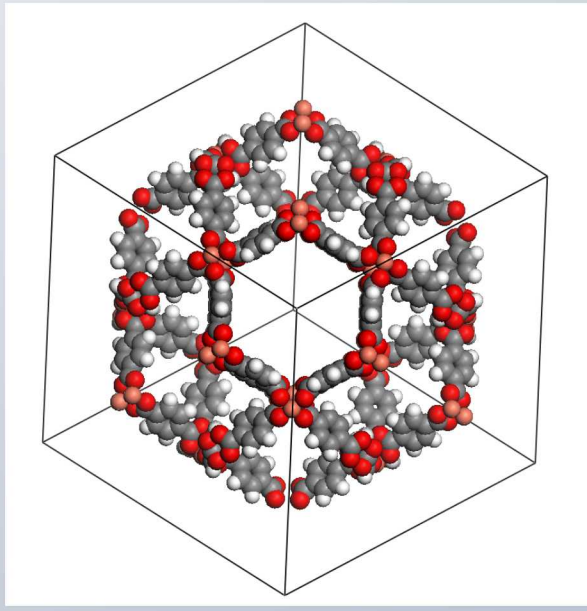
30,000 Pa

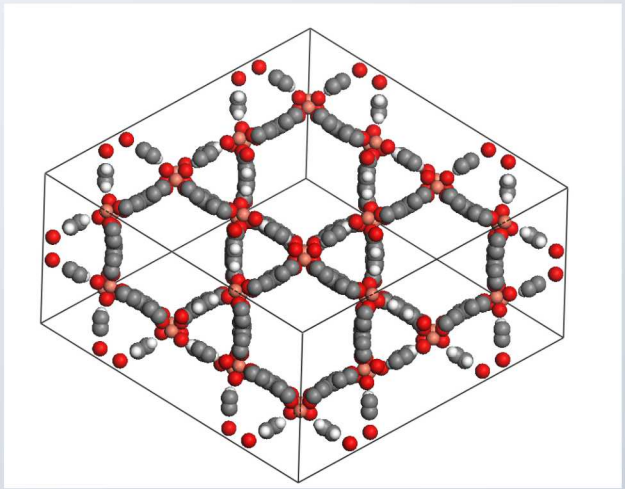




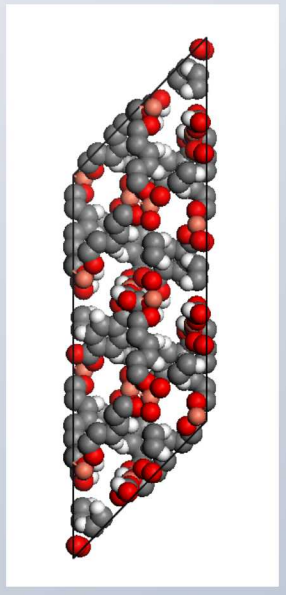


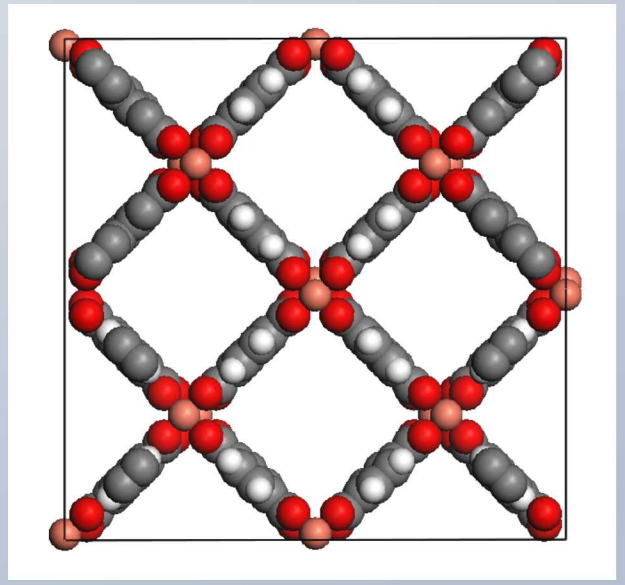
rhr_L12



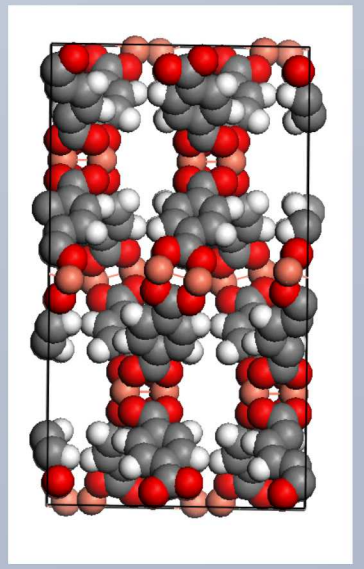


nbo_L12

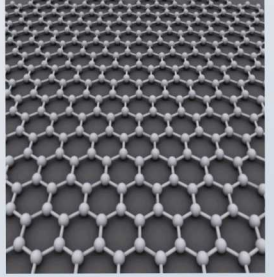




lvt_L12



Changing the topology can invert the selectivity from favoring linear to favoring branched isomers



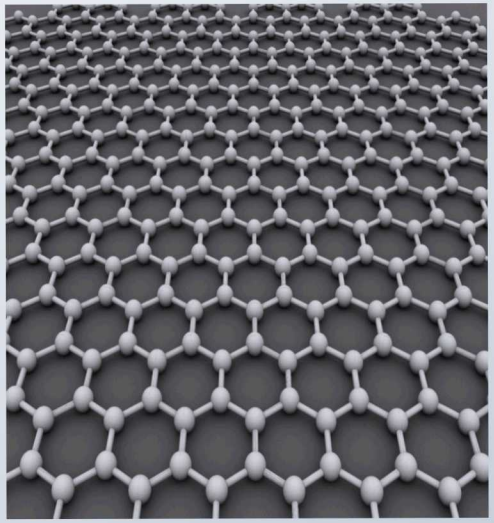
	Graphene	lvtb_L_12	nbob_L_12	rhrb_L_12
[C1]	-10.8	-11.1	-18.2	-7.8
[C2]	-17.1	-16.3	-29.8	-15.5
[C3]	-20.9	-20.0	-37.9	-22.0
[C4]	-23.4	-21.4	-45.3	-23.5
[isobutane]	-22.2	-22.3	-16.5	-27.0
[C5]	-26.9	-23.4	-52.4	-24.2
[2-methylbutane]	-24.7	-24.9	-18.1	-28.9
[neopentane]	-16.7	-21.8	-12.6	-27.5
[C6]	-30.2	-25.5	-60.0	-24.3
[2-methylpentane]	-29.9	-27.7	-25.1	-29.5
[3-methylpentane]	-27.7	-27.8	-20.9	-30.4
[22-dimethylbutane]	-20.8	-27.1	-14.9	-32.7
[23-dimethylbutane]	-24.2	-26.8	-14.9	-31.7

	lvtb_L_12	nbob_L_12	rhrb_L_12
[C1]	-0.2	-7.4	3.0
[C2]	0.8	-12.7	1.6
[C3]	1.0	-16.9	-1.1
[C4]	2.0	-21.9	-0.1
[isobutane]	-0.1	5.6	-4.8
[C5]	3.4	-25.5	2.6
[2-methylbutane]	-0.2	6.5	-4.3
[neopentane]	-5.1	4.1	-10.8
[C6]	4.7	-29.8	6.0
[2-methylpentane]	2.1	4.7	0.4
[3-methylpentane]	-0.2	6.8	-2.8
[22-dimethylbutane]	-6.3	6.0	-11.8
[23-dimethylbutane]	-2.6	9.3	-7.5

RED: stronger than graphene

BLUE: weaker than graphene

Changing the topology can invert the selectivity from favoring linear to favoring branched isomers

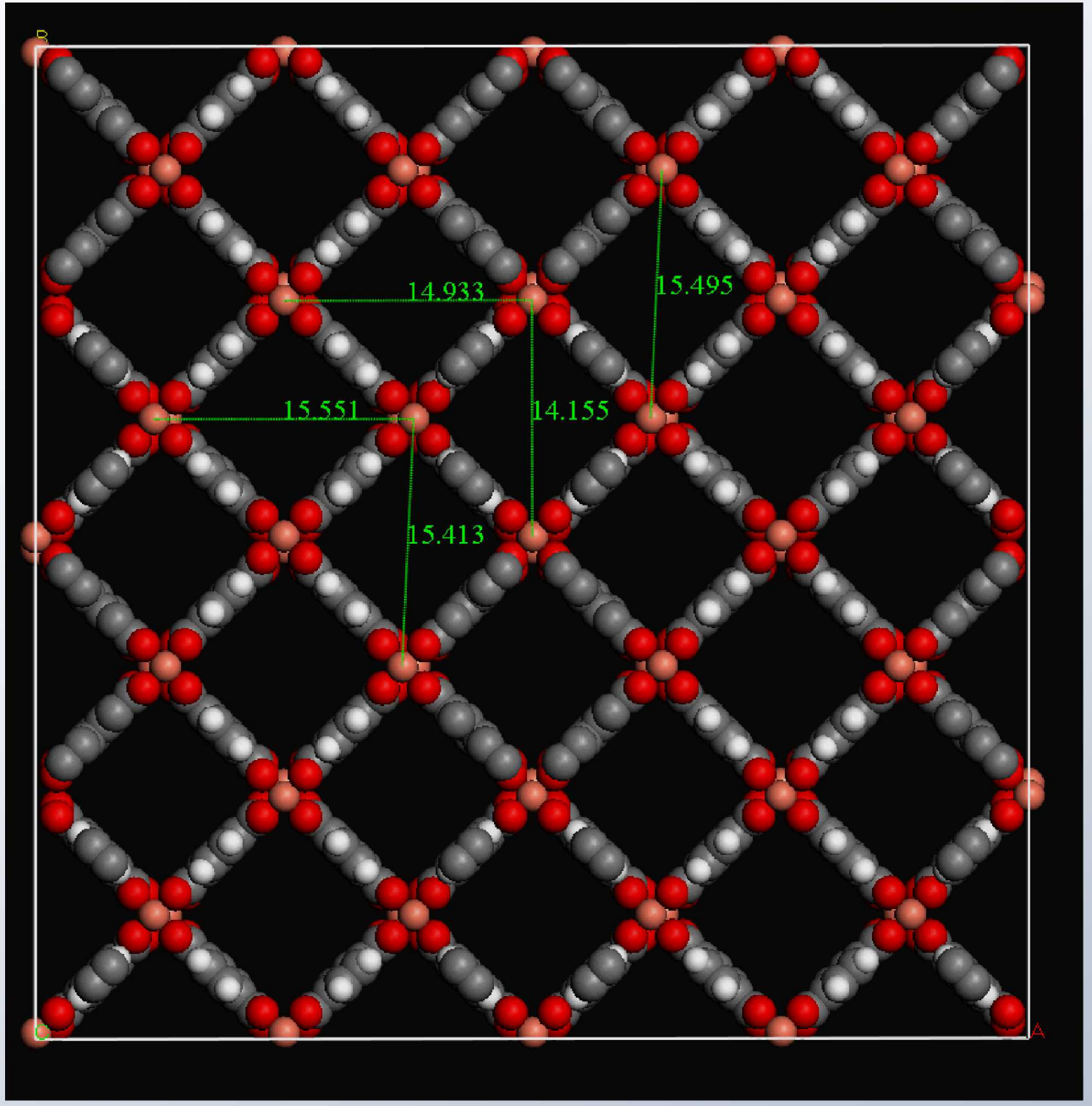


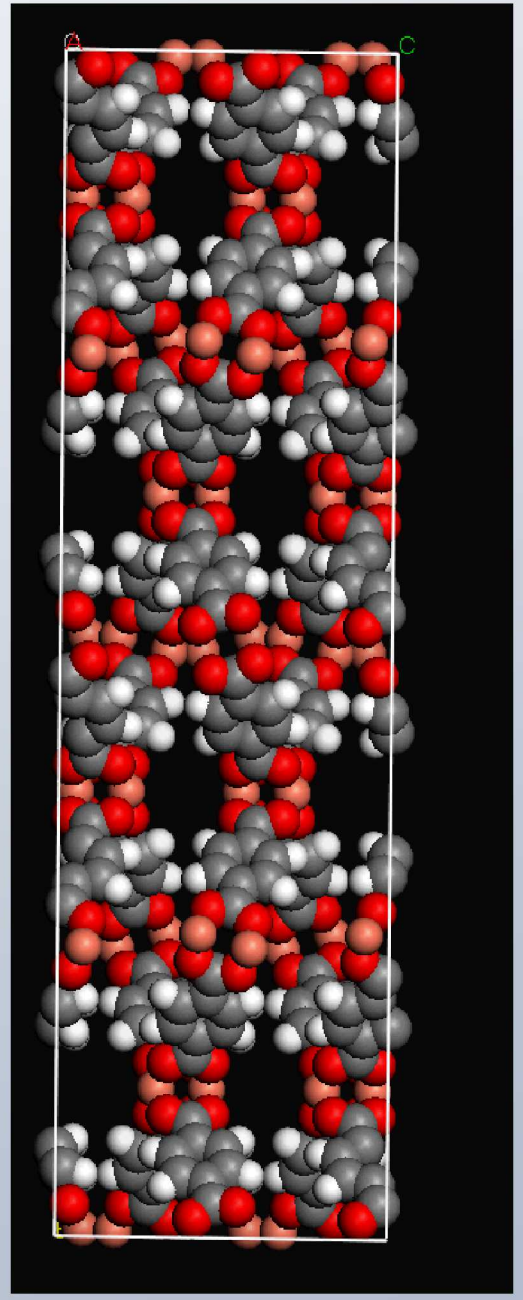
	Graphene	lvtb_L_12	nbob_L_12	rhrb_L_12
[C1]	0.0	0.0	0.0	0.0
[C2]	6.3	5.2	11.6	7.6
[C3]	10.1	8.9	19.6	14.2
[C4]	0.0	0.0	0.0	0.0
[isobutane]	1.2	-0.9	28.7	-3.5
[C5]	0.0	0.0	0.0	0.0
[2-methylbutane]	2.2	-1.4	34.3	-4.7
[neopentane]	10.2	1.6	39.8	-3.2
[C6]	0.0	0.0	0.0	0.0
[2-methylpentane]	0.3	-2.2	34.8	-5.2
[3-methylpentane]	2.5	-2.3	39.1	-6.2
[22-dimethylbutane]	9.4	-1.6	45.1	-8.4
[23-dimethylbutane]	6.0	-1.3	45.1	-7.5

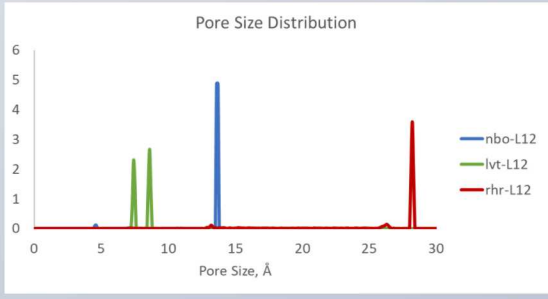
RED: more favored than linear alkane

BLUE: less favored than linear alkane

Lvt_L12


B
C
A
14.933
15.495
15.551
14.155
15.413


B
A
C


Pore Size Distribution
6
5
4
3
2
1
0
0
5
10
15
20
25
30
Pore Size, Å
nbo-L12
lvt-L12
rhr-L12