

Supporting Information

**Metal-Organic Frameworks (MOFs) beyond Crystallinity:
Amorphous MOFs, MOF Liquids and MOF Glasses**

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Table S1. Reported aMOFs.

aMOF	Stress	Highlights & Considerations	Ref.
a _p -ZIF-8	Pressure 0.34 GPa	<ul style="list-style-type: none"> • First porous aMOF • Accessible pore volume • Storage of I₂ 	[3,18]
a _p -ZIF-8	Pressure 1.9 GPa	<ul style="list-style-type: none"> • <i>In situ</i> TEM study • Irreversible morphological transition • BET surface area: 253 m²/g 	[17]
a _p -ZIF-8	Pressure 1.6 GPa Pressure 39 GPa	<ul style="list-style-type: none"> • Reversible amorphization • Irreversible amorphization 	[19]
a _p -ZIF-4	Pressure 0.35-0.98 GPa	<ul style="list-style-type: none"> • Reversible amorphization 	[28]
a _p -MOF-5	Pressure 3.5 MPa	<ul style="list-style-type: none"> • Irreversible amorphization • BET surface area: 6 m²/g 	[21]
a _p -MOF-5	Pressure 3.2 GPa	<ul style="list-style-type: none"> • DEF as pressure-transmitting medium 	[29]
a _p -MOF-177	Mechanical compression	<ul style="list-style-type: none"> • The maximum excess volumetric hydrogen storage uptake increased 80% 	[22]
a _p -Zn(CN) ₂	Pressure and X-ray 3.2 GPa	<ul style="list-style-type: none"> • Isopropanol as pressure-transmitting medium • No accessible pore volume 	[32]
a _p -Sc ₂ BDC ₃	Pressure 0.4 GPa Pressure 3 GPa	<ul style="list-style-type: none"> • Fluorinert-77 as pressure-transmitting medium • Reversible amorphization • Methanol as pressure-transmitting medium 	[33]
a _p -Sc ₂ (NO ₂ BDC) ₃	Pressure 2.6 GPa Pressure 3.3 GPa	<ul style="list-style-type: none"> • Fluorinert-77 as pressure-transmitting medium • Irreversible amorphization • Methanol as pressure-transmitting medium 	[33]
a _p -NH ₂ -MIL-53(In)	Pressure 23 GPa	<ul style="list-style-type: none"> • Irreversible amorphization 	[34]
a _p -UiO-66	Pressure 1.9 GPa	<ul style="list-style-type: none"> • Irreversible amorphization • BET surface area: 76 m²/g • Mechanical energy absorber 	[23]
a _p -Cu-BTC	Pressure 3.8 GPa	<ul style="list-style-type: none"> • Shock compression • Mechanical energy absorber 	[25]
a _p -Cu-BTC-Fc	Pressure 5.8 GPa	<ul style="list-style-type: none"> • Shock compression • Mechanical energy absorber 	[25]
a _p -Co ₂ (BDC) ₂ DABCO-4DMF·H ₂ O	Pressure 1.9 GPa	<ul style="list-style-type: none"> • Piezochromic aMOF 	[35]
a _T -Cu(Im) ₂	Temperature 110 °C	<ul style="list-style-type: none"> • Decomposition of Cu(Him)₂(CO₃)₂·H₂O 	[2]
a _T -Ni(Im) ₂	Temperature 130 °C Temperature 260 °C	<ul style="list-style-type: none"> • Decomposition of [Ni(Him)₂(Im)(CH₃COO)] • Decomposition of Ni(acac)₂(Him)₂ 	[36]
a _T -Pd(Im) ₂	Temperature 150 °C	<ul style="list-style-type: none"> • Decomposition of Pd(Him)₂(im)₂ 	[36]
a _T -Pt(Im) ₂	Temperature 250 °C	<ul style="list-style-type: none"> • Decomposition of Pt(Him)₂(im)₂ 	[36]
a _T -[ZnI ₂] ₃ (TPT) ₂	Temperature 277 °C	<ul style="list-style-type: none"> • Precursor of a crystalline phase 	[37,38]
a _T -[(ZnBr ₂) ₃ (TPT) ₂] _n ·(H ₂ O) a _T -[(ZnCl ₂)(μ-Cl)(ZnCl)(TPT)] _n	Temperature 277 °C	<ul style="list-style-type: none"> • Precursor of a crystalline phase 	[41]
a _T -ZIF-4	Temperature 300 °C	<ul style="list-style-type: none"> • Precursor of a crystalline phase • Reversible amorphization 	[42,45]

a _T -ZIF-1 a _T -ZIF-3 a _T -Co-ZIF-4	Temperature 300 °C	<ul style="list-style-type: none"> • Precursor of a crystalline phase 	[45]
a _T -ZIF-8	Temperature 250-350 °C	<ul style="list-style-type: none"> • Amorphization thanks to the protection of surrounding Matrimid® • Resulting to MMMs with the highest CO₂/CH₄ selectivities ever reported • BET surface area: 6 m²/g 	[50]
a _T -Cu ^I [Cu ^{III} (pdt) ₂]	Temperature 120 °C	<ul style="list-style-type: none"> • Reversible amorphization • Conductive aMOF 	[51]
a _T -Mn-250	Temperature 250 °C	<ul style="list-style-type: none"> • Catalytic activity for CO oxidation 	[52]
a _T -HS-ZIF-67	Temperature 260 °C	<ul style="list-style-type: none"> • OER catalytic activity • Catalytic activity for nonenzymatic glucose sensing 	[53]
a _m -CAU-7 a _T -CAU-7	Mechanical Ball milling Temperature 180 °C	<ul style="list-style-type: none"> • Controlled release of drugs 	[59]
a _m -ZIF-1	Ball milling 25 Hz, 30 min	<ul style="list-style-type: none"> • Irreversible amorphization • BET surface area: 11 m²g⁻¹ 	[62]
a _m -ZIF-3	Ball milling 25 Hz, 30 min	<ul style="list-style-type: none"> • Irreversible amorphization • BET surface area: 21 m²g⁻¹ 	[62]
a _m -ZIF-4	Ball milling 25 Hz, 30 min	<ul style="list-style-type: none"> • Irreversible amorphization • BET surface area: 10 m²g⁻¹ 	[62]
a _m -ZIF-8	Ball milling 25 Hz, 30 min	<ul style="list-style-type: none"> • Irreversible amorphization • BET surface area: 56 m²g⁻¹ 	[62]
a _m -ZIF-8	Ball milling 30 Hz, 30 min	<ul style="list-style-type: none"> • Irreversible amorphization • BET surface area: 55 m²g⁻¹ 	[48]
a _m -ZIF-8 a _m -ZIF-69 a _m -ZIF-mnlm	Ball milling 25 Hz, 30 min	<ul style="list-style-type: none"> • Storage of I₂ 	[66]
a _m -UiO-66	Ball milling 20 Hz, 30 min	<ul style="list-style-type: none"> • Storage and controlled release of calcein 	[67]
a _m -UiO-66 a _m -UiO-66-Br a _m -UiO-66-NO ₂ a _m -UiO-66-NH ₂ a _m -MIL-140B a _m -MIL-140C a _m -[Zr ₆ O ₄ (OH) ₄ L ₇] _n a _m -[Zr ₆ O ₄ (OH) ₄ L ₈] _n	Ball milling 20 Hz, 30 min	<ul style="list-style-type: none"> • Storage and controlled release of calcein • Storage and controlled release of α-CHC 	[68]
a _m -HKUST-1 a _m -Al-ndc a _m -ZIF-8 a _m -Zn-MOF-74	Ball milling 400 rpm, 60 min	<ul style="list-style-type: none"> • Identical EXAFS profile in pristine and amorphous phase • Reversible amorphization 	[70]
a _m -UiO-66	Ball milling 400 rpm, 60 min	<ul style="list-style-type: none"> • Distinct EXAFS profile in pristine and amorphous phase • Irreversible amorphization 	[70]
a _m -Ni-MOF-74	Ball milling 30 Hz, 90 min	<ul style="list-style-type: none"> • Irreversible amorphization • Significant decrease of the bulk magnetization 	[71]
a _m -Zn-MOF-74	Ball milling 30 Hz, 90 min	<ul style="list-style-type: none"> • Reversible amorphization 	[71]
a _m (S-IL@ZIF-8)	Ball milling 650 rpm, 15-30 min	<ul style="list-style-type: none"> • Partial amorphization • Storage of IL 	[72]
a _T -ZIF-L	Scanning electron beam 1.1·10 ⁻² C cm ⁻² , 5 kV	<ul style="list-style-type: none"> • Stabilization in water 	[75]
a _T -ZIF-4	Synchrotron X-ray radiation 6 min	<ul style="list-style-type: none"> • 50% of amorphization 	[76]
a _T -ZIF-62	Synchrotron X-ray radiation 14 min	<ul style="list-style-type: none"> • 50% of amorphization 	[76]
a _T -ZIF-zni	Synchrotron X-ray radiation 81 min	<ul style="list-style-type: none"> • 50% of amorphization 	[76]
a _e -MOF-5	Electrical discharge	<ul style="list-style-type: none"> • --- 	[78]
a _{CT} -LnHPA-I Ln = La, Ce, Pr, Sm, Eu, Gd, Tb, Dy	Dehydration Vacuum, RT	<ul style="list-style-type: none"> • Precursor of a crystalline phase 	[79]

$a_{CT}-Cu^{1.8}(ttc)_{0.6}(ttcH_3)_{0.4}$	Dechlorination $NH_3(aq)$ Vacuum, 130 °C	<ul style="list-style-type: none"> • Semiconductive aMOF • No accessible pore volume 	[80]
$a_{CT}-Co^{II}-MOF$	Desolvation Vacuum, 320 °C	<ul style="list-style-type: none"> • Improved magnetic properties 	[81]
$a_{CT}-PCMOF-17$	$\geq 60\%$ RH	<ul style="list-style-type: none"> • --- 	[83]
$a_{CT}-MOF-5$	$\geq 40\%$ RH, RT	<ul style="list-style-type: none"> • --- 	[84]
$a_{CT}-Ni-MOF-74$	$\geq 40\%$ RH, RT	<ul style="list-style-type: none"> • --- 	[84]
$a_{CT}-MIL-88B$	Ligand competition <i>BDC-NH₂</i> & MeIm ligands	<ul style="list-style-type: none"> • BET surface area: 60.2 m²g⁻¹ 	[85]
$a_{CT}-UiO-66$	Loading $\geq 6, 10$ and 14 wt%	<ul style="list-style-type: none"> • BET surface area: ~11 m²/g • Catalytic activity for toluene oxidation 	[86]
$a-Fe^{(III)}_3O(C_6H_3(COO)_3)_2NO_3$	Direct synthesis Sol-gel approach	<ul style="list-style-type: none"> • BET surface area: ~1618 m²/g 	[88]
$a-Nd-TP$	Direct synthesis ALD/MLD	<ul style="list-style-type: none"> • Precursor of a crystalline phase • Photoluminescence switching upon the amorphous-to-crystalline transition 	[90]
$a-ZIF-8$	Direct synthesis Supersonic cold spraying	<ul style="list-style-type: none"> • --- 	[92]
$a-Zn(ICA)-2$	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 251 m²g⁻¹ 	[93]
$a-UiO-66$	Direct synthesis	<ul style="list-style-type: none"> • Supercapacitor electrode material 	[94]
$a-UiO-66-SO_3H$	Direct synthesis	<ul style="list-style-type: none"> • Precursor of a crystalline phase • BET surface area: 11 m²g⁻¹ 	[95]
$a-Fe_1Ni_2(BDC-NH_2)$	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 142.2 m²g⁻¹ • Catalytic activity for OER 	[96]
$a-FMM-120$	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 42.6 m²g⁻¹ • Adsorbent for arsenic-contaminated water remediation 	[97]
CPPs-1	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 584.7 m²g⁻¹ • Adsorbent for dyes-contaminated water remediation 	[98]
CPPs-3	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 522.4 m²g⁻¹ • Adsorbent for dyes-contaminated water remediation 	[98]
CPPs-5	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 587.5 m²g⁻¹ • Adsorbent for dyes-contaminated water remediation 	[98]
$a-Ni-pPD$	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 35 m²g⁻¹ 	[99]
$a-Zr-MOF(23)$	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 410 m²g⁻¹ 	[100]
$a-Zr-MOF(24)$	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 955 m²g⁻¹ 	[100]
$a-Zr-MOF(25)$	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 535 m²g⁻¹ 	[100]
$a-Zr-MOF(26)$	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 546 m²g⁻¹ 	[100]
$a-[Ru(C_5H_5)\{B(CN)_4\}]_n$	Direct synthesis UV irradiation	<ul style="list-style-type: none"> • BET surface area: 80-202 m²g⁻¹ 	[101]
$a-Ni^{2+}-1,3,5$ -tribenzyl-1,3,5-triazine-2,4,6-(1H,3H,5H)-trione	Direct synthesis Microwave-assisted hydrothermal technique	<ul style="list-style-type: none"> • BET surface area: 3181 m²g⁻¹ • CO₂ adsorbent 	[102]
$a-Co(MeIm)_2$	Direct synthesis	<ul style="list-style-type: none"> • Precursor of a catalytic material 	[104]
$a-Fe-ZIF-67$	Direct synthesis	<ul style="list-style-type: none"> • Precursor of a catalytic material 	[105]
$a-ZIF$	Direct synthesis	<ul style="list-style-type: none"> • Enzyme encapsulation 	[106]
NEU-2	Direct synthesis Liquid-liquid interface strategy	<ul style="list-style-type: none"> • BET surface area: 220.9 m²g⁻¹ • CO₂ adsorbent • Adsorptive separations 	[110]
NEU-3	Direct synthesis Liquid-liquid interface strategy	<ul style="list-style-type: none"> • BET surface area: 94.2 m²g⁻¹ • Adsorptive separations 	[111]
NEU-4	Direct synthesis Liquid-liquid interface strategy	<ul style="list-style-type: none"> • BET surface area: 293.8 m²g⁻¹ • Adsorptive separations 	[111]
NEU-5	Direct synthesis Liquid-liquid interface strategy & metalloligand approach	<ul style="list-style-type: none"> • BET surface area: 10.2 m²g⁻¹ • Precursor of a catalytic material 	[112]
NEU-6	Direct synthesis Liquid-liquid interface strategy & metalloligand approach	<ul style="list-style-type: none"> • BET surface area: 18.2 m²g⁻¹ • Precursor of a catalytic material 	[112]

NEU-7	Direct synthesis Liquid-liquid interface strategy & metalloligand approach	<ul style="list-style-type: none"> • BET surface area: 174.2 m²g⁻¹ • Precursor of a catalytic material 	[112]
NEU-8	Direct synthesis Liquid-liquid interface strategy & metalloligand approach	<ul style="list-style-type: none"> • BET surface area: 244.2 m²g⁻¹ • Precursor of a catalytic material 	[112]
a-CoL NL	Direct synthesis	<ul style="list-style-type: none"> • BET surface area: 105.7 m²g⁻¹ 	[115]
a-Co-MOF-74-NP	Direct synthesis	<ul style="list-style-type: none"> • Precursor of a catalytic material 	[117]
a-Ni-MOF-74.	Direct synthesis	<ul style="list-style-type: none"> • Nanoparticles encapsulation • Precursor of a catalytic material 	[118]

Notation: amorphization by pressure (a_p); amorphization by heating (a_T); mechanical milling amorphization (a_m); amorphization by irradiation (a_i); amorphization by electrical discharge (a_e); amorphization by chemical treatment (a_{CT}); amorphous (a); 1,4-benzenedicarboxylate (BDC); 1,3,5-benzenetricarboxylate (BTC); ferrocene (Fc); 1,4-diazabicyclo[2.2.2]octane (DABCO); dimethylformamide (DMF); imidazole (Im); 2,4,6-tris(4-pyridyl)-1,3,5-triazine (TPT); 2,3-pyrazinedithiolate (pdt); hollow sphere (HS); 4-methyl-5-nitroimidazolone (mnlm); 4,4'-(1,2-diazenediyl)bis-benzoic acid (L7); 4,4'-(1,2-ethenediyl)bis-benzoic acid (L8); 2,6-naphthalenedicarboxylic acid (ndc); ionic liquid (IL); 2-hydroxyphosphonoacetate (HPA); trithiocyanurate (ttc); trithiocyanuric acid (ttcH₃); imidazole-2-carboxyaldehyde (ICA); p-phenylenediamine (pPD); 2-methylimidazolone (Melm).

Table S2. Reported MOF liquids and MOF glasses.

MOF	Formula	T _m (°C)	MOF glasses	T _g (°C)	Highlights & Considerations	Ref.
ZIF-4	Zn(Im) ₂	583	a _g -ZIF-4	292-316	<ul style="list-style-type: none"> Glass formation: melt-quenching ZIF-4 recrystallizes to ZIF-zni at 500 °C before melting ZIF-4 liquid is a type 1 porous liquid 	[5,120,124]
ZIF-GIS	Zn(Im) ₂	590	a _g -ZIF-GIS	292	<ul style="list-style-type: none"> Glass formation: melt-quenching The addition of successively larger organic ligands to the MOF structures increases T_g and reduces T_m 	[120,151]
TIF-4	Zn(Im) _{1.5} (mPhIm) _{0.5}	467	a _g -TIF-4	318		
ZIF-62	Zn(Im) _{1.75} (PhIm) _{0.25}	437	a _g -ZIF-62	333	<ul style="list-style-type: none"> a_g-ZIF-62 exhibits permanent accessible porosity 	
ZIF-UC-1a	Zn(Im) _{1.74} (PhIm) _{0.17} (mPhIm) _{0.09}	433	a _g -ZIF-UC-1a	315	<ul style="list-style-type: none"> Glass formation: melt-quenching The inclusion of small amounts of PhIm decreases T_m in non-stoichiometric variants of ZIF-62 Linear increment of T_g as the amount of mPhIm increased in the frameworks (a_g-ZIF-UC-1b-e) a_g-ZIF-UC-1a and a_g-ZIF-UC-1e exhibit permanent accessible porosity 	[144]
ZIF-UC-1b	Zn(Im) _{1.66} (PhIm) _{0.22} (mPhIm) _{0.12}	418	a _g -ZIF-UC-1b	305		
ZIF-UC-1c	Zn(Im) _{1.68} (PhIm) _{0.17} (mPhIm) _{0.15}	425	a _g -ZIF-UC-1c	311		
ZIF-UC-1d	Zn(Im) _{1.59} (PhIm) _{0.27} (mPhIm) _{0.14}	420	a _g -ZIF-UC-1d	310		
ZIF-UC-1e	Zn(Im) _{1.61} (PhIm) _{0.18} (mPhIm) _{0.21}	430	a _g -ZIF-UC-1e	316		
TIF-4	Zn(Im) _{1.8} (mPhIm) _{0.2}	440	a _g -TIF-4	350	<ul style="list-style-type: none"> Glass formation: melt-quenching The lower T_m of ZIF-UC-5 compared to TIF-4 was attributed to the electron withdrawing chlorine ligand which weakened the Zn-N coordination bond 	[147]
ZIF-UC-5	Zn(Im) _{1.8} (cPhIm) _{0.2}	428	a _g -ZIF-UC-5	336		
ZIF-UC-2	Zn(Im) _{1.87} (6-Cl-5-FPhIm) _{0.13}	406	a _g -ZIF-UC-2	250	<ul style="list-style-type: none"> Glass formation: melt-quenching The electron-withdrawing groups on the PhIm ligand yield a reduction in the T_m 	[148]
ZIF-UC-3	Zn(Im) _{1.81} (5-Cl-2-mPhIm) _{0.19}	390	a _g -ZIF-UC-3	336		
ZIF-UC-4	Zn(Im) _{1.66} (FPhIm) _{0.34}	421	a _g -ZIF-UC-4	290		
ZIF-UC-5	Zn(Im) _{1.72} (cPhIm) _{0.28}	432	a _g -ZIF-UC-5	320		
ZIF-76-mPhIm	Zn(Im)(mPhIm)	No melt	---	---		
ZIF-76	Zn(Im)(cPhIm)	No melt	---	---		[147]
ZIF-76-mPhIm	Zn(Im) _{1.33} (mPhIm) _{0.67}	471	a _g -ZIF-76-mPhIm	317	<ul style="list-style-type: none"> Glass formation: melt-quenching The melting was assisted by impurities acting as a flux a_g-ZIF-76 and a_g-ZIF-76-mPhIm exhibit permanent accessible porosity 	[149]
ZIF-76	Zn(Im) _{1.62} (cPhIm) _{0.38}	451	a _g -ZIF-76	310		
ZIF-62(Co)	Co(Im) _{1.70} (PhIm) _{0.30}	432	a _g -ZIF-62(Co)	290	<ul style="list-style-type: none"> Glass formation: melt-quenching a_g-ZIF-62(Co) exhibits permanent accessible porosity 	[174]

Zn(Im) _{1.98} (PhIm) _{0.02}	Zn(Im) _{1.98} (PhIm) _{0.02}	No melt	---	---	
Zn(Im) _{1.97} (PhIm) _{0.03}	Zn(Im) _{1.97} (PhIm) _{0.03}	370	a _g -[Zn(Im) _{1.97} (PhIm) _{0.03}]	292	
Zn(Im) _{1.95} (PhIm) _{0.05}	Zn(Im) _{1.95} (PhIm) _{0.05}	372	a _g -[Zn(Im) _{1.95} (PhIm) _{0.05}]	298	<ul style="list-style-type: none"> Glass formation: melt-quenching The limit of melting and formation of stable liquids that did not recrystallize to ZIF-zn(M) lied around x = 0.05 a_g-ZIF-62 exhibits permanent accessible porosity The specific surface areas for a_g-ZIF-62 materials (determined by NLDFT fitting of the CO₂ adsorption isotherms) range between 235 and 269 m² g⁻¹
Zn(Im) _{1.94} (PhIm) _{0.06}	Zn(Im) _{1.94} (PhIm) _{0.06}	380	a _g -[Zn(Im) _{1.94} (PhIm) _{0.06}]	298	
Zn(Im) _{1.88} (PhIm) _{0.12}	Zn(Im) _{1.88} (PhIm) _{0.12}	395	a _g -[Zn(Im) _{1.88} (PhIm) _{0.12}]	303	
Zn(Im) _{1.83} (PhIm) _{0.17}	Zn(Im) _{1.83} (PhIm) _{0.17}	409	a _g -[Zn(Im) _{1.83} (PhIm) _{0.17}]	303	
Zn(Im) _{1.76} (PhIm) _{0.24}	Zn(Im) _{1.76} (PhIm) _{0.24}	423	a _g -[Zn(Im) _{1.76} (PhIm) _{0.24}]	305	
Zn(Im) _{1.65} (PhIm) _{0.35}	Zn(Im) _{1.65} (PhIm) _{0.35}	441	a _g -[Zn(Im) _{1.65} (PhIm) _{0.35}]	320	
Co(Im) _{1.90} (PhIm) _{0.10}	Co(Im) _{1.90} (PhIm) _{0.10}	386	a _g -[Co(Im) _{1.90} (PhIm) _{0.10}]	260	
Co(Im) _{1.83} (PhIm) _{0.17}	Co(Im) _{1.83} (PhIm) _{0.17}	410	a _g -[Co(Im) _{1.83} (PhIm) _{0.17}]	278	
Co(Im) _{1.70} (PhIm) _{0.30}	Co(Im) _{1.70} (PhIm) _{0.30}	432	a _g -[Co(Im) _{1.70} (PhIm) _{0.30}]	290	
Zn(Im) _{1.7} (PhIm) _{0.3}	Zn(Im) _{1.7} (PhIm) _{0.3}	447	a _g -[Zn(Im) _{1.7} (PhIm) _{0.3}]	328	
Zn _{0.9} Co _{0.1} (Im) _{1.7} (PhIm) _{0.3}	Zn _{0.9} Co _{0.1} (Im) _{1.7} (PhIm) _{0.3}	446	a _g -[Zn _{0.9} Co _{0.1} (Im) _{1.7} (PhIm) _{0.3}]	331	<ul style="list-style-type: none"> Glass formation: melt-quenching The luminescence enhances by increasing the concentration level of Co ions
Zn _{0.5} Co _{0.5} (Im) _{1.7} (PhIm) _{0.3}	Zn _{0.5} Co _{0.5} (Im) _{1.7} (PhIm) _{0.3}	452	a _g -[Zn _{0.5} Co _{0.5} (Im) _{1.7} (PhIm) _{0.3}]	334	
Zn(Im) _{1.95} (cPhIm) _{0.05}	Zn(Im) _{1.95} (cPhIm) _{0.05}	366	a _g -[Zn(Im) _{1.95} (cPhIm) _{0.05}]	296	<ul style="list-style-type: none"> Glass formation: melt-quenching
Zn(Im) _{1.95} (PhIm) _{0.025} (cPhIm) _{0.025}	Zn(Im) _{1.95} (PhIm) _{0.025} (cPhIm) _{0.025}	356	a _g -[Zn(Im) _{1.95} (PhIm) _{0.025} (cPhIm) _{0.025}]	296	
Zn _{0.8} Co _{0.2} (Im) _{1.95} (PhIm) _{0.025} (cPhIm) _{0.025}	Zn _{0.8} Co _{0.2} (Im) _{1.95} (PhIm) _{0.025} (cPhIm) _{0.025}	310	a _g -[Zn _{0.8} Co _{0.2} (Im) _{1.95} (PhIm) _{0.025} (cPhIm) _{0.025}]	288	
[[Co(μ-NCS) ₂ (C ₅ H ₅ N ₃ O) ₂] ₂ ·C ₅ H ₅ N ₃ O] _n	---	197	a _g -[[Co(μ-NCS) ₂ (C ₅ H ₅ N ₃ O) ₂] ₂ ·C ₅ H ₅ N ₃ O] _n	78	<ul style="list-style-type: none"> Glass formation: melt-quenching Glass formation: melt-quenching [Cd(μ-NCS)₂(C₅H₅N₃O)₂]_n recrystallizes at 144 °C upon cooling from 220 °C
[Cd(μ-NCS) ₂ (C ₅ H ₅ N ₃ O) ₂] _n	---	205	a _g -[Cd(μ-NCS) ₂ (C ₅ H ₅ N ₃ O) ₂] _n	---	
Cu ₂ (SCN) ₃ (C ₂ bpy)	---	187	a _g -[Cu ₂ (SCN) ₃ (C ₂ bpy)]	68	<ul style="list-style-type: none"> Glass formation: melt-quenching T_m is influenced by the substituent group on pyridinium ligands
Cu ₂ (SCN) ₃ (C ₄ bpy)	---	138	a _g -[Cu ₂ (SCN) ₃ (C ₄ bpy)]	59	
Cu ₈ (SCN) ₁₂ (Phbpy) ₄	---	217	a _g -[Cu ₈ (SCN) ₁₂ (Phbpy) ₄]	71	
Cu(SCN) ₂ (3-Pybpy)	---	203	a _g -[Cu(SCN) ₂ (3-Pybpy)]	72	
[[Zn ₂ (HPO ₄) ₂ (H ₂ PO ₄)](cPhImH ⁺) ₂ ·(H ₂ PO ₄) ⁻ ·(MeOH)] _n	---	148	a _g -[[Zn ₂ (HPO ₄) ₂ (H ₂ PO ₄)](cPhImH ⁺) ₂ ·(H ₂ PO ₄) ⁻ ·(MeOH)] _n	72	<ul style="list-style-type: none"> Glass formation: melt-quenching
[Cd(sdc)(L4)]·DMF	---	---	---	---	
[Cd(DMF)(sdc)(L6)]·DMF	---	---	---	---	<ul style="list-style-type: none"> Glass formation: ultrafast melting <i>via</i> femtosecond IR laser pulses
[Zn(sdc)(L6)]·DMF	---	---	---	---	

[145]

[178]

[152]

[176]

[177]

[153]

CdTz-40	---	---	a_g -CdTz-40	79	<ul style="list-style-type: none"> Glass formation: Solvent-free mechanical milling under an Ar atmosphere Milling times are 40, 240, and 500 minutes, yielding a_g-CdTz-40, a_g-CdTz-240, and a_g-CdTz-500, respectively 	[189]
CdTz-240	---	---	a_g -CdTz-240	85		
CdTz-500	---	---	a_g -CdTz-500	90		
CrTz	---	---	a_g -CrTz	62	<ul style="list-style-type: none"> Glass formation: Solvent-free mechanical milling under an Ar atmosphere 	[190]
MnTz	---	---	a_g -MnTz	81		
Ti-BPA	---	---	a_g -Ti-BPA	---	<ul style="list-style-type: none"> Glass formation: assembling metal nodes and organic struts and subsequent evaporation of a plasticizer-modulator solvent. ^(#)Surface area of a_g-Ti-BPA 330 m² g⁻¹ ^(#)Surface area of a_g-Ti-BPP 267 m² g⁻¹ 	[191]
Ti-BPP	---	---	a_g -Ti-BPP	---		
[Ag(<i>p</i> L2)(CF ₃ SO ₃)]·2C ₆ H ₆	---	271	a_g -[Ag(<i>p</i> L2)(CF ₃ SO ₃)]·2C ₆ H ₆	161	<ul style="list-style-type: none"> Glass formation: melt-quenching Glass formation: mechanical hand-grinding in Ar atmosphere a_g-[Ag(<i>p</i>L2)(CF₃SO₃)]·2C₆H₆ exhibits permanent accessible porosity 	[193]
[Ag(<i>p</i> L2)(CF ₃ SO ₃)]·2C ₆ H ₆	---	---	a_g -[[Ag(<i>p</i> L2)(CF ₃ SO ₃)]·2C ₆ H ₆]	107		
[Ag(<i>m</i> L1)(CF ₃ SO ₃)]·2C ₆ H ₆	---	169	a_g -[[Ag(<i>m</i> L1)(CF ₃ SO ₃)]·2C ₆ H ₆]	68		
[Ag(<i>m</i> L1)(CF ₃ SO ₃)]·2C ₆ H ₆	---	---	a_g -[[Ag(<i>m</i> L1)(CF ₃ SO ₃)]·2C ₆ H ₆]	68	<ul style="list-style-type: none"> Glass formation: melt-quenching Glass formation: mechanical hand-grinding in Ar atmosphere a_g-[[Ag(<i>m</i>L1)(CF₃SO₃)]·2C₆H₆] exhibits permanent accessible porosity 	[194]
(ZIF-4-Zn) _{0.5} (ZIF-62) _{0.5}	---	590	(ZIF-4-Zn) _{0.5} (ZIF-62) _{0.5} blend	306		
(ZIF-4-Co) _{0.5} (ZIF-62) _{0.5}	---	425	(ZIF-4-Co) _{0.5} (ZIF-62) _{0.5} blend	300	<ul style="list-style-type: none"> Glass formation: melt-quenching 	[6]
(ZIF-8)(ZIF-62)(20/80)	---	500	a_g -[(ZIF-8) _{0.2} (ZIF-62) _{0.8}]	334	<ul style="list-style-type: none"> Flux melted MOF glass Surface area of a_g-[(ZIF-8)_{0.2}(ZIF-62)_{0.8}]^(*)202 m² g⁻¹ Surface area of a_g-[(ZIF-67)_{0.2}(ZIF-62)_{0.8}]^(*)194 m² g⁻¹ 	[7]
(ZIF-67)(ZIF-62)(20/80)	---	497	a_g -[(ZIF-67) _{0.2} (ZIF-62) _{0.8}]	---		
(MIL-53) _{0.25} (ZIF-62) _{0.75}	---	---	(MIL-53) _{0.25} (a_g -ZIF-62) _{0.75}	310	<ul style="list-style-type: none"> Flux melted MOF glass 	[8]
(UiO-66) _{0.25} (ZIF-62) _{0.75}	---	---	(UiO-66) _{0.25} (a_g -ZIF-62) _{0.75}	---	<ul style="list-style-type: none"> Flux melted MOF glass 	

Notation: imidazole (Im); glass phase (a_g); 5-methylbenzimidazole (mPhIm); benzimidazole (PhIm); *chlorobenzimidazole* (cPhIm); 4,4'-stilbenedicarbonylate (sdc); dimethylformamide (DMF); 1,4-di(1H-imidazol-1-yl)butane (L4); 1,6-di(1H-imidazol-1-yl)hexane (L6); 6-chloro-5-fluorobenzimidazole (6-Cl-5-FPhIm); 5-chloro-2-methylbenzimidazole (5-Cl-2-mPhIm); 5-fluorobenzimidazole (FPhIm); 1-ethyl-[4,4'-bipyridin]-1-ium thiocyanate (C₂bpySCN); 1-butyl-[4,4'-bipyridin]-1-ium thiocyanate (C₄bpySCN); 1-phenyl-[4,4'-bipyridin]-1-ium thiocyanate (PhbpySCN); 3,1':4',4''-terpyridin]-1'-ium thiocyanate (3-PybySCN); 1,2,4-triazole (Tr); 2,2-Bis(4-hydroxyphenyl)propane (BPA); 2,2-Bis(4-hydroxyphenyl)propane (BPP); 1,3,5-tris(4-ethynylbenzotrile)benzene (*p*L2); 1,3,5-tris(3-cyanophenylethynyl)benzene (*m*L1).

^(#) Analysis of the N₂ isotherm
^(*) Analysis of the CO₂ isotherm