Constraining the coordination geometries of lanthanide centers and magnetic building blocks in frameworks: a new strategy for molecular nanomagnets

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Table S1 The structures and abbreviations of the ligands in this review

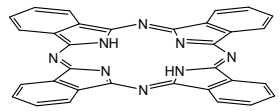
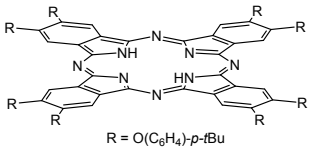
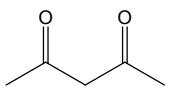
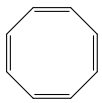
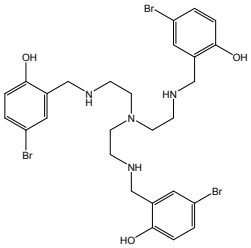
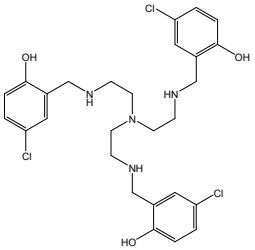
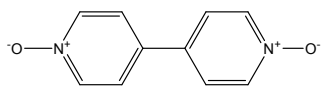
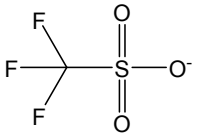
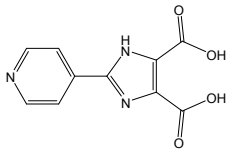
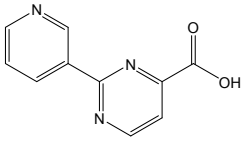
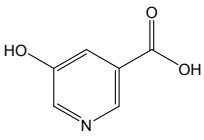
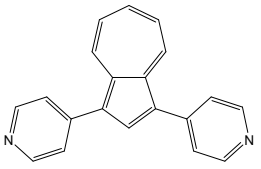
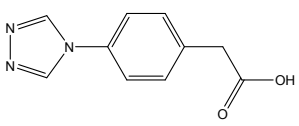
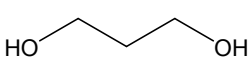
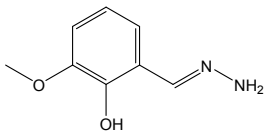
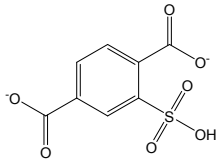
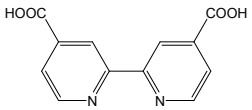
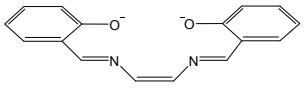
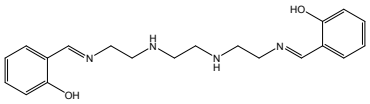
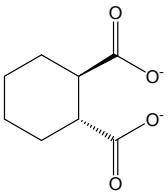
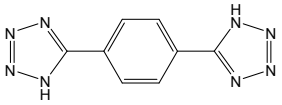
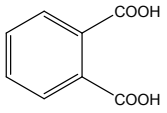
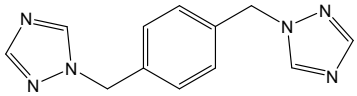
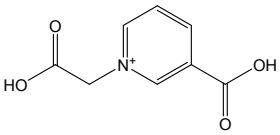
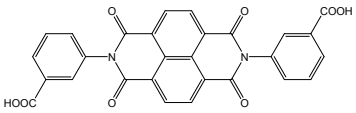
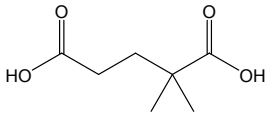
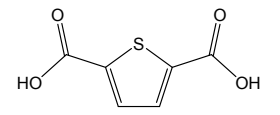
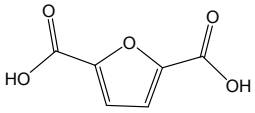
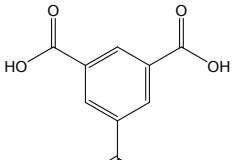
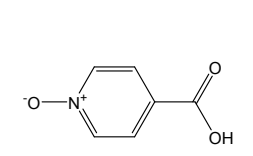
		
Pc	Pc’	acac
		
COT	L^1	L^2
		
bipyNO	TfO	H_3PIDC
		
H3-py-4-pmc	H_2NA	azbbpy
		
		H_2hmi

Table S1 The structures and abbreviations of the ligands in this review

HTPA	H_2pd	
		
stp	H_2BPDC	$salen^{2-}$
	trans-chdc  1,2-	
H_2sal_2-trien		H_2bdt
		
H_2pta	btx	H_2L
		
H_2L	H_2L	H_2TDA
		
H_2FDA	H_3BTC	HINO

Pc [21] = phthalocyanine

Pc' [6] = octa(tert-butylphenoxy)-phthalocyanine

acac [22] = acetylacetone

COT [2] = cyclooctatetraene

L^1 [24] = 2,2',2''-(((nitrilotris(ethane-2,1-diyl))tris(azanediyl))tris(methylene))tris(4-bromophenol)

L^2 [25] = 2,2',2''-(((nitrilotris(ethane-2,1-diyl))tris(azanediyl))tris(methylene))tris(4-chlorophenol)

bipyNO [29]= 4,4'-bypyridyl-N,N'-dioxide

TfO [29] = triflate

H_3PIDC [30] = 2-(Pyridin-4-yl)-1H-imidazole-4,5-dicarboxylic acid

H3-py-4-pmc [31] = 2-(3-pyridyl)pyrimidine-4-carboxylic acid

H_2NA [28] = 5-hydroxynicotinic acid

Azbbpy [32] = 1,3-bis(4-pyridyl)azulene

Table S1 The structures and abbreviations of the ligands in this review

HTPA [33] =4-(1,2,4-triazol-4-yl)-phenylacetic acid
H_2pd [34] = 1,3-propanediol
H_2hmi [35] =(2-hydroxy-3-methoxyphenyl)methylene(isonicotino)hydrazine
stp [36] = m_5-2-sulfoterephthalate
H_2BPDC [37] = 4,4'-dicarboxylate-2, 2'-dipyridine
salen^{2-} [40] = N,N'-ethylene-bis-(salicylideneiminate)
H_2sal$_2$-trien [40] = N,N'-disalicylidenetriethylenetetramine
trans-1,2-chdc [41] = trans-1,2-cyclohexanedicarboxylate
H_2bdt [42] = 5,5'-(1,4-phenylene)bis(1H-tetrazole)
H_2pta [43] = phthalic acid
btx [43] = 1,4-bis(1,2,4-triazole-1-ylmethyl)benzene
H_2L [44] = 1-carboxymethylpyridinium-3-carboxylate acid
H_2L[45]=3,3'-(1,3,6,8-tetraoxobenzol[lmn][3,8]-phenanthroline-2,7(1H,3H,6H,8H)diyl)-di-benzoic acid
H_2L [46] = 2,2-dimethylglutaric acid
H_2TDA [47] = thiophene-2,5-dicarboxylic acid
H_2FDA [49] = furan-2,5-dicarboxylic acid
H_3BTC [50] = 1,3,5-benzenetricarboxylate acid
HINO [48] = isonicotinic acid N-oxide