CheckMyMetal(CMM) report for PDB code: 7kol

PDB title: The crystal structure of papain-like protease of sars cov-2 in complex with plp_snyder496 inhibitor (2.6Å)

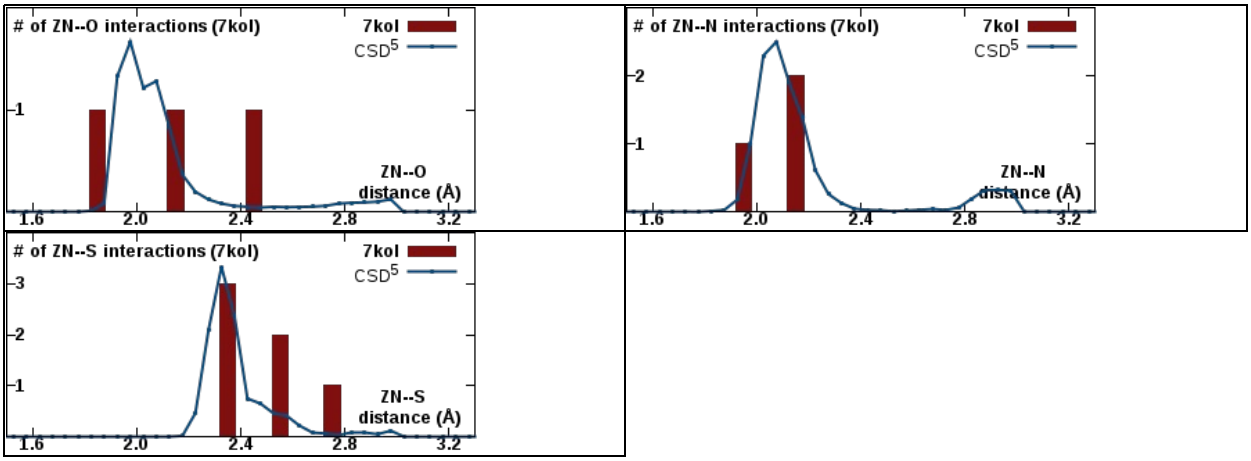
Warning: Coordinating ligands by symmetry operation are labeled with prefix 'sym-'

ID	Res.	Metal	*Occupancy*	*B factor (env.)*[1]	*Ligands*	*Valence*[2]	*nVECSUM*[3]	*Geometry*[1,4]	*gRMSD*(°)[1]	*Vacancy*[1]	Bidentate	Alt. metal
A:502	ZN	Zn	1	154.3 (150.2)	S_4	*1.2*	*0.22*	Tetrahedral	**23.7°**	0	0	Cu
A:503	ZN	Zn	1	105 (96.8)	$O_1N_1S_1$	*1.2*	**0.3**	Tetrahedral	*14.2°*	*25%*	0	
A:504	ZN	Zn	1	90.6 (94.3)	*O_1N_1*	1.6	*0.19*	Tetrahedral	3.4°	**50%**	0	Cu, Co
A:505	ZN	Zn	1	91.6 (104.3)	$O_1N_1S_1$	*1.4*	**0.35**	Tetrahedral	11°	*25%*	0	Cu, Co
A:506	CL	Cl	1	93.5 (104.6)		N/A	N/A	**Free**	N/A	N/A	N/A	
A:507	CL	Cl	1	100.5 (93.4)		N/A	N/A	**Free**	N/A	N/A	N/A	
A:508	CL	Cl	1	**118.8 (0)**		N/A	N/A	**Free**	N/A	N/A	N/A	
A:509	CL	Cl	1	**107.9 (0)**		N/A	N/A	**Free**	N/A	N/A	N/A	

Legend: Not applicable | **Outlier** | *Borderline* | Acceptable

Column	Description
Occupancy	Occupancy of ion under consideration
B factor (env.)[1]	Metal ion B factor, with valence-weighted environmental average B factor in parenthesis
Ligands	Elemental composition of the coordination sphere
Valence[2]	Summation of bond valence values for an ion binding site. *Valence* accounts for metal-ligand distances
nVECSUM[3]	Summation of ligand vectors, weighted by bond valence values and normalized by overall valence. Increase when the coordination sphere is not symmetrical due to incompleteness.
Geometry[1,4]	Arrangement of ligands around the ion, as defined by the *NEIGHBORHOOD* algorithm
gRMSD(°)[1]	R.M.S. Deviation of observed geometry angles (L-M-L angles) compared to ideal geometry, in degrees
Vacancy[1]	Percentage of unoccupied sites in the coordination sphere for the given geometry
Bidentate	Number of residues that form a bidentate interaction instead of being considered as multiple ligands
Alt. metal	A list of alternative metal(s) is proposed in descending order of confidency, assuming metal environment is accurately determined. This feature is still experimental. It requires user discrimination and cannot be blindly accepted

Metal-ligand distance distributions for pdb7kol.ent in comparison with CSD



(1) Zheng H, Chordia MD, Cooper DR, Chruszcz M, Müller P, Sheldrick GM, Minor W (2014) *Nature Protocols, 9(1),* 156-70.
(2) Brown ID (2009) *Chem. Rev., 109,* 6858-6919.
(3) Müller P, Köpke S, Sheldrick GM (2003) *Acta Crystallogr. D Biol. Crystallogr., 59,* 32-37.
(4) Kuppuraj G, Dudev M, Lim C (2009) *J. Phys. Chem. B, 113,* 2952-2960.
(5) *CSD: Cambridge Structural Database*
Maintained by: Heping Zheng <dust@iwonka.med.virginia.edu>

Citing CheckMyMetal (CMM):

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Validation of metal-binding sites in macromolecular structures with the CheckMyMetal web server. Zheng,H., Chordia,M.D., Cooper,D.R., Chruszcz,M., Müller,P., Sheldrick,G.M., Minor,W. (2014) *Nature Protocols, 9(1),* 156-70.