



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 10:51 AM BST

PDB ID : 4DCX
Title : X-ray structure of NikA in complex with Fe(1R,2R)-N,N'-Bis(2-pyridylmethyl)-N,N'-dicarboxymethyl-1,2-cyclohexanediamine
Authors : Cherrier, M.V.; Girgenti, E.; Amara, P.; Iannello, M.; Marchi-Delapierre, C.; Fontecilla-Camps, J.C.; Menage, S.; Cavazza, C.
Deposited on : 2012-01-18
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

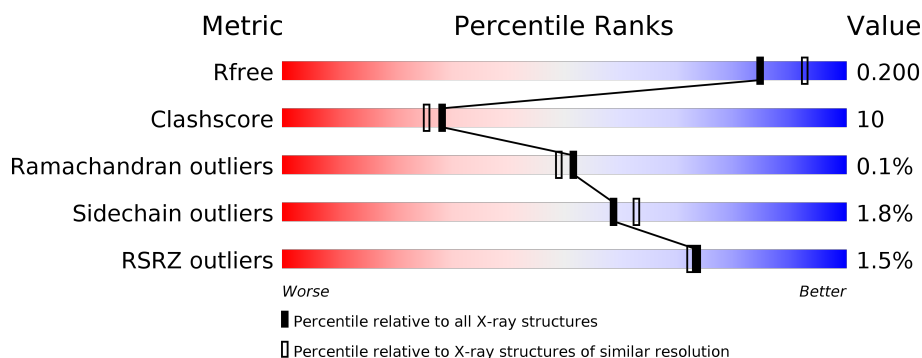
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	502	
1	B	502	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ACT	A	615	-	-	X	-
3	ACT	A	616	-	-	X	X
3	ACT	A	617	-	-	X	-
3	ACT	B	606	-	-	X	-
4	GOL	A	626	-	-	-	X
4	GOL	A	627	-	-	X	-
4	GOL	A	628	-	-	-	X
4	GOL	B	612	-	-	X	-
4	GOL	B	618	-	-	X	X

2 Entry composition [i](#)

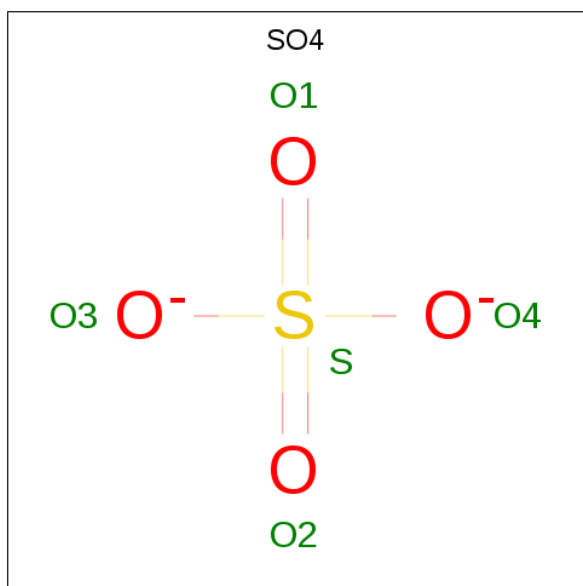
There are 6 unique types of molecules in this entry. The entry contains 9194 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Nickel-binding periplasmic protein.

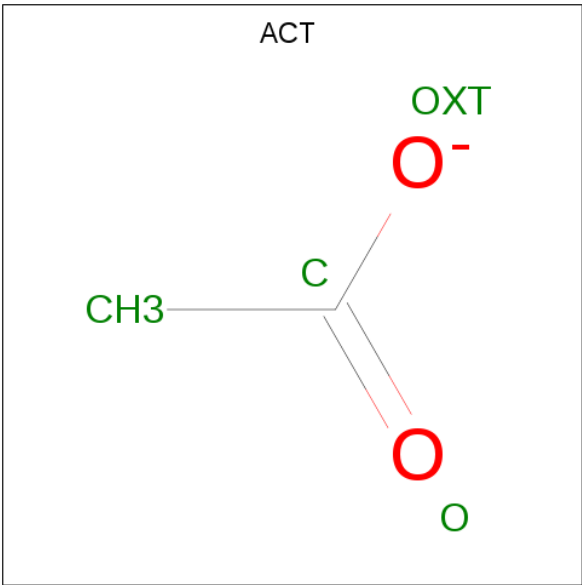
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	497	Total	C	N	O	S	9	20	0
			4050	2606	676	755	13			
1	B	497	Total	C	N	O	S	11	9	0
			3992	2562	671	748	11			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



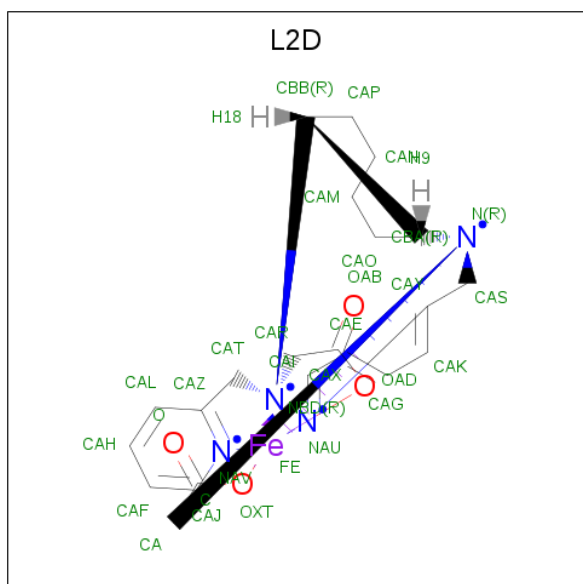
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is {2,2'-[(1R,2R)-cyclohexane-1,2-diylbis{[(pyridin-2-yl-kappaN)methyl]imino-kappaN}]]diacetato-kappaO(2-)}iron (three-letter code: L2D) (formula: C₂₂H₂₆FeN₄O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	Fe	N	O	0	0
			31	22	1	4	4		
5	B	1	Total	C	Fe	N	O	0	0
			31	22	1	4	4		

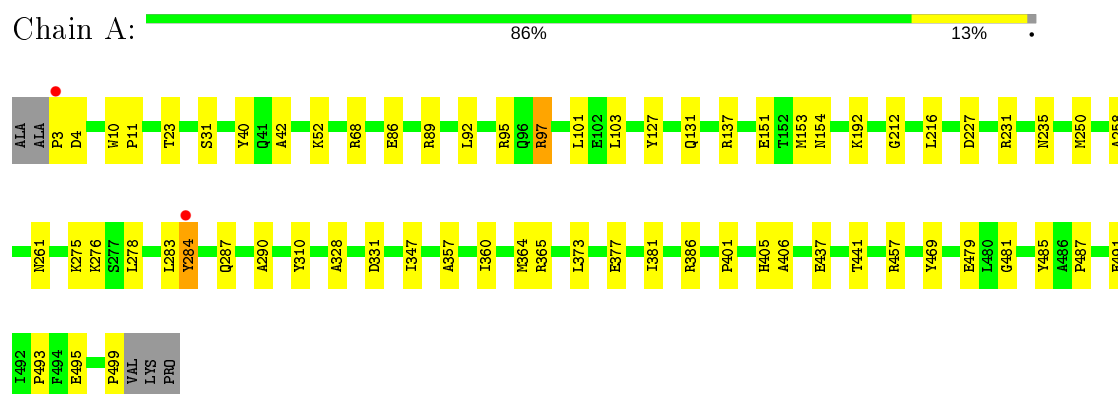
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	511	Total	O	2	0
			511	511		
6	B	347	Total	O	0	1
			348	348		

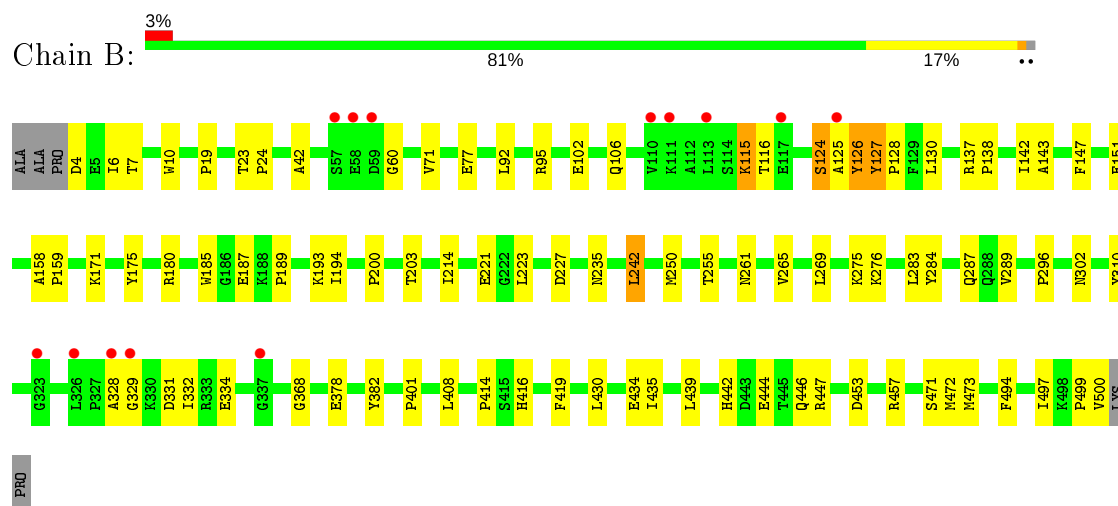
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Nickel-binding periplasmic protein



- Molecule 1: Nickel-binding periplasmic protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.35Å 95.64Å 125.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.82 – 2.00 47.82 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.82-2.00) 99.6 (47.82-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.64 (at 2.00Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.155 , 0.203 0.153 , 0.200	Depositor DCC
R_{free} test set	3519 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	27.5	Xtriage
Anisotropy	0.346	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 65.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9194	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4, L2D, ACT

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.37	0/4219	0.52	0/5743
1	B	0.34	0/4122	0.52	0/5616
All	All	0.36	0/8341	0.52	0/11359

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4050	0	4059	64	0
1	B	3992	0	3967	84	0
2	A	10	0	0	0	0
2	B	5	0	0	0	0
3	A	68	0	51	13	0
3	B	40	0	30	13	0
4	A	60	0	80	16	0
4	B	48	0	64	21	0
5	A	31	0	26	0	0
5	B	31	0	26	2	0
6	A	511	0	0	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	348	0	0	15	0
All	All	9194	0	8303	162	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (162) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:331:ASP:HB2	3:B:606:ACT:H3	1.32	1.06
1:B:235:ASN:HD21	4:B:616:GOL:H2	1.19	1.02
3:B:606:ACT:C	4:B:618:GOL:H12	1.94	0.97
1:A:52:LYS:HE3	1:A:68:ARG:HA	1.47	0.93
3:B:606:ACT:H2	4:B:618:GOL:H31	1.49	0.93
1:B:310:TYR:HB2	4:B:612:GOL:H12	1.51	0.90
1:B:6:ILE:HD11	1:B:194:ILE:HG12	1.50	0.90
1:B:331:ASP:HB2	3:B:606:ACT:CH3	2.07	0.84
1:B:331:ASP:CB	3:B:606:ACT:H3	2.07	0.84
1:B:10:TRP:HZ2	4:B:615:GOL:H11	1.40	0.83
1:A:499:PRO:HB2	6:A:1101:HOH:O	1.80	0.82
1:A:227:ASP:HB3	1:A:284[B]:TYR:CZ	2.15	0.81
3:B:606:ACT:CH3	4:B:618:GOL:H12	2.09	0.81
1:A:437:GLU:O	1:A:441[B]:THR:HG23	1.81	0.80
1:B:10:TRP:CZ2	4:B:615:GOL:H11	2.19	0.77
1:B:235:ASN:HD21	4:B:616:GOL:C2	1.96	0.77
1:A:365:ARG:HH22	4:A:622:GOL:H32	1.53	0.74
1:A:52:LYS:CE	1:A:68:ARG:HA	2.18	0.73
1:A:151[B]:GLU:HG3	6:A:953:HOH:O	1.88	0.72
1:B:414:PRO:HA	1:B:419:PHE:CD1	2.25	0.72
3:B:605:ACT:H1	6:B:1008:HOH:O	1.89	0.72
1:A:331:ASP:H	4:A:622:GOL:H11	1.55	0.72
1:B:332:ILE:HG12	3:B:606:ACT:H1	1.70	0.71
1:A:328:ALA:HB2	3:A:615:ACT:H2	1.71	0.71
1:A:441[B]:THR:HG22	6:A:1096:HOH:O	1.90	0.70
1:B:7:THR:HG22	1:B:214:ILE:HG22	1.73	0.69
1:B:499:PRO:O	1:B:500:VAL:HB	1.91	0.68
1:B:227:ASP:HB3	1:B:284:TYR:CZ	2.30	0.66
1:B:19:PRO:HG3	1:B:142:ILE:HB	1.78	0.66
3:B:606:ACT:H2	4:B:618:GOL:H12	1.77	0.65
1:A:92:LEU:HD12	4:A:621:GOL:H31	1.79	0.64
1:B:378:GLU:HG2	1:B:382:TYR:CE2	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:491:GLU:OE1	4:A:627:GOL:H32	1.99	0.62
1:B:6:ILE:CD1	1:B:194:ILE:HG12	2.25	0.62
1:A:328:ALA:HB2	3:A:615:ACT:CH3	2.29	0.61
1:B:310:TYR:CB	4:B:612:GOL:H12	2.26	0.61
1:B:444:GLU:HG3	6:B:928:HOH:O	1.99	0.61
1:A:89:ARG:HG3	4:A:626:GOL:H12	1.82	0.61
1:B:221:GLU:OE2	4:B:613:GOL:H31	2.01	0.61
1:A:275:LYS:HB2	4:A:625:GOL:H32	1.83	0.60
1:B:143:ALA:HB1	6:B:1020:HOH:O	2.00	0.60
1:B:416:HIS:HB2	6:B:964:HOH:O	2.01	0.60
1:B:442:HIS:CD2	4:B:617:GOL:H32	2.38	0.59
1:B:115:LYS:H	1:B:115:LYS:HD2	1.67	0.59
1:A:276:LYS:HD2	6:A:1167:HOH:O	2.01	0.59
1:A:231[A]:ARG:NH2	6:A:1162:HOH:O	2.37	0.58
1:A:360:ILE:O	1:A:364[B]:MET:HG2	2.03	0.58
1:B:331:ASP:H	3:B:606:ACT:C	2.16	0.58
4:A:623:GOL:H11	6:A:876:HOH:O	2.02	0.58
1:B:275:LYS:HB2	4:B:612:GOL:H2	1.86	0.58
1:A:347:ILE:HD13	1:A:381:ILE:HG13	1.86	0.58
1:A:86:GLU:HG2	6:A:1117:HOH:O	2.04	0.58
1:A:377:GLU:HG2	6:A:1168:HOH:O	2.03	0.57
3:B:609:ACT:H1	6:B:956:HOH:O	2.03	0.57
1:A:52:LYS:NZ	6:A:869:HOH:O	2.30	0.57
1:B:125:ALA:HB3	1:B:442:HIS:CE1	2.40	0.56
1:B:276:LYS:H	4:B:612:GOL:H31	1.70	0.56
1:B:151:GLU:HG3	6:B:735:HOH:O	2.05	0.56
1:A:495:GLU:CD	1:A:495:GLU:H	2.09	0.56
1:B:180:ARG:HG3	1:B:189:PRO:HG2	1.88	0.56
1:A:89:ARG:HD3	4:A:626:GOL:H31	1.87	0.55
1:B:193:LYS:HD2	6:B:1032:HOH:O	2.05	0.55
1:B:71:VAL:HG22	1:B:185:TRP:CG	2.42	0.55
3:A:616:ACT:H1	6:A:917:HOH:O	2.06	0.55
1:B:171:LYS:HD3	1:B:175:TYR:OH	2.07	0.54
1:B:414:PRO:HA	1:B:419:PHE:CG	2.41	0.54
1:A:153[B]:MET:HE2	6:A:1206:HOH:O	2.08	0.54
1:A:235:ASN:HD21	3:A:618:ACT:C	2.21	0.54
1:A:479:GLU:HG2	6:A:1092:HOH:O	2.08	0.53
1:B:106:GLN:NE2	1:B:124:SER:H	2.07	0.53
1:A:365:ARG:NH2	4:A:622:GOL:H32	2.23	0.53
1:B:242:LEU:HD22	1:B:471:SER:HB2	1.91	0.53
1:A:92:LEU:O	1:A:95[A]:ARG:HB2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:LEU:C	1:A:278:LEU:HD23	2.29	0.52
1:B:223:LEU:HA	4:B:619:GOL:H12	1.92	0.52
1:A:491:GLU:HB3	4:A:627:GOL:H11	1.91	0.52
1:B:296:PRO:HB3	1:B:302:ASN:HD22	1.74	0.52
1:A:258:ALA:HA	1:A:261:ASN:OD1	2.10	0.51
1:A:485:TYR:CD1	3:A:614:ACT:H2	2.45	0.51
1:B:171:LYS:HD3	1:B:175:TYR:CZ	2.45	0.51
1:A:52:LYS:HE3	1:A:68:ARG:CA	2.32	0.50
1:A:31:SER:HB3	4:A:627:GOL:H12	1.93	0.50
1:B:124:SER:HB3	6:B:985:HOH:O	2.11	0.49
1:B:200:PRO:HD2	6:B:969:HOH:O	2.12	0.49
1:B:276:LYS:HD3	1:B:289:VAL:HG21	1.93	0.49
1:A:101[B]:LEU:HD21	1:A:103:LEU:HB3	1.94	0.48
1:B:102:GLU:HB3	1:B:126:TYR:OH	2.13	0.48
1:A:52:LYS:HG3	1:A:68:ARG:HG2	1.95	0.48
1:A:42:ALA:HB3	1:B:42:ALA:HB3	1.96	0.48
1:A:10:TRP:CD1	1:A:11:PRO:HD2	2.49	0.48
1:B:275:LYS:HG3	4:B:612:GOL:H11	1.95	0.48
1:A:457:ARG:HH22	3:A:609:ACT:H3	1.78	0.48
1:A:310:TYR:H	4:A:625:GOL:H2	1.78	0.48
4:A:624:GOL:H32	6:A:1197:HOH:O	2.13	0.48
1:A:101[B]:LEU:CD2	1:A:103:LEU:HB3	2.43	0.47
1:A:52:LYS:HE2	3:A:617:ACT:CH3	2.44	0.47
4:A:623:GOL:H12	6:A:909:HOH:O	2.15	0.47
1:B:106:GLN:HE22	1:B:124:SER:H	1.60	0.47
1:A:137:ARG:HA	6:A:1153:HOH:O	2.14	0.47
1:B:331:ASP:O	1:B:368:GLY:HA2	2.14	0.47
1:A:386:ARG:HE	3:A:616:ACT:CH3	2.28	0.47
1:B:283:LEU:HD13	1:B:287[B]:GLN:OE1	2.15	0.47
5:B:620:L2D:H7	5:B:620:L2D:H11	1.69	0.46
1:A:3:PRO:HG2	1:A:4:ASP:H	1.80	0.46
1:B:127:TYR:HA	4:B:617:GOL:H31	1.97	0.46
1:B:457:ARG:HG3	6:B:972:HOH:O	2.15	0.46
1:A:491:GLU:CD	4:A:627:GOL:H32	2.36	0.46
1:B:127:TYR:CD2	1:B:447:ARG:NH1	2.84	0.45
1:B:446:GLN:HA	1:B:446:GLN:OE1	2.16	0.45
1:A:328:ALA:CB	3:A:615:ACT:H2	2.42	0.45
1:B:408:LEU:O	1:B:435:ILE:HD13	2.17	0.45
1:B:494:PHE:HA	1:B:497:ILE:HD13	1.98	0.45
1:A:386:ARG:HE	3:A:616:ACT:H2	1.82	0.45
1:A:52:LYS:HE2	3:A:617:ACT:H1	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:142:ILE:HG23	1:B:147:PHE:HE2	1.83	0.44
5:B:620:L2D:H16	5:B:620:L2D:H21	1.68	0.44
1:B:457:ARG:HD2	6:B:996:HOH:O	2.17	0.44
1:A:154:ASN:HA	6:A:1165:HOH:O	2.16	0.44
1:B:430:LEU:HD23	1:B:430:LEU:C	2.38	0.44
1:B:126:TYR:CD1	1:B:128:PRO:HD2	2.53	0.44
1:A:23[A]:THR:HG22	6:A:1080:HOH:O	2.18	0.43
1:A:405:HIS:CG	1:A:406:ALA:N	2.86	0.43
1:A:493:PRO:HA	1:A:495:GLU:OE2	2.18	0.43
4:B:614:GOL:H2	6:B:977:HOH:O	2.19	0.43
1:A:131:GLN:HG2	6:A:1079:HOH:O	2.19	0.43
3:A:615:ACT:H3	6:A:1012:HOH:O	2.18	0.43
1:B:265:VAL:O	1:B:269:LEU:HG	2.18	0.43
1:B:328:ALA:HA	1:B:329:GLY:HA2	1.71	0.43
1:B:401:PRO:HB2	6:B:806:HOH:O	2.18	0.43
1:B:92:LEU:O	1:B:95:ARG:HB2	2.18	0.43
1:B:125:ALA:O	1:B:126:TYR:C	2.57	0.43
1:B:223:LEU:HB3	1:B:473:MET:CE	2.49	0.43
3:A:613:ACT:H2	3:A:617:ACT:H3	2.00	0.43
1:B:203:THR:HG22	3:B:610:ACT:H2	2.01	0.43
1:B:124:SER:C	1:B:126:TYR:H	2.23	0.42
1:A:40:TYR:CE1	1:A:401:PRO:HB3	2.53	0.42
1:B:158:ALA:HA	1:B:159:PRO:HD3	1.84	0.42
1:B:60:GLY:O	1:B:125:ALA:HA	2.19	0.42
1:B:127:TYR:HA	4:B:617:GOL:H11	2.00	0.42
1:B:23[B]:THR:HG23	1:B:24:PRO:HA	2.02	0.42
3:B:606:ACT:H2	4:B:618:GOL:C3	2.35	0.42
4:A:627:GOL:H31	6:A:1050:HOH:O	2.19	0.42
1:B:137:ARG:HB2	1:B:138:PRO:HA	2.02	0.42
1:B:115:LYS:H	1:B:115:LYS:CD	2.28	0.41
1:A:290:ALA:HA	1:A:469:TYR:CE2	2.55	0.41
1:B:275:LYS:H	4:B:612:GOL:H11	1.85	0.41
1:A:487:PRO:HD2	6:A:803:HOH:O	2.20	0.41
1:A:97:ARG:HG2	6:A:1087:HOH:O	2.19	0.41
1:A:481:GLY:HA2	1:B:444:GLU:CG	2.51	0.41
1:B:434:GLU:OE1	1:B:453[B]:ASP:OD2	2.38	0.41
1:B:71:VAL:HG22	1:B:185:TRP:CB	2.50	0.41
1:B:116:THR:HG22	6:B:1024:HOH:O	2.20	0.41
1:B:127:TYR:N	1:B:128:PRO:CD	2.83	0.41
1:B:447:ARG:HG3	1:B:447:ARG:HH11	1.86	0.41
1:B:4:ASP:HA	1:B:499:PRO:CB	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:472:MET:HA	1:B:472:MET:HE2	2.03	0.41
1:A:283:LEU:HD13	1:A:287[B]:GLN:OE1	2.21	0.41
1:A:481:GLY:CA	1:B:444:GLU:HG2	2.50	0.41
1:B:255:THR:O	1:B:261:ASN:HA	2.20	0.40
1:B:77:GLU:OE1	6:B:1020:HOH:O	2.22	0.40
1:A:357:ALA:HB1	1:A:373:LEU:HD22	2.04	0.40
1:A:212:GLY:HA2	6:A:937:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	516/502 (103%)	499 (97%)	17 (3%)	0	100	100
1	B	504/502 (100%)	486 (96%)	17 (3%)	1 (0%)	47	44
All	All	1020/1004 (102%)	985 (97%)	34 (3%)	1 (0%)	51	49

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	126	TYR

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	443/425 (104%)	435 (98%)	8 (2%)	59	63
1	B	431/425 (101%)	421 (98%)	10 (2%)	50	53
All	All	874/850 (103%)	856 (98%)	18 (2%)	59	57

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	97	ARG
1	A	127	TYR
1	A	192	LYS
1	A	216	LEU
1	A	250[A]	MET
1	A	250[B]	MET
1	A	284[A]	TYR
1	A	284[B]	TYR
1	B	115	LYS
1	B	124	SER
1	B	127	TYR
1	B	130	LEU
1	B	187	GLU
1	B	242	LEU
1	B	250[A]	MET
1	B	250[B]	MET
1	B	334	GLU
1	B	439	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	174	GLN
1	A	288	GLN
1	A	416	HIS
1	B	106	GLN
1	B	288	GLN
1	B	302	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

50 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	GOL	B	613	-	5,5,5	0.40	0	5,5,5	0.42	0
3	ACT	A	607	-	1,3,3	1.42	0	0,3,3	0.00	-
3	ACT	A	613	-	1,3,3	1.52	0	0,3,3	0.00	-
4	GOL	A	625	-	5,5,5	0.45	0	5,5,5	0.26	0
3	ACT	A	610	-	1,3,3	1.61	0	0,3,3	0.00	-
3	ACT	A	609	-	1,3,3	1.46	0	0,3,3	0.00	-
4	GOL	B	615	-	5,5,5	0.35	0	5,5,5	0.61	0
3	ACT	B	607	-	1,3,3	1.58	0	0,3,3	0.00	-
3	ACT	B	602	-	1,3,3	1.36	0	0,3,3	0.00	-
4	GOL	A	622	-	5,5,5	0.43	0	5,5,5	0.25	0
5	L2D	B	620	-	28,38,38	4.17	15 (53%)	28,69,69	2.43	10 (35%)
3	ACT	A	612	-	1,3,3	1.36	0	0,3,3	0.00	-
4	GOL	A	626	-	5,5,5	0.36	0	5,5,5	0.33	0
4	GOL	B	618	-	5,5,5	0.37	0	5,5,5	0.38	0
4	GOL	A	629	-	5,5,5	0.44	0	5,5,5	0.21	0
3	ACT	A	619	-	1,3,3	1.62	0	0,3,3	0.00	-
4	GOL	A	620	-	5,5,5	0.33	0	5,5,5	0.40	0
4	GOL	A	623	-	5,5,5	0.42	0	5,5,5	0.37	0
4	GOL	A	624	-	5,5,5	0.35	0	5,5,5	0.28	0
3	ACT	B	605	-	1,3,3	1.38	0	0,3,3	0.00	-
3	ACT	B	611	-	1,3,3	1.16	0	0,3,3	0.00	-
3	ACT	B	608	-	1,3,3	1.25	0	0,3,3	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ACT	A	605	-	1,3,3	1.29	0	0,3,3	0.00	-
3	ACT	B	610	-	1,3,3	1.41	0	0,3,3	0.00	-
3	ACT	A	608	-	1,3,3	1.56	0	0,3,3	0.00	-
4	GOL	A	621	-	5,5,5	0.29	0	5,5,5	0.48	0
3	ACT	A	618	-	1,3,3	1.24	0	0,3,3	0.00	-
2	SO4	A	602	-	4,4,4	0.13	0	6,6,6	0.12	0
4	GOL	B	617	-	5,5,5	0.36	0	5,5,5	0.21	0
3	ACT	A	614	-	1,3,3	1.21	0	0,3,3	0.00	-
3	ACT	B	604	-	1,3,3	1.51	0	0,3,3	0.00	-
2	SO4	B	601	-	4,4,4	0.12	0	6,6,6	0.13	0
4	GOL	A	627	-	5,5,5	0.35	0	5,5,5	0.37	0
3	ACT	B	609	-	1,3,3	1.39	0	0,3,3	0.00	-
4	GOL	B	619	-	5,5,5	0.30	0	5,5,5	0.35	0
3	ACT	A	616	-	1,3,3	0.54	0	0,3,3	0.00	-
3	ACT	B	606	-	1,3,3	1.28	0	0,3,3	0.00	-
4	GOL	B	614	-	5,5,5	0.40	0	5,5,5	0.25	0
2	SO4	A	601	-	4,4,4	0.18	0	6,6,6	0.13	0
3	ACT	A	606	-	1,3,3	0.85	0	0,3,3	0.00	-
4	GOL	B	612	-	5,5,5	0.41	0	5,5,5	0.31	0
5	L2D	A	630	-	28,38,38	3.71	11 (39%)	28,69,69	2.81	13 (46%)
4	GOL	B	616	-	5,5,5	0.34	0	5,5,5	0.13	0
3	ACT	A	611	-	1,3,3	1.44	0	0,3,3	0.00	-
4	GOL	A	628	-	5,5,5	0.35	0	5,5,5	0.39	0
3	ACT	A	603	-	1,3,3	1.64	0	0,3,3	0.00	-
3	ACT	B	603	-	1,3,3	1.40	0	0,3,3	0.00	-
3	ACT	A	604	-	1,3,3	1.34	0	0,3,3	0.00	-
3	ACT	A	615	-	1,3,3	0.93	0	0,3,3	0.00	-
3	ACT	A	617	-	1,3,3	0.92	0	0,3,3	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	613	-	-	4/4/4/4	-
4	GOL	A	621	-	-	2/4/4/4	-
4	GOL	A	620	-	-	3/4/4/4	-
4	GOL	A	625	-	-	1/4/4/4	-
4	GOL	B	618	-	-	3/4/4/4	-
4	GOL	A	629	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	612	-	-	4/4/4/4	-
4	GOL	B	615	-	-	4/4/4/4	-
4	GOL	A	623	-	-	4/4/4/4	-
4	GOL	A	624	-	-	2/4/4/4	-
4	GOL	A	627	-	-	2/4/4/4	-
4	GOL	A	628	-	-	2/4/4/4	-
4	GOL	B	619	-	-	2/4/4/4	-
4	GOL	A	622	-	-	2/4/4/4	-
4	GOL	A	626	-	-	2/4/4/4	-
4	GOL	B	617	-	-	4/4/4/4	-
4	GOL	B	616	-	-	2/4/4/4	-
4	GOL	B	614	-	-	1/4/4/4	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	620	L2D	CAZ-NAV	-12.14	1.31	1.48
5	B	620	L2D	CAY-NAU	-11.77	1.32	1.48
5	A	630	L2D	CAZ-NAV	-10.61	1.34	1.48
5	A	630	L2D	CAY-NAU	-10.05	1.34	1.48
5	A	630	L2D	CAG-CAK	-5.51	1.38	1.53
5	B	620	L2D	CAH-CAL	-5.46	1.38	1.53
5	A	630	L2D	CAH-CAL	-5.27	1.39	1.53
5	B	620	L2D	CAL-CAZ	-5.19	1.38	1.52
5	B	620	L2D	CAK-CAY	-5.12	1.38	1.52
5	A	630	L2D	CAK-CAY	-5.10	1.38	1.52
5	B	620	L2D	CAG-CAK	-5.04	1.39	1.53
5	A	630	L2D	CAL-CAZ	-4.95	1.39	1.52
5	B	620	L2D	CAF-CAH	-3.62	1.37	1.51
5	A	630	L2D	CAI-CAE	-3.50	1.38	1.51
5	A	630	L2D	CAJ-CAF	-3.37	1.39	1.51
5	B	620	L2D	CAJ-CAF	-3.28	1.39	1.51
5	B	620	L2D	CAR-NBD	-3.28	1.44	1.48
5	B	620	L2D	CAI-CAE	-3.27	1.39	1.51
5	A	630	L2D	CAE-CAG	-3.11	1.39	1.51
5	B	620	L2D	CBA-N	-3.05	1.46	1.50
5	A	630	L2D	CAF-CAH	-3.00	1.39	1.51
5	B	620	L2D	CAE-CAG	-2.98	1.39	1.51
5	B	620	L2D	CA-N	-2.58	1.45	1.48
5	B	620	L2D	OXT-C	-2.42	1.23	1.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	620	L2D	OAD-CAX	-2.19	1.23	1.28
5	A	630	L2D	CA-C	2.03	1.56	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	630	L2D	CAP-CBB-NBD	-5.54	108.45	114.64
5	B	620	L2D	CAH-CAL-CAZ	4.87	121.19	110.83
5	A	630	L2D	CAJ-CAF-CAH	4.78	119.53	111.44
5	A	630	L2D	CAI-CAE-CAG	4.67	119.35	111.44
5	B	620	L2D	CAI-CAE-CAG	4.41	118.91	111.44
5	B	620	L2D	CAP-CBB-NBD	-4.24	109.90	114.64
5	A	630	L2D	CAE-CAG-CAK	4.19	119.95	111.42
5	B	620	L2D	CAG-CAK-CAY	4.17	119.70	110.83
5	A	630	L2D	CAS-N-CA	-3.92	105.91	111.76
5	B	620	L2D	CAO-CBA-N	-3.84	110.35	114.64
5	A	630	L2D	CAO-CBA-N	-3.82	110.37	114.64
5	B	620	L2D	CAT-NBD-CAR	-3.81	106.07	111.76
5	A	630	L2D	CAF-CAH-CAL	3.78	119.13	111.42
5	A	630	L2D	CAH-CAL-CAZ	3.66	118.61	110.83
5	B	620	L2D	CAJ-CAF-CAH	3.55	117.45	111.44
5	A	630	L2D	CAG-CAK-CAY	3.50	118.27	110.83
5	B	620	L2D	CAF-CAH-CAL	3.48	118.50	111.42
5	A	630	L2D	C-CA-N	3.41	115.03	111.32
5	A	630	L2D	CA-N-CBA	3.28	118.35	112.44
5	B	620	L2D	CAE-CAG-CAK	3.05	117.62	111.42
5	B	620	L2D	CA-N-CBA	2.93	117.72	112.44
5	A	630	L2D	CAT-NBD-CAR	-2.44	108.11	111.76
5	A	630	L2D	CAR-NBD-CBB	-2.42	108.07	112.44

There are no chirality outliers.

All (46) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	613	GOL	O1-C1-C2-C3
4	B	615	GOL	O1-C1-C2-C3
4	A	622	GOL	O1-C1-C2-O2
4	A	622	GOL	O1-C1-C2-C3
4	A	626	GOL	O1-C1-C2-C3
4	A	629	GOL	O1-C1-C2-O2
4	A	629	GOL	O1-C1-C2-C3
4	A	623	GOL	C1-C2-C3-O3

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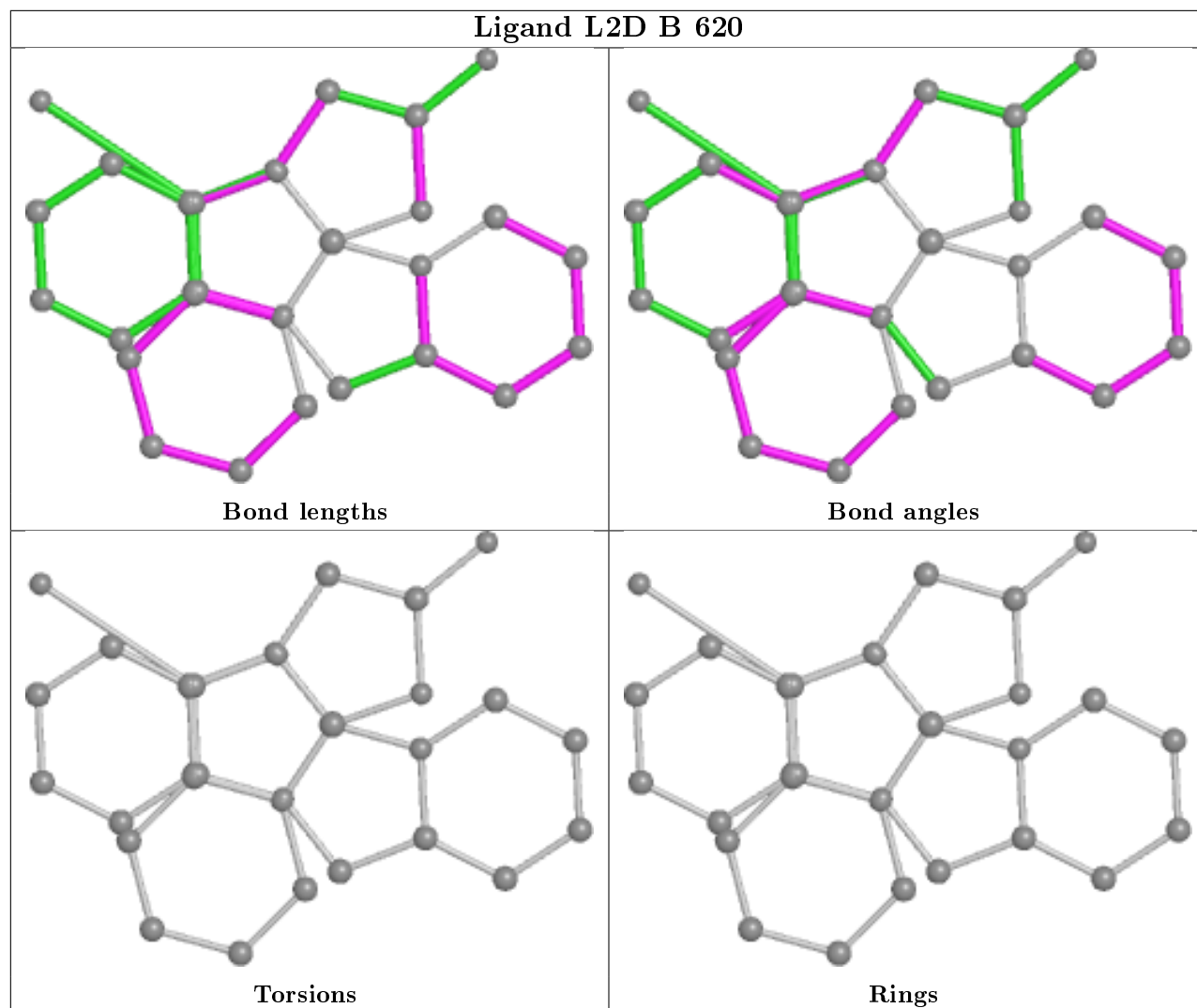
Mol	Chain	Res	Type	Atoms
4	A	624	GOL	O1-C1-C2-C3
4	B	618	GOL	O1-C1-C2-C3
4	B	612	GOL	O1-C1-C2-C3
4	B	612	GOL	C1-C2-C3-O3
4	B	616	GOL	O1-C1-C2-C3
4	A	628	GOL	O1-C1-C2-C3
4	A	626	GOL	O1-C1-C2-O2
4	A	620	GOL	O1-C1-C2-O2
4	B	612	GOL	O2-C2-C3-O3
4	B	613	GOL	C1-C2-C3-O3
4	B	615	GOL	C1-C2-C3-O3
4	A	620	GOL	O1-C1-C2-C3
4	A	623	GOL	O1-C1-C2-C3
4	A	621	GOL	O1-C1-C2-C3
4	B	617	GOL	O1-C1-C2-C3
4	B	617	GOL	C1-C2-C3-O3
4	B	618	GOL	C1-C2-C3-O3
4	B	619	GOL	O1-C1-C2-C3
4	B	613	GOL	O1-C1-C2-O2
4	B	615	GOL	O1-C1-C2-O2
4	B	615	GOL	O2-C2-C3-O3
4	A	621	GOL	O1-C1-C2-O2
4	B	617	GOL	O1-C1-C2-O2
4	B	617	GOL	O2-C2-C3-O3
4	B	619	GOL	O1-C1-C2-O2
4	B	616	GOL	O1-C1-C2-O2
4	A	623	GOL	O1-C1-C2-O2
4	A	623	GOL	O2-C2-C3-O3
4	A	624	GOL	O1-C1-C2-O2
4	B	612	GOL	O1-C1-C2-O2
4	A	628	GOL	O1-C1-C2-O2
4	B	614	GOL	O1-C1-C2-C3
4	B	613	GOL	O2-C2-C3-O3
4	B	618	GOL	O1-C1-C2-O2
4	A	625	GOL	O1-C1-C2-O2
4	A	627	GOL	O1-C1-C2-C3
4	A	627	GOL	O1-C1-C2-O2
4	A	620	GOL	C1-C2-C3-O3

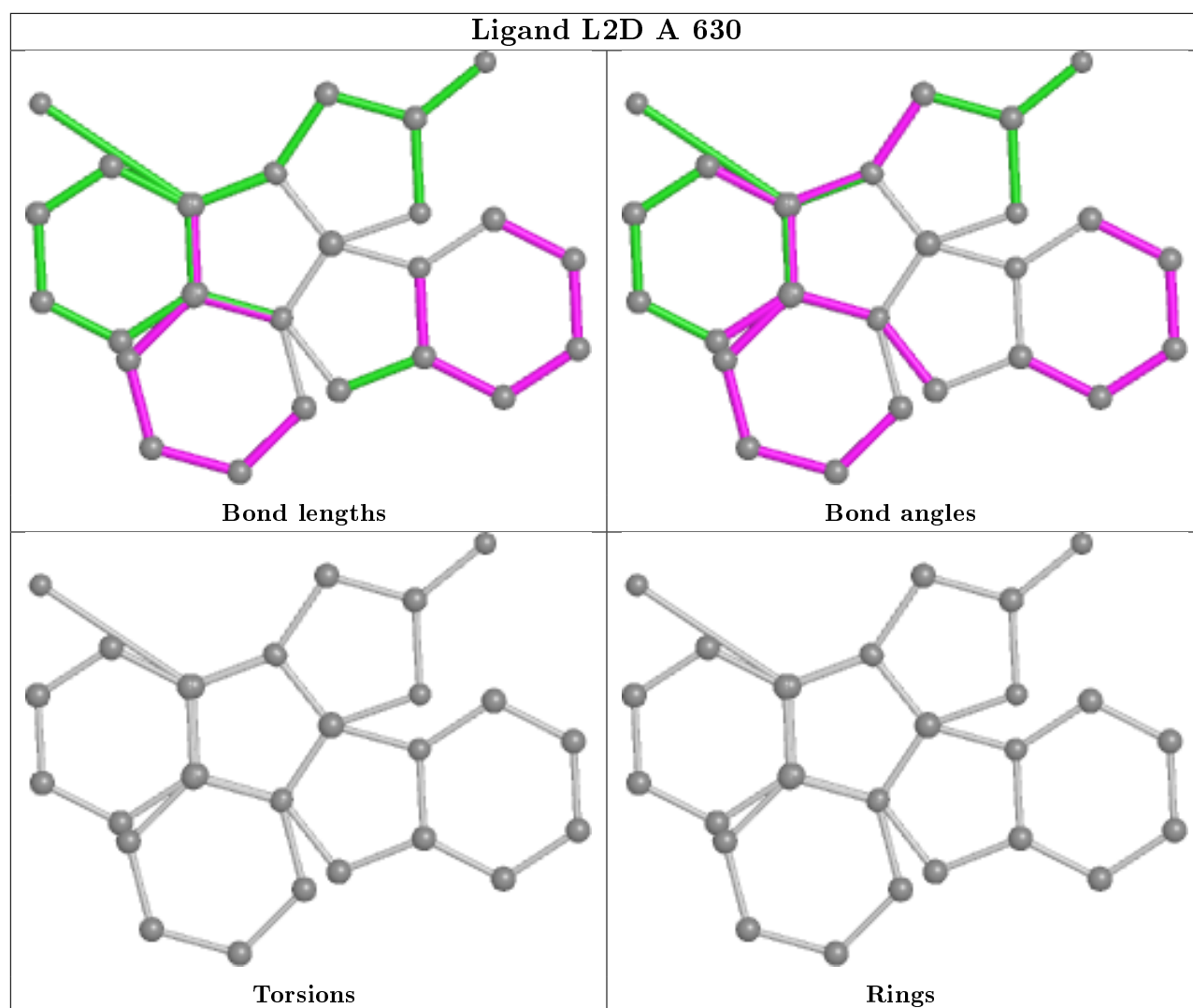
There are no ring outliers.

27 monomers are involved in 60 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	613	GOL	1	0
3	A	613	ACT	1	0
4	A	625	GOL	2	0
3	A	609	ACT	1	0
4	B	615	GOL	2	0
4	A	622	GOL	3	0
5	B	620	L2D	2	0
4	A	626	GOL	2	0
4	B	618	GOL	5	0
4	A	623	GOL	2	0
4	A	624	GOL	1	0
3	B	605	ACT	1	0
3	B	610	ACT	1	0
4	A	621	GOL	1	0
3	A	618	ACT	1	0
4	B	617	GOL	3	0
3	A	614	ACT	1	0
4	A	627	GOL	5	0
3	B	609	ACT	1	0
4	B	619	GOL	1	0
3	A	616	ACT	3	0
3	B	606	ACT	10	0
4	B	614	GOL	1	0
4	B	612	GOL	6	0
4	B	616	GOL	2	0
3	A	615	ACT	4	0
3	A	617	ACT	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	497/502 (99%)	-0.47	2 (0%) 92 92	17, 27, 49, 85	1 (0%)
1	B	497/502 (99%)	-0.19	13 (2%) 56 54	18, 36, 73, 113	3 (0%)
All	All	994/1004 (99%)	-0.33	15 (1%) 73 72	17, 30, 64, 113	4 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	329	GLY	3.5
1	B	337	GLY	3.4
1	B	125	ALA	3.1
1	B	328	ALA	3.0
1	B	323	GLY	2.9
1	B	113	LEU	2.9
1	B	59	ASP	2.8
1	A	3	PRO	2.8
1	B	117	GLU	2.7
1	B	57	SER	2.6
1	B	110	VAL	2.4
1	B	58	GLU	2.4
1	A	284[A]	TYR	2.1
1	B	326	LEU	2.1
1	B	111	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	B	617	6/6	0.63	0.39	90,91,92,95	0
3	ACT	A	617	4/4	0.66	0.26	63,64,70,79	0
4	GOL	B	616	6/6	0.67	0.39	75,78,79,80	0
4	GOL	B	614	6/6	0.67	0.39	80,81,82,84	0
4	GOL	A	626	6/6	0.68	0.65	120,122,123,125	0
3	ACT	A	608	4/4	0.68	0.36	48,59,64,68	4
3	ACT	B	603	4/4	0.69	0.24	77,82,82,84	0
4	GOL	A	628	6/6	0.70	0.42	74,77,82,85	0
3	ACT	A	616	4/4	0.71	0.40	35,39,40,43	4
3	ACT	A	607	4/4	0.73	0.31	44,54,56,57	4
4	GOL	B	618	6/6	0.76	0.41	60,70,75,83	0
3	ACT	A	615	4/4	0.77	0.32	34,38,39,40	4
3	ACT	A	619	4/4	0.78	0.21	64,72,72,78	0
4	GOL	A	627	6/6	0.78	0.22	33,45,56,58	0
3	ACT	B	604	4/4	0.80	0.21	55,59,62,64	0
4	GOL	A	629	6/6	0.81	0.23	52,67,74,77	0
4	GOL	B	612	6/6	0.81	0.23	61,69,71,78	0
3	ACT	A	603	4/4	0.81	0.16	63,69,69,70	0
4	GOL	A	624	6/6	0.82	0.17	59,79,85,87	0
3	ACT	A	605	4/4	0.84	0.27	74,82,83,86	0
3	ACT	A	609	4/4	0.84	0.25	79,83,83,83	0
3	ACT	B	610	4/4	0.84	0.25	38,63,69,72	0
4	GOL	A	620	6/6	0.85	0.20	51,65,68,70	0
4	GOL	A	623	6/6	0.85	0.21	29,54,61,61	0
4	GOL	B	613	6/6	0.86	0.20	38,52,61,61	0
2	SO4	A	602	5/5	0.87	0.25	91,96,101,101	0
3	ACT	A	611	4/4	0.87	0.37	77,81,81,82	0
3	ACT	A	613	4/4	0.87	0.14	64,68,69,69	0
4	GOL	B	619	6/6	0.88	0.14	48,59,60,63	0
3	ACT	A	614	4/4	0.89	0.31	55,58,60,66	4
3	ACT	A	612	4/4	0.89	0.15	72,74,74,74	0
3	ACT	A	618	4/4	0.89	0.38	68,69,71,73	0
3	ACT	B	606	4/4	0.89	0.34	59,60,61,61	4
3	ACT	A	604	4/4	0.90	0.15	45,54,54,60	0
3	ACT	B	611	4/4	0.90	0.58	44,45,49,55	4
4	GOL	A	621	6/6	0.90	0.15	35,44,59,63	0
3	ACT	B	602	4/4	0.91	0.13	68,69,71,75	0

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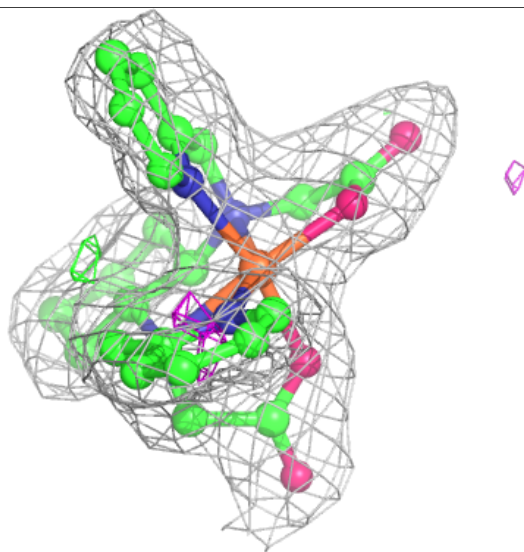
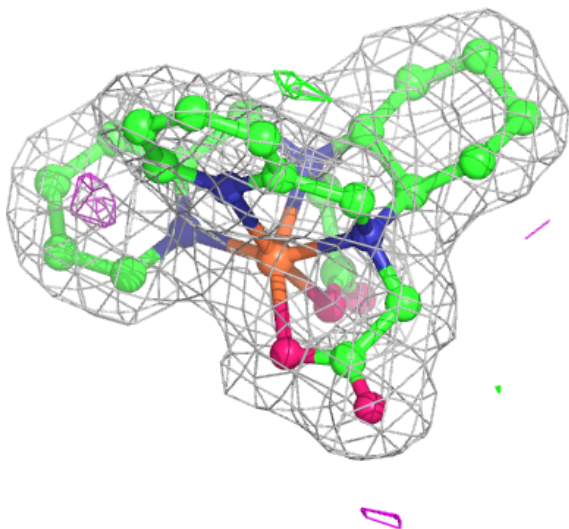
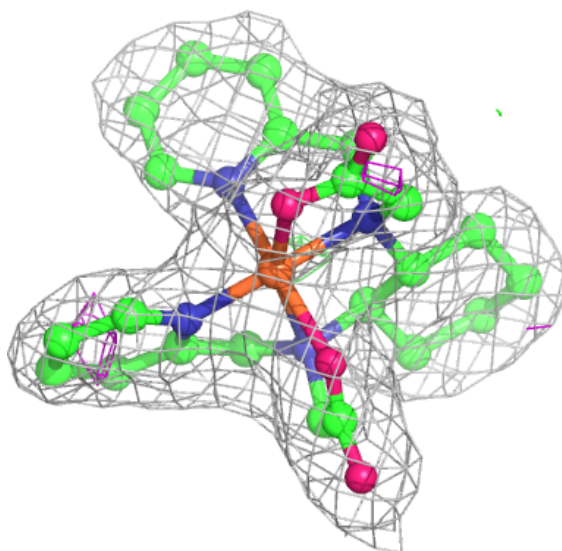
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	GOL	A	625	6/6	0.91	0.25	45,52,55,58	0
3	ACT	B	607	4/4	0.91	0.25	51,60,61,64	0
4	GOL	B	615	6/6	0.92	0.16	33,44,48,60	0
4	GOL	A	622	6/6	0.92	0.12	51,56,58,65	0
3	ACT	B	605	4/4	0.92	0.11	46,54,56,58	0
3	ACT	B	608	4/4	0.93	0.13	81,82,82,83	0
3	ACT	A	610	4/4	0.94	0.15	54,55,55,59	0
3	ACT	B	609	4/4	0.94	0.24	66,71,73,75	0
3	ACT	A	606	4/4	0.95	0.15	37,38,41,52	0
5	L2D	B	620	31/31	0.97	0.09	24,33,49,53	0
2	SO4	B	601	5/5	0.97	0.15	63,65,74,78	0
2	SO4	A	601	5/5	0.97	0.16	60,61,66,72	0
5	L2D	A	630	31/31	0.98	0.11	19,26,37,46	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

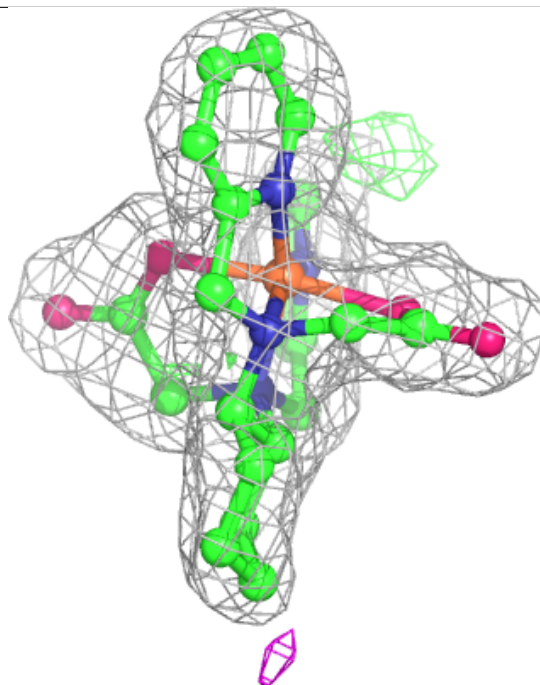
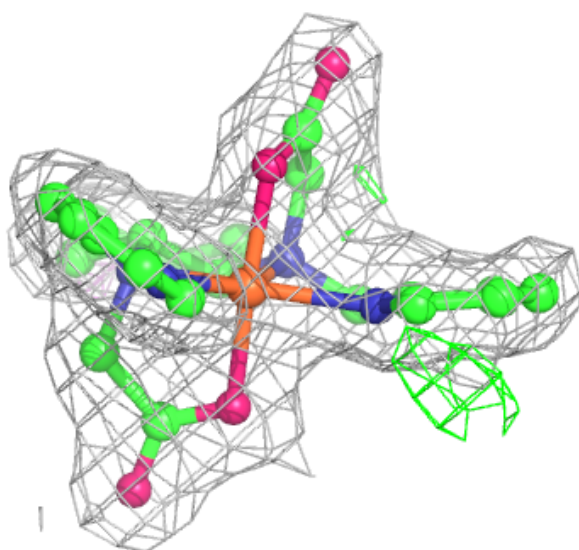
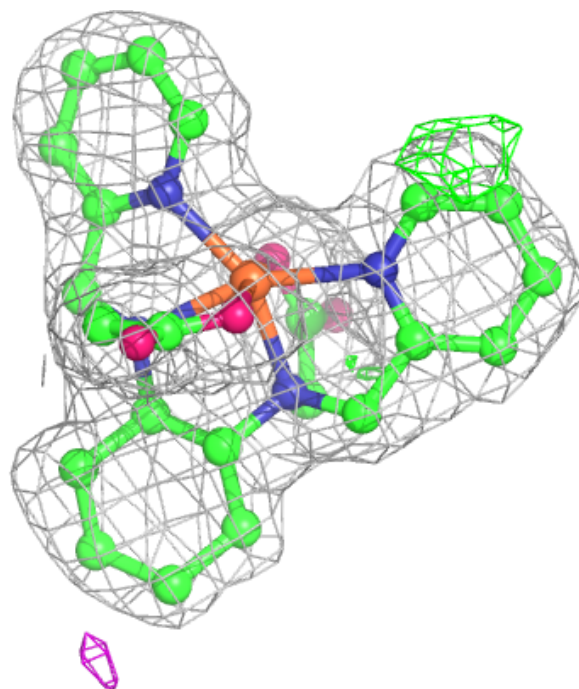
Electron density around L2D B 620:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around L2D A 630:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.