



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2026 – 11:34 pm BST

PDB ID : 4XE5 / pdb_00004xe5
Title : Crystal structure of the Na,K-ATPase from bovine
Authors : Gregersen, J.L.; Mattle, D.; Fedosova, N.U.; Nissen, P.; Reinhard, L.
Deposited on : 2014-12-22
Resolution : 3.90 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

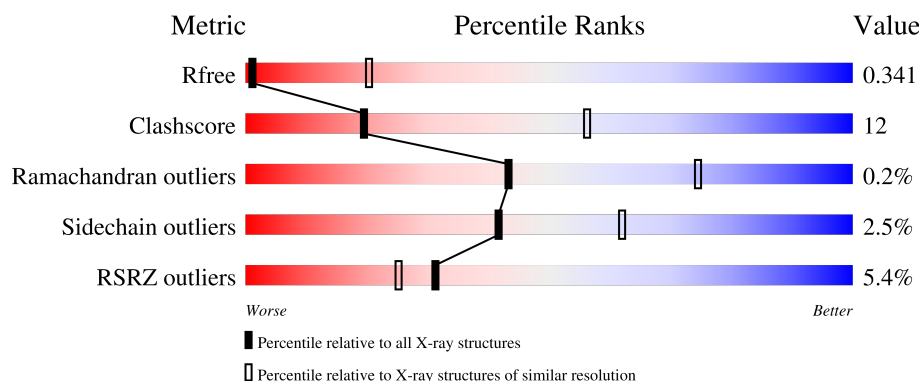
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.90 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}	180053	1270 (4.10-3.70)
Clashscore	190562	1034 (4.08-3.72)
Ramachandran outliers	187476	1251 (4.10-3.70)
Sidechain outliers	187428	1243 (4.10-3.70)
RSRZ outliers	180081	1269 (4.10-3.70)

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 of 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	1021	
2	B	303	
3	G	58	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 10362 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 Sodium/potassium-transporting ATPase subunit alpha-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	995	Total	C	N	O	S	0	0	0
			7713	4909	1303	1453	48			

- Molecule 2 is a protein called Sodium/potassium-transporting ATPase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	281	Total	C	N	O	S	0	0	0
			2286	1480	372	422	12			

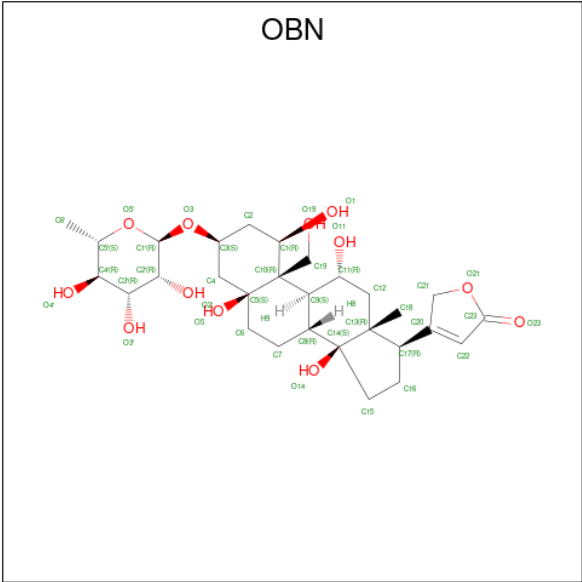
- Molecule 3 is a protein called Sodium/potassium-transporting ATPase subunit gamma.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	32	Total	C	N	O	0	0	0
			254	173	37	44			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

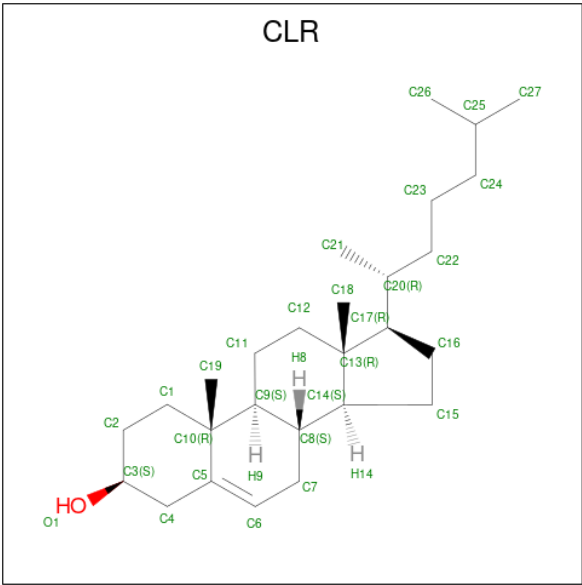
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Mg	0	0
			3	3		

- Molecule 5 is OUABAIN (CCD ID: OBN) (formula: C₂₉H₄₄O₁₂).



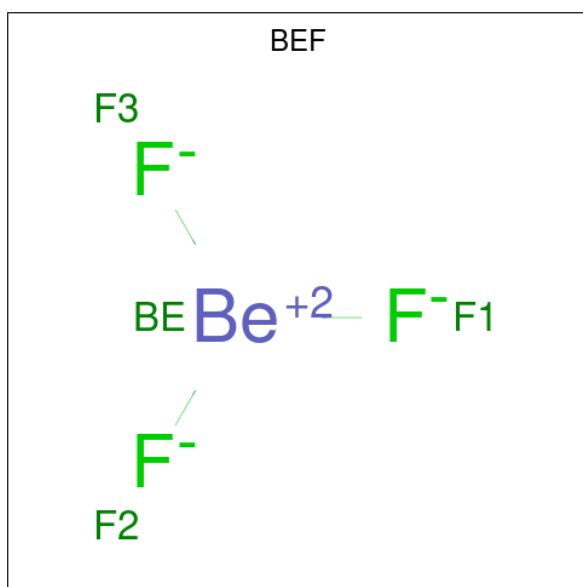
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			41	29	12		

- Molecule 6 is CHOLESTEROL (CCD ID: CLR) (formula: C₂₇H₄₆O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			28	27	1		
6	A	1	Total	C	O	0	0
			28	27	1		

- Molecule 7 is BERYLLIUM TRIFLUORIDE ION (CCD ID: BEF) (formula: BeF₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	Be	F	0	0
			4	1	3		

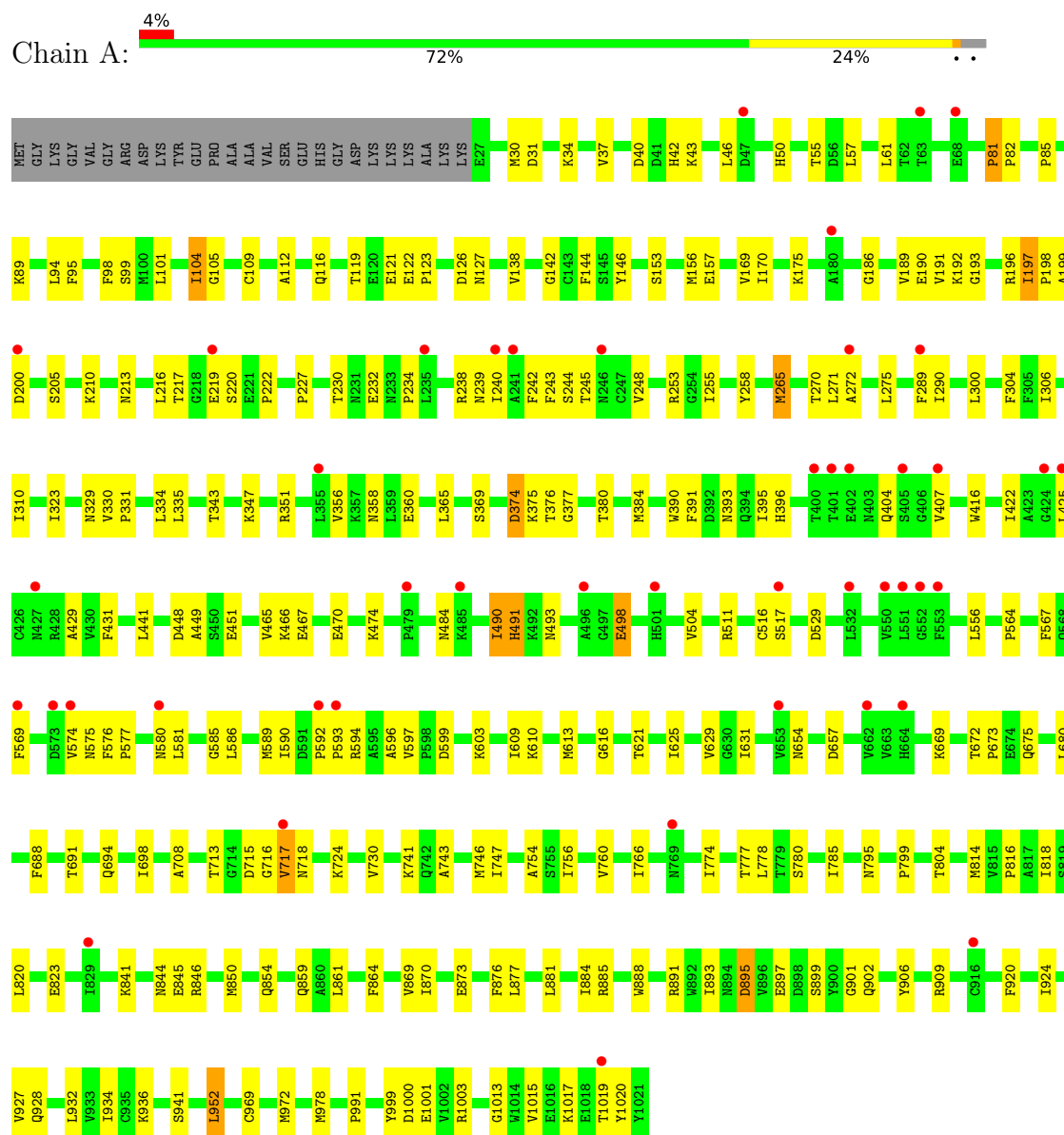
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	5	Total	O	0	0
			5	5		

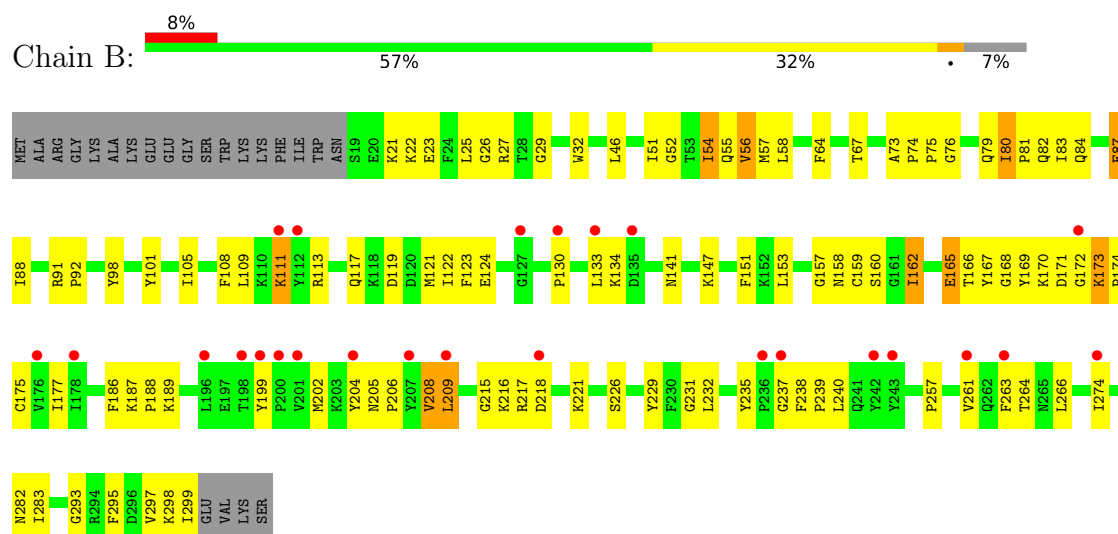
3 Residue-property plots

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.

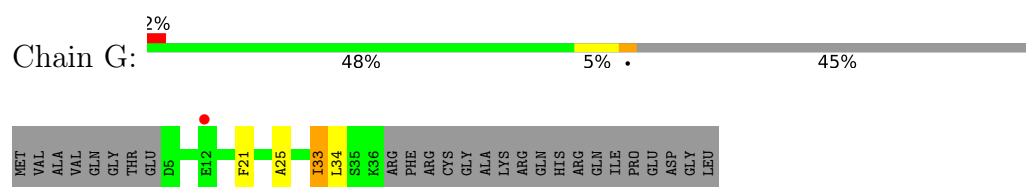
- Molecule 1: Sodium/potassium-transporting ATPase subunit alpha-1



- Molecule 2: Sodium/potassium-transporting ATPase subunit beta



• Molecule 3: Sodium/potassium-transporting ATPase subunit gamma



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	64.75Å 301.07Å 242.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.74 – 3.90 29.74 – 3.90	Depositor EDS
% Data completeness (in resolution range)	64.9 (29.74-3.90) 67.0 (29.74-3.90)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 3.75Å)	Xtriage
Refinement program	PHENIX dev_1888	Depositor
R, R_{free}	0.322 , 0.336 0.341 , 0.341	Depositor DCC
R_{free} test set	762 reflections (2.73%)	wwPDB-VP
Wilson B-factor (Å ²)	91.1	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 61.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.78	EDS
Total number of atoms	10362	wwPDB-VP
Average B, all atoms (Å ²)	111.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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: OBN, CLR, MG, BEF

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	0.34	0/7861	0.79	6/10668 (0.1%)
2	B	0.38	0/2344	0.83	5/3163 (0.2%)
3	G	0.39	0/260	0.84	0/353
All	All	0.35	0/10465	0.80	11/14184 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	197	ILE	CA-C-N	5.91	125.11	118.97
1	A	197	ILE	C-N-CA	5.91	125.11	118.97
2	B	87	GLU	N-CA-C	5.68	117.69	108.26
1	A	81	PRO	CA-C-N	5.68	125.36	119.56
1	A	81	PRO	C-N-CA	5.68	125.36	119.56
2	B	209	LEU	CA-C-N	5.56	125.76	120.31
2	B	209	LEU	C-N-CA	5.56	125.76	120.31
1	A	374	ASP	N-CA-C	-5.38	102.56	110.52
2	B	173	LYS	CA-C-N	5.13	125.04	119.76
2	B	173	LYS	C-N-CA	5.13	125.04	119.76
1	A	40	ASP	N-CA-C	5.08	115.23	108.38

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	165	GLU	Peptide

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	7713	0	7755	161	0
2	B	2286	0	2268	79	0
3	G	254	0	257	2	0
4	A	3	0	0	0	0
5	A	41	0	44	3	0
6	A	56	0	92	3	0
7	A	4	0	0	0	0
8	A	5	0	0	1	0
All	All	10362	0	10416	241	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 12.

All (241) 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)
2:B:80:ILE:HG13	2:B:81:PRO:HD3	1.58	0.85
1:A:123:PRO:HA	1:A:126:ASP:HB2	1.64	0.77
2:B:130:PRO:HD3	2:B:232:LEU:HD12	1.68	0.75
1:A:404:GLN:HE21	1:A:441:LEU:HD11	1.51	0.74
1:A:893:ILE:O	1:A:909:ARG:NH2	2.22	0.73
1:A:390:TRP:HB3	1:A:586:LEU:H	1.54	0.72
1:A:885:ARG:HA	1:A:888:TRP:HB3	1.73	0.71
1:A:112:ALA:HB2	1:A:323:ILE:HG21	1.73	0.71
1:A:609:ILE:HD11	1:A:760:VAL:HG21	1.74	0.69
1:A:200:ASP:HB2	1:A:258:TYR:HB2	1.74	0.69
2:B:80:ILE:HD11	2:B:177:ILE:H	1.58	0.69
2:B:165:GLU:HB3	2:B:166:THR:HA	1.74	0.68
1:A:493:ASN:ND2	1:A:498:GLU:O	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:170:LYS:HB2	2:B:174:PRO:HA	1.76	0.67
2:B:202:MET:HB2	2:B:204:TYR:CE2	2.30	0.65
1:A:876:PHE:HB3	1:A:881:LEU:HD21	1.78	0.65
1:A:85:PRO:HB2	1:A:89:LYS:HD3	1.78	0.63
1:A:869:VAL:HG22	2:B:57:MET:HG3	1.79	0.63
2:B:202:MET:HB2	2:B:204:TYR:HE2	1.63	0.63
2:B:217:ARG:NH1	2:B:218:ASP:OD2	2.33	0.62
2:B:121:MET:O	2:B:147:LYS:NZ	2.32	0.62
1:A:122:GLU:OE2	5:A:1104:OBN:O5	2.17	0.62
2:B:29:GLY:HA2	2:B:32:TRP:CD1	2.35	0.62
2:B:170:LYS:HB2	2:B:175:CYS:H	1.63	0.62
1:A:854:GLN:OE1	1:A:999:TYR:OH	2.17	0.61
1:A:884:ILE:HG22	1:A:888:TRP:HB2	1.82	0.61
1:A:380:THR:HA	1:A:593:PRO:HA	1.81	0.61
1:A:34:LYS:HE2	1:A:270:THR:HB	1.83	0.61
1:A:517:SER:HB3	1:A:580:ASN:HA	1.81	0.61
2:B:216:LYS:HB2	2:B:221:LYS:HG2	1.83	0.61
1:A:1015:VAL:O	1:A:1019:THR:OG1	2.14	0.60
1:A:895:ASP:OD1	2:B:82:GLN:NE2	2.35	0.60
2:B:189:LYS:H	2:B:282:ASN:HB2	1.67	0.60
1:A:109:CYS:O	1:A:127:ASN:ND2	2.35	0.60
2:B:168:GLY:HA3	2:B:171:ASP:HB3	1.84	0.60
1:A:190:GLU:HB2	1:A:253:ARG:HD3	1.83	0.59
1:A:242:PHE:O	1:A:245:THR:OG1	2.19	0.59
1:A:873:GLU:O	2:B:67:THR:OG1	2.21	0.59
2:B:133:LEU:HG	2:B:240:LEU:HB3	1.84	0.59
1:A:804:THR:HG22	1:A:978:MET:HE2	1.85	0.59
1:A:196:ARG:HE	1:A:244:SER:HA	1.68	0.58
1:A:186:GLY:HA2	1:A:255:ILE:HG23	1.85	0.58
2:B:187:LYS:O	2:B:282:ASN:ND2	2.36	0.58
1:A:795:ASN:HA	1:A:885:ARG:HE	1.69	0.57
2:B:160:SER:HB2	2:B:169:TYR:CE1	2.40	0.57
1:A:55:THR:HG23	1:A:61:LEU:HG	1.86	0.56
2:B:238:PHE:HD2	2:B:257:PRO:HB2	1.69	0.56
1:A:936:LYS:HE2	1:A:952:LEU:HD13	1.88	0.56
1:A:730:VAL:HG11	1:A:756:ILE:HD11	1.86	0.56
1:A:191:VAL:HG11	1:A:197:ILE:HD13	1.88	0.56
1:A:42:HIS:HB3	1:A:240:ILE:HD11	1.88	0.55
2:B:73:ALA:O	2:B:75:PRO:HD3	2.07	0.55
1:A:50:HIS:ND1	1:A:57:LEU:HD21	2.20	0.55
2:B:79:GLN:HB3	2:B:295:PHE:CZ	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:616:GLY:HA2	1:A:691:THR:H	1.72	0.54
1:A:567:PHE:CE2	1:A:569:PHE:HB3	2.43	0.54
1:A:42:HIS:O	1:A:239:ASN:ND2	2.40	0.54
1:A:425:LEU:HB3	1:A:491:HIS:HE1	1.73	0.54
1:A:594:ARG:HB2	1:A:597:VAL:HG23	1.91	0.53
1:A:895:ASP:OD1	1:A:895:ASP:N	2.41	0.53
1:A:230:THR:HG21	1:A:238:ARG:HD2	1.91	0.53
1:A:846:ARG:HD3	1:A:1020:TYR:O	2.08	0.53
2:B:88:ILE:HB	2:B:299:ILE:HG22	1.90	0.53
2:B:83:ILE:HG22	2:B:84:GLN:N	2.23	0.53
2:B:170:LYS:HB2	2:B:175:CYS:N	2.23	0.53
1:A:99:SER:HB3	1:A:138:VAL:HG13	1.91	0.53
2:B:215:GLY:HA2	2:B:274:ILE:HD13	1.91	0.53
2:B:209:LEU:HD21	2:B:283:ILE:HD11	1.91	0.53
1:A:98:PHE:CE2	1:A:335:LEU:HB2	2.44	0.53
1:A:123:PRO:C	1:A:126:ASP:H	2.18	0.52
1:A:42:HIS:CE1	1:A:234:PRO:HG3	2.45	0.52
1:A:713:THR:HG22	1:A:730:VAL:HB	1.91	0.52
1:A:275:LEU:HB2	1:A:724:LYS:HG2	1.92	0.52
1:A:688:PHE:HB3	1:A:691:THR:HG21	1.92	0.52
1:A:610:LYS:NZ	1:A:680:LEU:O	2.42	0.52
2:B:188:PRO:HB3	2:B:209:LEU:HD22	1.92	0.52
2:B:113:ARG:HA	2:B:153:LEU:HD11	1.92	0.51
2:B:186:PHE:CZ	2:B:282:ASN:HB3	2.45	0.51
1:A:205:SER:HA	1:A:227:PRO:HG3	1.91	0.51
2:B:109:LEU:HD23	2:B:153:LEU:HD23	1.93	0.51
1:A:467:GLU:HA	1:A:470:GLU:HG2	1.93	0.51
1:A:741:LYS:HG3	1:A:747:ILE:HD12	1.92	0.51
1:A:516:CYS:SG	1:A:581:LEU:HB2	2.51	0.51
2:B:171:ASP:OD1	2:B:172:GLY:N	2.44	0.51
2:B:174:PRO:HG2	2:B:263:PHE:HB2	1.93	0.51
1:A:217:THR:HG23	1:A:219:GLU:H	1.75	0.51
1:A:306:ILE:O	1:A:310:ILE:HG12	2.11	0.51
2:B:229:TYR:CD1	2:B:261:VAL:HG22	2.46	0.50
1:A:694:GLN:O	1:A:698:ILE:HG12	2.12	0.50
1:A:46:LEU:HD12	1:A:46:LEU:H	1.75	0.50
1:A:104:ILE:HG13	1:A:105:GLY:N	2.27	0.50
2:B:124:GLU:OE1	2:B:134:LYS:NZ	2.45	0.50
1:A:422:ILE:HG23	1:A:504:VAL:HB	1.93	0.50
1:A:850:MET:HA	1:A:854:GLN:HE21	1.77	0.50
1:A:220:SER:HB2	1:A:376:THR:HG21	1.92	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:208:VAL:HG22	2:B:239:PRO:HA	1.93	0.49
1:A:272:ALA:O	1:A:724:LYS:NZ	2.34	0.49
1:A:391:PHE:CE1	1:A:416:TRP:HA	2.47	0.49
1:A:599:ASP:O	1:A:603:LYS:HG2	2.13	0.49
1:A:816:PRO:HB3	1:A:932:LEU:HD22	1.94	0.49
2:B:80:ILE:HD12	2:B:105:ILE:HD12	1.93	0.49
1:A:616:GLY:HA2	1:A:691:THR:N	2.28	0.49
2:B:73:ALA:HB3	2:B:74:PRO:HD3	1.95	0.49
1:A:189:VAL:HG11	1:A:198:PRO:HG2	1.93	0.48
1:A:216:LEU:O	1:A:718:ASN:HB3	2.13	0.48
2:B:119:ASP:HB3	2:B:122:ILE:HG12	1.94	0.48
1:A:30:MET:O	1:A:34:LYS:HG3	2.13	0.48
1:A:112:ALA:HB3	1:A:127:ASN:HD21	1.77	0.48
1:A:265:MET:HE2	1:A:717:VAL:HG11	1.94	0.48
1:A:377:GLY:HA3	1:A:715:ASP:OD2	2.12	0.48
1:A:814:MET:O	1:A:818:ILE:HG12	2.13	0.48
1:A:897:GLU:HG3	1:A:901:GLY:HA2	1.95	0.48
1:A:613:MET:HE3	1:A:631:ILE:HD12	1.94	0.48
1:A:390:TRP:CZ2	1:A:393:ASN:HA	2.48	0.48
1:A:431:PHE:HB2	1:A:466:LYS:NZ	2.28	0.48
1:A:850:MET:SD	1:A:854:GLN:NE2	2.86	0.48
1:A:170:ILE:HG12	1:A:175:LYS:HG2	1.96	0.48
1:A:243:PHE:HD2	1:A:265:MET:HG2	1.78	0.48
1:A:569:PHE:HB2	1:A:575:ASN:CG	2.38	0.48
2:B:25:LEU:H	2:B:29:GLY:HA3	1.78	0.48
1:A:213:ASN:HB3	1:A:217:THR:HG22	1.96	0.47
2:B:166:THR:N	2:B:167:TYR:HA	2.29	0.47
1:A:116:GLN:HB3	1:A:121:GLU:HG2	1.95	0.47
3:G:21:PHE:CE1	3:G:25:ALA:HB2	2.49	0.47
1:A:941:SER:OG	1:A:1001:GLU:OE2	2.20	0.47
2:B:119:ASP:O	2:B:123:PHE:HB2	2.15	0.47
1:A:777:THR:O	1:A:780:SER:OG	2.30	0.47
1:A:934:ILE:HB	1:A:1000:ASP:OD2	2.15	0.47
1:A:1003:ARG:HE	1:A:1019:THR:HB	1.80	0.47
2:B:98:TYR:HA	2:B:101:TYR:HD2	1.79	0.47
1:A:425:LEU:HB3	1:A:491:HIS:CE1	2.50	0.46
1:A:780:SER:OG	8:A:1201:HOH:O	2.21	0.46
1:A:358:ASN:HD21	1:A:360:GLU:HG2	1.79	0.46
2:B:235:TYR:O	2:B:237:GLY:N	2.47	0.46
1:A:156:MET:HE2	1:A:156:MET:HA	1.97	0.46
1:A:891:ARG:HA	1:A:906:TYR:CD1	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:820:LEU:HD12	1:A:820:LEU:HA	1.81	0.46
2:B:26:GLY:HA2	2:B:27:ARG:HA	1.70	0.46
1:A:820:LEU:HD11	1:A:936:LYS:HA	1.97	0.46
1:A:841:LYS:HB2	1:A:844:ASN:HB3	1.98	0.46
2:B:76:GLY:HA2	2:B:293:GLY:H	1.81	0.46
2:B:133:LEU:HD12	2:B:133:LEU:H	1.81	0.46
1:A:43:LYS:NZ	1:A:232:GLU:OE2	2.32	0.46
1:A:390:TRP:HD1	1:A:395:ILE:HG12	1.81	0.45
1:A:654:ASN:HB3	1:A:657:ASP:OD2	2.15	0.45
2:B:204:TYR:C	2:B:206:PRO:HD2	2.42	0.45
1:A:490:ILE:HG12	1:A:569:PHE:HZ	1.81	0.45
2:B:22:LYS:HA	2:B:23:GLU:HA	1.36	0.45
2:B:101:TYR:O	2:B:105:ILE:HG12	2.16	0.45
1:A:861:LEU:HD11	2:B:46:LEU:HB3	1.99	0.45
5:A:1104:OBN:H8	5:A:1104:OBN:H191	1.72	0.45
2:B:151:PHE:HE1	2:B:232:LEU:HD23	1.81	0.45
1:A:192:LYS:HG2	1:A:193:GLY:H	1.82	0.45
1:A:920:PHE:O	1:A:924:ILE:HG12	2.16	0.45
1:A:390:TRP:HE3	1:A:585:GLY:HA2	1.82	0.45
2:B:80:ILE:HG22	2:B:108:PHE:CG	2.52	0.45
2:B:111:LYS:H	2:B:111:LYS:HG3	1.62	0.45
2:B:87:GLU:HA	2:B:298:LYS:O	2.17	0.45
1:A:556:LEU:HD22	1:A:581:LEU:HD23	1.98	0.45
1:A:596:ALA:HB1	1:A:754:ALA:HB2	1.98	0.44
1:A:672:THR:N	1:A:675:GLN:OE1	2.44	0.44
1:A:799:PRO:HG2	1:A:864:PHE:HE1	1.82	0.44
1:A:877:LEU:HD23	1:A:899:SER:HB2	1.99	0.44
1:A:969:CYS:HB3	1:A:972:MET:HG3	1.99	0.44
1:A:94:LEU:HD12	1:A:142:GLY:HA3	1.99	0.44
2:B:208:VAL:HG22	2:B:238:PHE:O	2.17	0.44
2:B:54:ILE:O	2:B:58:LEU:HG	2.17	0.44
1:A:216:LEU:HD11	1:A:243:PHE:O	2.18	0.44
1:A:448:ASP:OD1	1:A:449:ALA:N	2.46	0.44
1:A:474:LYS:HA	1:A:491:HIS:HD2	1.83	0.43
1:A:669:LYS:HE2	1:A:669:LYS:HB3	1.86	0.43
2:B:158:ASN:O	2:B:166:THR:OG1	2.36	0.43
1:A:199:ALA:HB1	1:A:258:TYR:O	2.18	0.43
1:A:300:LEU:O	1:A:304:PHE:HB2	2.19	0.43
2:B:174:PRO:HG3	2:B:266:LEU:HD12	2.00	0.43
1:A:101:LEU:HD22	1:A:290:ILE:HG23	2.00	0.43
1:A:153:SER:O	1:A:157:GLU:HG2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:64:PHE:CZ	2:B:141:ASN:HB2	2.53	0.43
2:B:21:LYS:HG3	2:B:22:LYS:H	1.84	0.43
1:A:210:LYS:HB3	1:A:222:PRO:HB2	1.99	0.43
1:A:823:GLU:OE2	1:A:936:LYS:NZ	2.42	0.43
1:A:1013:GLY:O	1:A:1017:LYS:HG2	2.19	0.43
1:A:369:SER:OG	1:A:708:ALA:HB1	2.19	0.43
1:A:629:VAL:HG23	1:A:631:ILE:HG13	2.01	0.43
1:A:50:HIS:HB2	1:A:55:THR:O	2.19	0.43
1:A:484:ASN:OD1	1:A:511:ARG:NH1	2.52	0.42
2:B:51:ILE:O	2:B:55:GLN:HG2	2.19	0.42
1:A:906:TYR:HA	1:A:909:ARG:HE	1.84	0.42
1:A:375:LYS:HG3	1:A:625:ILE:HG21	2.01	0.42
1:A:123:PRO:HG3	5:A:1104:OBN:O19	2.19	0.42
1:A:564:PRO:HG2	1:A:567:PHE:HB2	2.00	0.42
1:A:574:VAL:HG13	1:A:576:PHE:O	2.19	0.42
1:A:289:PHE:HE1	1:A:778:LEU:HD11	1.84	0.42
1:A:724:LYS:HB2	1:A:743:ALA:HB1	2.00	0.42
2:B:52:GLY:O	2:B:56:VAL:HG23	2.19	0.42
1:A:715:ASP:CG	1:A:716:GLY:N	2.78	0.42
2:B:80:ILE:HG12	2:B:177:ILE:HB	2.02	0.42
2:B:130:PRO:HB3	2:B:208:VAL:HG23	2.02	0.42
1:A:888:TRP:HA	1:A:909:ARG:NH1	2.35	0.42
1:A:391:PHE:HD2	1:A:396:HIS:CD2	2.38	0.42
6:A:1105:CLR:H162	6:A:1105:CLR:H221	1.69	0.42
2:B:134:LYS:HE2	2:B:134:LYS:HB3	1.88	0.41
2:B:157:GLY:C	2:B:159:CYS:H	2.28	0.41
1:A:81:PRO:HA	1:A:82:PRO:HD3	1.77	0.41
1:A:845:GLU:CD	1:A:845:GLU:H	2.28	0.41
1:A:144:PHE:CE2	1:A:343:THR:HG21	2.56	0.41
1:A:271:LEU:O	1:A:275:LEU:HG	2.19	0.41
1:A:730:VAL:HG22	1:A:746:MET:HE2	2.03	0.41
1:A:144:PHE:CZ	1:A:343:THR:HG21	2.55	0.41
1:A:330:VAL:HA	1:A:331:PRO:HD3	1.74	0.41
3:G:33:ILE:HG22	3:G:34:LEU:HD23	2.02	0.41
1:A:347:LYS:O	1:A:351:ARG:HG3	2.21	0.41
1:A:897:GLU:HA	1:A:902:GLN:O	2.21	0.41
2:B:80:ILE:CD1	2:B:177:ILE:H	2.30	0.41
2:B:205:ASN:N	2:B:206:PRO:HD2	2.36	0.41
1:A:592:PRO:HA	1:A:593:PRO:HD3	1.94	0.41
2:B:204:TYR:CE1	2:B:235:TYR:HB3	2.56	0.41
1:A:89:LYS:HG3	1:A:146:TYR:HE1	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:PRO:HB3	1:A:248:VAL:HG21	2.02	0.41
1:A:329:ASN:HA	1:A:785:ILE:HD11	2.02	0.41
1:A:991:PRO:HG3	6:A:1105:CLR:C15	2.51	0.41
1:A:1000:ASP:OD1	1:A:1003:ARG:NH1	2.54	0.41
2:B:226:SER:OG	2:B:264:THR:OG1	2.14	0.41
1:A:429:ALA:HB2	1:A:451:GLU:HB3	2.02	0.41
2:B:172:GLY:C	2:B:173:LYS:HG3	2.46	0.41
2:B:231:GLY:HA3	2:B:235:TYR:O	2.20	0.41
1:A:590:ILE:HD11	1:A:621:THR:HG23	2.02	0.40
1:A:672:THR:HB	1:A:673:PRO:HD2	2.02	0.40
2:B:91:ARG:HA	2:B:92:PRO:HD3	1.95	0.40
1:A:95:PHE:O	1:A:99:SER:HB2	2.21	0.40
1:A:859:GLN:HG2	1:A:927:VAL:HB	2.04	0.40
1:A:924:ILE:O	1:A:928:GLN:HG2	2.21	0.40
2:B:21:LYS:HD2	2:B:21:LYS:HA	1.70	0.40
1:A:334:LEU:HD11	1:A:774:ILE:HA	2.03	0.40
6:A:1105:CLR:H183	6:A:1105:CLR:H20	1.98	0.40
1:A:384:MET:HE3	1:A:589:MET:HG3	2.02	0.40
2:B:204:TYR:H	2:B:204:TYR:HD2	1.70	0.40
1:A:529:ASP:OD1	1:A:529:ASP:N	2.54	0.40
1:A:556:LEU:CD2	1:A:577:PRO:HG2	2.52	0.40
2:B:83:ILE:HD11	2:B:88:ILE:HG13	2.03	0.40
2:B:208:VAL:HA	2:B:240:LEU:HD13	2.03	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	993/1021 (97%)	949 (96%)	44 (4%)	0	100	100
2	B	279/303 (92%)	257 (92%)	20 (7%)	2 (1%)	18	53

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	G	30/58 (52%)	30 (100%)	0	0	100	100
All	All	1302/1382 (94%)	1236 (95%)	64 (5%)	2 (0%)	43	74

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	162	ILE
2	B	199	TYR

5.3.2 Protein sidechains ⓘ

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	843/862 (98%)	824 (98%)	19 (2%)	44	63
2	B	250/268 (93%)	242 (97%)	8 (3%)	34	56
3	G	26/47 (55%)	25 (96%)	1 (4%)	29	52
All	All	1119/1177 (95%)	1091 (98%)	28 (2%)	42	62

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ASP
1	A	37	VAL
1	A	104	ILE
1	A	119	THR
1	A	169	VAL
1	A	265	MET
1	A	356	VAL
1	A	365	LEU
1	A	374	ASP
1	A	407	VAL
1	A	465	VAL
1	A	490	ILE
1	A	491	HIS

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Mol	Chain	Res	Type
1	A	498	GLU
1	A	717	VAL
1	A	766	ILE
1	A	870	ILE
1	A	895	ASP
1	A	952	LEU
2	B	54	ILE
2	B	56	VAL
2	B	80	ILE
2	B	111	LYS
2	B	117	GLN
2	B	162	ILE
2	B	208	VAL
2	B	297	VAL
3	G	33	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	GLN
1	A	127	ASN
1	A	404	GLN
1	A	487	GLN
1	A	491	HIS
1	A	903	GLN
2	B	241	GLN
2	B	262	GLN
3	G	16	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 for Mogul analysis.

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
6	CLR	A	1105	-	31,31,31	2.41	8 (25%)	48,48,48	2.87	17 (35%)
5	OBN	A	1104	-	44,46,46	1.07	4 (9%)	66,76,76	1.64	13 (19%)
6	CLR	A	1106	-	31,31,31	2.34	8 (25%)	48,48,48	3.21	17 (35%)
7	BEF	A	1107	1	0,3,3	-	-	-	-	-

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
6	CLR	A	1105	-	-	8/10/68/68	0/4/4/4
5	OBN	A	1104	-	-	3/11/116/116	0/6/6/6
6	CLR	A	1106	-	-	2/10/68/68	0/4/4/4

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1106	CLR	C6-C5	9.19	1.53	1.33
6	A	1105	CLR	C6-C5	9.09	1.53	1.33
6	A	1105	CLR	C11-C9	5.05	1.62	1.53
6	A	1106	CLR	C11-C9	4.67	1.61	1.53
6	A	1105	CLR	C8-C14	3.80	1.60	1.53
6	A	1106	CLR	C8-C14	3.63	1.60	1.53
6	A	1105	CLR	C12-C13	3.23	1.59	1.54
6	A	1106	CLR	C15-C14	-3.16	1.47	1.54
6	A	1105	CLR	C15-C14	-3.02	1.48	1.54
6	A	1105	CLR	C20-C17	-2.55	1.49	1.54
6	A	1106	CLR	C16-C17	2.51	1.59	1.54
6	A	1106	CLR	C20-C17	-2.48	1.50	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1104	OBN	C4-C5	2.48	1.56	1.53
6	A	1105	CLR	C13-C17	-2.48	1.50	1.55
5	A	1104	OBN	O14-C14	-2.35	1.40	1.44
6	A	1106	CLR	C12-C13	2.25	1.58	1.54
6	A	1105	CLR	C16-C17	2.24	1.59	1.54
5	A	1104	OBN	O5-C5	-2.16	1.40	1.44
5	A	1104	OBN	C4'-C5'	2.13	1.57	1.52
6	A	1106	CLR	C13-C17	-2.05	1.51	1.55

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1106	CLR	C7-C6-C5	-8.44	109.49	125.06
6	A	1105	CLR	C4-C5-C6	-7.90	109.22	120.61
6	A	1106	CLR	C10-C5-C6	-6.88	112.37	122.90
6	A	1105	CLR	C7-C6-C5	-6.86	112.41	125.06
6	A	1106	CLR	C19-C10-C9	-6.78	103.59	111.68
6	A	1106	CLR	C12-C13-C17	-6.51	106.82	116.57
6	A	1105	CLR	C12-C13-C17	-6.40	106.99	116.57
6	A	1106	CLR	C21-C20-C17	6.35	122.65	112.92
6	A	1106	CLR	C18-C13-C14	6.13	123.15	111.71
6	A	1105	CLR	C1-C10-C9	6.04	117.16	108.73
5	A	1104	OBN	C1'-O5'-C5'	5.98	123.95	113.67
6	A	1106	CLR	C1-C10-C9	5.96	117.06	108.73
6	A	1105	CLR	C19-C10-C9	-5.80	104.76	111.68
6	A	1105	CLR	C10-C5-C6	-5.76	114.09	122.90
6	A	1106	CLR	C8-C7-C6	-5.70	104.54	112.73
6	A	1105	CLR	C7-C8-C9	4.59	115.28	109.71
6	A	1106	CLR	C7-C8-C14	4.47	117.38	110.91
6	A	1105	CLR	C7-C8-C14	4.37	117.24	110.91
5	A	1104	OBN	O5'-C5'-C4'	4.26	117.16	109.52
6	A	1106	CLR	C4-C5-C6	-3.98	114.88	120.61
6	A	1106	CLR	C21-C20-C22	-3.64	104.66	110.36
6	A	1105	CLR	C21-C20-C17	3.60	118.43	112.92
5	A	1104	OBN	O5-C5-C10	-3.59	103.97	109.54
6	A	1105	CLR	C16-C17-C13	-3.49	99.63	103.84
6	A	1106	CLR	C15-C14-C13	-3.42	99.72	103.84
6	A	1105	CLR	C11-C9-C10	3.16	117.25	113.08
6	A	1106	CLR	C18-C13-C12	-3.10	105.70	110.59
6	A	1106	CLR	C7-C8-C9	3.04	113.39	109.71
6	A	1105	CLR	C12-C11-C9	2.89	118.12	113.11
5	A	1104	OBN	C6'-C5'-C4'	-2.87	107.77	113.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1104	OBN	O5'-C1'-C2'	2.84	116.37	110.35
5	A	1104	OBN	O1-C1-C10	-2.79	105.31	110.35
6	A	1106	CLR	C11-C9-C8	2.74	115.70	111.75
6	A	1105	CLR	C1-C2-C3	2.68	113.91	110.47
5	A	1104	OBN	C14-C13-C17	2.67	106.62	103.56
6	A	1105	CLR	C17-C13-C14	2.62	103.18	100.07
6	A	1105	CLR	C19-C10-C1	-2.61	105.31	109.43
5	A	1104	OBN	C4-C5-C10	2.57	115.00	111.58
5	A	1104	OBN	O2'-C2'-C1'	-2.50	103.96	110.05
6	A	1106	CLR	C16-C15-C14	-2.43	100.31	105.13
5	A	1104	OBN	O3'-C3'-C2'	-2.32	104.98	110.35
5	A	1104	OBN	C7-C8-C14	-2.17	109.26	111.39
6	A	1106	CLR	C11-C12-C13	-2.14	109.10	112.78
5	A	1104	OBN	C9-C10-C1	2.13	115.45	110.96
6	A	1105	CLR	C11-C9-C8	2.12	114.80	111.75
5	A	1104	OBN	O4'-C4'-C5'	-2.02	105.18	109.67
6	A	1105	CLR	C21-C20-C22	-2.00	107.22	110.36

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1104	OBN	O5'-C1'-O3-C3
6	A	1105	CLR	C13-C17-C20-C22
6	A	1105	CLR	C13-C17-C20-C21
6	A	1105	CLR	C20-C22-C23-C24
6	A	1106	CLR	C22-C23-C24-C25
6	A	1105	CLR	C17-C20-C22-C23
5	A	1104	OBN	C1-C10-C19-O19
5	A	1104	OBN	C9-C10-C19-O19
6	A	1105	CLR	C16-C17-C20-C21
6	A	1105	CLR	C16-C17-C20-C22
6	A	1105	CLR	C23-C24-C25-C26
6	A	1105	CLR	C23-C24-C25-C27
6	A	1106	CLR	C20-C22-C23-C24

There are no ring outliers.

2 monomers are involved in 6 short contacts:

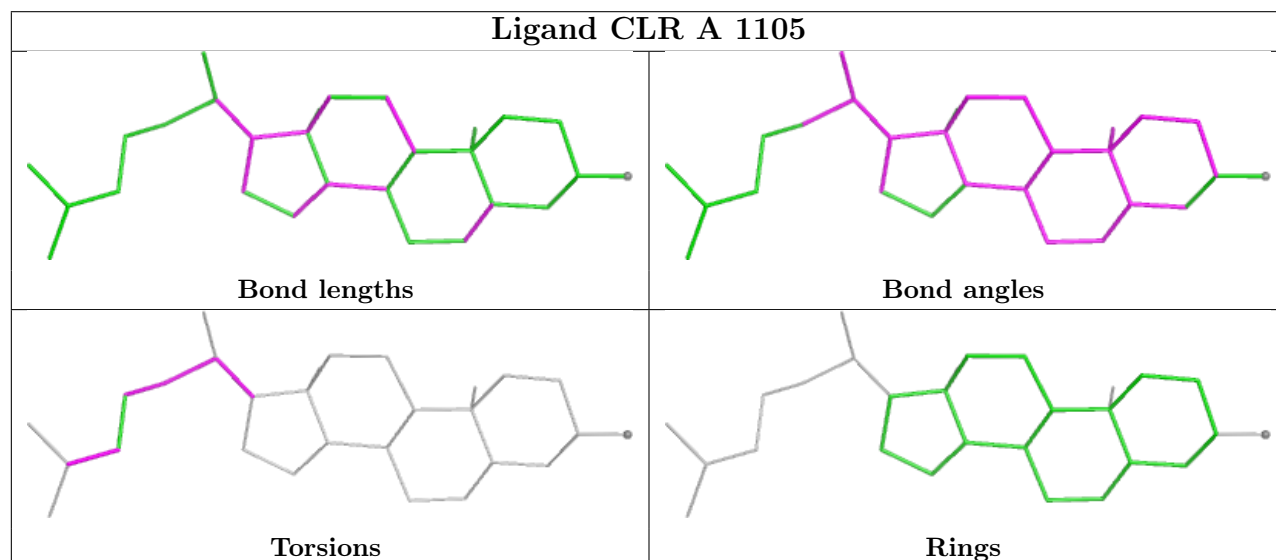
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1105	CLR	3	0

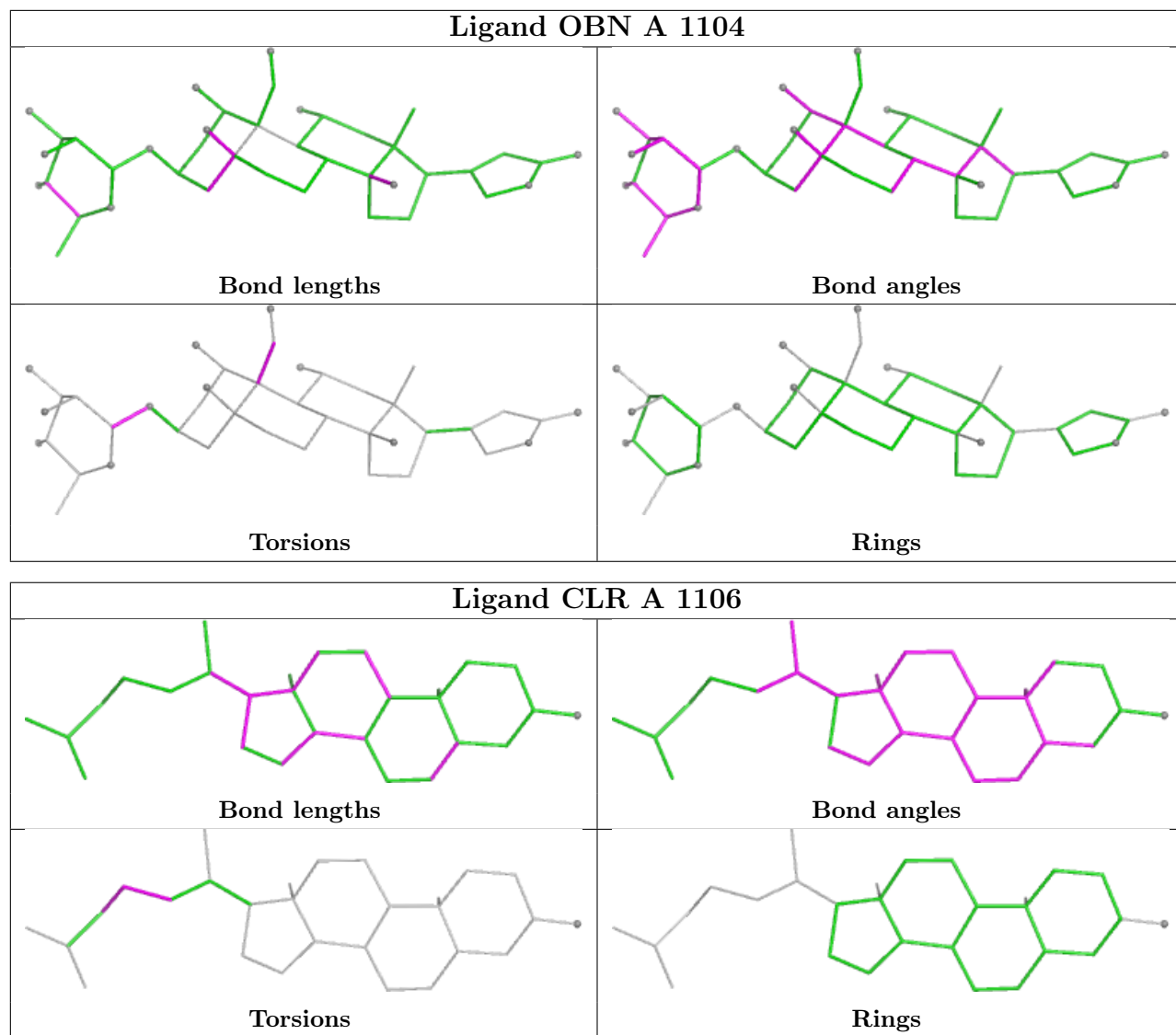
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1104	OBN	3	0

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

6.1 Protein, DNA and RNA chains

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	995/1021 (97%)	0.58	45 (4%) 38 28	61, 110, 163, 195	0
2	B	281/303 (92%)	0.81	25 (8%) 15 16	70, 114, 153, 167	0
3	G	32/58 (55%)	0.07	1 (3%) 51 36	61, 71, 91, 92	0
All	All	1308/1382 (94%)	0.62	71 (5%) 31 25	61, 112, 160, 195	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	263	PHE	3.9
1	A	240	ILE	3.7
1	A	496	ALA	3.7
2	B	218	ASP	3.7
1	A	241	ALA	3.5
2	B	127	GLY	3.3
2	B	178	ILE	3.3
2	B	198	THR	3.2
1	A	769	ASN	3.1
1	A	593	PRO	3.1
1	A	402	GLU	3.0
2	B	242	TYR	3.0
1	A	552	GLY	3.0
2	B	199	TYR	2.9
1	A	485	LYS	2.8
1	A	355	LEU	2.8
2	B	135	ASP	2.7
1	A	1019	THR	2.6
1	A	235	LEU	2.6
2	B	236	PRO	2.6
2	B	237	GLY	2.6
1	A	272	ALA	2.6
2	B	130	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	68	GLU	2.6
1	A	653	VAL	2.6
2	B	112	TYR	2.5
1	A	425	LEU	2.5
2	B	172	GLY	2.5
2	B	133	LEU	2.5
1	A	289	PHE	2.5
1	A	829	ILE	2.4
2	B	204	TYR	2.4
1	A	580	ASN	2.4
1	A	550	VAL	2.4
2	B	274	ILE	2.4
1	A	405	SER	2.4
2	B	111	LYS	2.4
2	B	209	LEU	2.4
2	B	176	VAL	2.4
1	A	532	LEU	2.3
1	A	501	HIS	2.3
2	B	196	LEU	2.3
1	A	63	THR	2.3
1	A	400	THR	2.3
2	B	200	PRO	2.3
2	B	201	VAL	2.2
1	A	246	ASN	2.2
1	A	592	PRO	2.2
1	A	180	ALA	2.2
1	A	664	HIS	2.2
1	A	574	VAL	2.2
1	A	427	ASN	2.2
1	A	424	GLY	2.1
1	A	517	SER	2.1
1	A	407	VAL	2.1
1	A	200	ASP	2.1
2	B	207	TYR	2.1
2	B	243	TYR	2.1
2	B	261	VAL	2.1
1	A	916	CYS	2.1
1	A	479	PRO	2.0
1	A	401	THR	2.0
1	A	573	ASP	2.0
1	A	717	VAL	2.0
3	G	12	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	662	VAL	2.0
1	A	219	GLU	2.0
1	A	47	ASP	2.0
1	A	551	LEU	2.0
1	A	553	PHE	2.0
1	A	569	PHE	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

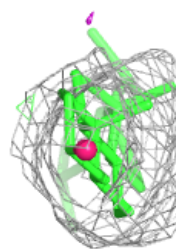
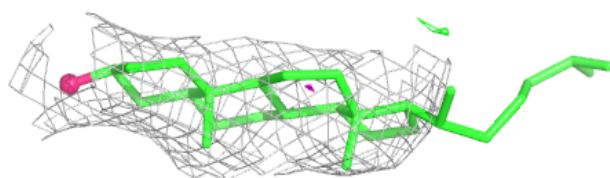
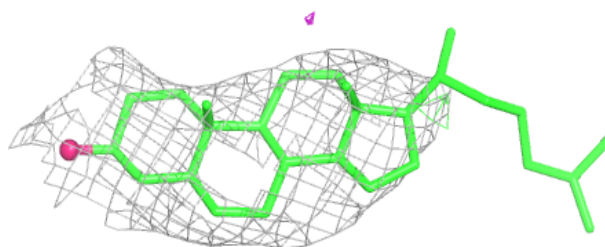
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	A	1102	1/1	0.68	0.28	85,85,85,85	0
6	CLR	A	1105	28/28	0.83	0.24	86,91,95,96	0
5	OBN	A	1104	41/41	0.86	0.13	86,91,95,97	0
6	CLR	A	1106	28/28	0.89	0.20	86,88,91,92	0
4	MG	A	1103	1/1	0.96	0.08	97,97,97,97	0
7	BEF	A	1107	4/4	0.97	0.04	106,108,108,110	0
4	MG	A	1101	1/1	0.98	0.03	89,89,89,89	0

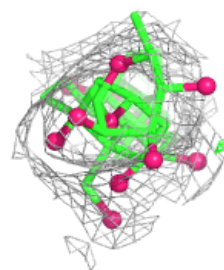
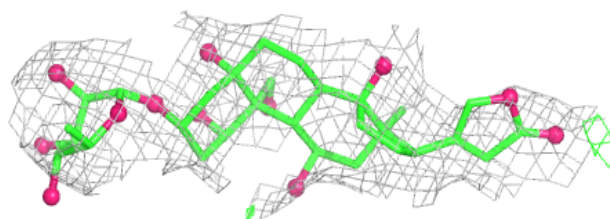
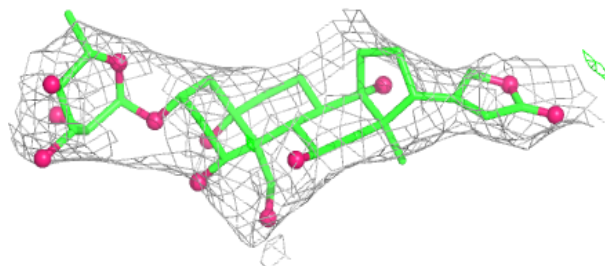
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

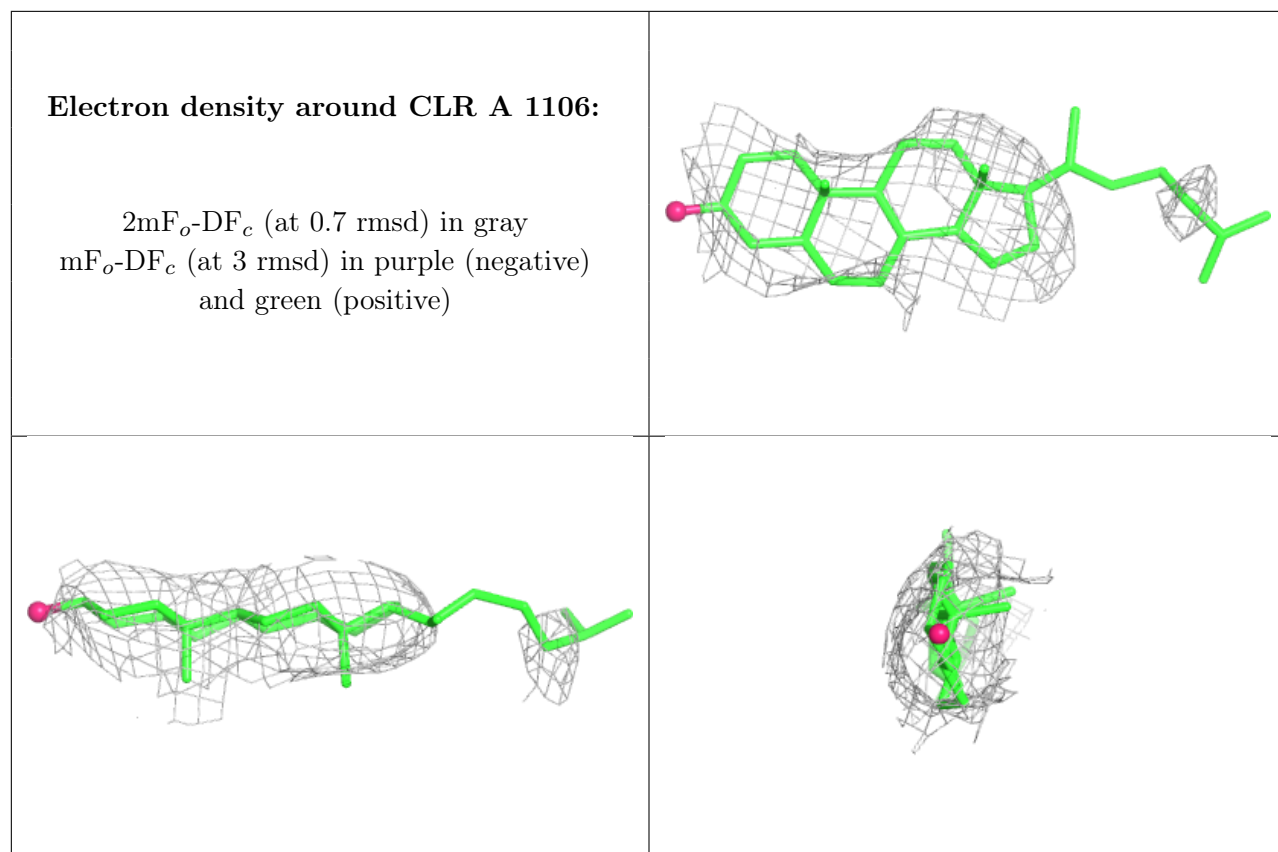
Electron density around CLR A 1105:

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

**Electron density around OBN A 1104:**

$2mF_o-DF_c$ (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.