



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:43 AM UTC

PDB ID : 5X9H / pdb_00005x9h
Title : Crystal structure of the Mg²⁺ channel MgtE in complex with ATP
Authors : Tomita, A.; Hattori, M.; Nureki, O.
Deposited on : 2017-03-07
Resolution : 3.60 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 : 2022.3.0, CSD as543be (2022)
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 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

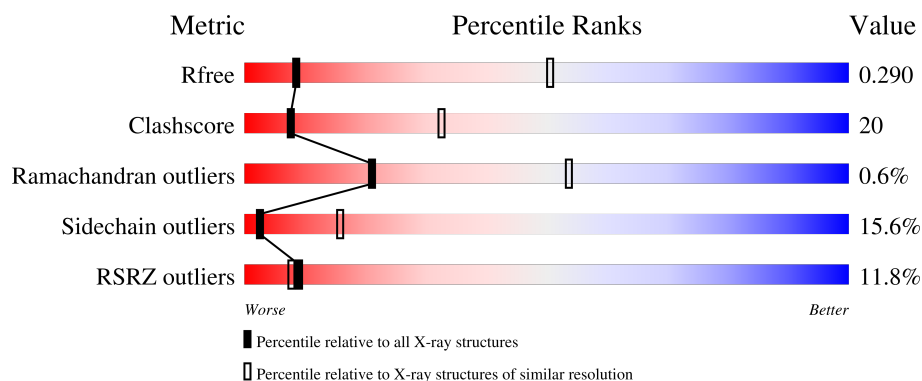
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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	473	10% 52% 32% 6% 10%
1	B	473	11% 52% 32% 6% 10%

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	MG	A	502	-	-	-	X
3	MG	B	503	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6522 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 Magnesium transporter MgtE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	427	Total	C	N	O	S	0	0	0
			3248	2108	528	605	7			
1	B	426	Total	C	N	O	S	0	0	0
			3199	2084	504	604	7			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	expression tag	UNP Q5SMG8
A	-21	GLY	-	expression tag	UNP Q5SMG8
A	-20	SER	-	expression tag	UNP Q5SMG8
A	-19	SER	-	expression tag	UNP Q5SMG8
A	-18	HIS	-	expression tag	UNP Q5SMG8
A	-17	HIS	-	expression tag	UNP Q5SMG8
A	-16	HIS	-	expression tag	UNP Q5SMG8
A	-15	HIS	-	expression tag	UNP Q5SMG8
A	-14	HIS	-	expression tag	UNP Q5SMG8
A	-13	HIS	-	expression tag	UNP Q5SMG8
A	-12	SER	-	expression tag	UNP Q5SMG8
A	-11	SER	-	expression tag	UNP Q5SMG8
A	-10	GLY	-	expression tag	UNP Q5SMG8
A	-9	LEU	-	expression tag	UNP Q5SMG8
A	-8	GLY	-	expression tag	UNP Q5SMG8
A	-7	VAL	-	expression tag	UNP Q5SMG8
A	-6	LEU	-	expression tag	UNP Q5SMG8
A	-5	PRO	-	expression tag	UNP Q5SMG8
A	-4	GLY	-	expression tag	UNP Q5SMG8
A	-3	GLY	-	expression tag	UNP Q5SMG8
A	-2	PRO	-	expression tag	UNP Q5SMG8
A	-1	LEU	-	expression tag	UNP Q5SMG8
A	0	HIS	-	expression tag	UNP Q5SMG8
B	-22	MET	-	expression tag	UNP Q5SMG8
B	-21	GLY	-	expression tag	UNP Q5SMG8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	SER	-	expression tag	UNP Q5SMG8
B	-19	SER	-	expression tag	UNP Q5SMG8
B	-18	HIS	-	expression tag	UNP Q5SMG8
B	-17	HIS	-	expression tag	UNP Q5SMG8
B	-16	HIS	-	expression tag	UNP Q5SMG8
B	-15	HIS	-	expression tag	UNP Q5SMG8
B	-14	HIS	-	expression tag	UNP Q5SMG8
B	-13	HIS	-	expression tag	UNP Q5SMG8
B	-12	SER	-	expression tag	UNP Q5SMG8
B	-11	SER	-	expression tag	UNP Q5SMG8
B	-10	GLY	-	expression tag	UNP Q5SMG8
B	-9	LEU	-	expression tag	UNP Q5SMG8
B	-8	GLY	-	expression tag	UNP Q5SMG8
B	-7	VAL	-	expression tag	UNP Q5SMG8
B	-6	LEU	-	expression tag	UNP Q5SMG8
B	-5	PRO	-	expression tag	UNP Q5SMG8
B	-4	GLY	-	expression tag	UNP Q5SMG8
B	-3	GLY	-	expression tag	UNP Q5SMG8
B	-2	PRO	-	expression tag	UNP Q5SMG8
B	-1	LEU	-	expression tag	UNP Q5SMG8
B	0	HIS	-	expression tag	UNP Q5SMG8

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- The diagram illustrates the chemical structure of Adenosine Triphosphate (ATP). It consists of an adenine base (a purine ring system with an amino group at N6) attached to a ribose sugar (a five-membered ring with hydroxyl groups at C2' and C3'). The ribose is linked to a chain of three phosphate groups (O3A, O3B, O3C) via phosphodiester bonds. The structure is color-coded: adenine is blue, ribose is green, and the phosphate chain is pink. Stereochemistry is indicated with wedges and dashes at the C1' and C5' positions of the ribose ring.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 31	C 10	N 5	O 13	P 3	0	0



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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

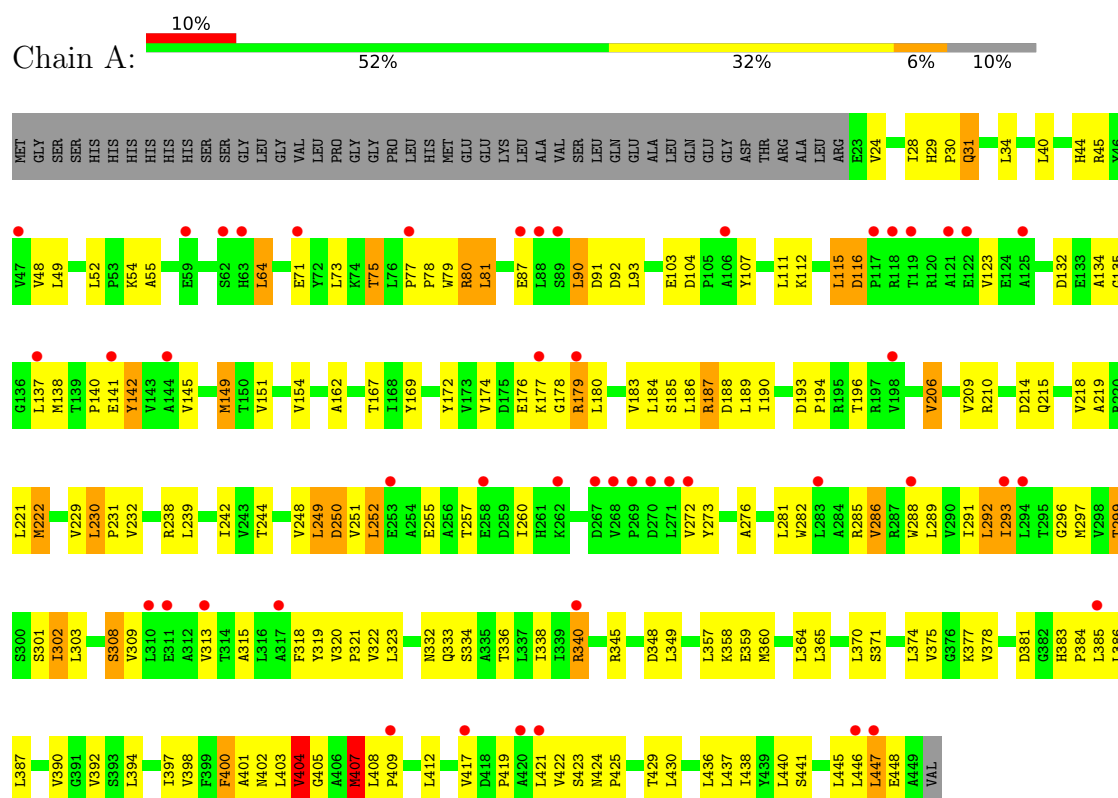
- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total	Mg	0	0
			8	8		
3	B	5	Total	Mg	0	0
			5	5		

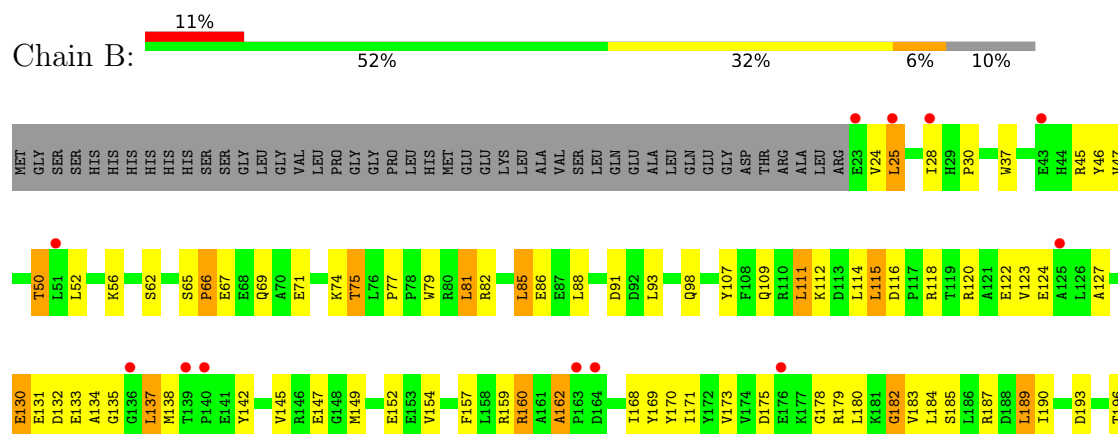
3 Residue-property plots [i](#)

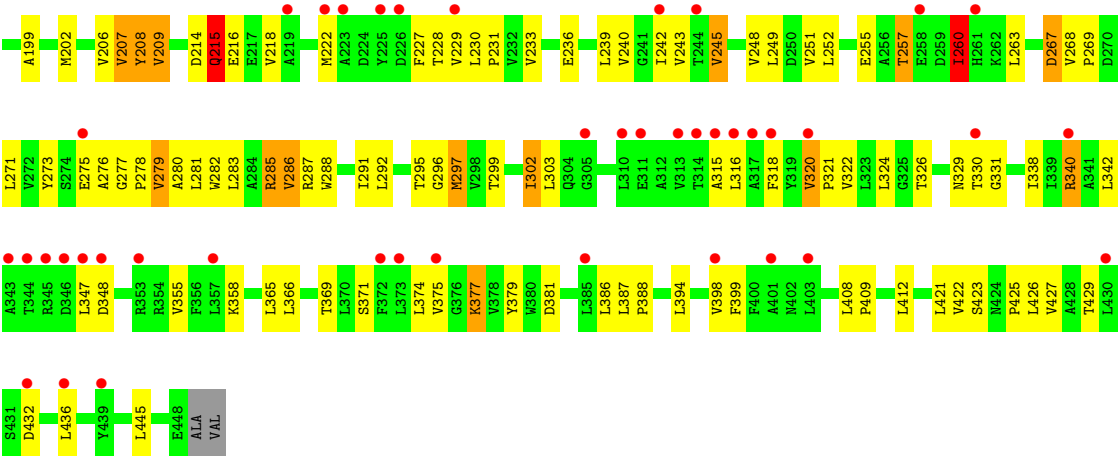
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Magnesium transporter MgtE



• Molecule 1: Magnesium transporter MgtE





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	135.33Å 85.65Å 156.56Å 90.00° 100.02° 90.00°	Depositor
Resolution (Å)	50.33 – 3.60 50.33 – 3.60	Depositor EDS
% Data completeness (in resolution range)	98.3 (50.33-3.60) 98.2 (50.33-3.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 3.57Å)	Xtriage
Refinement program	PHENIX (1.10_2155)	Depositor
R, R_{free}	0.250 , 0.283 0.257 , 0.290	Depositor DCC
R_{free} test set	2000 reflections (9.65%)	wwPDB-VP
Wilson B-factor (Å ²)	152.5	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 81.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	6522	wwPDB-VP
Average B, all atoms (Å ²)	169.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ATP, MG

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	0.64	1/3308 (0.0%)	1.01	7/4534 (0.2%)
1	B	0.61	1/3259 (0.0%)	1.04	16/4478 (0.4%)
All	All	0.63	2/6567 (0.0%)	1.02	23/9012 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	178	GLY	C-O	5.75	1.31	1.23
1	B	66	PRO	N-CD	5.09	1.54	1.47

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	277	GLY	CA-C-N	8.13	128.59	119.32
1	B	277	GLY	C-N-CA	8.13	128.59	119.32
1	B	162	ALA	CA-C-N	-7.54	112.03	119.87
1	B	162	ALA	C-N-CA	-7.54	112.03	119.87
1	B	52	LEU	CA-C-N	7.35	127.77	119.83
1	B	52	LEU	C-N-CA	7.35	127.77	119.83
1	A	407	MET	N-CA-C	7.33	119.35	111.36
1	A	104	ASP	CA-C-N	7.18	126.43	118.97
1	A	104	ASP	C-N-CA	7.18	126.43	118.97
1	A	308	SER	N-CA-C	6.59	118.26	111.14
1	B	182	GLY	N-CA-C	6.54	120.16	110.42
1	B	65	SER	CA-C-N	-5.97	112.42	119.05
1	B	65	SER	C-N-CA	-5.97	112.42	119.05
1	B	215	GLN	N-CA-C	-5.97	104.78	111.28
1	A	404	VAL	N-CA-C	-5.91	104.59	110.62
1	B	260	ILE	N-CA-C	5.64	116.42	110.72
1	B	267	ASP	N-CA-C	5.49	117.06	109.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	116	ASP	CA-C-N	5.33	126.50	119.84
1	A	116	ASP	C-N-CA	5.33	126.50	119.84
1	B	297	MET	N-CA-C	-5.31	105.49	111.28
1	B	116	ASP	CA-C-N	5.24	124.86	119.05
1	B	116	ASP	C-N-CA	5.24	124.86	119.05
1	B	28	ILE	N-CA-C	5.00	116.03	109.58

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	3248	0	3273	154	0
1	B	3199	0	3187	147	0
2	A	31	0	12	1	0
2	B	31	0	12	2	0
3	A	8	0	0	0	0
3	B	5	0	0	0	0
All	All	6522	0	6484	262	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 20.

All (262) 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:A:29:HIS:ND1	1:A:30:PRO:HD2	1.28	1.39
1:A:29:HIS:ND1	1:A:30:PRO:CD	1.96	1.26
1:B:132:ASP:HA	1:B:215:GLN:OE1	1.07	1.23
1:B:132:ASP:CA	1:B:215:GLN:OE1	1.96	1.11
1:B:216:GLU:HG2	1:B:255:GLU:HG2	1.40	1.03
1:B:132:ASP:O	1:B:215:GLN:CG	2.14	0.95
1:B:132:ASP:O	1:B:215:GLN:HG3	1.74	0.87
1:A:29:HIS:ND1	1:A:30:PRO:HD3	1.91	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:ASP:O	1:B:215:GLN:HG2	1.78	0.82
1:A:93:LEU:HD23	1:A:123:VAL:HG21	1.61	0.81
1:B:132:ASP:C	1:B:215:GLN:HG2	2.08	0.79
1:A:31:GLN:HG3	1:A:34:LEU:HD12	1.65	0.79
1:B:216:GLU:HG2	1:B:255:GLU:CG	2.13	0.79
1:A:400:PHE:O	1:A:404:VAL:HG23	1.83	0.77
1:A:29:HIS:CE1	1:A:30:PRO:CD	2.69	0.75
1:B:93:LEU:HD21	1:B:115:LEU:HD21	1.67	0.75
1:B:216:GLU:HG3	1:B:255:GLU:OE2	1.86	0.75
1:B:132:ASP:HA	1:B:215:GLN:CD	2.08	0.74
1:A:73:LEU:O	1:A:107:TYR:OH	2.05	0.74
1:B:147:GLU:HB2	1:B:199:ALA:HB2	1.70	0.74
1:B:324:LEU:HD12	1:B:432:ASP:HA	1.70	0.73
1:A:29:HIS:CE1	1:A:30:PRO:HD2	2.20	0.72
1:B:81:LEU:HD11	1:B:114:LEU:HD13	1.71	0.72
1:B:379:TYR:HB2	1:B:387:LEU:HD21	1.72	0.71
1:B:24:VAL:HG13	1:B:25:LEU:HD23	1.72	0.70
1:B:185:SER:HB3	2:B:501:ATP:H3'	1.73	0.69
1:B:216:GLU:CG	1:B:255:GLU:OE2	2.40	0.69
1:A:315:ALA:HA	1:A:318:PHE:CD2	2.29	0.68
1:A:332:ASN:O	1:A:336:THR:HG23	1.94	0.67
1:A:215:GLN:HE21	1:A:251:VAL:HG13	1.59	0.66
1:B:184:LEU:HD22	1:B:189:LEU:HD21	1.78	0.66
1:B:208:TYR:HD1	1:B:208:TYR:H	1.43	0.65
1:A:91:ASP:OD1	1:A:92:ASP:N	2.29	0.65
1:A:359:GLU:OE1	1:B:285:ARG:NH2	2.29	0.65
1:B:180:LEU:HD13	1:B:239:LEU:HD13	1.80	0.64
1:A:371:SER:HA	1:A:394:LEU:HD23	1.80	0.64
1:B:132:ASP:C	1:B:215:GLN:CG	2.69	0.63
1:A:31:GLN:HA	1:A:34:LEU:HB2	1.81	0.63
1:B:280:ALA:HA	1:B:283:LEU:HD12	1.78	0.63
1:A:78:PRO:HA	1:A:81:LEU:HB2	1.80	0.63
1:B:173:VAL:HG21	1:B:202:MET:SD	2.38	0.63
1:A:219:ALA:HB1	1:A:252:LEU:HD12	1.82	0.61
1:A:423:SER:HB2	1:A:425:PRO:HD2	1.83	0.61
1:B:209:VAL:HG21	1:B:218:VAL:HG22	1.82	0.61
1:A:378:VAL:HG11	1:A:387:LEU:HA	1.83	0.60
1:A:176:GLU:HG2	1:A:177:LYS:HG3	1.81	0.60
1:A:319:TYR:C	1:A:321:PRO:HD2	2.27	0.60
1:B:112:LYS:NZ	1:B:124:GLU:OE2	2.28	0.59
1:A:30:PRO:HD2	1:A:31:GLN:H	1.65	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:293:ILE:HD11	1:B:326:THR:HG23	1.83	0.59
1:A:336:THR:HB	1:A:340:ARG:HH22	1.68	0.59
1:A:30:PRO:CD	1:A:31:GLN:H	2.16	0.58
1:A:303:LEU:HD23	1:A:440:LEU:HD11	1.85	0.58
1:A:348:ASP:OD1	1:A:349:LEU:N	2.36	0.58
1:A:318:PHE:HA	1:B:303:LEU:HD13	1.85	0.58
1:A:377:LYS:HB3	1:B:297:MET:HE2	1.85	0.58
1:A:358:LYS:NZ	1:B:276:ALA:O	2.37	0.57
1:A:360:MET:HA	1:A:403:LEU:HD11	1.86	0.57
1:A:417:VAL:O	1:A:419:PRO:HD3	2.04	0.57
1:B:71:GLU:O	1:B:75:THR:OG1	2.23	0.57
1:A:54:LYS:HA	1:A:80:ARG:HH12	1.69	0.57
1:B:222:MET:HE2	1:B:230:LEU:HG	1.87	0.57
1:A:29:HIS:CG	1:A:30:PRO:CD	2.83	0.57
1:B:175:ASP:OD1	1:B:178:GLY:N	2.37	0.57
1:A:323:LEU:HD23	1:A:397:ILE:HD12	1.86	0.57
1:B:132:ASP:CB	1:B:215:GLN:OE1	2.53	0.56
1:A:407:MET:O	1:A:408:LEU:C	2.49	0.56
1:B:288:TRP:HZ3	1:B:426:LEU:HD12	1.70	0.56
1:A:322:VAL:HG22	1:B:296:GLY:C	2.30	0.56
1:A:134:ALA:O	1:A:135:GLY:C	2.47	0.56
1:A:320:VAL:N	1:A:321:PRO:CD	2.69	0.55
1:A:29:HIS:O	1:A:30:PRO:C	2.46	0.55
1:B:109:GLN:HA	1:B:112:LYS:HG2	1.87	0.55
1:B:66:PRO:O	1:B:69:GLN:HB2	2.07	0.55
1:B:142:TYR:HB3	1:B:242:ILE:HG12	1.89	0.55
1:B:145:VAL:HG22	1:B:149:MET:HG3	1.90	0.54
1:A:222:MET:HE3	1:A:230:LEU:HD22	1.90	0.54
1:B:408:LEU:N	1:B:409:PRO:HD2	2.22	0.54
1:A:309:VAL:HG12	1:A:447:LEU:HD12	1.90	0.54
1:A:340:ARG:CZ	1:B:260:ILE:HB	2.38	0.54
1:A:24:VAL:HG12	1:A:28:ILE:HG23	1.88	0.54
1:A:282:TRP:O	1:A:286:VAL:HB	2.08	0.54
1:B:81:LEU:O	1:B:85:LEU:HB2	2.07	0.54
1:B:145:VAL:O	1:B:173:VAL:HA	2.08	0.53
1:A:184:LEU:HD21	1:A:189:LEU:HD21	1.91	0.53
1:A:285:ARG:NH2	1:B:330:THR:O	2.41	0.53
1:B:371:SER:HA	1:B:394:LEU:HD23	1.90	0.53
1:B:214:ASP:O	1:B:215:GLN:C	2.49	0.52
1:A:132:ASP:O	1:A:215:GLN:HB2	2.09	0.52
1:B:37:TRP:CE2	1:B:45:ARG:HD2	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:PRO:CD	1:A:322:VAL:H	2.22	0.52
1:A:436:LEU:HD11	1:B:321:PRO:HG2	1.92	0.52
1:B:281:LEU:O	1:B:285:ARG:HG3	2.10	0.52
1:A:185:SER:HB2	2:A:501:ATP:H3'	1.91	0.52
1:A:319:TYR:OH	1:A:378:VAL:HA	2.10	0.52
1:B:279:VAL:O	1:B:282:TRP:HB3	2.09	0.52
1:A:55:ALA:HB2	1:A:87:GLU:OE2	2.10	0.51
1:A:179:ARG:HA	1:A:238:ARG:HA	1.93	0.51
1:B:216:GLU:CG	1:B:255:GLU:CD	2.84	0.51
1:A:142:TYR:CE2	1:A:239:LEU:HG	2.45	0.51
1:A:297:MET:HE2	1:B:374:LEU:HA	1.93	0.51
1:A:377:LYS:O	1:A:381:ASP:N	2.44	0.51
1:A:383:HIS:C	1:A:385:LEU:H	2.19	0.51
1:B:208:TYR:HA	1:B:231:PRO:HD2	1.92	0.51
1:A:140:PRO:HD2	1:A:141:GLU:OE1	2.10	0.51
1:B:387:LEU:N	1:B:388:PRO:HD2	2.26	0.51
1:A:292:LEU:HD23	1:B:329:ASN:ND2	2.25	0.51
1:B:315:ALA:HA	1:B:318:PHE:CD2	2.45	0.51
1:B:287:ARG:O	1:B:291:ILE:HG12	2.11	0.51
1:B:296:GLY:O	1:B:297:MET:C	2.53	0.50
1:A:77:PRO:HG2	1:A:80:ARG:HB2	1.94	0.50
1:A:321:PRO:CG	1:A:322:VAL:N	2.73	0.50
1:A:250:ASP:OD1	1:A:250:ASP:N	2.40	0.50
1:A:333:GLN:OE1	1:B:285:ARG:HD2	2.11	0.50
1:B:145:VAL:HG21	1:B:154:VAL:HG22	1.94	0.50
1:B:222:MET:SD	1:B:245:VAL:HG23	2.52	0.50
1:A:288:TRP:HA	1:A:291:ILE:HD12	1.94	0.50
1:B:222:MET:HE1	1:B:243:VAL:HG12	1.95	0.49
1:A:249:LEU:HD22	1:B:249:LEU:HD13	1.95	0.49
1:A:79:TRP:CG	1:A:80:ARG:N	2.81	0.49
1:A:142:TYR:HE2	1:A:239:LEU:HG	1.76	0.49
1:A:276:ALA:O	1:B:358:LYS:NZ	2.24	0.49
1:B:47:VAL:HA	1:B:50:THR:HG22	1.95	0.49
1:B:134:ALA:HA	1:B:137:LEU:HD11	1.94	0.49
1:B:282:TRP:CE2	1:B:286:VAL:HG21	2.48	0.49
1:A:29:HIS:CE1	1:A:30:PRO:HD3	2.41	0.48
1:A:77:PRO:HB2	1:A:79:TRP:HD1	1.78	0.48
1:A:167:THR:HG21	1:A:229:VAL:HG21	1.94	0.48
1:A:359:GLU:HG3	1:A:402:ASN:HD21	1.78	0.48
1:B:74:LYS:HA	1:B:107:TYR:CE2	2.48	0.48
1:A:400:PHE:O	1:A:404:VAL:CG2	2.58	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:184:LEU:HD22	1:B:189:LEU:CD2	2.44	0.48
1:B:216:GLU:HG2	1:B:255:GLU:CD	2.39	0.48
1:A:336:THR:HB	1:A:340:ARG:NH2	2.28	0.48
1:A:176:GLU:CG	1:A:177:LYS:HG3	2.44	0.48
1:A:289:LEU:HD12	1:B:366:LEU:HD22	1.95	0.47
1:B:183:VAL:HG12	1:B:206:VAL:HB	1.96	0.47
1:B:193:ASP:O	1:B:196:THR:OG1	2.25	0.47
1:A:365:LEU:HB2	1:B:282:TRP:CD1	2.49	0.47
1:A:30:PRO:CD	1:A:31:GLN:N	2.77	0.47
1:A:193:ASP:HB3	1:A:196:THR:HG23	1.96	0.47
1:A:296:GLY:O	1:B:322:VAL:HG22	2.15	0.47
1:A:282:TRP:HZ2	1:B:369:THR:HG21	1.78	0.47
1:A:321:PRO:CG	1:A:322:VAL:H	2.27	0.47
1:A:322:VAL:HG23	1:B:299:THR:OG1	2.14	0.47
1:A:370:LEU:HD12	1:A:398:VAL:HG21	1.96	0.47
1:A:52:LEU:O	1:A:80:ARG:NH2	2.46	0.47
1:A:273:TYR:CD2	1:B:347:LEU:HD11	2.50	0.47
1:B:112:LYS:NZ	1:B:120:ARG:HH21	2.13	0.46
1:B:112:LYS:HZ2	1:B:120:ARG:HE	1.63	0.46
1:B:296:GLY:O	1:B:299:THR:N	2.48	0.46
1:A:187:ARG:HB2	1:B:168:ILE:HD11	1.96	0.46
1:B:377:LYS:HG3	1:B:381:ASP:HB2	1.97	0.46
1:A:303:LEU:HD13	1:B:318:PHE:HA	1.97	0.46
1:A:319:TYR:OH	1:A:381:ASP:OD2	2.24	0.46
1:A:44:HIS:O	1:A:48:VAL:HG23	2.15	0.46
1:A:115:LEU:HD12	1:A:115:LEU:HA	1.78	0.46
1:B:257:THR:HA	1:B:260:ILE:HG13	1.98	0.46
1:B:216:GLU:HG2	1:B:255:GLU:OE2	2.16	0.46
1:A:169:TYR:HB3	1:B:169:TYR:HE2	1.81	0.46
1:A:383:HIS:O	1:A:385:LEU:N	2.49	0.46
1:A:424:ASN:HB2	1:A:425:PRO:HD3	1.98	0.46
1:A:138:MET:SD	1:A:232:VAL:HG11	2.56	0.46
1:A:219:ALA:HB1	1:A:252:LEU:CD1	2.46	0.45
1:B:118:ARG:O	1:B:122:GLU:HG2	2.16	0.45
1:A:134:ALA:O	1:A:137:LEU:N	2.43	0.45
1:A:142:TYR:HD1	1:A:142:TYR:H	1.62	0.45
1:A:321:PRO:HG2	1:A:322:VAL:N	2.30	0.45
1:B:320:VAL:HG13	1:B:324:LEU:HD23	1.98	0.45
1:A:71:GLU:O	1:A:75:THR:OG1	2.33	0.45
1:A:90:LEU:HD13	1:A:141:GLU:OE1	2.17	0.45
1:B:208:TYR:CD2	1:B:233:VAL:HG21	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:423:SER:HB2	1:B:425:PRO:HD2	1.99	0.45
1:A:142:TYR:N	1:A:142:TYR:CD1	2.81	0.45
1:A:162:ALA:HB1	1:B:187:ARG:NH1	2.32	0.45
1:A:183:VAL:HB	1:A:206:VAL:HG13	1.99	0.45
1:A:321:PRO:HD2	1:A:322:VAL:H	1.81	0.45
1:B:86:GLU:C	1:B:88:LEU:H	2.24	0.45
1:B:145:VAL:HG22	1:B:149:MET:HE3	1.99	0.45
1:B:282:TRP:CZ2	1:B:286:VAL:HG21	2.52	0.45
1:B:37:TRP:CZ2	1:B:45:ARG:HD2	2.52	0.45
1:A:322:VAL:CG2	1:B:296:GLY:O	2.65	0.45
1:A:340:ARG:CD	1:B:260:ILE:HD12	2.46	0.45
1:B:30:PRO:HG3	1:B:56:LYS:HD2	1.99	0.45
1:B:91:ASP:N	1:B:91:ASP:OD1	2.48	0.45
1:B:112:LYS:HZ2	1:B:120:ARG:HH21	1.64	0.45
1:A:180:LEU:HD11	1:A:231:PRO:HB2	1.99	0.44
1:B:170:TYR:HE2	2:B:501:ATP:H8	1.64	0.44
1:A:190:ILE:O	1:B:159:ARG:HG2	2.17	0.44
1:B:207:VAL:CG1	1:B:227:PHE:HE2	2.31	0.44
1:A:221:LEU:HD23	1:A:221:LEU:HA	1.85	0.44
1:A:321:PRO:HG2	1:A:322:VAL:H	1.82	0.44
1:A:301:SER:HB2	1:B:377:LYS:NZ	2.33	0.44
1:B:222:MET:HE3	1:B:222:MET:HB2	1.71	0.44
1:B:331:GLY:C	1:B:427:VAL:HG11	2.42	0.44
1:A:334:SER:OG	1:A:405:GLY:HA3	2.18	0.44
1:A:340:ARG:HD2	1:B:260:ILE:HD12	1.99	0.44
1:B:199:ALA:HA	1:B:202:MET:HG2	1.99	0.44
1:A:142:TYR:HD1	1:A:142:TYR:N	2.16	0.44
1:B:175:ASP:OD1	1:B:179:ARG:N	2.50	0.43
1:A:151:VAL:HB	1:A:194:PRO:HA	2.01	0.43
1:A:446:LEU:HD23	1:A:446:LEU:HA	1.86	0.43
1:B:278:PRO:O	1:B:282:TRP:HB2	2.18	0.43
1:A:383:HIS:HB3	1:A:386:LEU:CD2	2.49	0.43
1:B:134:ALA:O	1:B:137:LEU:HD12	2.18	0.43
1:B:375:VAL:HB	1:B:387:LEU:HD23	1.99	0.43
1:A:273:TYR:HE2	1:B:355:VAL:HG22	1.82	0.43
1:A:338:ILE:HD11	1:B:273:TYR:CZ	2.54	0.43
1:B:338:ILE:O	1:B:342:LEU:HG	2.18	0.43
1:A:333:GLN:OE1	1:B:285:ARG:HB3	2.19	0.42
1:B:77:PRO:HB2	1:B:79:TRP:NE1	2.35	0.42
1:B:98:GLN:NE2	1:B:127:ALA:HA	2.34	0.42
1:B:436:LEU:HD23	1:B:436:LEU:HA	1.80	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:LEU:HD12	1:A:64:LEU:HA	1.81	0.42
1:A:402:ASN:ND2	1:A:403:LEU:HD12	2.35	0.42
1:B:228:THR:HG23	1:B:229:VAL:HG23	2.01	0.42
1:A:222:MET:CE	1:A:230:LEU:HB2	2.50	0.42
1:A:401:ALA:HA	1:A:404:VAL:HG23	2.02	0.42
1:B:445:LEU:HD23	1:B:445:LEU:HA	1.81	0.42
1:A:187:ARG:NH2	1:B:162:ALA:O	2.49	0.42
1:B:107:TYR:O	1:B:111:LEU:HB2	2.19	0.42
1:A:321:PRO:CD	1:A:322:VAL:N	2.83	0.42
1:A:374:LEU:O	1:A:378:VAL:HG23	2.20	0.42
1:B:299:THR:O	1:B:302:ILE:HG22	2.19	0.42
1:A:340:ARG:N	1:A:340:ARG:HD3	2.32	0.42
1:A:209:VAL:HG11	1:A:218:VAL:HG22	2.02	0.42
1:B:216:GLU:CG	1:B:255:GLU:CG	2.92	0.42
1:A:49:LEU:HD23	1:A:49:LEU:HA	1.93	0.41
1:B:132:ASP:CA	1:B:215:GLN:CD	2.82	0.41
1:B:182:GLY:HA3	1:B:202:MET:HE3	2.01	0.41
1:B:295:THR:HB	1:B:429:THR:HG21	2.02	0.41
1:A:303:LEU:HD23	1:A:303:LEU:HA	1.83	0.41
1:A:408:LEU:O	1:A:412:LEU:HG	2.20	0.41
1:A:172:TYR:OH	1:A:229:VAL:HG12	2.20	0.41
1:A:289:LEU:O	1:A:293:ILE:HD13	2.20	0.41
1:B:320:VAL:O	1:B:324:LEU:HD23	2.21	0.41
1:A:186:LEU:HD21	1:B:190:ILE:HD12	2.00	0.41
1:A:214:ASP:OD1	1:A:215:GLN:N	2.53	0.41
1:B:98:GLN:HE21	1:B:127:ALA:HA	1.85	0.41
1:A:185:SER:HB3	1:A:188:ASP:CG	2.46	0.41
1:A:273:TYR:CZ	1:B:338:ILE:HD11	2.55	0.41
1:A:425:PRO:HB3	1:B:329:ASN:CG	2.46	0.41
1:B:157:PHE:HA	1:B:160:ARG:HG2	2.01	0.41
1:B:230:LEU:HA	1:B:231:PRO:HD3	1.74	0.41
1:B:46:TYR:CD1	1:B:46:TYR:C	2.99	0.41
1:B:412:LEU:HD21	1:B:421:LEU:HB2	2.02	0.41
1:A:79:TRP:CD1	1:A:80:ARG:H	2.39	0.41
1:A:296:GLY:O	1:A:299:THR:HG23	2.20	0.41
1:A:302:ILE:HD11	1:A:437:LEU:HD21	2.03	0.41
1:A:340:ARG:NE	1:B:260:ILE:HB	2.36	0.41
1:A:180:LEU:HA	1:A:180:LEU:HD23	1.75	0.41
1:A:222:MET:HE1	1:A:230:LEU:HB2	2.02	0.41
1:A:249:LEU:CD2	1:B:249:LEU:HD13	2.50	0.41
1:B:135:GLY:HA2	1:B:138:MET:SD	2.61	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:LEU:N	1:A:409:PRO:HD2	2.37	0.40
1:A:149:MET:HE2	1:A:154:VAL:HG22	2.02	0.40
1:A:401:ALA:CA	1:A:404:VAL:HG23	2.51	0.40
1:A:179:ARG:CA	1:A:238:ARG:HA	2.50	0.40
1:B:145:VAL:CG2	1:B:149:MET:HE3	2.52	0.40
1:A:257:THR:HG23	1:B:340:ARG:HG3	2.03	0.40
1:B:130:GLU:HB2	1:B:133:GLU:HG3	2.04	0.40
1:B:263:LEU:HD23	1:B:263:LEU:HA	1.83	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	425/473 (90%)	392 (92%)	30 (7%)	3 (1%)	18	51
1	B	424/473 (90%)	395 (93%)	27 (6%)	2 (0%)	24	57
All	All	849/946 (90%)	787 (93%)	57 (7%)	5 (1%)	21	54

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	269	PRO
1	B	422	VAL
1	A	422	VAL
1	A	384	PRO
1	A	390	VAL

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	340/400 (85%)	283 (83%)	57 (17%)	2	13
1	B	333/400 (83%)	285 (86%)	48 (14%)	3	18
All	All	673/800 (84%)	568 (84%)	105 (16%)	2	16

All (105) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	40	LEU
1	A	45	ARG
1	A	64	LEU
1	A	75	THR
1	A	80	ARG
1	A	81	LEU
1	A	90	LEU
1	A	103	GLU
1	A	111	LEU
1	A	112	LYS
1	A	115	LEU
1	A	116	ASP
1	A	142	TYR
1	A	145	VAL
1	A	149	MET
1	A	174	VAL
1	A	179	ARG
1	A	187	ARG
1	A	206	VAL
1	A	210	ARG
1	A	222	MET
1	A	230	LEU
1	A	242	ILE
1	A	244	THR
1	A	248	VAL
1	A	249	LEU

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Mol	Chain	Res	Type
1	A	250	ASP
1	A	252	LEU
1	A	255	GLU
1	A	260	ILE
1	A	272	VAL
1	A	281	LEU
1	A	286	VAL
1	A	292	LEU
1	A	293	ILE
1	A	299	THR
1	A	302	ILE
1	A	308	SER
1	A	313	VAL
1	A	340	ARG
1	A	345	ARG
1	A	357	LEU
1	A	364	LEU
1	A	375	VAL
1	A	392	VAL
1	A	400	PHE
1	A	404	VAL
1	A	407	MET
1	A	421	LEU
1	A	429	THR
1	A	430	LEU
1	A	438	ILE
1	A	441	SER
1	A	445	LEU
1	A	447	LEU
1	A	448	GLU
1	B	25	LEU
1	B	50	THR
1	B	62	SER
1	B	67	GLU
1	B	75	THR
1	B	81	LEU
1	B	82	ARG
1	B	85	LEU
1	B	111	LEU
1	B	115	LEU
1	B	123	VAL
1	B	130	GLU

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Mol	Chain	Res	Type
1	B	131	GLU
1	B	137	LEU
1	B	152	GLU
1	B	160	ARG
1	B	171	ILE
1	B	189	LEU
1	B	207	VAL
1	B	208	TYR
1	B	209	VAL
1	B	215	GLN
1	B	236	GLU
1	B	240	VAL
1	B	245	VAL
1	B	248	VAL
1	B	251	VAL
1	B	252	LEU
1	B	257	THR
1	B	260	ILE
1	B	267	ASP
1	B	268	VAL
1	B	271	LEU
1	B	275	GLU
1	B	279	VAL
1	B	285	ARG
1	B	286	VAL
1	B	292	LEU
1	B	302	ILE
1	B	316	LEU
1	B	320	VAL
1	B	340	ARG
1	B	348	ASP
1	B	365	LEU
1	B	377	LYS
1	B	386	LEU
1	B	398	VAL
1	B	399	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 13 are monoatomic - leaving 2 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	$\# Z > 2$
2	ATP	A	501	-	32,33,33	1.49	7 (21%)	48,52,52	1.73	10 (20%)
2	ATP	B	501	-	32,33,33	1.56	7 (21%)	48,52,52	1.78	9 (18%)

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
2	ATP	A	501	-	-	0/22/38/38	0/3/3/3
2	ATP	B	501	-	-	1/22/38/38	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	ATP	C5-C4	5.05	1.48	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	ATP	C5-C4	4.51	1.47	1.39
2	B	501	ATP	PB-O3B	3.29	1.63	1.59
2	A	501	ATP	PB-O3B	3.13	1.62	1.59
2	B	501	ATP	C8-N7	2.53	1.36	1.31
2	A	501	ATP	C8-N7	2.52	1.36	1.31
2	A	501	ATP	C4-N9	-2.49	1.32	1.37
2	B	501	ATP	C5-C6	2.47	1.47	1.41
2	B	501	ATP	C5-N7	-2.43	1.34	1.39
2	A	501	ATP	C5-N7	-2.27	1.34	1.39
2	B	501	ATP	PA-O3A	2.26	1.61	1.59
2	A	501	ATP	PB-O3A	2.24	1.61	1.59
2	B	501	ATP	PB-O3A	2.22	1.61	1.59
2	A	501	ATP	C5-C6	2.03	1.46	1.41

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	ATP	C5-C4-N3	-5.61	118.99	126.72
2	B	501	ATP	N3-C4-N9	4.65	135.07	127.17
2	A	501	ATP	C5-C4-N3	-4.45	120.59	126.72
2	A	501	ATP	N3-C4-N9	3.97	133.91	127.17
2	B	501	ATP	C2-N3-C4	3.53	120.44	111.83
2	A	501	ATP	N3-C2-N1	-3.52	123.25	128.58
2	A	501	ATP	C2-N3-C4	3.31	119.92	111.83
2	B	501	ATP	N3-C2-N1	-3.31	123.57	128.58
2	B	501	ATP	C2'-C1'-N9	-3.27	105.18	113.30
2	A	501	ATP	C4-N9-C8	2.89	108.77	105.74
2	B	501	ATP	C4-C5-N7	-2.71	107.49	110.58
2	A	501	ATP	O5'-C5'-C4'	-2.53	100.37	108.99
2	B	501	ATP	C4-N9-C8	2.28	108.13	105.74
2	A	501	ATP	N6-C6-N1	2.22	123.33	118.38
2	A	501	ATP	O2A-PA-O1A	2.13	122.36	112.44
2	B	501	ATP	C2-N1-C6	2.12	122.21	118.73
2	A	501	ATP	C5-C6-N6	-2.10	118.08	123.29
2	B	501	ATP	C5'-C4'-C3'	-2.10	107.65	115.21
2	A	501	ATP	O3G-PG-O2G	2.02	115.38	107.80

There are no chirality outliers.

All (1) torsion outliers are listed below:

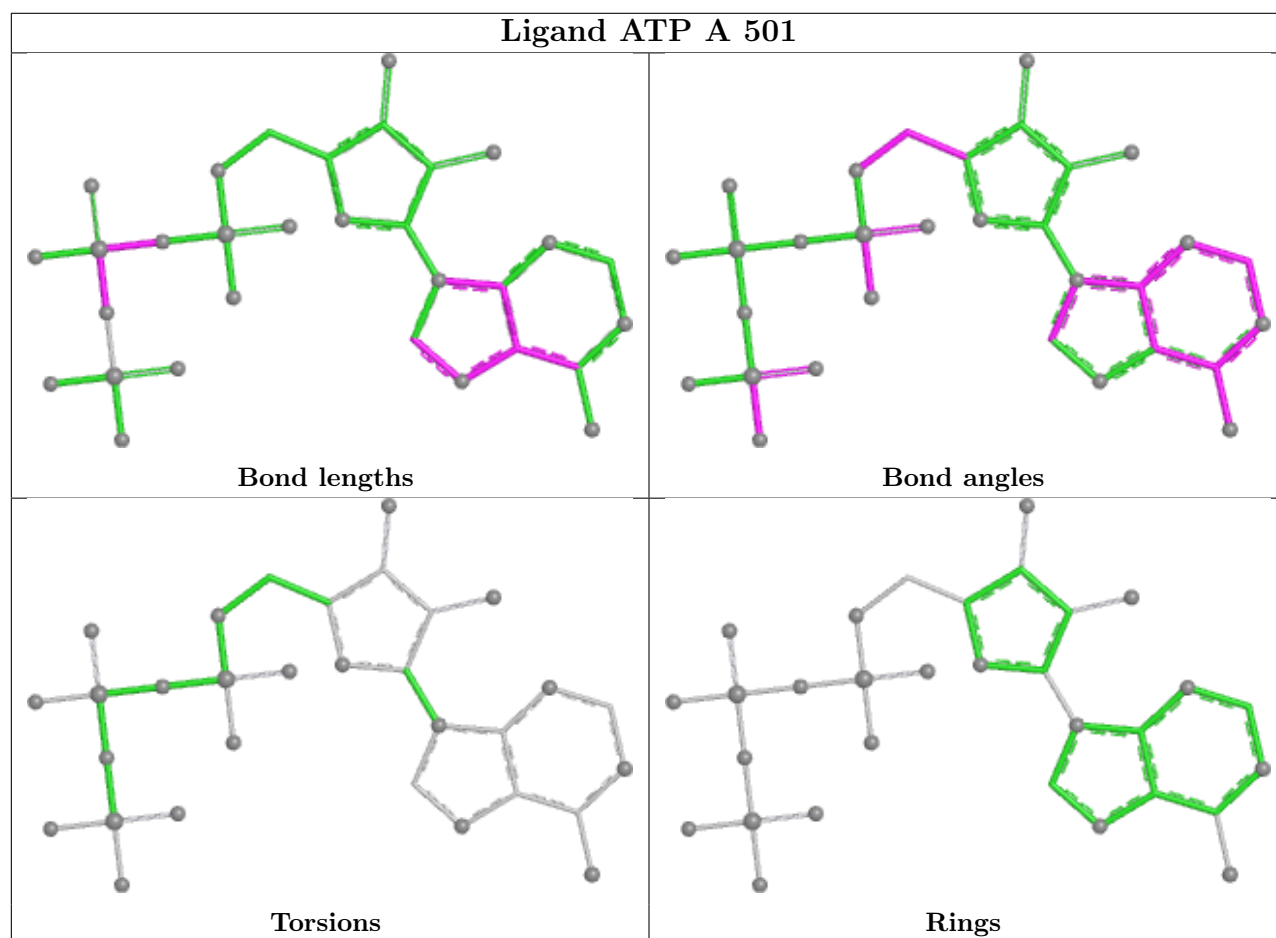
Mol	Chain	Res	Type	Atoms
2	B	501	ATP	O4'-C4'-C5'-O5'

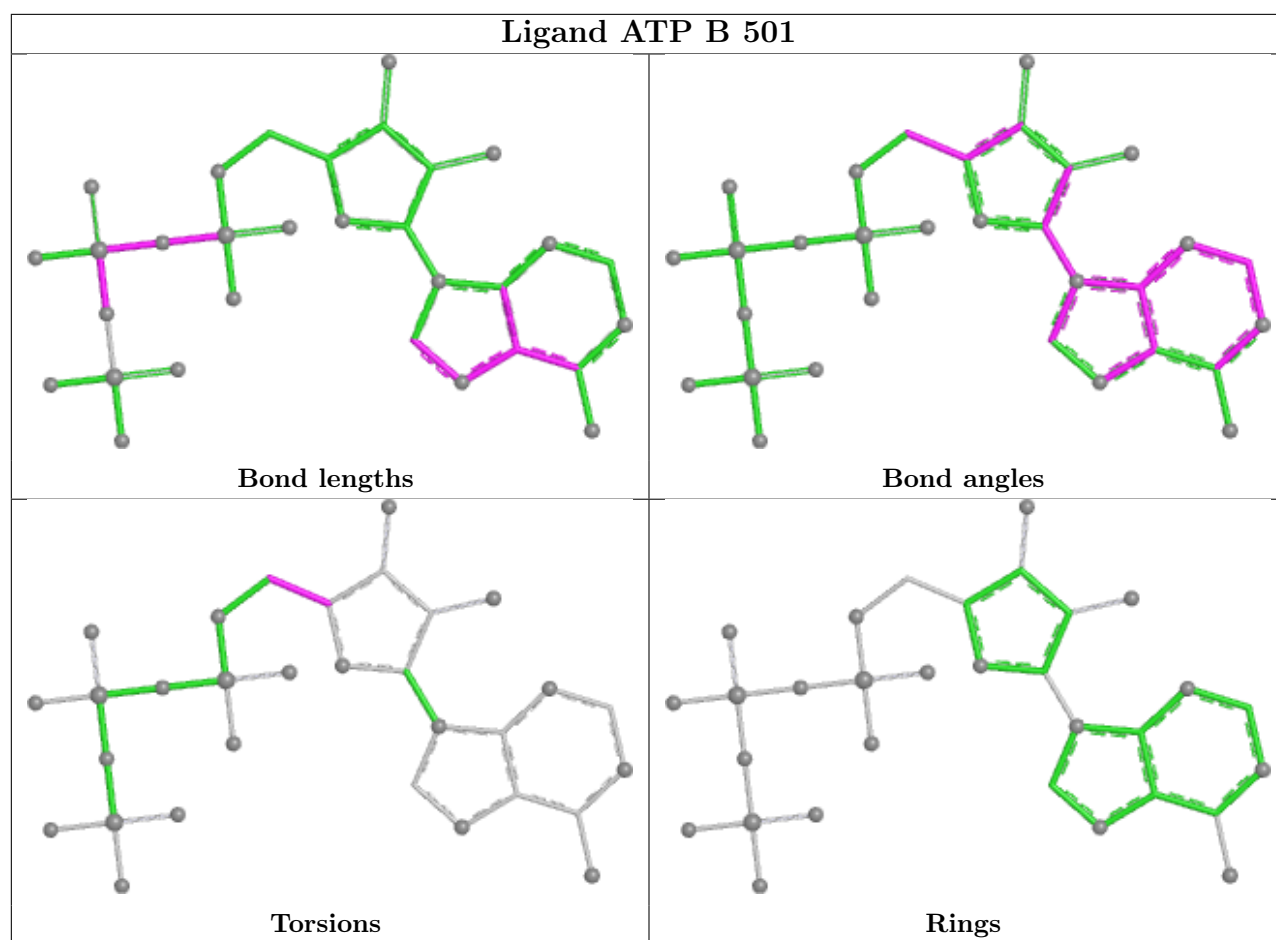
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	ATP	1	0
2	B	501	ATP	2	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	427/473 (90%)	0.63	47 (11%) 10 9	110, 165, 220, 264	0
1	B	426/473 (90%)	0.69	54 (12%) 8 7	115, 165, 235, 283	0
All	All	853/946 (90%)	0.66	101 (11%) 9 8	110, 165, 227, 283	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	267	ASP	7.6
1	A	59	GLU	6.7
1	B	317	ALA	6.7
1	A	63	HIS	6.7
1	A	421	LEU	5.8
1	B	316	LEU	5.8
1	A	117	PRO	5.5
1	B	222	MET	5.4
1	B	347	LEU	5.3
1	A	268	VAL	4.9
1	A	270	ASP	4.8
1	A	271	LEU	4.8
1	B	315	ALA	4.7
1	A	447	LEU	4.6
1	A	385	LEU	4.4
1	B	345	ARG	4.4
1	B	25	LEU	4.3
1	A	62	SER	4.2
1	B	344	THR	4.2
1	B	164	ASP	4.1
1	B	310	LEU	4.1
1	A	340	ARG	4.0
1	B	223	ALA	4.0
1	A	141	GLU	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	253	GLU	3.9
1	B	28	ILE	3.8
1	A	198	VAL	3.8
1	B	372	PHE	3.6
1	A	269	PRO	3.5
1	B	436	LEU	3.5
1	B	125	ALA	3.5
1	A	317	ALA	3.4
1	B	348	ASP	3.3
1	B	163	PRO	3.3
1	A	283	LEU	3.3
1	B	226	ASP	3.2
1	B	244	THR	3.1
1	B	242	ILE	3.1
1	B	51	LEU	3.1
1	A	288	TRP	3.1
1	B	261	HIS	3.1
1	B	375	VAL	3.0
1	B	258	GLU	3.0
1	A	121	ALA	3.0
1	B	343	ALA	3.0
1	A	446	LEU	3.0
1	A	313	VAL	2.9
1	A	77	PRO	2.9
1	B	346	ASP	2.9
1	B	140	PRO	2.9
1	B	403	LEU	2.8
1	B	373	LEU	2.8
1	B	176	GLU	2.8
1	A	122	GLU	2.7
1	B	353	ARG	2.7
1	A	258	GLU	2.7
1	A	177	LYS	2.7
1	A	88	LEU	2.7
1	A	420	ALA	2.7
1	B	219	ALA	2.6
1	B	311	GLU	2.6
1	B	313	VAL	2.6
1	B	340	ARG	2.6
1	B	330	THR	2.6
1	A	293	ILE	2.5
1	B	229	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	314	THR	2.5
1	A	272	VAL	2.5
1	B	139	THR	2.5
1	B	385	LEU	2.5
1	B	225	TYR	2.5
1	A	262	LYS	2.5
1	B	320	VAL	2.5
1	A	106	ALA	2.4
1	B	136	GLY	2.4
1	B	432	ASP	2.4
1	B	318	PHE	2.4
1	B	439	TYR	2.4
1	B	275	GLU	2.3
1	A	179	ARG	2.3
1	A	71	GLU	2.2
1	B	23	GLU	2.2
1	A	294	LEU	2.2
1	A	119	THR	2.2
1	A	89	SER	2.2
1	A	118	ARG	2.2
1	B	430	LEU	2.2
1	A	311	GLU	2.1
1	A	409	PRO	2.1
1	A	137	LEU	2.1
1	A	125	ALA	2.1
1	A	87	GLU	2.1
1	A	417	VAL	2.1
1	B	398	VAL	2.1
1	B	305	GLY	2.1
1	B	401	ALA	2.1
1	B	357	LEU	2.0
1	A	47	VAL	2.0
1	B	43	GLU	2.0
1	A	310	LEU	2.0
1	A	144	ALA	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

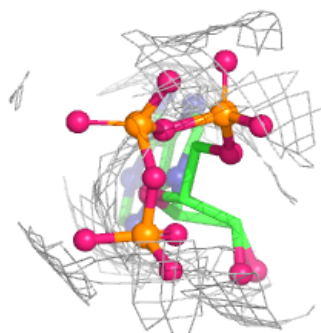
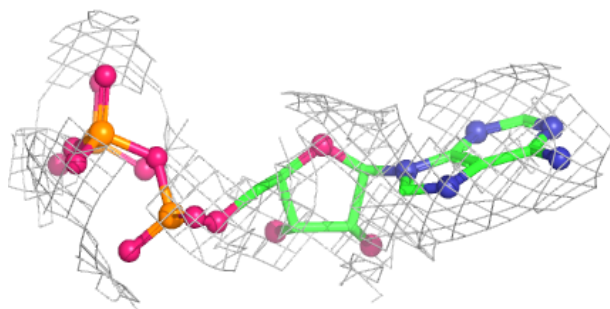
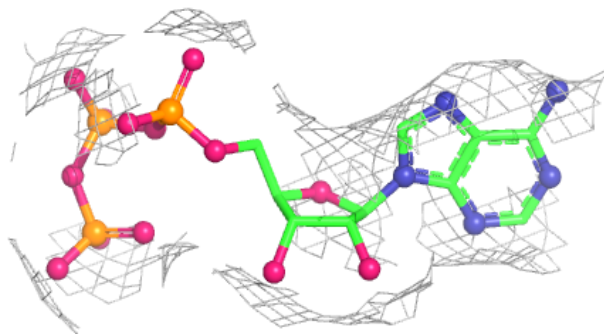
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MG	A	502	1/1	0.35	0.59	157,157,157,157	0
3	MG	B	503	1/1	0.49	0.43	137,137,137,137	0
3	MG	B	504	1/1	0.60	0.33	150,150,150,150	0
3	MG	A	504	1/1	0.66	0.35	136,136,136,136	0
3	MG	B	505	1/1	0.74	0.33	100,100,100,100	0
3	MG	A	505	1/1	0.78	0.32	117,117,117,117	0
3	MG	A	503	1/1	0.80	0.47	129,129,129,129	0
3	MG	A	506	1/1	0.82	0.36	104,104,104,104	0
3	MG	A	507	1/1	0.83	0.26	131,131,131,131	0
3	MG	A	508	1/1	0.87	0.27	143,143,143,143	0
3	MG	B	506	1/1	0.89	0.18	111,111,111,111	0
2	ATP	B	501	31/31	0.90	0.10	140,155,218,226	0
2	ATP	A	501	31/31	0.91	0.09	107,127,203,214	0
3	MG	A	509	1/1	0.91	0.16	106,106,106,106	0
3	MG	B	502	1/1	0.93	0.16	133,133,133,133	0

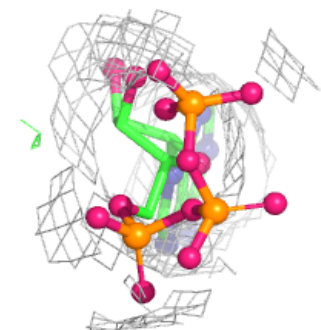
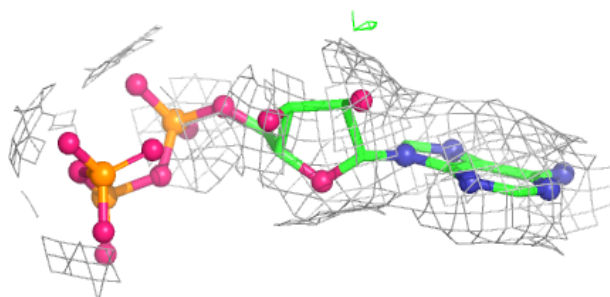
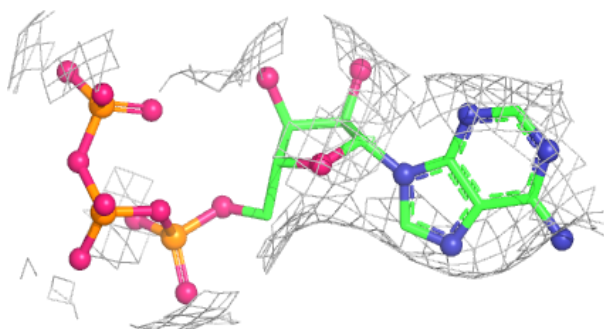
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ATP B 501:

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

**Electron density around ATP A 501:**

$2mF_o-DF_c$ (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.