



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 06:32 am GMT

PDB ID : 2DPY
Title : Crystal structure of the flagellar type III ATPase FliI
Authors : Imada, K.; Namba, K.; Minamino, T.
Deposited on : 2006-05-18
Resolution : 2.40 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

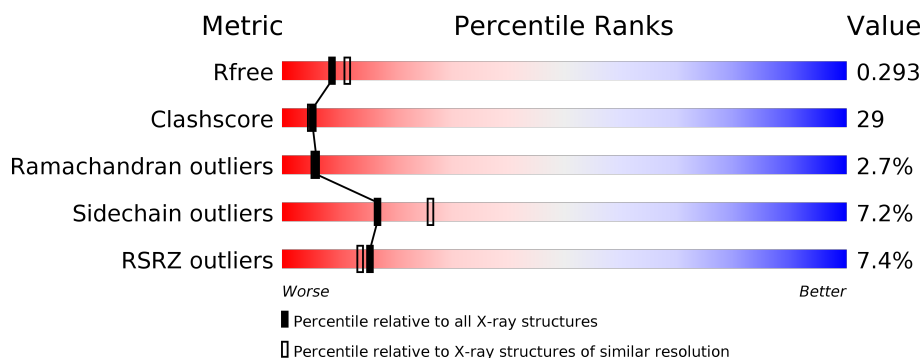
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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. 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	438	
1	B	438	

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
2	ADP	B	600	-	-	-	X

2 Entry composition [i](#)

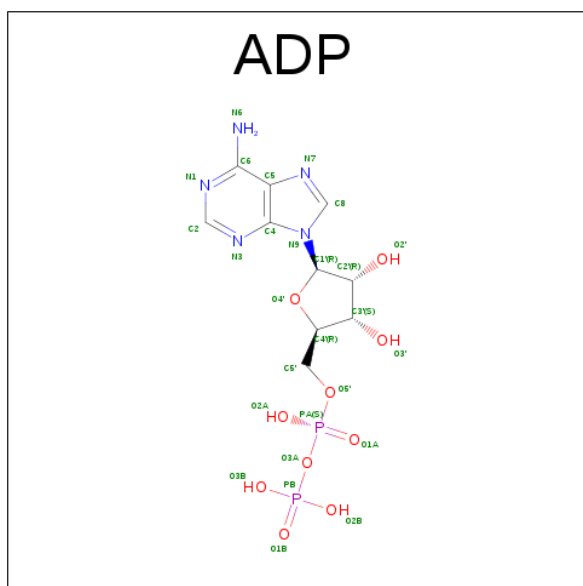
There are 3 unique types of molecules in this entry. The entry contains 6604 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Flagellum-specific ATP synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	0	0	0
			3211	2032	575	594	10			
1	B	426	Total	C	N	O	S	0	0	0
			3233	2045	579	599	10			

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



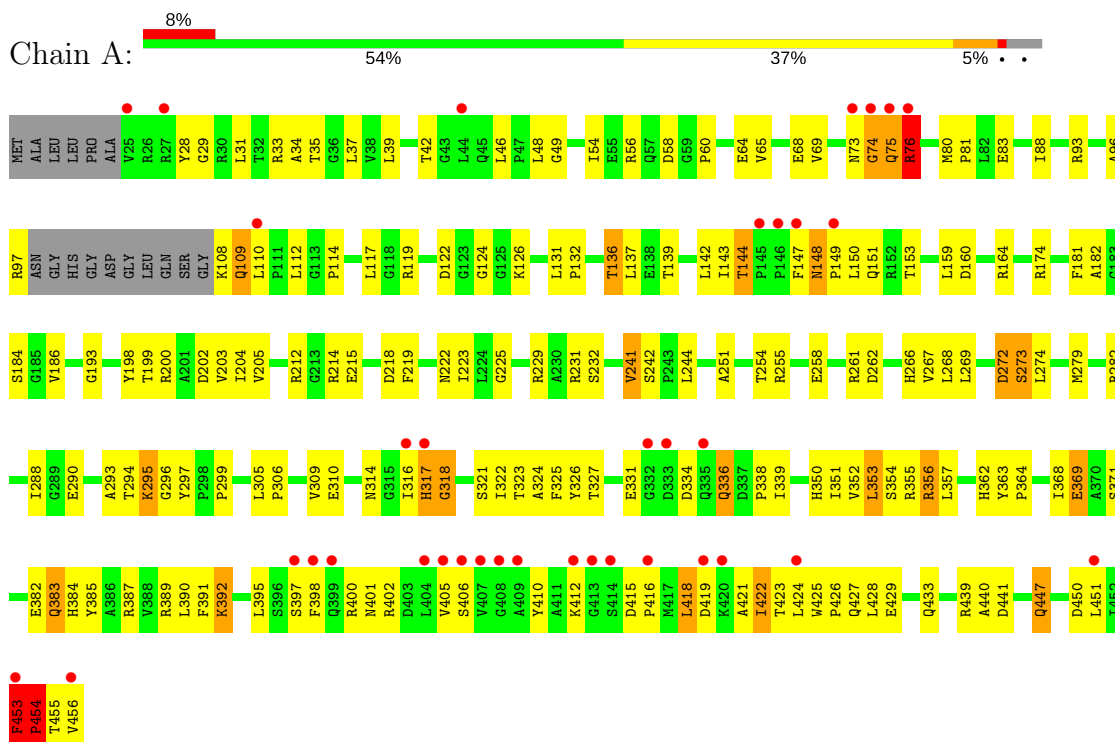
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	62	Total	O	0	0
			62	62		

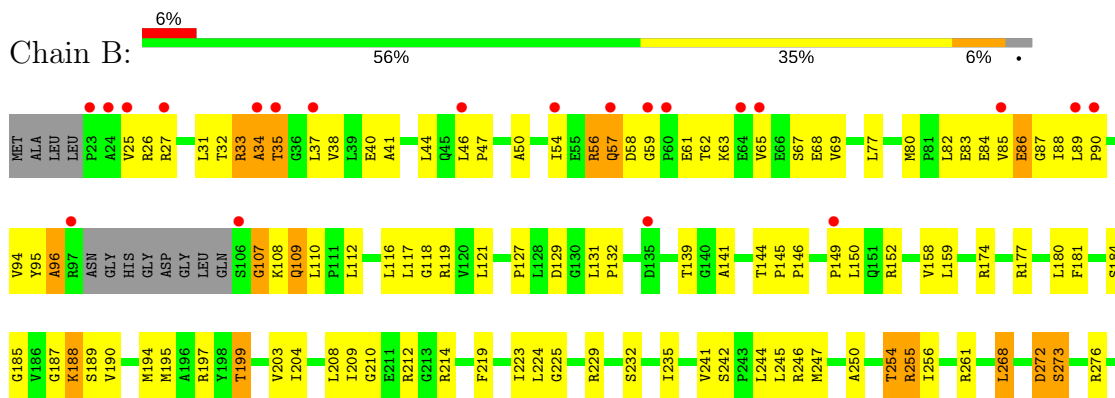
3 Residue-property plots

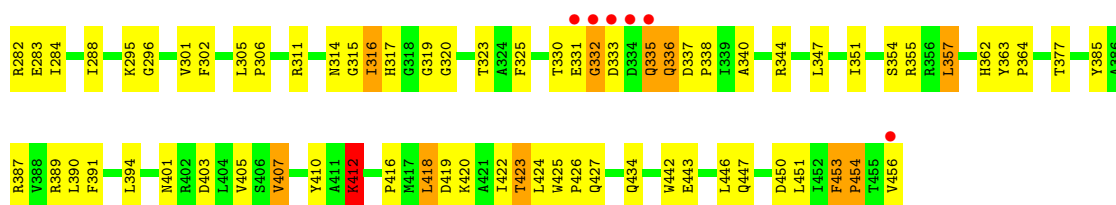
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Flagellum-specific ATP synthase



• Molecule 1: Flagellum-specific ATP synthase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.16Å 72.75Å 125.74Å 90.00° 94.13° 90.00°	Depositor
Resolution (Å)	45.98 – 2.40 45.98 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.4 (45.98-2.40) 99.6 (45.98-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.85 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.244 , 0.295 0.243 , 0.293	Depositor DCC
R_{free} test set	1680 reflections (4.95%)	DCC
Wilson B-factor (Å ²)	43.1	Xtriage
Anisotropy	0.407	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6604	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.88% 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: ADP

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	0.40	0/3268	0.71	3/4431 (0.1%)
1	B	0.39	0/3291	0.68	1/4462 (0.0%)
All	All	0.39	0/6559	0.69	4/8893 (0.0%)

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
1	A	0	2
1	B	0	2
All	All	0	4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	273	SER	N-CA-CB	9.16	124.24	110.50
1	A	318	GLY	N-CA-C	-9.03	90.53	113.10
1	B	273	SER	N-CA-CB	7.54	121.81	110.50
1	A	29	GLY	N-CA-C	-5.93	98.28	113.10

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	272	ASP	Mainchain,Peptide
1	B	272	ASP	Mainchain,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	3211	0	3277	186	0
1	B	3233	0	3298	191	0
2	B	27	0	12	8	0
3	A	71	0	0	3	0
3	B	62	0	0	16	0
All	All	6604	0	6587	376	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 29.

All (376) 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:412:LYS:HD3	1:B:412:LYS:H	1.30	0.97
1:B:450:ASP:HA	1:B:454:PRO:HG3	1.46	0.97
1:A:356:ARG:HD3	1:A:356:ARG:H	1.27	0.96
1:B:54:ILE:HG23	1:B:65:VAL:HB	1.50	0.92
1:A:148:ASN:ND2	1:A:151:GLN:HE21	1.67	0.91
1:A:261:ARG:HH11	1:A:261:ARG:HG2	1.35	0.90
1:B:38:VAL:HG21	1:B:288:ILE:HD11	1.57	0.87
1:A:148:ASN:HD21	1:A:151:GLN:HE21	1.22	0.86
1:B:109:GLN:HG3	1:B:139:THR:HG22	1.55	0.86
1:A:33:ARG:HH11	1:A:35:THR:HG21	1.45	0.82
1:A:142:LEU:HD21	1:A:251:ALA:HB1	1.59	0.82
1:B:450:ASP:HA	1:B:454:PRO:CG	2.10	0.81
1:A:425:TRP:CE3	1:A:428:LEU:HD11	2.16	0.81
1:B:158:VAL:HG12	3:B:656:HOH:O	1.80	0.79
1:B:209:ILE:HG22	1:B:276:ARG:HB3	1.64	0.79
1:B:295:LYS:HD3	1:B:338:PRO:CD	2.13	0.79
1:B:419:ASP:O	1:B:423:THR:HG23	1.83	0.79
1:A:387:ARG:HD2	1:A:454:PRO:HG3	1.66	0.78
1:B:107:GLY:HA2	3:B:628:HOH:O	1.85	0.77
1:A:153:THR:HG21	1:A:317:HIS:HB2	1.66	0.77
1:B:242:SER:OG	1:B:245:LEU:HD23	1.84	0.77
1:A:279:MET:HG2	1:A:282:ARG:HH22	1.49	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:GLN:NE2	1:B:336:GLN:H	1.83	0.76
1:B:27:ARG:HG3	3:B:640:HOH:O	1.86	0.75
1:B:204:ILE:HG23	1:B:268:LEU:HD12	1.69	0.75
1:A:136:THR:O	1:A:137:LEU:HB2	1.84	0.75
1:A:282:ARG:HH11	1:A:282:ARG:HG2	1.50	0.74
1:A:447:GLN:O	1:A:451:LEU:HG	1.86	0.74
1:A:453:PHE:HB3	1:A:454:PRO:HD3	1.68	0.74
1:A:33:ARG:HD3	1:A:35:THR:OG1	1.87	0.73
1:B:187:GLY:HA2	2:B:600:ADP:H5'1	1.71	0.73
1:B:412:LYS:CD	1:B:412:LYS:H	2.02	0.73
1:B:56:ARG:HD2	1:B:63:LYS:HD2	1.70	0.73
1:B:85:VAL:HG22	1:B:88:ILE:HD12	1.71	0.72
1:B:141:ALA:O	1:B:255:ARG:HD3	1.89	0.72
1:B:447:GLN:O	1:B:451:LEU:HG	1.90	0.72
1:B:412:LYS:N	1:B:412:LYS:HD3	2.04	0.71
1:B:453:PHE:HB2	1:B:454:PRO:CD	2.21	0.71
1:B:85:VAL:HG13	1:B:88:ILE:HD12	1.73	0.71
1:A:182:ALA:HB2	1:A:353:LEU:HB2	1.72	0.71
1:A:314:ASN:HD22	1:A:321:SER:HB2	1.56	0.70
1:A:305:LEU:HB2	1:A:306:PRO:HD3	1.74	0.69
1:A:136:THR:HG23	1:A:137:LEU:H	1.57	0.69
1:A:428:LEU:HD12	1:A:429:GLU:N	2.07	0.69
1:B:453:PHE:H	1:B:454:PRO:HD2	1.57	0.69
1:A:33:ARG:NH1	1:A:35:THR:HG21	2.07	0.69
1:A:415:ASP:OD2	1:A:418:LEU:HB2	1.92	0.69
1:A:109:GLN:HG3	1:A:139:THR:HB	1.75	0.69
1:B:174:ARG:HA	1:B:323:THR:OG1	1.93	0.68
1:B:295:LYS:HD3	1:B:338:PRO:HD3	1.74	0.67
1:A:279:MET:HG2	1:A:282:ARG:NH2	2.10	0.67
1:A:424:LEU:HD23	1:A:427:GLN:NE2	2.08	0.67
1:B:302:PHE:HZ	1:B:338:PRO:HB3	1.60	0.67
1:B:416:PRO:O	1:B:420:LYS:HG3	1.95	0.67
1:B:282:ARG:HH21	1:B:296:GLY:HA3	1.59	0.67
1:A:147:PHE:HE2	1:A:316:ILE:H	1.41	0.67
1:A:96:ALA:O	1:A:97:ARG:HB2	1.96	0.66
1:B:302:PHE:CZ	1:B:338:PRO:HB3	2.30	0.66
1:A:383:GLN:HB3	3:A:509:HOH:O	1.95	0.66
1:B:212:ARG:HD3	1:B:214:ARG:NH1	2.10	0.66
1:B:316:ILE:HG13	1:B:317:HIS:H	1.60	0.65
1:B:44:LEU:HG	1:B:46:LEU:HD23	1.79	0.65
1:B:225:GLY:O	1:B:229:ARG:HG3	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:GLN:HA	1:B:57:GLN:HE21	1.60	0.65
1:A:148:ASN:C	1:A:150:LEU:H	2.00	0.65
1:A:200:ARG:HE	1:A:439:ARG:HH21	1.45	0.65
1:A:33:ARG:HH11	1:A:35:THR:CG2	2.09	0.65
1:A:153:THR:CG2	1:A:317:HIS:HB2	2.27	0.65
1:B:58:ASP:HB2	1:B:63:LYS:HE2	1.78	0.65
1:B:187:GLY:HA2	2:B:600:ADP:H8	1.62	0.64
1:B:31:LEU:HD13	1:B:54:ILE:HD13	1.78	0.64
1:A:261:ARG:HG2	1:A:261:ARG:NH1	2.10	0.64
1:B:121:LEU:HD23	1:B:127:PRO:HA	1.80	0.64
1:B:190:VAL:HG12	1:B:194:MET:HE2	1.79	0.63
1:A:317:HIS:HD2	1:A:318:GLY:H	1.47	0.63
1:A:453:PHE:CB	1:A:454:PRO:HD3	2.29	0.63
1:A:108:LYS:HE2	1:A:110:LEU:HD21	1.80	0.63
1:B:47:PRO:O	1:B:69:VAL:HG11	1.99	0.63
1:A:148:ASN:HD21	1:A:151:GLN:NE2	1.95	0.62
1:B:119:ARG:NH1	1:B:119:ARG:HB3	2.15	0.62
1:B:25:VAL:HG22	1:B:26:ARG:H	1.63	0.62
1:B:82:LEU:HA	1:B:245:LEU:HD21	1.80	0.62
1:A:425:TRP:HA	1:A:428:LEU:HG	1.81	0.62
1:A:439:ARG:HG3	1:A:439:ARG:HH11	1.64	0.62
1:B:56:ARG:HD2	1:B:63:LYS:CD	2.30	0.62
1:A:200:ARG:HE	1:A:439:ARG:NH2	1.96	0.61
1:A:401:ASN:CG	1:A:418:LEU:HD11	2.21	0.61
1:A:114:PRO:O	1:A:117:LEU:HD23	2.01	0.61
1:A:288:ILE:HD12	1:A:290:GLU:H	1.65	0.61
1:A:73:ASN:O	1:A:75:GLN:N	2.34	0.61
1:A:159:LEU:HD22	1:A:323:THR:HG21	1.83	0.61
1:B:363:TYR:HA	1:B:364:PRO:C	2.21	0.60
1:B:56:ARG:HD3	1:B:57:GLN:N	2.16	0.60
1:B:84:GLU:HG3	3:B:611:HOH:O	2.00	0.60
1:A:387:ARG:HD2	1:A:454:PRO:CG	2.31	0.60
1:B:119:ARG:HH11	1:B:119:ARG:HB3	1.65	0.60
1:B:184:SER:O	1:B:355:ARG:NH2	2.34	0.60
1:B:57:GLN:HA	1:B:61:GLU:O	2.02	0.60
1:A:142:LEU:HD21	1:A:251:ALA:CB	2.30	0.60
1:B:194:MET:HE1	2:B:600:ADP:N6	2.16	0.60
1:B:50:ALA:H	1:B:69:VAL:CG1	2.14	0.60
1:B:250:ALA:O	1:B:254:THR:HG22	2.00	0.60
1:A:42:THR:OG1	1:A:76:ARG:NE	2.36	0.59
1:A:261:ARG:HH11	1:A:261:ARG:CG	2.13	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:ASN:O	1:A:150:LEU:N	2.34	0.59
1:A:200:ARG:NE	1:A:439:ARG:HH21	2.00	0.59
1:A:109:GLN:HE21	1:A:109:GLN:CA	2.16	0.58
1:B:235:ILE:N	1:B:235:ILE:HD12	2.18	0.58
1:B:390:LEU:HD13	1:B:453:PHE:CD2	2.38	0.58
1:B:456:VAL:OXT	1:B:456:VAL:HG12	2.04	0.58
1:A:383:GLN:HG2	1:A:387:ARG:NH1	2.19	0.58
1:A:109:GLN:HG3	1:A:139:THR:CB	2.34	0.58
1:A:31:LEU:HD11	1:A:39:LEU:HB3	1.84	0.58
1:A:261:ARG:NH1	1:A:262:ASP:OD1	2.35	0.57
1:A:405:VAL:HG22	1:A:410:TYR:HB2	1.85	0.57
1:B:57:GLN:NE2	1:B:61:GLU:O	2.36	0.57
1:B:424:LEU:O	1:B:427:GLN:HB2	2.05	0.57
1:A:174:ARG:HH11	1:A:317:HIS:HE1	1.51	0.57
1:B:208:LEU:C	1:B:209:ILE:HD12	2.25	0.57
1:B:241:VAL:HG12	1:B:245:LEU:HB2	1.87	0.56
1:A:356:ARG:N	1:A:356:ARG:HD3	2.10	0.56
1:A:164:ARG:HB2	1:A:440:ALA:CB	2.36	0.56
1:B:387:ARG:HD3	1:B:454:PRO:CG	2.35	0.56
1:A:336:GLN:NE2	1:A:336:GLN:N	2.54	0.56
1:A:385:TYR:O	1:A:389:ARG:HG3	2.05	0.56
1:B:336:GLN:HE21	1:B:336:GLN:H	1.54	0.56
1:A:122:ASP:OD2	1:A:126:LYS:HB2	2.06	0.55
1:B:340:ALA:O	1:B:344:ARG:HG3	2.05	0.55
1:B:442:TRP:CE2	1:B:446:LEU:HD11	2.41	0.55
1:A:297:TYR:OH	1:A:339:ILE:HD11	2.07	0.55
1:B:185:GLY:HA2	3:B:633:HOH:O	2.07	0.55
1:B:254:THR:HG23	1:B:311:ARG:NH2	2.21	0.55
1:B:385:TYR:HE2	1:B:389:ARG:HH21	1.54	0.55
1:B:110:LEU:O	1:B:139:THR:HG23	2.06	0.55
1:B:387:ARG:HD3	1:B:454:PRO:HD3	1.89	0.55
1:A:114:PRO:HA	1:A:117:LEU:HD23	1.89	0.55
1:A:164:ARG:NH2	1:A:433:GLN:O	2.40	0.55
1:A:306:PRO:O	1:A:310:GLU:HB2	2.05	0.55
1:A:64:GLU:OE1	1:A:97:ARG:NH2	2.40	0.55
1:B:25:VAL:HG22	1:B:26:ARG:N	2.22	0.55
1:A:119:ARG:NH2	1:A:132:PRO:O	2.40	0.54
1:B:96:ALA:N	3:B:640:HOH:O	2.40	0.54
1:A:317:HIS:CD2	1:A:318:GLY:H	2.24	0.54
1:A:416:PRO:HA	1:A:419:ASP:OD2	2.07	0.54
1:B:189:SER:HB2	2:B:600:ADP:O2A	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:453:PHE:HB2	1:B:454:PRO:HD3	1.89	0.54
1:B:197:ARG:HD3	1:B:224:LEU:O	2.08	0.54
1:B:210:GLY:HA3	1:B:276:ARG:CG	2.37	0.54
1:B:194:MET:SD	3:B:625:HOH:O	2.58	0.54
1:A:383:GLN:OE1	1:A:384:HIS:HB2	2.07	0.54
1:A:412:LYS:NZ	1:A:419:ASP:HA	2.23	0.54
1:A:450:ASP:OD1	1:A:456:VAL:HG22	2.09	0.54
1:B:50:ALA:O	1:B:68:GLU:HA	2.08	0.54
1:A:398:PHE:O	1:A:402:ARG:HB2	2.07	0.53
1:B:333:ASP:HB2	3:B:648:HOH:O	2.07	0.53
1:A:401:ASN:ND2	1:A:418:LEU:HD11	2.23	0.53
1:A:142:LEU:HD11	1:A:251:ALA:HB3	1.91	0.53
1:A:109:GLN:HA	1:A:109:GLN:NE2	2.24	0.53
1:B:131:LEU:HD12	1:B:131:LEU:N	2.22	0.53
1:B:301:VAL:O	1:B:305:LEU:HD13	2.08	0.53
1:A:297:TYR:CZ	1:A:339:ILE:HD11	2.44	0.53
1:B:235:ILE:HD11	1:B:256:ILE:HD12	1.90	0.53
1:B:442:TRP:NE1	1:B:446:LEU:HD11	2.23	0.53
1:B:41:ALA:HB1	1:B:94:VAL:HG21	1.91	0.53
1:B:152:ARG:HH12	1:B:177:ARG:HH12	1.55	0.53
1:A:453:PHE:HB3	1:A:454:PRO:CD	2.39	0.52
1:A:117:LEU:HD11	1:A:203:VAL:HG13	1.90	0.52
1:A:336:GLN:H	1:A:336:GLN:HE21	1.57	0.52
1:A:74:GLY:HA2	3:A:521:HOH:O	2.08	0.52
1:B:69:VAL:HG21	1:B:77:LEU:HD11	1.91	0.52
1:B:453:PHE:N	1:B:454:PRO:HD2	2.18	0.52
1:A:109:GLN:HE21	1:A:109:GLN:HA	1.75	0.52
1:A:295:LYS:HB2	1:A:338:PRO:HG3	1.91	0.52
1:B:129:ASP:HB2	3:B:612:HOH:O	2.09	0.52
1:A:164:ARG:HB2	1:A:440:ALA:HB1	1.91	0.52
1:B:195:MET:O	1:B:199:THR:HB	2.10	0.52
1:B:387:ARG:HD3	1:B:454:PRO:HG3	1.91	0.52
1:A:205:VAL:CG2	1:A:269:LEU:CD1	2.88	0.51
1:A:65:VAL:HG13	1:A:83:GLU:HB2	1.92	0.51
1:A:54:ILE:HG21	1:A:88:ILE:HD13	1.92	0.51
1:B:109:GLN:HE21	1:B:109:GLN:N	2.08	0.51
1:A:205:VAL:HG23	1:A:269:LEU:HD12	1.92	0.51
1:A:387:ARG:HG2	1:A:454:PRO:HD3	1.92	0.51
1:B:241:VAL:HG12	1:B:242:SER:N	2.25	0.51
1:B:443:GLU:HG2	1:B:447:GLN:HE21	1.76	0.51
1:B:335:GLN:NE2	1:B:335:GLN:N	2.58	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:377:THR:HG23	3:B:631:HOH:O	2.11	0.51
1:B:401:ASN:O	1:B:405:VAL:HG23	2.10	0.51
1:B:59:GLY:C	1:B:61:GLU:H	2.13	0.51
1:B:112:LEU:HA	1:B:116:LEU:CD2	2.41	0.51
1:B:407:VAL:O	1:B:407:VAL:HG13	2.10	0.51
1:B:85:VAL:HG22	1:B:88:ILE:CD1	2.40	0.51
1:A:391:PHE:CZ	1:A:428:LEU:HB2	2.45	0.51
1:B:204:ILE:O	1:B:232:SER:HA	2.11	0.50
1:B:335:GLN:O	1:B:337:ASP:N	2.44	0.50
1:A:450:ASP:CG	1:A:456:VAL:HG22	2.31	0.50
1:A:109:GLN:O	1:A:110:LEU:HD23	2.11	0.50
1:B:152:ARG:HH12	1:B:177:ARG:NH1	2.08	0.50
1:A:225:GLY:O	1:A:229:ARG:HG3	2.11	0.50
1:A:331:GLU:HG3	1:A:331:GLU:O	2.11	0.50
1:A:422:ILE:HG13	1:A:423:THR:N	2.26	0.50
1:B:336:GLN:NE2	1:B:336:GLN:N	2.58	0.50
1:A:142:LEU:O	1:A:142:LEU:HD23	2.11	0.50
1:B:108:LYS:O	1:B:141:ALA:HA	2.11	0.50
1:B:110:LEU:C	1:B:139:THR:HG23	2.32	0.50
1:A:357:LEU:HD13	1:A:362:HIS:CD2	2.47	0.50
1:A:46:LEU:HB2	1:A:69:VAL:HG21	1.92	0.50
1:A:160:ASP:O	1:A:199:THR:HA	2.12	0.49
1:A:397:SER:O	1:A:401:ASN:HB3	2.12	0.49
1:B:144:THR:HB	1:B:145:PRO:HD2	1.94	0.49
1:A:363:TYR:HA	1:A:364:PRO:C	2.32	0.49
1:B:214:ARG:HG3	1:B:214:ARG:HH11	1.77	0.49
1:B:131:LEU:HB3	1:B:132:PRO:HD2	1.94	0.49
1:B:410:TYR:HH	1:B:419:ASP:CG	2.16	0.49
1:B:54:ILE:HA	1:B:94:VAL:HG12	1.94	0.49
1:A:218:ASP:OD1	1:A:222:ASN:ND2	2.36	0.49
1:A:455:THR:HG23	1:A:455:THR:O	2.12	0.49
1:B:112:LEU:HA	1:B:116:LEU:HD22	1.94	0.49
1:A:354:SER:OG	1:A:356:ARG:NE	2.45	0.49
1:B:62:THR:HG21	3:B:613:HOH:O	2.12	0.49
1:A:193:GLY:HA3	1:A:223:ILE:HG22	1.95	0.49
1:A:28:TYR:HD2	1:A:93:ARG:HB3	1.78	0.49
1:B:295:LYS:HD3	1:B:338:PRO:HD2	1.91	0.49
1:A:352:VAL:H	1:A:371:SER:HB3	1.78	0.48
1:B:118:GLY:HA2	1:B:229:ARG:O	2.12	0.48
1:B:241:VAL:CG1	1:B:245:LEU:HB2	2.42	0.48
1:A:148:ASN:C	1:A:150:LEU:N	2.66	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:LEU:HD23	1:B:159:LEU:C	2.33	0.48
1:B:330:THR:C	1:B:332:GLY:H	2.16	0.48
1:B:325:PHE:HE2	1:B:351:ILE:HD12	1.79	0.48
1:B:67:SER:HB2	1:B:80:MET:O	2.14	0.48
1:A:148:ASN:ND2	1:A:151:GLN:HG3	2.28	0.48
1:A:451:LEU:N	1:A:451:LEU:HD23	2.27	0.48
1:B:203:VAL:HG12	1:B:204:ILE:N	2.28	0.48
1:B:158:VAL:HG21	1:B:442:TRP:CZ3	2.48	0.48
1:A:204:ILE:HB	1:A:232:SER:HB3	1.96	0.48
1:A:368:ILE:HD12	1:A:368:ILE:N	2.29	0.48
1:B:109:GLN:HG3	1:B:139:THR:CG2	2.37	0.48
1:B:394:LEU:HD11	1:B:453:PHE:HE2	1.79	0.48
1:B:419:ASP:O	1:B:423:THR:CG2	2.57	0.47
1:A:136:THR:OG1	1:A:137:LEU:N	2.47	0.47
1:A:153:THR:HG22	1:A:314:ASN:HB3	1.97	0.47
1:A:204:ILE:HD11	1:A:231:ARG:NH1	2.29	0.47
1:A:455:THR:O	1:A:456:VAL:C	2.53	0.47
1:A:119:ARG:NH2	1:A:131:LEU:HB2	2.30	0.47
1:A:148:ASN:ND2	1:A:151:GLN:NE2	2.50	0.47
1:B:357:LEU:HG	1:B:362:HIS:CD2	2.50	0.47
1:A:205:VAL:CG2	1:A:269:LEU:HD12	2.44	0.47
1:B:109:GLN:HE21	1:B:109:GLN:CA	2.26	0.47
1:A:425:TRP:N	1:A:426:PRO:HD2	2.30	0.47
1:B:194:MET:HE1	2:B:600:ADP:HN61	1.79	0.47
1:A:400:ARG:HH11	1:A:400:ARG:HG2	1.79	0.47
1:B:188:LYS:HE2	1:B:272:ASP:OD2	2.15	0.47
1:A:336:GLN:N	1:A:336:GLN:HE21	2.12	0.46
1:B:204:ILE:HG23	1:B:268:LEU:CD1	2.44	0.46
1:B:424:LEU:O	1:B:427:GLN:N	2.48	0.46
1:A:75:GLN:O	1:A:76:ARG:HB2	2.14	0.46
1:B:332:GLY:O	1:B:333:ASP:HB2	2.15	0.46
1:A:204:ILE:O	1:A:232:SER:HB2	2.15	0.46
1:A:241:VAL:HG13	1:A:242:SER:N	2.30	0.46
1:A:274:LEU:HD22	1:A:305:LEU:CD2	2.45	0.46
1:A:269:LEU:HB3	1:A:324:ALA:HB2	1.95	0.46
1:B:387:ARG:HD3	1:B:454:PRO:CD	2.45	0.46
1:A:122:ASP:OD1	1:A:124:GLY:N	2.45	0.46
1:A:33:ARG:HG2	1:A:34:ALA:N	2.30	0.46
1:B:219:PHE:O	1:B:223:ILE:HB	2.15	0.46
1:A:282:ARG:NH1	1:A:282:ARG:HG2	2.26	0.46
1:B:434:GLN:NE2	3:B:625:HOH:O	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305:LEU:CB	1:A:306:PRO:HD3	2.45	0.46
1:A:391:PHE:O	1:A:395:LEU:HB2	2.16	0.46
1:A:68:GLU:OE1	1:A:244:LEU:HD22	2.15	0.46
1:A:282:ARG:CG	1:A:282:ARG:HH11	2.24	0.46
1:B:109:GLN:CG	1:B:139:THR:HG22	2.38	0.46
1:A:241:VAL:CG1	1:A:242:SER:N	2.78	0.46
1:B:149:PRO:HG2	3:B:662:HOH:O	2.14	0.46
1:B:204:ILE:HD13	1:B:268:LEU:HB3	1.97	0.46
1:B:245:LEU:HD13	3:B:659:HOH:O	2.16	0.46
1:B:56:ARG:HD3	1:B:57:GLN:C	2.36	0.45
1:A:200:ARG:NE	1:A:439:ARG:NH2	2.60	0.45
1:A:212:ARG:HB2	1:A:215:GLU:HG3	1.98	0.45
1:B:261:ARG:HD2	1:B:315:GLY:CA	2.45	0.45
1:A:294:THR:C	1:A:296:GLY:H	2.20	0.45
1:B:305:LEU:N	1:B:305:LEU:HD12	2.32	0.45
1:B:272:ASP:HA	1:B:273:SER:HA	1.69	0.45
1:B:89:LEU:HB3	1:B:90:PRO:HD2	1.99	0.45
1:B:159:LEU:HD13	1:B:323:THR:HG21	1.98	0.45
1:B:80:MET:SD	1:B:244:LEU:HG	2.56	0.45
1:B:288:ILE:O	1:B:288:ILE:HG23	2.16	0.45
1:A:117:LEU:N	1:A:117:LEU:HD22	2.31	0.45
1:A:336:GLN:NE2	1:A:336:GLN:H	2.13	0.45
1:B:131:LEU:CD1	1:B:131:LEU:N	2.79	0.45
1:B:95:TYR:O	1:B:96:ALA:O	2.35	0.45
1:B:32:THR:HG21	1:B:40:GLU:OE2	2.17	0.45
1:B:209:ILE:N	1:B:209:ILE:HD12	2.32	0.44
1:B:261:ARG:HD2	1:B:315:GLY:HA3	1.99	0.44
1:B:54:ILE:CG2	1:B:65:VAL:HB	2.35	0.44
1:A:149:PRO:O	1:A:150:LEU:HD23	2.15	0.44
1:A:293:ALA:HB2	1:A:299:PRO:HG3	1.98	0.44
1:A:96:ALA:O	1:A:97:ARG:CB	2.65	0.44
1:A:219:PHE:HA	1:A:223:ILE:CG1	2.47	0.44
1:B:418:LEU:O	1:B:422:ILE:HG12	2.17	0.44
1:B:85:VAL:CG2	1:B:88:ILE:HD12	2.46	0.44
1:A:453:PHE:CB	1:A:454:PRO:CD	2.92	0.44
1:B:210:GLY:HA3	1:B:276:ARG:HG2	1.99	0.44
1:B:246:ARG:HD2	3:B:634:HOH:O	2.17	0.44
1:B:247:MET:CE	1:B:284:ILE:HD12	2.48	0.44
1:B:145:PRO:HA	1:B:146:PRO:HD2	1.92	0.44
1:A:425:TRP:O	1:A:428:LEU:HG	2.18	0.44
1:B:188:LYS:HD2	2:B:600:ADP:O2B	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:ARG:HG3	1:B:214:ARG:NH1	2.33	0.43
1:A:288:ILE:HD12	1:A:288:ILE:C	2.38	0.43
1:A:182:ALA:HB1	1:A:186:VAL:CG2	2.48	0.43
1:B:305:LEU:N	1:B:306:PRO:HD2	2.33	0.43
1:B:330:THR:O	1:B:332:GLY:N	2.50	0.43
1:B:34:ALA:O	1:B:35:THR:C	2.55	0.43
1:A:267:VAL:HG12	1:A:268:LEU:N	2.33	0.43
1:A:184:SER:O	1:A:355:ARG:NH2	2.51	0.43
1:A:56:ARG:HG3	1:A:56:ARG:O	2.18	0.43
1:A:258:GLU:HA	1:A:322:ILE:HD11	1.98	0.43
1:A:398:PHE:CD1	1:A:422:ILE:HG22	2.54	0.43
1:B:58:ASP:O	1:B:61:GLU:HB3	2.18	0.43
1:B:261:ARG:HG3	1:B:320:GLY:C	2.39	0.43
1:A:119:ARG:CZ	1:A:131:LEU:HB2	2.49	0.43
1:A:294:THR:O	1:A:296:GLY:N	2.46	0.43
1:B:37:LEU:HD11	1:B:283:GLU:HG2	2.01	0.43
1:B:314:ASN:HA	1:B:314:ASN:HD22	1.68	0.43
1:A:148:ASN:HD22	1:A:151:GLN:HG3	1.84	0.43
1:A:261:ARG:HD3	1:A:322:ILE:HG13	2.01	0.43
1:A:418:LEU:O	1:A:421:ALA:HB3	2.19	0.43
1:A:142:LEU:CD2	1:A:251:ALA:HB1	2.40	0.42
1:A:273:SER:HA	1:A:327:THR:OG1	2.19	0.42
1:B:241:VAL:CG1	1:B:242:SER:N	2.81	0.42
1:B:181:PHE:N	1:B:181:PHE:CD2	2.87	0.42
1:B:32:THR:HG23	1:B:33:ARG:N	2.34	0.42
1:A:254:THR:O	1:A:258:GLU:HG3	2.19	0.42
1:B:288:ILE:O	1:B:288:ILE:CG2	2.68	0.42
1:B:187:GLY:CA	2:B:600:ADP:H8	2.30	0.42
1:B:119:ARG:NH1	1:B:129:ASP:OD1	2.53	0.42
1:B:330:THR:C	1:B:332:GLY:N	2.72	0.42
1:A:205:VAL:HG23	1:A:269:LEU:CD1	2.50	0.42
1:A:325:PHE:HE2	1:A:351:ILE:HD12	1.85	0.42
1:A:181:PHE:CD2	1:A:181:PHE:N	2.87	0.42
1:A:205:VAL:CG2	1:A:269:LEU:HD13	2.50	0.42
1:B:319:GLY:N	3:B:605:HOH:O	2.53	0.42
1:B:85:VAL:CG1	1:B:88:ILE:HD12	2.47	0.42
1:A:369:GLU:HG3	1:A:392:LYS:NZ	2.34	0.41
1:A:401:ASN:OD1	1:A:418:LEU:HD11	2.20	0.41
1:B:117:LEU:O	1:B:119:ARG:HG2	2.20	0.41
1:A:202:ASP:OD1	1:A:266:HIS:HB2	2.20	0.41
1:B:425:TRP:N	1:B:426:PRO:HD2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:LEU:C	1:A:142:LEU:HD23	2.39	0.41
1:A:144:THR:HG21	1:A:255:ARG:HE	1.85	0.41
1:A:153:THR:CB	1:A:317:HIS:HB2	2.50	0.41
1:A:272:ASP:HA	1:A:273:SER:HA	1.66	0.41
1:B:109:GLN:NE2	1:B:109:GLN:CA	2.84	0.41
1:B:247:MET:HE1	1:B:284:ILE:HD12	2.01	0.41
1:A:350:HIS:O	1:A:371:SER:HB2	2.21	0.41
1:A:198:TYR:HB3	1:A:439:ARG:NH1	2.35	0.41
1:A:75:GLN:O	1:A:76:ARG:HD3	2.21	0.41
1:B:57:GLN:O	1:B:59:GLY:N	2.54	0.41
1:A:406:SER:HB2	3:A:510:HOH:O	2.19	0.41
1:A:182:ALA:HB1	1:A:186:VAL:HG21	2.03	0.41
1:B:335:GLN:H	1:B:335:GLN:NE2	2.18	0.41
1:A:425:TRP:CZ3	1:A:428:LEU:HD11	2.54	0.41
1:A:200:ARG:HH21	1:A:439:ARG:HH21	1.69	0.41
1:A:200:ARG:NH2	1:A:439:ARG:HH21	2.18	0.41
1:A:424:LEU:HD23	1:A:427:GLN:HE22	1.85	0.40
1:B:407:VAL:O	1:B:407:VAL:CG1	2.69	0.40
1:B:390:LEU:O	1:B:394:LEU:HG	2.20	0.40
1:A:80:MET:HA	1:A:81:PRO:HD3	1.83	0.40
1:B:188:LYS:HB3	2:B:600:ADP:O1B	2.21	0.40
1:A:274:LEU:HG	1:A:326:TYR:HB3	2.03	0.40
1:A:137:LEU:CD2	1:B:26:ARG:HB3	2.52	0.40
1:B:86:GLU:CD	1:B:86:GLU:N	2.75	0.40
1:B:209:ILE:CG2	1:B:210:GLY:N	2.85	0.40
1:B:86:GLU:OE1	1:B:87:GLY:N	2.46	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	418/438 (95%)	379 (91%)	28 (7%)	11 (3%)	6	6
1	B	422/438 (96%)	368 (87%)	42 (10%)	12 (3%)	6	5
All	All	840/876 (96%)	747 (89%)	70 (8%)	23 (3%)	6	6

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	74	GLY
1	A	136	THR
1	A	144	THR
1	A	454	PRO
1	B	34	ALA
1	B	35	THR
1	B	96	ALA
1	B	316	ILE
1	B	336	GLN
1	B	412	LYS
1	B	453	PHE
1	A	453	PHE
1	B	454	PRO
1	B	331	GLU
1	B	332	GLY
1	A	76	ARG
1	A	295	LYS
1	A	60	PRO
1	A	75	GLN
1	B	83	GLU
1	B	107	GLY
1	A	49	GLY
1	A	143	ILE

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	332/342 (97%)	306 (92%)	26 (8%)	15	23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	334/342 (98%)	312 (93%)	22 (7%)	19	30
All	All	666/684 (97%)	618 (93%)	48 (7%)	17	26

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	LEU
1	A	48	LEU
1	A	58	ASP
1	A	76	ARG
1	A	109	GLN
1	A	112	LEU
1	A	148	ASN
1	A	214	ARG
1	A	241	VAL
1	A	309	VAL
1	A	317	HIS
1	A	334	ASP
1	A	336	GLN
1	A	353	LEU
1	A	356	ARG
1	A	369	GLU
1	A	382	GLU
1	A	383	GLN
1	A	390	LEU
1	A	392	LYS
1	A	418	LEU
1	A	422	ILE
1	A	441	ASP
1	A	447	GLN
1	A	453	PHE
1	A	454	PRO
1	B	33	ARG
1	B	56	ARG
1	B	57	GLN
1	B	86	GLU
1	B	109	GLN
1	B	150	LEU
1	B	180	LEU
1	B	188	LYS
1	B	199	THR
1	B	254	THR

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Mol	Chain	Res	Type
1	B	255	ARG
1	B	268	LEU
1	B	335	GLN
1	B	347	LEU
1	B	354	SER
1	B	357	LEU
1	B	391	PHE
1	B	403	ASP
1	B	407	VAL
1	B	412	LYS
1	B	418	LEU
1	B	423	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	GLN
1	A	73	ASN
1	A	109	GLN
1	A	148	ASN
1	A	281	GLN
1	A	314	ASN
1	A	317	HIS
1	A	335	GLN
1	A	336	GLN
1	A	399	GLN
1	A	427	GLN
1	B	57	GLN
1	B	109	GLN
1	B	248	GLN
1	B	314	ASN
1	B	317	HIS
1	B	335	GLN
1	B	336	GLN
1	B	393	GLN
1	B	433	GLN
1	B	447	GLN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z > 2$
2	ADP	B	600	-	25,29,29	2.07	7 (28%)	24,45,45	2.19	3 (12%)

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	ADP	B	600	-	-	0/12/32/32	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	600	ADP	PB-O3A	-4.44	1.52	1.60
2	B	600	ADP	PA-O2A	-3.79	1.36	1.55
2	B	600	ADP	C5-N7	-2.54	1.30	1.39
2	B	600	ADP	C2-N1	2.52	1.38	1.33
2	B	600	ADP	PB-O2B	2.97	1.67	1.54
2	B	600	ADP	C2-N3	3.60	1.38	1.32
2	B	600	ADP	O4'-C1'	4.72	1.47	1.41

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	600	ADP	N3-C2-N1	-9.82	120.31	128.86
2	B	600	ADP	C4-C5-N7	-2.57	106.93	109.41
2	B	600	ADP	O2A-PA-O5'	2.01	117.61	108.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	600	ADP	8	0

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	422/438 (96%)	0.30	36 (8%) 11 10	22, 47, 96, 111	0
1	B	426/438 (97%)	0.37	27 (6%) 21 19	23, 48, 93, 113	0
All	All	848/876 (96%)	0.34	63 (7%) 15 14	22, 48, 94, 113	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	25	VAL	11.2
1	B	85	VAL	9.9
1	B	60	PRO	8.1
1	A	407	VAL	7.9
1	B	59	GLY	7.3
1	B	456	VAL	6.8
1	B	24	ALA	6.1
1	B	333	ASP	5.9
1	A	404	LEU	5.6
1	A	408	GLY	5.1
1	B	332	GLY	5.0
1	B	334	ASP	4.9
1	B	23	PRO	4.7
1	A	456	VAL	4.6
1	A	316	ILE	4.5
1	B	54	ILE	4.3
1	A	76	ARG	4.3
1	A	332	GLY	4.2
1	B	65	VAL	4.2
1	A	74	GLY	4.1
1	A	405	VAL	4.0
1	B	57	GLN	4.0
1	B	34	ALA	3.9
1	A	406	SER	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	397	SER	3.9
1	A	317	HIS	3.8
1	A	416	PRO	3.8
1	B	331	GLU	3.8
1	A	333	ASP	3.7
1	A	414	SER	3.7
1	A	147	PHE	3.6
1	A	409	ALA	3.5
1	A	73	ASN	3.3
1	B	135	ASP	3.3
1	B	46	LEU	3.3
1	B	149	PRO	3.2
1	B	90	PRO	3.2
1	A	424	LEU	3.1
1	B	335	GLN	2.9
1	A	453	PHE	2.8
1	A	412	LYS	2.8
1	A	145	PRO	2.8
1	A	399	GLN	2.7
1	A	27	ARG	2.6
1	B	106	SER	2.6
1	B	97	ARG	2.5
1	A	75	GLN	2.5
1	A	420	LYS	2.4
1	B	89	LEU	2.4
1	A	398	PHE	2.3
1	A	413	GLY	2.3
1	A	110	LEU	2.3
1	A	146	PRO	2.3
1	B	64	GLU	2.3
1	A	451	LEU	2.2
1	A	25	VAL	2.2
1	A	335	GLN	2.2
1	A	419	ASP	2.1
1	B	35	THR	2.1
1	A	44	LEU	2.1
1	B	27	ARG	2.0
1	B	37	LEU	2.0
1	A	149	PRO	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	ADP	B	600	27/27	0.59	0.50	3.78	127,130,133,133	0

6.5 Other polymers [i](#)

There are no such residues in this entry.