



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 14, 2016 – 08:22 PM EDT

PDB ID : 4ZHI
Title : Eg5 motor domain mutant E162S
Authors : Luo, M.; Park, C.; Worthylake, D.; Kim, S.
Deposited on : 2015-04-24
Resolution : 2.30 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027790
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : rb-20027790

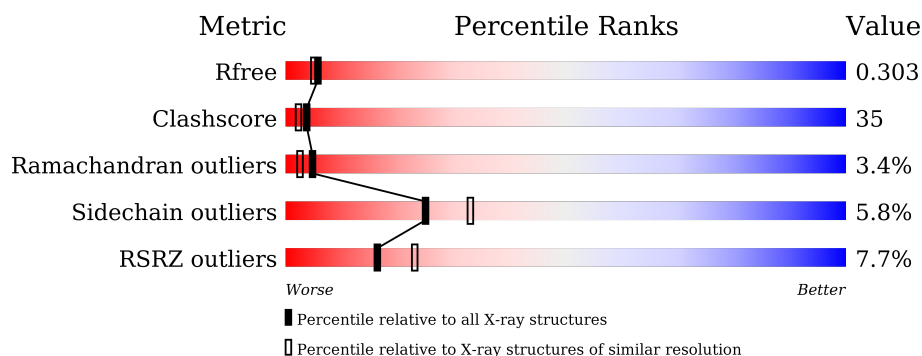
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	369	
1	B	369	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5527 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.

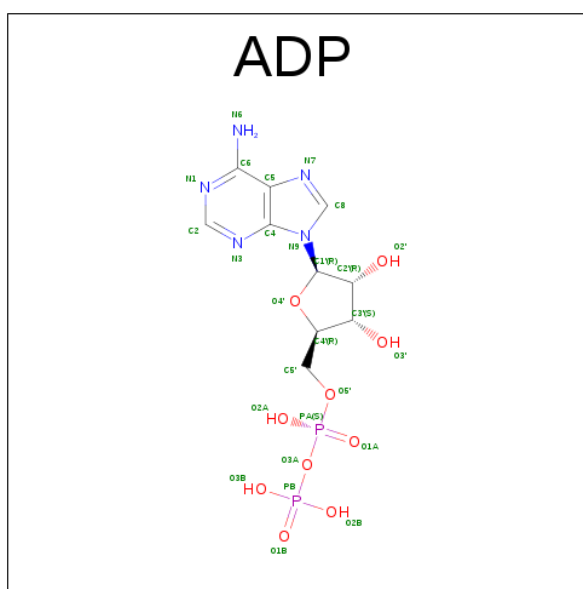
- Molecule 1 is a protein called Kinesin-like protein KIF11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	0	0
			2577	1614	450	503	10			
1	B	330	Total	C	N	O	S	0	0	0
			2588	1622	450	506	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	162	SER	GLU	engineered mutation	UNP P52732
B	162	SER	GLU	engineered mutation	UNP P52732

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

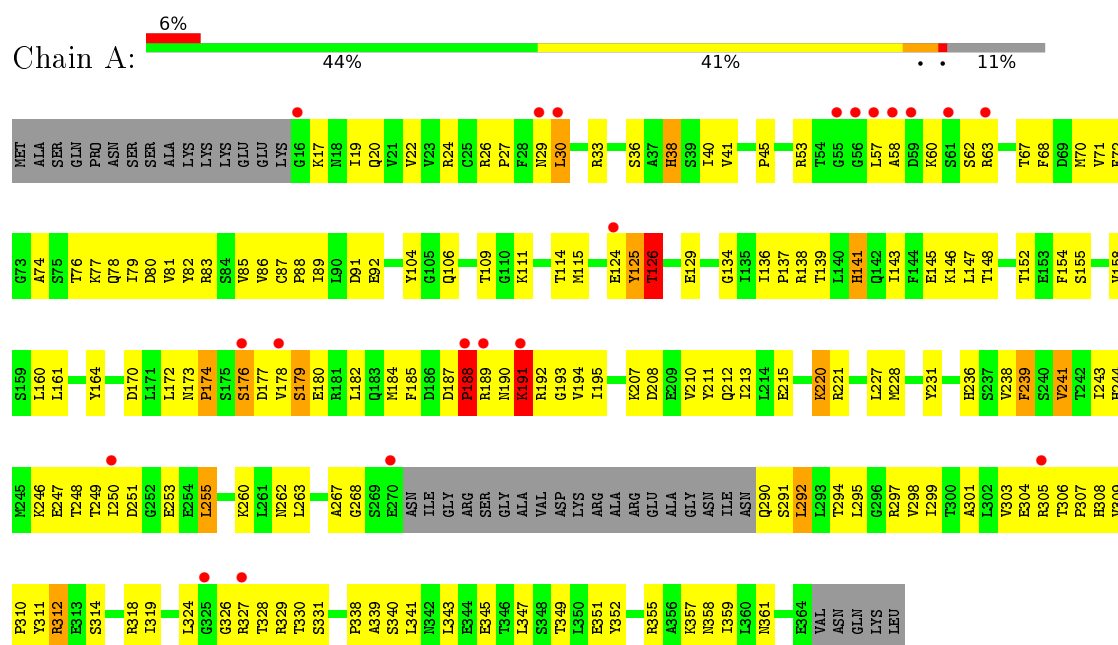
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	184	Total	O	0	0
			184	184		
3	B	124	Total	O	0	0
			124	124		

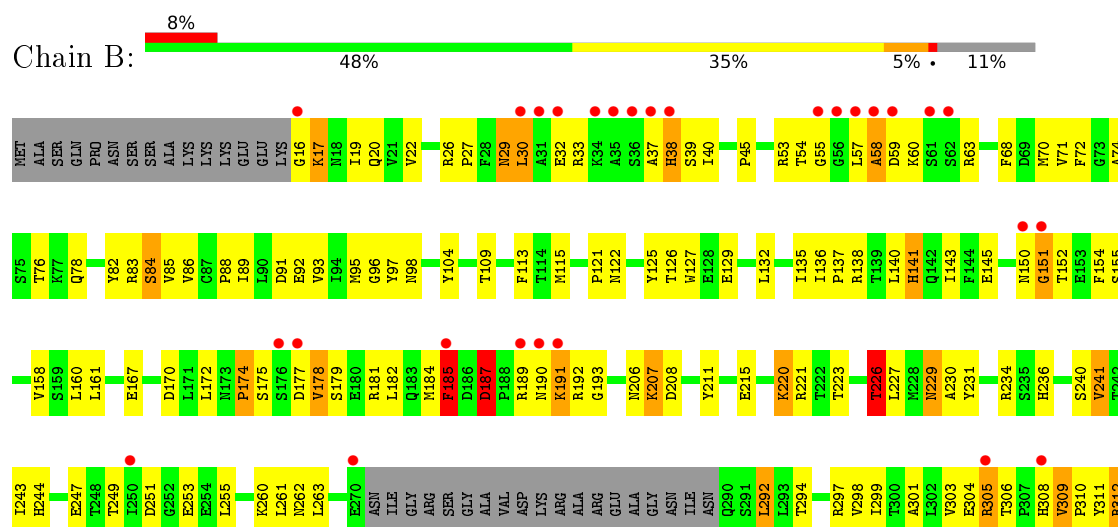
3 Residue-property plots

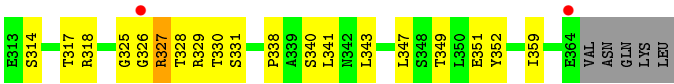
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Kinesin-like protein KIF11



• Molecule 1: Kinesin-like protein KIF11





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.72Å 81.22Å 82.60Å 90.00° 94.31° 90.00°	Depositor
Resolution (Å)	25.00 – 2.30 24.79 – 2.22	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.30) 96.9 (24.79-2.22)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.22Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.254 , 0.303 0.253 , 0.303	Depositor DCC
R_{free} test set	1517 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	24.0	Xtriage
Anisotropy	0.528	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 19.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5527	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.83% 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	0.42	0/2615	0.68	2/3535 (0.1%)
1	B	0.47	1/2627 (0.0%)	0.89	7/3551 (0.2%)
All	All	0.45	1/5242 (0.0%)	0.79	9/7086 (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
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	185	PHE	CB-CG	-7.63	1.38	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	185	PHE	CB-CG-CD2	-19.60	107.08	120.80
1	B	185	PHE	CB-CG-CD1	17.04	132.73	120.80
1	B	185	PHE	N-CA-CB	-17.03	79.94	110.60
1	B	187	ASP	CB-CG-OD2	-8.43	110.71	118.30
1	B	185	PHE	CB-CA-C	7.21	124.82	110.40
1	B	187	ASP	CB-CG-OD1	5.64	123.37	118.30
1	A	126	THR	N-CA-C	-5.36	96.52	111.00
1	B	189	ARG	NE-CZ-NH2	-5.23	117.69	120.30
1	A	324	LEU	N-CA-C	-5.10	97.22	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	185	PHE	Sidechain

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	2577	0	2606	193	0
1	B	2588	0	2617	181	0
2	A	27	0	12	3	0
2	B	27	0	12	1	0
3	A	184	0	0	70	0
3	B	124	0	0	33	0
All	All	5527	0	5247	368	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 35.

All (368) 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:249:THR:HG22	1:B:253:GLU:H	1.13	1.09
1:A:78:GLN:HB2	3:A:729:HOH:O	1.56	1.06
1:A:40:ILE:HG12	3:A:702:HOH:O	1.57	1.04
1:A:329:ARG:HD3	3:A:825:HOH:O	1.63	0.98
1:A:294:THR:O	3:A:701:HOH:O	1.85	0.94
1:B:206:ASN:HD21	1:B:208:ASP:HB2	1.33	0.94
1:A:249:THR:HG23	1:A:251:ASP:HB2	1.51	0.93
1:B:40:ILE:HD13	1:B:340:SER:HA	1.50	0.93
1:A:188:PRO:HD2	3:A:847:HOH:O	1.68	0.93
1:A:213:ILE:HG13	3:A:856:HOH:O	1.68	0.92
1:A:76:THR:HG22	3:A:703:HOH:O	1.69	0.91
1:B:249:THR:HG23	1:B:251:ASP:H	1.36	0.91
1:B:329:ARG:HD3	3:B:782:HOH:O	1.73	0.89
1:B:85:VAL:N	3:B:703:HOH:O	2.05	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:314:SER:O	3:B:701:HOH:O	1.89	0.89
1:B:187:ASP:HB2	1:B:190:ASN:HB2	1.51	0.88
1:B:178:VAL:HG12	1:B:179:SER:H	1.38	0.88
1:B:317:THR:N	3:B:701:HOH:O	2.07	0.88
1:B:347:LEU:O	1:B:351:GLU:HG3	1.74	0.88
1:A:187:ASP:HB3	1:A:195:ILE:HG12	1.58	0.86
1:A:17:LYS:HE3	3:A:750:HOH:O	1.77	0.85
1:B:32:GLU:O	1:B:37:ALA:HB3	1.77	0.85
1:A:249:THR:HG21	1:A:253:GLU:OE1	1.76	0.85
1:A:30:LEU:HA	1:A:33:ARG:HG3	1.58	0.84
1:B:249:THR:HG22	1:B:253:GLU:N	1.93	0.84
1:B:29:ASN:HB3	3:B:810:HOH:O	1.78	0.83
1:B:74:ALA:HB2	3:B:730:HOH:O	1.76	0.83
1:A:312:ARG:HB3	1:A:312:ARG:HH11	1.44	0.83
1:B:247:GLU:HB3	3:B:705:HOH:O	1.77	0.83
1:B:312:ARG:HB3	1:B:312:ARG:HH11	1.43	0.82
1:B:92:GLU:O	3:B:702:HOH:O	1.95	0.82
1:B:230:ALA:O	1:B:234:ARG:HG2	1.79	0.81
1:A:297:ARG:N	3:A:701:HOH:O	2.11	0.81
1:B:327:ARG:HE	1:B:327:ARG:HA	1.45	0.81
1:B:37:ALA:HA	3:B:745:HOH:O	1.79	0.80
1:A:328:THR:HB	3:A:846:HOH:O	1.80	0.80
1:B:97:TYR:N	3:B:702:HOH:O	2.14	0.80
1:B:187:ASP:CG	1:B:190:ASN:HD22	1.84	0.80
1:A:338:PRO:HA	3:A:735:HOH:O	1.79	0.80
1:B:192:ARG:HD3	3:B:800:HOH:O	1.83	0.79
1:A:136:ILE:HG12	1:A:263:LEU:HD13	1.63	0.79
1:A:347:LEU:O	1:A:351:GLU:HG2	1.83	0.79
1:A:295:LEU:C	3:A:701:HOH:O	2.22	0.78
1:B:167:GLU:OE1	1:B:181:ARG:HD3	1.82	0.77
1:B:305:ARG:HG2	1:B:305:ARG:HH11	1.49	0.77
1:B:249:THR:HG23	1:B:251:ASP:N	2.00	0.77
1:A:194:VAL:HB	3:A:716:HOH:O	1.85	0.77
1:A:330:THR:HG23	3:A:711:HOH:O	1.84	0.77
1:A:92:GLU:HG2	3:A:865:HOH:O	1.83	0.76
1:B:187:ASP:CB	1:B:190:ASN:HB2	2.15	0.76
1:A:299:ILE:O	1:A:303:VAL:HG23	1.85	0.76
1:A:312:ARG:NH1	1:A:312:ARG:H	1.83	0.76
1:A:329:ARG:HG2	3:A:753:HOH:O	1.86	0.75
1:B:38:HIS:N	1:B:38:HIS:CD2	2.55	0.75
1:A:40:ILE:HD13	1:A:340:SER:HA	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:PRO:O	3:A:702:HOH:O	2.06	0.74
1:B:82:TYR:O	3:B:703:HOH:O	2.04	0.74
1:A:190:ASN:ND2	1:A:193:GLY:HA3	2.02	0.74
1:A:158:VAL:HG12	1:A:241:VAL:HG23	1.69	0.73
1:A:126:THR:OG1	1:A:129:GLU:HG3	1.87	0.73
1:B:158:VAL:HG12	1:B:241:VAL:HG23	1.72	0.72
1:B:312:ARG:H	1:B:312:ARG:NH1	1.87	0.72
1:A:81:VAL:HG23	3:A:703:HOH:O	1.88	0.72
1:A:38:HIS:CD2	1:A:38:HIS:N	2.58	0.71
1:B:226:THR:HG22	1:B:227:LEU:HD23	1.71	0.71
1:A:41:VAL:HG21	3:A:735:HOH:O	1.91	0.71
1:B:244:HIS:HD2	3:B:814:HOH:O	1.72	0.71
1:B:59:ASP:O	1:B:60:LYS:HD3	1.90	0.71
1:B:328:THR:HB	3:B:789:HOH:O	1.91	0.71
1:A:80:ASP:HB2	3:A:703:HOH:O	1.90	0.71
1:A:239:PHE:N	3:A:709:HOH:O	2.23	0.71
1:B:160:LEU:HB3	1:B:172:LEU:HD13	1.70	0.70
1:B:190:ASN:HB3	1:B:193:GLY:H	1.55	0.70
1:A:111:LYS:HE2	3:A:704:HOH:O	1.91	0.70
1:A:305:ARG:HH11	1:A:305:ARG:HG2	1.53	0.70
1:B:299:ILE:O	1:B:303:VAL:HG23	1.92	0.70
1:A:138:ARG:NH1	3:A:705:HOH:O	2.18	0.70
1:A:160:LEU:HB3	1:A:172:LEU:HD13	1.73	0.70
1:A:305:ARG:HH12	1:B:30:LEU:HD23	1.56	0.69
1:A:106:GLN:O	3:A:704:HOH:O	2.11	0.69
1:A:136:ILE:HG12	1:A:263:LEU:CD1	2.22	0.69
1:A:77:LYS:O	3:A:703:HOH:O	2.10	0.69
1:A:267:ALA:O	1:A:292:LEU:HD11	1.93	0.69
1:B:184:MET:O	1:B:185:PHE:HB2	1.90	0.69
1:B:27:PRO:HB3	1:B:74:ALA:HB1	1.74	0.69
1:A:53:ARG:NH1	1:A:63:ARG:HD3	2.09	0.68
1:B:229:ASN:HD22	1:B:230:ALA:H	1.41	0.68
1:B:40:ILE:CD1	1:B:340:SER:HA	2.24	0.67
1:A:53:ARG:NH1	1:A:63:ARG:HB3	2.09	0.67
1:B:318:ARG:N	3:B:701:HOH:O	2.17	0.66
1:A:211:TYR:HB2	3:A:783:HOH:O	1.94	0.66
1:B:325:GLY:C	1:B:327:ARG:H	1.99	0.66
1:B:206:ASN:ND2	1:B:208:ASP:HB2	2.07	0.66
1:A:106:GLN:NE2	1:A:268:GLY:HA3	2.10	0.66
1:A:82:TYR:CG	1:A:138:ARG:HD2	2.31	0.65
1:B:167:GLU:HB3	1:B:181:ARG:HD3	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:TYR:CD2	3:B:702:HOH:O	2.49	0.65
1:A:143:ILE:HD13	1:A:243:ILE:HD11	1.77	0.65
1:B:53:ARG:NH1	1:B:63:ARG:HB3	2.11	0.65
1:A:211:TYR:O	1:A:215:GLU:HG3	1.95	0.65
1:B:292:LEU:HD23	1:B:352:TYR:CE2	2.32	0.65
1:A:62:SER:HA	3:A:730:HOH:O	1.95	0.64
1:A:303:VAL:CG2	3:A:875:HOH:O	2.45	0.64
1:A:148:THR:N	3:A:708:HOH:O	2.21	0.64
1:A:341:LEU:HB2	3:A:764:HOH:O	1.96	0.64
1:A:236:HIS:HE1	1:A:292:LEU:HD12	1.62	0.64
1:A:305:ARG:NH1	1:A:305:ARG:HG2	2.13	0.64
1:A:250:ILE:HG22	1:A:250:ILE:O	1.98	0.63
1:A:301:ALA:HB3	1:A:309:VAL:HG22	1.80	0.63
1:B:249:THR:HG21	1:B:253:GLU:OE1	1.98	0.63
1:A:22:VAL:HG12	1:A:70:MET:HB2	1.81	0.63
1:A:310:PRO:HB3	3:A:738:HOH:O	1.99	0.63
1:B:305:ARG:HG2	1:B:305:ARG:NH1	2.12	0.63
1:B:220:LYS:O	1:B:220:LYS:HE2	1.99	0.63
1:B:327:ARG:HA	1:B:327:ARG:NE	2.12	0.63
1:B:143:ILE:HD13	1:B:243:ILE:HD11	1.81	0.62
1:A:192:ARG:HD3	3:A:850:HOH:O	1.99	0.62
1:A:134:GLY:O	1:A:137:PRO:HD2	1.99	0.62
1:B:229:ASN:HD22	1:B:230:ALA:N	1.97	0.62
1:A:305:ARG:NH1	1:B:30:LEU:HD23	2.13	0.62
1:A:83:ARG:HD2	3:A:769:HOH:O	1.99	0.62
1:A:38:HIS:H	1:A:38:HIS:HD2	1.46	0.62
1:B:84:SER:N	3:B:703:HOH:O	2.32	0.62
1:B:178:VAL:HG12	1:B:179:SER:N	2.13	0.61
1:A:26:ARG:HB2	1:A:27:PRO:HD2	1.83	0.61
1:A:184:MET:HG3	3:A:716:HOH:O	2.00	0.61
1:A:155:SER:OG	1:A:244:HIS:HB2	2.01	0.61
1:A:190:ASN:HD22	1:A:193:GLY:HA3	1.65	0.61
1:A:248:THR:HG22	1:A:249:THR:H	1.65	0.60
1:A:38:HIS:CD2	1:A:38:HIS:H	2.17	0.60
1:B:167:GLU:HB3	1:B:181:ARG:CD	2.30	0.60
1:B:38:HIS:HD2	1:B:38:HIS:H	1.48	0.60
1:A:81:VAL:N	3:A:703:HOH:O	2.35	0.60
1:A:267:ALA:HB2	3:A:786:HOH:O	2.02	0.60
1:A:60:LYS:O	1:A:60:LYS:HG3	2.01	0.60
1:B:82:TYR:CG	1:B:138:ARG:HD2	2.37	0.60
1:B:59:ASP:C	1:B:60:LYS:HD3	2.23	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:VAL:HG12	1:A:241:VAL:CG2	2.32	0.59
1:B:158:VAL:HG12	1:B:241:VAL:CG2	2.31	0.59
1:A:136:ILE:HB	1:A:137:PRO:HD3	1.84	0.59
1:A:143:ILE:HD13	1:A:243:ILE:CD1	2.32	0.59
1:A:312:ARG:CB	1:A:312:ARG:HH11	2.15	0.59
1:A:40:ILE:N	3:A:702:HOH:O	2.34	0.59
1:B:329:ARG:HG2	3:B:771:HOH:O	2.03	0.59
1:A:20:GLN:HE22	1:A:85:VAL:HA	1.68	0.59
1:B:22:VAL:HG12	1:B:70:MET:HB2	1.85	0.58
1:B:38:HIS:CD2	1:B:38:HIS:H	2.19	0.58
1:B:211:TYR:O	1:B:215:GLU:HG3	2.02	0.58
1:B:249:THR:CG2	1:B:253:GLU:H	2.01	0.58
1:A:221:ARG:HG3	1:A:231:TYR:HE2	1.68	0.58
1:B:96:GLY:N	3:B:702:HOH:O	2.35	0.58
1:B:136:ILE:HB	1:B:137:PRO:HD3	1.85	0.58
1:B:236:HIS:HE1	1:B:292:LEU:HD12	1.69	0.58
1:B:341:LEU:HD13	3:B:731:HOH:O	2.04	0.57
1:A:220:LYS:O	1:A:220:LYS:HE2	2.04	0.57
1:A:357:LYS:HB2	3:A:875:HOH:O	2.04	0.57
1:B:301:ALA:HB3	1:B:309:VAL:HG22	1.84	0.57
1:B:326:GLY:C	1:B:328:THR:H	2.06	0.57
1:B:312:ARG:CB	1:B:312:ARG:HH11	2.16	0.57
1:A:238:VAL:HG12	1:A:238:VAL:O	2.05	0.57
1:B:53:ARG:NH1	1:B:63:ARG:HD3	2.20	0.57
1:B:150:ASN:ND2	1:B:152:THR:OG1	2.37	0.57
1:A:111:LYS:N	2:A:601:ADP:O2B	2.38	0.56
1:B:260:LYS:HE2	1:B:262:ASN:HD21	1.69	0.56
1:B:57:LEU:O	1:B:58:ALA:HB3	2.04	0.56
1:B:26:ARG:HB2	1:B:27:PRO:HD2	1.88	0.56
1:B:135:ILE:HD11	3:B:811:HOH:O	2.06	0.56
1:B:38:HIS:N	1:B:38:HIS:HD2	1.99	0.56
1:A:191:LYS:HA	1:A:191:LYS:HZ3	1.71	0.55
1:B:32:GLU:HB3	1:B:37:ALA:CB	2.37	0.55
1:A:345:GLU:HG3	3:A:849:HOH:O	2.06	0.55
1:A:57:LEU:O	1:A:58:ALA:HB3	2.06	0.55
1:B:22:VAL:CG1	1:B:70:MET:HB2	2.37	0.55
1:A:185:PHE:N	3:A:716:HOH:O	2.33	0.55
1:A:294:THR:O	1:A:298:VAL:HG23	2.06	0.55
1:B:236:HIS:CE1	1:B:292:LEU:HD12	2.42	0.54
1:B:230:ALA:HB3	1:B:234:ARG:HD2	1.90	0.54
1:A:187:ASP:OD2	1:A:190:ASN:OD1	2.24	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:HIS:CE1	1:A:292:LEU:HD12	2.42	0.54
1:A:309:VAL:HG11	1:A:311:TYR:CZ	2.42	0.54
1:B:17:LYS:HE3	1:B:17:LYS:HA	1.89	0.54
1:B:20:GLN:HE22	1:B:85:VAL:HA	1.72	0.54
1:A:154:PHE:HA	1:A:244:HIS:O	2.08	0.54
1:B:16:GLY:O	1:B:17:LYS:HG2	2.07	0.54
1:B:223:THR:O	1:B:226:THR:HB	2.08	0.54
1:B:184:MET:HG2	1:B:185:PHE:N	2.23	0.54
1:B:30:LEU:C	1:B:32:GLU:H	2.10	0.54
1:B:86:VAL:N	3:B:703:HOH:O	2.41	0.54
1:A:147:LEU:N	3:A:708:HOH:O	2.41	0.53
1:A:161:LEU:HD12	1:A:161:LEU:C	2.28	0.53
1:A:19:ILE:HD13	1:A:330:THR:HB	1.90	0.53
1:B:127:TRP:HA	1:B:141:HIS:CD2	2.43	0.53
1:A:303:VAL:HG22	3:A:875:HOH:O	2.05	0.53
1:A:303:VAL:HG21	3:A:875:HOH:O	2.06	0.53
1:B:207:LYS:HD2	1:B:208:ASP:N	2.24	0.53
1:B:85:VAL:HG13	1:B:331:SER:OG	2.09	0.53
1:A:184:MET:HE3	1:A:318:ARG:CZ	2.39	0.53
1:B:109:THR:HG21	3:B:750:HOH:O	2.09	0.53
1:A:187:ASP:HB3	1:A:195:ILE:CG1	2.33	0.52
1:A:319:ILE:HG23	3:A:868:HOH:O	2.08	0.52
1:A:249:THR:CG2	1:A:253:GLU:HB2	2.38	0.52
1:A:85:VAL:HG13	1:A:331:SER:OG	2.10	0.52
1:A:125:TYR:O	1:A:129:GLU:OE1	2.27	0.52
1:A:22:VAL:CG1	1:A:70:MET:HB2	2.39	0.52
1:B:240:SER:OG	1:B:262:ASN:ND2	2.42	0.52
1:B:309:VAL:HG11	1:B:311:TYR:CZ	2.44	0.52
1:B:229:ASN:ND2	1:B:230:ALA:N	2.57	0.52
1:A:125:TYR:O	1:A:126:THR:HG23	2.10	0.52
1:A:191:LYS:HZ2	1:A:191:LYS:HB3	1.74	0.51
1:B:150:ASN:O	1:B:152:THR:N	2.39	0.51
1:A:355:ARG:HA	1:B:121:PRO:HG3	1.92	0.51
1:B:93:VAL:HG21	1:B:261:LEU:HB2	1.92	0.51
1:B:29:ASN:ND2	3:B:711:HOH:O	2.37	0.51
1:A:179:SER:HA	3:A:790:HOH:O	2.09	0.51
1:A:347:LEU:O	1:A:351:GLU:CG	2.58	0.51
1:B:184:MET:HG2	1:B:185:PHE:H	1.75	0.51
1:A:146:LYS:HE3	3:A:808:HOH:O	2.11	0.50
1:A:67:THR:OG1	1:A:361:ASN:ND2	2.43	0.50
1:A:248:THR:HG22	1:A:249:THR:N	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:LEU:HD23	1:A:352:TYR:CE2	2.46	0.50
1:A:88:PRO:O	1:A:91:ASP:HB2	2.10	0.50
1:B:136:ILE:HG12	1:B:263:LEU:HD13	1.93	0.50
1:B:326:GLY:C	1:B:328:THR:N	2.64	0.50
1:B:343:LEU:HD11	1:B:347:LEU:HD11	1.94	0.50
1:B:155:SER:OG	1:B:244:HIS:HB2	2.12	0.50
1:A:109:THR:HG21	3:A:787:HOH:O	2.11	0.50
1:B:126:THR:HG23	1:B:129:GLU:OE1	2.12	0.50
1:B:53:ARG:HH12	1:B:63:ARG:HB3	1.74	0.50
1:A:68:PHE:HA	1:A:359:ILE:HD12	1.94	0.49
1:B:113:PHE:CD1	2:B:601:ADP:H2'	2.47	0.49
1:A:330:THR:HA	3:A:711:HOH:O	2.11	0.49
1:B:88:PRO:O	1:B:91:ASP:HB2	2.12	0.49
1:A:195:ILE:N	3:A:716:HOH:O	2.46	0.49
1:A:221:ARG:HG3	1:A:231:TYR:CE2	2.47	0.49
1:B:174:PRO:O	1:B:177:ASP:HB2	2.12	0.49
1:A:79:ILE:N	3:A:729:HOH:O	2.45	0.49
1:A:221:ARG:NH2	3:A:731:HOH:O	2.46	0.49
1:A:72:PHE:CD1	1:A:76:THR:HG21	2.47	0.49
1:A:358:ASN:ND2	1:B:121:PRO:CA	2.76	0.49
1:A:213:ILE:HD11	3:A:871:HOH:O	2.13	0.49
1:A:301:ALA:HB3	1:A:309:VAL:CG2	2.43	0.49
1:A:246:LYS:HA	1:A:255:LEU:O	2.13	0.48
1:B:309:VAL:O	1:B:312:ARG:NH2	2.45	0.48
1:A:249:THR:CG2	1:A:251:ASP:HB2	2.34	0.48
1:A:173:ASN:HB3	3:A:727:HOH:O	2.13	0.48
1:B:83:ARG:HD2	3:B:741:HOH:O	2.13	0.48
1:A:210:VAL:HA	3:A:856:HOH:O	2.13	0.48
1:B:294:THR:O	1:B:298:VAL:HG23	2.13	0.48
1:B:310:PRO:HA	1:B:312:ARG:NH2	2.28	0.48
1:B:141:HIS:HE1	3:B:719:HOH:O	1.96	0.48
1:B:82:TYR:C	3:B:703:HOH:O	2.47	0.48
1:A:176:SER:HB3	1:A:180:GLU:OE2	2.13	0.48
1:A:310:PRO:HA	1:A:312:ARG:NH2	2.29	0.48
1:A:355:ARG:HA	1:B:121:PRO:CG	2.44	0.48
1:B:136:ILE:HG12	1:B:263:LEU:CD1	2.44	0.48
1:A:339:ALA:HB1	3:A:764:HOH:O	2.14	0.47
1:A:38:HIS:N	1:A:38:HIS:HD2	2.04	0.47
1:B:167:GLU:HG2	1:B:181:ARG:HH11	1.79	0.47
1:B:314:SER:O	1:B:318:ARG:HG3	2.14	0.47
1:A:111:LYS:NZ	3:A:706:HOH:O	2.20	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:241:VAL:O	1:B:241:VAL:HG12	2.14	0.47
1:A:125:TYR:O	1:A:129:GLU:HB2	2.15	0.47
1:A:187:ASP:HB2	3:A:847:HOH:O	2.14	0.47
1:A:241:VAL:HG12	1:A:241:VAL:O	2.13	0.47
1:A:89:ILE:HG12	1:A:329:ARG:HB3	1.95	0.47
1:A:86:VAL:HB	3:A:741:HOH:O	2.14	0.47
1:B:127:TRP:N	1:B:145:GLU:OE1	2.45	0.47
1:B:181:ARG:HD2	3:B:743:HOH:O	2.14	0.47
1:A:212:GLN:HG3	3:A:783:HOH:O	2.15	0.47
1:A:53:ARG:HH12	1:A:63:ARG:HB3	1.77	0.47
1:B:167:GLU:HG2	1:B:181:ARG:NH1	2.30	0.47
1:B:150:ASN:C	1:B:152:THR:H	2.18	0.47
1:B:226:THR:HG22	1:B:227:LEU:CD2	2.42	0.47
1:B:57:LEU:O	1:B:58:ALA:CB	2.62	0.46
1:B:84:SER:C	3:B:703:HOH:O	2.47	0.46
1:A:314:SER:O	1:A:318:ARG:HG3	2.16	0.46
1:B:54:THR:O	1:B:59:ASP:HB3	2.15	0.46
1:B:95:MET:C	3:B:702:HOH:O	2.54	0.46
1:A:249:THR:HG21	1:A:253:GLU:HB2	1.98	0.46
1:B:122:ASN:HB3	1:B:125:TYR:CD2	2.51	0.46
1:A:309:VAL:O	1:A:312:ARG:NH2	2.46	0.46
1:B:78:GLN:HG3	1:B:132:LEU:HA	1.99	0.45
1:A:86:VAL:HG12	3:A:741:HOH:O	2.15	0.45
1:B:104:TYR:HE1	1:B:349:THR:HA	1.82	0.45
1:A:238:VAL:C	3:A:709:HOH:O	2.55	0.45
1:B:249:THR:CG2	1:B:253:GLU:HB2	2.46	0.45
1:B:325:GLY:C	1:B:327:ARG:N	2.65	0.45
1:B:154:PHE:HA	1:B:244:HIS:O	2.17	0.45
1:B:19:ILE:CD1	1:B:330:THR:HB	2.47	0.45
1:A:357:LYS:HD2	3:A:875:HOH:O	2.17	0.45
2:A:601:ADP:O2A	3:A:707:HOH:O	2.20	0.45
1:B:187:ASP:HB3	1:B:190:ASN:H	1.80	0.45
1:A:158:VAL:CG1	1:A:241:VAL:HG23	2.43	0.45
1:B:317:THR:CA	3:B:701:HOH:O	2.58	0.45
1:A:247:GLU:HG2	3:A:775:HOH:O	2.17	0.44
1:A:308:HIS:CG	1:A:308:HIS:O	2.67	0.44
1:A:312:ARG:N	1:A:312:ARG:NH1	2.61	0.44
1:A:238:VAL:CG1	1:A:238:VAL:O	2.65	0.44
1:A:238:VAL:O	1:A:239:PHE:C	2.55	0.44
1:B:98:ASN:HA	1:B:260:LYS:O	2.17	0.44
1:B:82:TYR:CD2	1:B:138:ARG:HD2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:ARG:HH11	1:A:63:ARG:HD3	1.82	0.44
1:B:151:GLY:O	3:B:704:HOH:O	2.21	0.44
1:B:167:GLU:CD	1:B:181:ARG:HD3	2.38	0.44
1:A:178:VAL:O	1:A:179:SER:CB	2.66	0.43
1:A:304:GLU:HB2	1:A:306:THR:HG23	2.00	0.43
1:A:309:VAL:HG12	1:A:309:VAL:O	2.17	0.43
1:B:297:ARG:HG2	1:B:297:ARG:HH11	1.83	0.43
1:A:241:VAL:O	1:A:241:VAL:CG1	2.66	0.43
1:B:58:ALA:HB1	1:B:60:LYS:CE	2.48	0.43
1:A:104:TYR:HE1	1:A:349:THR:HA	1.83	0.43
1:A:207:LYS:HG3	1:A:208:ASP:N	2.33	0.43
1:A:40:ILE:HD12	1:A:343:LEU:HD13	1.99	0.43
1:B:304:GLU:HB2	1:B:306:THR:HG23	2.01	0.43
1:A:27:PRO:HB3	1:A:74:ALA:HB1	2.01	0.43
1:A:38:HIS:CD2	1:A:340:SER:HG	2.37	0.43
1:B:191:LYS:HG3	1:B:192:ARG:HG3	1.99	0.43
1:B:33:ARG:HH11	1:B:33:ARG:HG2	1.84	0.43
1:B:83:ARG:C	3:B:703:HOH:O	2.56	0.43
1:B:68:PHE:HA	1:B:359:ILE:HD12	2.00	0.43
1:A:297:ARG:HH11	1:A:297:ARG:HG2	1.84	0.43
1:A:45:PRO:HB3	1:A:71:VAL:O	2.18	0.43
1:B:143:ILE:HD13	1:B:243:ILE:CD1	2.47	0.43
1:A:291:SER:HA	1:A:314:SER:HB2	2.01	0.43
1:B:309:VAL:HG12	1:B:311:TYR:CD2	2.54	0.43
1:B:312:ARG:NH1	1:B:312:ARG:N	2.63	0.43
1:A:309:VAL:HG12	1:A:311:TYR:CD2	2.54	0.42
1:B:72:PHE:CD1	1:B:76:THR:HG21	2.54	0.42
1:A:138:ARG:NH2	3:A:729:HOH:O	2.52	0.42
1:A:40:ILE:CD1	1:A:340:SER:HA	2.42	0.42
1:B:161:LEU:C	1:B:161:LEU:HD12	2.39	0.42
1:A:177:ASP:HB2	3:A:754:HOH:O	2.18	0.42
1:A:307:PRO:O	1:A:308:HIS:HB3	2.20	0.42
1:B:170:ASP:HB2	1:B:182:LEU:HD21	2.02	0.42
1:B:309:VAL:O	1:B:309:VAL:HG12	2.20	0.42
1:A:298:VAL:N	3:A:701:HOH:O	1.88	0.42
1:A:355:ARG:HA	1:B:121:PRO:CB	2.49	0.42
1:B:298:VAL:HG21	1:B:317:THR:HG21	2.01	0.42
1:A:24:ARG:NH1	1:A:114:THR:OG1	2.50	0.42
1:A:260:LYS:HE2	1:A:262:ASN:HD21	1.83	0.42
1:B:167:GLU:HB3	1:B:181:ARG:HD2	2.02	0.42
1:A:141:HIS:CE1	1:A:145:GLU:OE1	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:LEU:O	1:A:33:ARG:HB2	2.20	0.41
1:B:175:SER:C	1:B:177:ASP:N	2.73	0.41
1:B:241:VAL:O	1:B:241:VAL:CG1	2.69	0.41
1:B:325:GLY:O	1:B:327:ARG:N	2.48	0.41
1:B:39:SER:HA	1:B:338:PRO:O	2.20	0.41
1:A:136:ILE:CG1	1:A:263:LEU:HD13	2.44	0.41
1:B:221:ARG:HG3	1:B:231:TYR:HE2	1.85	0.41
1:A:82:TYR:CD2	1:A:138:ARG:HD2	2.55	0.41
1:A:87:CYS:N	3:A:741:HOH:O	2.53	0.41
1:B:45:PRO:HB3	1:B:71:VAL:O	2.19	0.41
1:A:164:TYR:CZ	1:A:228:MET:HB3	2.55	0.41
1:B:89:ILE:HG12	1:B:329:ARG:HB3	2.02	0.41
1:A:173:ASN:HA	1:A:174:PRO:HD3	1.89	0.41
1:B:126:THR:OG1	1:B:129:GLU:HG3	2.21	0.41
1:B:29:ASN:OD1	1:B:32:GLU:HG3	2.21	0.41
1:B:190:ASN:HB3	1:B:193:GLY:N	2.28	0.41
1:A:82:TYR:OH	1:A:139:THR:HA	2.20	0.41
1:A:338:PRO:CA	3:A:735:HOH:O	2.55	0.40
1:B:178:VAL:CG1	1:B:179:SER:H	2.20	0.40
1:A:170:ASP:HB2	1:A:182:LEU:HD21	2.03	0.40
1:A:298:VAL:HG13	1:A:309:VAL:CG1	2.51	0.40
1:A:358:ASN:ND2	1:B:121:PRO:HA	2.36	0.40
1:A:147:LEU:HB2	3:A:708:HOH:O	2.22	0.40
1:B:308:HIS:HE1	1:B:312:ARG:HE	1.68	0.40
1:A:111:LYS:HB2	2:A:601:ADP:O2B	2.20	0.40
1:B:140:LEU:HA	1:B:140:LEU:HD13	1.92	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	326/369 (88%)	290 (89%)	23 (7%)	13 (4%)	4	2
1	B	326/369 (88%)	296 (91%)	21 (6%)	9 (3%)	6	4
All	All	652/738 (88%)	586 (90%)	44 (7%)	22 (3%)	5	2

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	179	SER
1	A	239	PHE
1	B	226	THR
1	A	326	GLY
1	B	178	VAL
1	A	30	LEU
1	A	36	SER
1	A	125	TYR
1	A	126	THR
1	A	174	PRO
1	A	176	SER
1	A	188	PRO
1	A	191	LYS
1	B	29	ASN
1	B	30	LEU
1	B	58	ALA
1	B	151	GLY
1	A	29	ASN
1	A	124	GLU
1	B	174	PRO
1	B	185	PHE
1	B	55	GLY

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	290/323 (90%)	275 (95%)	15 (5%)	29	38

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	292/323 (90%)	273 (94%)	19 (6%)	21	27
All	All	582/646 (90%)	548 (94%)	34 (6%)	25	33

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	HIS
1	A	115	MET
1	A	141	HIS
1	A	152	THR
1	A	188	PRO
1	A	189	ARG
1	A	191	LYS
1	A	220	LYS
1	A	227	LEU
1	A	241	VAL
1	A	255	LEU
1	A	290	GLN
1	A	292	LEU
1	A	312	ARG
1	A	327	ARG
1	B	17	LYS
1	B	38	HIS
1	B	84	SER
1	B	115	MET
1	B	141	HIS
1	B	185	PHE
1	B	187	ASP
1	B	191	LYS
1	B	207	LYS
1	B	220	LYS
1	B	226	THR
1	B	229	ASN
1	B	241	VAL
1	B	255	LEU
1	B	292	LEU
1	B	305	ARG
1	B	309	VAL
1	B	312	ARG
1	B	327	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	GLN
1	A	38	HIS
1	A	106	GLN
1	A	165	ASN
1	A	183	GLN
1	A	229	ASN
1	A	262	ASN
1	A	290	GLN
1	A	321	GLN
1	A	358	ASN
1	A	361	ASN
1	B	20	GLN
1	B	38	HIS
1	B	106	GLN
1	B	141	HIS
1	B	183	GLN
1	B	190	ASN
1	B	206	ASN
1	B	229	ASN
1	B	244	HIS
1	B	262	ASN
1	B	321	GLN
1	B	361	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ADP	A	601	-	24,29,29	1.67	4 (16%)	23,45,45	3.15	3 (13%)
2	ADP	B	601	-	24,29,29	1.66	4 (16%)	23,45,45	3.23	3 (13%)

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	A	601	-	-	0/12/32/32	0/3/3/3
2	ADP	B	601	-	-	0/12/32/32	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	ADP	C5-C4	-2.99	1.33	1.40
2	B	601	ADP	C5-C4	-2.90	1.34	1.40
2	B	601	ADP	O4'-C1'	2.08	1.44	1.41
2	A	601	ADP	O4'-C1'	2.12	1.44	1.41
2	A	601	ADP	C2-N1	3.37	1.40	1.33
2	B	601	ADP	C2-N1	3.52	1.40	1.33
2	B	601	ADP	C2-N3	4.55	1.40	1.32
2	A	601	ADP	C2-N3	4.68	1.40	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	ADP	N3-C2-N1	-14.39	117.57	128.87
2	A	601	ADP	N3-C2-N1	-14.02	117.86	128.87
2	A	601	ADP	C4'-O4'-C1'	-2.28	107.23	109.64
2	A	601	ADP	C2'-C3'-C4'	-2.18	98.17	102.64
2	B	601	ADP	C4'-O4'-C1'	-2.16	107.36	109.64
2	B	601	ADP	O4'-C1'-N9	2.69	113.18	108.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	ADP	3	0
2	B	601	ADP	1	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	330/369 (89%)	0.02	21 (6%)	23 31	9, 28, 96, 200	0
1	B	330/369 (89%)	0.23	30 (9%)	11 17	10, 29, 109, 188	0
All	All	660/738 (89%)	0.13	51 (7%)	16 23	9, 28, 105, 200	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	57	LEU	18.0
1	B	56	GLY	17.7
1	A	58	ALA	10.2
1	B	58	ALA	9.7
1	A	57	LEU	8.8
1	A	56	GLY	8.7
1	B	36	SER	8.3
1	B	34	LYS	7.4
1	B	16	GLY	6.9
1	A	176	SER	6.9
1	B	59	ASP	6.2
1	B	31	ALA	6.1
1	A	59	ASP	5.9
1	B	35	ALA	5.5
1	A	327	ARG	5.5
1	B	30	LEU	5.4
1	A	178	VAL	4.7
1	B	191	LYS	4.5
1	A	16	GLY	4.2
1	B	62	SER	4.1
1	B	37	ALA	4.0
1	B	177	ASP	3.6
1	A	325	GLY	3.6
1	A	61	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	188	PRO	3.3
1	B	308	HIS	3.3
1	B	305	ARG	3.3
1	A	305	ARG	3.3
1	B	55	GLY	3.0
1	A	250	ILE	3.0
1	B	189	ARG	3.0
1	A	55	GLY	2.9
1	A	30	LEU	2.8
1	A	29	ASN	2.8
1	A	124	GLU	2.8
1	B	150	ASN	2.6
1	B	32	GLU	2.6
1	A	191	LYS	2.6
1	B	190	ASN	2.5
1	A	270	GLU	2.5
1	B	364	GLU	2.3
1	B	38	HIS	2.3
1	B	61	SER	2.2
1	B	270	GLU	2.2
1	B	185	PHE	2.2
1	A	63	ARG	2.1
1	A	189	ARG	2.1
1	B	176	SER	2.0
1	B	250	ILE	2.0
1	B	151	GLY	2.0
1	B	326	GLY	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	A	601	27/27	0.89	0.16	1.96	32,38,43,49	0
2	ADP	B	601	27/27	0.92	0.13	-0.00	37,43,49,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.