Full wwPDB X-ray Structure Validation Report (i)

Feb 4, 2024 – 02:43 AM EST

PDB ID	:	1Q1E
Title	:	The ATPase component of E. coli maltose transporter (MalK) in the nucleotide-free form
Authors	:	Chen, J.; Lu, G.; Lin, J.; Davidson, A.L.; Quiocho, F.A.
Deposited on	:	2003-07-19
Resolution	:	2.90 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

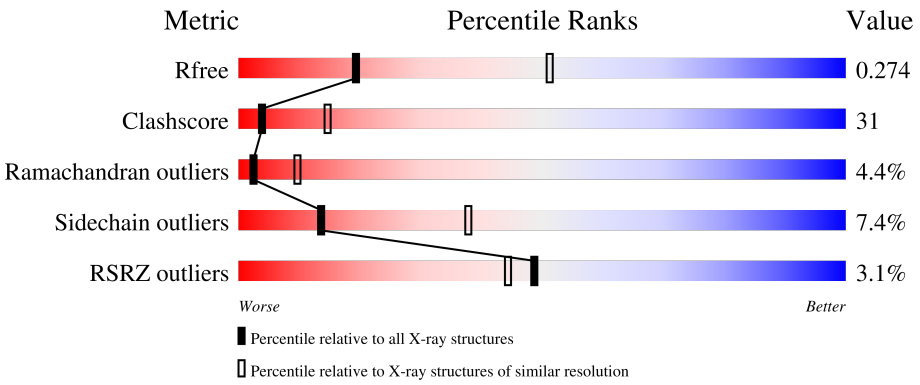
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 <=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	381	3% 52% 35% 7% • •
1	B	381	3% 54% 35% 6% • •

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 5706 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 Maltose/maltodextrin transport ATP-binding protein malK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	367	2853	1805	511	524	13	0	0	0
1	B	367	2853	1805	511	524	13	0	0	0

There are 20 discrepancies between the modelled and reference sequences:

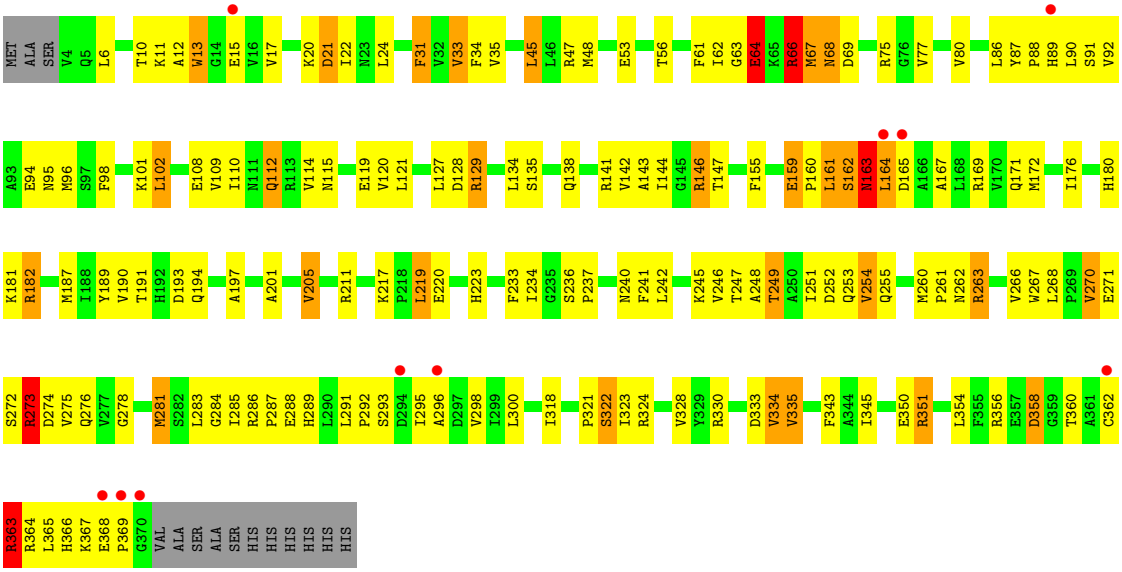
Chain	Residue	Modelled	Actual	Comment	Reference
A	372	ALA	-	expression tag	UNP P68187
A	373	SER	-	expression tag	UNP P68187
A	374	ALA	-	expression tag	UNP P68187
A	375	SER	-	expression tag	UNP P68187
A	376	HIS	-	expression tag	UNP P68187
A	377	HIS	-	expression tag	UNP P68187
A	378	HIS	-	expression tag	UNP P68187
A	379	HIS	-	expression tag	UNP P68187
A	380	HIS	-	expression tag	UNP P68187
A	381	HIS	-	expression tag	UNP P68187
B	372	ALA	-	expression tag	UNP P68187
B	373	SER	-	expression tag	UNP P68187
B	374	ALA	-	expression tag	UNP P68187
B	375	SER	-	expression tag	UNP P68187
B	376	HIS	-	expression tag	UNP P68187
B	377	HIS	-	expression tag	UNP P68187
B	378	HIS	-	expression tag	UNP P68187
B	379	HIS	-	expression tag	UNP P68187
B	380	HIS	-	expression tag	UNP P68187
B	381	HIS	-	expression tag	UNP P68187

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.

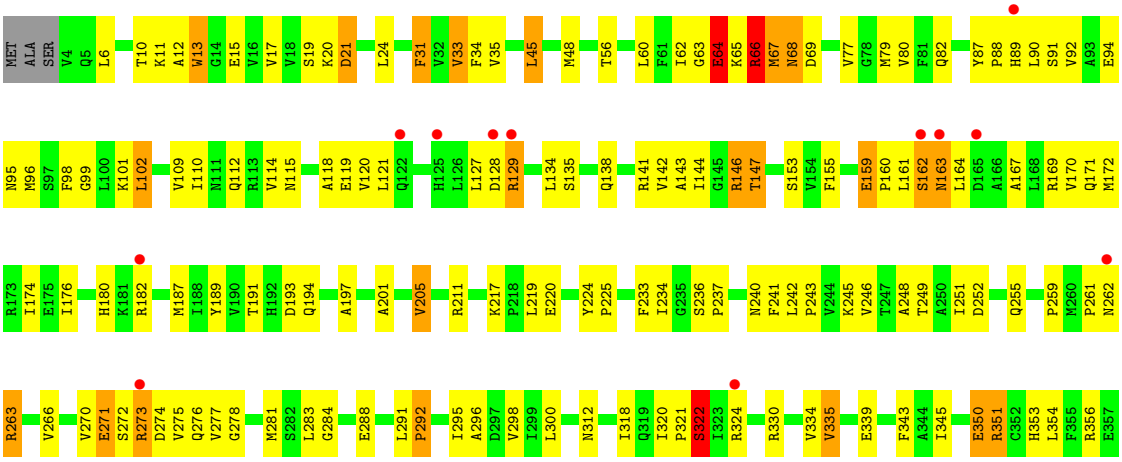
• Molecule 1: Maltose/maltodextrin transport ATP-binding protein malK

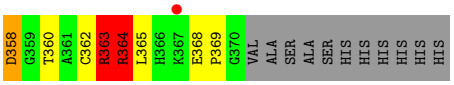
Chain A:



• Molecule 1: Maltose/maltodextrin transport ATP-binding protein malK

Chain B:




D358
G359
T360
A361
C362
R363
R364
L365
H366
K367
E368
P369
G370
VAL
ALA
SER
ALA
SER
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	64.61Å 102.04Å 99.29Å 90.00° 107.10° 90.00°	Depositor
Resolution (Å)	15.78 – 2.90 15.78 – 2.86	Depositor EDS
% Data completeness (in resolution range)	98.2 (15.78-2.90) 96.0 (15.78-2.86)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.76 (at 2.86Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.244 , 0.276 0.244 , 0.274	Depositor DCC
R_{free} test set	2174 reflections (7.79%)	wwPDB-VP
Wilson B-factor (Å^2)	44.3	Xtriage
Anisotropy	0.342	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 47.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	5706	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.68% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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)

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	0.59	0/2903	0.89	11/3936 (0.3%)
1	B	0.47	0/2903	0.79	7/3936 (0.2%)
All	All	0.54	0/5806	0.84	18/7872 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	363	ARG	O-C-N	8.22	135.85	122.70
1	A	364	ARG	NE-CZ-NH2	7.67	124.13	120.30
1	A	363	ARG	NE-CZ-NH2	7.60	124.10	120.30
1	B	363	ARG	NE-CZ-NH2	7.49	124.05	120.30
1	A	356	ARG	NE-CZ-NH2	7.42	124.01	120.30
1	A	324	ARG	NE-CZ-NH2	7.30	123.95	120.30
1	A	351	ARG	NE-CZ-NH2	7.21	123.90	120.30
1	B	364	ARG	NE-CZ-NH2	6.94	123.77	120.30
1	B	324	ARG	NE-CZ-NH2	6.93	123.77	120.30
1	A	66	ARG	NE-CZ-NH2	6.67	123.63	120.30
1	A	273	ARG	NE-CZ-NH2	6.65	123.63	120.30
1	A	330	ARG	NE-CZ-NH2	6.63	123.61	120.30
1	A	286	ARG	NE-CZ-NH2	6.62	123.61	120.30
1	B	66	ARG	NE-CZ-NH2	6.62	123.61	120.30
1	B	363	ARG	CA-C-N	-6.37	103.20	117.20
1	B	67	MET	CG-SD-CE	6.11	109.98	100.20
1	A	281	MET	CG-SD-CE	5.78	109.45	100.20
1	A	67	MET	CG-SD-CE	5.69	109.30	100.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	2853	0	2918	184	0
1	B	2853	0	2918	179	0
All	All	5706	0	5836	354	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 31.

All (354) 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:102:LEU:H	1:B:102:LEU:CD2	1.56	1.19
1:A:367:LYS:O	1:A:368:GLU:HG2	1.50	1.09
1:A:67:MET:CE	1:A:75:ARG:HA	1.83	1.08
1:A:369:PRO:CG	1:B:334:VAL:HG11	1.83	1.07
1:A:281:MET:HE3	1:A:354:LEU:HD21	1.30	1.06
1:B:321:PRO:O	1:B:322:SER:OG	1.76	1.04
1:A:354:LEU:HB3	1:A:362:CYS:SG	2.01	1.01
1:A:368:GLU:HB3	1:A:369:PRO:HD2	1.40	0.99
1:A:369:PRO:CB	1:B:334:VAL:HG11	1.94	0.98
1:B:102:LEU:CD2	1:B:102:LEU:N	2.27	0.97
1:A:368:GLU:HB3	1:A:369:PRO:CD	1.97	0.94
1:B:102:LEU:H	1:B:102:LEU:HD22	1.30	0.94
1:B:102:LEU:N	1:B:102:LEU:HD23	1.80	0.94
1:A:87:TYR:H	1:A:95:ASN:HD21	1.14	0.93
1:B:87:TYR:H	1:B:95:ASN:HD21	1.14	0.93
1:B:164:LEU:HD23	1:B:169:ARG:NE	1.83	0.93
1:B:102:LEU:H	1:B:102:LEU:HD23	1.34	0.92
1:B:68:ASN:H	1:B:68:ASN:ND2	1.67	0.90
1:A:67:MET:HE1	1:A:75:ARG:HA	1.54	0.90
1:B:68:ASN:H	1:B:68:ASN:HD22	0.90	0.88
1:A:270:VAL:HG13	1:A:362:CYS:O	1.72	0.88
1:B:68:ASN:HD22	1:B:68:ASN:N	1.73	0.85
1:B:356:ARG:CZ	1:B:362:CYS:HB3	2.06	0.85
1:B:164:LEU:HD11	1:B:193:ASP:CB	2.07	0.84
1:A:289:HIS:CE1	1:A:351:ARG:HH21	1.96	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:362:CYS:O	1:B:362:CYS:SG	2.37	0.83
1:B:33:VAL:CG1	1:B:201:ALA:HB2	2.09	0.82
1:A:164:LEU:HD11	1:A:193:ASP:HB2	1.62	0.81
1:B:10:THR:HG22	1:B:21:ASP:H	1.46	0.81
1:B:68:ASN:O	1:B:69:ASP:HB2	1.81	0.81
1:A:10:THR:HG22	1:A:21:ASP:H	1.46	0.80
1:A:366:HIS:ND1	1:A:367:LYS:N	2.30	0.80
1:B:164:LEU:CD2	1:B:169:ARG:NE	2.45	0.80
1:A:33:VAL:CG1	1:A:201:ALA:HB2	2.11	0.80
1:A:321:PRO:O	1:A:322:SER:OG	1.97	0.80
1:B:270:VAL:HG22	1:B:363:ARG:CB	2.11	0.79
1:A:246:VAL:CG2	1:A:281:MET:CE	2.61	0.79
1:A:369:PRO:CG	1:B:334:VAL:CG1	2.61	0.79
1:A:246:VAL:CG2	1:A:281:MET:HE2	2.12	0.78
1:B:164:LEU:HD11	1:B:193:ASP:HB2	1.65	0.78
1:A:246:VAL:HG23	1:A:281:MET:HE2	1.65	0.78
1:A:252:ASP:HB3	1:A:365:LEU:CD1	2.12	0.78
1:A:80:VAL:HA	1:A:147:THR:HG21	1.65	0.78
1:A:35:VAL:HG11	1:A:197:ALA:HB2	1.65	0.77
1:A:369:PRO:HG2	1:B:334:VAL:CG1	2.14	0.77
1:A:12:ALA:H	1:A:56:THR:HG21	1.49	0.77
1:A:62:ILE:HB	1:A:67:MET:HG3	1.65	0.77
1:A:366:HIS:CE1	1:A:367:LYS:O	2.38	0.77
1:A:120:VAL:HG23	1:A:121:LEU:HD13	1.68	0.76
1:A:67:MET:CE	1:A:75:ARG:CA	2.63	0.75
1:B:87:TYR:H	1:B:95:ASN:ND2	1.83	0.75
1:B:87:TYR:N	1:B:95:ASN:HD21	1.83	0.75
1:B:281:MET:HE1	1:B:354:LEU:HD21	1.69	0.74
1:B:120:VAL:HG23	1:B:121:LEU:HD13	1.69	0.74
1:B:35:VAL:HG11	1:B:197:ALA:HB2	1.69	0.74
1:B:80:VAL:HA	1:B:147:THR:HG21	1.69	0.74
1:A:33:VAL:HG11	1:A:201:ALA:HB2	1.71	0.73
1:B:12:ALA:H	1:B:56:THR:HG21	1.54	0.73
1:A:295:ILE:HD12	1:A:295:ILE:N	2.03	0.73
1:B:164:LEU:HD23	1:B:169:ARG:CD	2.18	0.73
1:A:87:TYR:N	1:A:95:ASN:HD21	1.86	0.73
1:B:295:ILE:N	1:B:295:ILE:HD12	2.04	0.73
1:A:281:MET:CE	1:A:354:LEU:HD21	2.13	0.72
1:B:354:LEU:H	1:B:363:ARG:HD3	1.55	0.72
1:A:369:PRO:HG2	1:B:334:VAL:HG11	1.65	0.72
1:B:194:GLN:HG3	1:B:234:ILE:HA	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:366:HIS:ND1	1:A:367:LYS:O	2.23	0.72
1:B:270:VAL:HG22	1:B:363:ARG:HB2	1.72	0.71
1:B:252:ASP:HB3	1:B:365:LEU:HD11	1.71	0.71
1:B:217:LYS:HB2	1:B:220:GLU:HG3	1.73	0.71
1:B:271:GLU:H	1:B:363:ARG:HA	1.54	0.71
1:B:225:PRO:HG2	1:B:363:ARG:HH21	1.56	0.70
1:B:164:LEU:CD2	1:B:169:ARG:CZ	2.70	0.70
1:A:164:LEU:HD11	1:A:193:ASP:CB	2.21	0.70
1:B:99:GLY:HA2	1:B:102:LEU:HD21	1.72	0.70
1:A:163:ASN:O	1:A:165:ASP:N	2.20	0.69
1:A:87:TYR:H	1:A:95:ASN:ND2	1.88	0.68
1:A:291:LEU:HB3	1:A:292:PRO:HD2	1.74	0.68
1:A:271:GLU:N	1:A:363:ARG:O	2.24	0.67
1:A:295:ILE:HD12	1:A:295:ILE:H	1.59	0.67
1:B:92:VAL:O	1:B:96:MET:HG3	1.95	0.66
1:B:358:ASP:OD2	1:B:358:ASP:N	2.24	0.66
1:A:68:ASN:O	1:A:69:ASP:HB2	1.95	0.66
1:A:271:GLU:HG3	1:A:273:ARG:HG3	1.78	0.66
1:B:33:VAL:HG11	1:B:201:ALA:HB2	1.77	0.66
1:A:68:ASN:OD1	1:A:68:ASN:N	2.26	0.65
1:A:266:VAL:HG21	1:A:298:VAL:HG21	1.79	0.64
1:B:10:THR:HG22	1:B:21:ASP:N	2.12	0.64
1:B:172:MET:O	1:B:176:ILE:HG13	1.96	0.64
1:B:281:MET:HE1	1:B:354:LEU:CD2	2.27	0.64
1:A:275:VAL:O	1:A:276:GLN:HB2	1.96	0.64
1:B:164:LEU:HD21	1:B:169:ARG:CZ	2.27	0.64
1:B:63:GLY:O	1:B:64:GLU:HB2	1.97	0.64
1:A:194:GLN:HG3	1:A:234:ILE:HA	1.79	0.63
1:B:110:ILE:O	1:B:114:VAL:HG23	1.96	0.63
1:A:236:SER:HA	1:A:237:PRO:C	2.18	0.63
1:A:271:GLU:HB2	1:A:363:ARG:HB3	1.80	0.63
1:B:295:ILE:HD12	1:B:295:ILE:H	1.62	0.63
1:B:135:SER:OG	1:B:138:GLN:HG2	1.97	0.63
1:B:164:LEU:HD23	1:B:169:ARG:HD3	1.80	0.62
1:B:134:LEU:HA	1:B:138:GLN:HE21	1.64	0.62
1:A:217:LYS:HB2	1:A:220:GLU:HG3	1.79	0.62
1:A:20:LYS:HG3	1:A:211:ARG:HG2	1.81	0.62
1:A:67:MET:HE2	1:A:75:ARG:CB	2.28	0.62
1:B:87:TYR:CB	1:B:90:LEU:HD12	2.30	0.62
1:A:246:VAL:CG2	1:A:281:MET:HE1	2.28	0.62
1:A:358:ASP:OD2	1:A:358:ASP:N	2.29	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:270:VAL:HG13	1:B:362:CYS:O	1.99	0.62
1:A:368:GLU:CB	1:A:369:PRO:HD2	2.25	0.62
1:B:60:LEU:H	1:B:68:ASN:HD21	1.47	0.62
1:B:236:SER:HA	1:B:237:PRO:C	2.20	0.62
1:A:10:THR:HG22	1:A:21:ASP:N	2.13	0.61
1:A:367:LYS:C	1:A:368:GLU:HG2	2.19	0.61
1:A:367:LYS:O	1:A:368:GLU:CG	2.39	0.61
1:B:161:LEU:H	1:B:161:LEU:HD12	1.66	0.61
1:A:334:VAL:O	1:A:335:VAL:HB	1.99	0.60
1:B:271:GLU:HG3	1:B:273:ARG:HG3	1.83	0.60
1:A:223:HIS:CE1	1:A:368:GLU:OE1	2.55	0.60
1:A:62:ILE:HD13	1:A:77:VAL:CG1	2.31	0.60
1:A:162:SER:HB2	1:A:169:ARG:HB2	1.83	0.60
1:B:164:LEU:HD23	1:B:169:ARG:CZ	2.30	0.60
1:B:62:ILE:HB	1:B:67:MET:HG3	1.83	0.60
1:B:164:LEU:CD1	1:B:193:ASP:HB2	2.31	0.60
1:B:270:VAL:HG22	1:B:363:ARG:HB3	1.82	0.59
1:B:248:ALA:HB3	1:B:255:GLN:HB3	1.85	0.59
1:A:63:GLY:O	1:A:64:GLU:HB2	2.01	0.59
1:B:252:ASP:HB3	1:B:365:LEU:CD1	2.33	0.59
1:A:11:LYS:HA	1:A:56:THR:HG23	1.85	0.59
1:A:253:GLN:O	1:A:268:LEU:O	2.21	0.59
1:A:322:SER:O	1:A:323:ILE:HG13	2.03	0.58
1:A:334:VAL:HG23	1:A:335:VAL:H	1.67	0.58
1:B:45:LEU:HA	1:B:48:MET:HE2	1.84	0.58
1:B:353:HIS:HD2	1:B:363:ARG:NH2	2.01	0.58
1:A:172:MET:O	1:A:176:ILE:HG13	2.03	0.58
1:A:161:LEU:HD12	1:A:161:LEU:H	1.67	0.58
1:A:366:HIS:CG	1:A:367:LYS:H	2.21	0.58
1:A:67:MET:HE2	1:A:75:ARG:CA	2.33	0.58
1:A:368:GLU:CB	1:A:369:PRO:CD	2.66	0.58
1:A:102:LEU:HD13	1:A:102:LEU:H	1.69	0.58
1:B:363:ARG:HG3	1:B:363:ARG:HH11	1.69	0.57
1:A:102:LEU:CD1	1:A:102:LEU:N	2.67	0.57
1:B:35:VAL:CG1	1:B:197:ALA:HB2	2.35	0.57
1:B:356:ARG:CZ	1:B:362:CYS:CB	2.81	0.57
1:A:120:VAL:HG12	1:A:182:ARG:HH11	1.69	0.57
1:B:164:LEU:CD1	1:B:193:ASP:CB	2.81	0.57
1:B:363:ARG:HG3	1:B:363:ARG:NH1	2.19	0.57
1:A:134:LEU:HA	1:A:138:GLN:HE21	1.69	0.57
1:A:155:PHE:HB3	1:A:187:MET:HG2	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:VAL:HG22	1:A:281:MET:CE	2.34	0.57
1:B:334:VAL:HG23	1:B:335:VAL:H	1.70	0.56
1:A:252:ASP:HB3	1:A:365:LEU:HD11	1.86	0.56
1:A:159:GLU:HG2	1:A:191:THR:HA	1.86	0.56
1:B:275:VAL:O	1:B:276:GLN:HB2	2.05	0.56
1:B:167:ALA:O	1:B:171:GLN:HG2	2.05	0.56
1:B:155:PHE:HB3	1:B:187:MET:HG2	1.87	0.56
1:A:91:SER:HB2	1:A:128:ASP:HA	1.87	0.56
1:B:283:LEU:C	1:B:283:LEU:HD23	2.25	0.56
1:B:68:ASN:O	1:B:69:ASP:CB	2.50	0.56
1:A:233:PHE:HD2	1:A:234:ILE:HD12	1.70	0.55
1:A:287:PRO:HB3	1:A:328:VAL:O	2.07	0.55
1:A:31:PHE:HB2	1:A:180:HIS:CE1	2.42	0.55
1:A:159:GLU:CG	1:A:191:THR:HA	2.37	0.55
1:A:12:ALA:H	1:A:56:THR:CG2	2.19	0.55
1:A:110:ILE:O	1:A:114:VAL:HG23	2.07	0.54
1:B:271:GLU:O	1:B:272:SER:HB3	2.06	0.54
1:B:291:LEU:HB3	1:B:292:PRO:HD2	1.88	0.54
1:B:12:ALA:O	1:B:13:TRP:HE3	1.91	0.54
1:B:88:PRO:O	1:B:89:HIS:HB2	2.08	0.54
1:B:91:SER:HB2	1:B:128:ASP:HA	1.90	0.54
1:A:292:PRO:O	1:A:293:SER:HB3	2.08	0.54
1:B:363:ARG:NH1	1:B:364:ARG:HB2	2.22	0.54
1:A:12:ALA:O	1:A:13:TRP:HE3	1.91	0.54
1:A:91:SER:OG	1:A:94:GLU:HG3	2.08	0.54
1:A:88:PRO:O	1:A:89:HIS:HB2	2.08	0.53
1:A:135:SER:OG	1:A:138:GLN:HG2	2.07	0.53
1:A:248:ALA:HB3	1:A:255:GLN:HB3	1.89	0.53
1:A:20:LYS:HB3	1:A:211:ARG:NH1	2.24	0.53
1:B:62:ILE:HD13	1:B:77:VAL:CG1	2.38	0.53
1:B:20:LYS:HG3	1:B:211:ARG:HG2	1.91	0.53
1:A:272:SER:O	1:A:273:ARG:C	2.46	0.53
1:A:289:HIS:CE1	1:A:351:ARG:NH2	2.71	0.53
1:B:162:SER:HB2	1:B:169:ARG:HB2	1.91	0.53
1:A:261:PRO:O	1:A:262:ASN:CG	2.46	0.53
1:A:366:HIS:CG	1:A:367:LYS:N	2.76	0.53
1:B:356:ARG:NH2	1:B:362:CYS:CB	2.72	0.53
1:A:162:SER:CB	1:A:169:ARG:HB2	2.38	0.52
1:A:17:VAL:HG11	1:A:20:LYS:HD3	1.91	0.52
1:A:354:LEU:CB	1:A:362:CYS:SG	2.87	0.52
1:A:367:LYS:C	1:A:368:GLU:CG	2.77	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:LEU:HD13	1:A:102:LEU:N	2.24	0.52
1:B:20:LYS:O	1:B:21:ASP:C	2.48	0.52
1:B:87:TYR:HB2	1:B:90:LEU:HD12	1.91	0.52
1:B:138:GLN:HB3	1:B:141:ARG:HH21	1.74	0.52
1:B:35:VAL:HG11	1:B:197:ALA:CB	2.39	0.52
1:A:295:ILE:H	1:A:295:ILE:CD1	2.22	0.52
1:B:170:VAL:O	1:B:174:ILE:HG13	2.10	0.52
1:B:163:ASN:O	1:B:169:ARG:NH1	2.43	0.52
1:B:350:GLU:CD	1:B:350:GLU:H	2.13	0.52
1:A:260:MET:HA	1:A:322:SER:OG	2.10	0.51
1:A:318:ILE:HD11	1:A:343:PHE:HD2	1.76	0.51
1:B:6:LEU:HD22	1:B:24:LEU:HD12	1.91	0.51
1:B:12:ALA:H	1:B:56:THR:CG2	2.22	0.51
1:B:79:MET:HG3	1:B:80:VAL:N	2.24	0.51
1:A:368:GLU:HB3	1:A:369:PRO:HD3	1.91	0.51
1:B:87:TYR:HB3	1:B:90:LEU:HD12	1.93	0.51
1:A:321:PRO:O	1:A:322:SER:CB	2.59	0.51
1:B:295:ILE:H	1:B:295:ILE:CD1	2.24	0.51
1:B:353:HIS:HD2	1:B:363:ARG:CZ	2.23	0.51
1:A:350:GLU:H	1:A:350:GLU:CD	2.14	0.51
1:B:33:VAL:HA	1:B:189:TYR:O	2.11	0.50
1:B:91:SER:HA	1:B:129:ARG:O	2.12	0.50
1:B:295:ILE:N	1:B:295:ILE:CD1	2.73	0.50
1:B:281:MET:CE	1:B:354:LEU:HD21	2.39	0.50
1:B:246:VAL:HG12	1:B:277:VAL:HA	1.94	0.50
1:B:11:LYS:HA	1:B:56:THR:HG23	1.93	0.50
1:B:164:LEU:HA	1:B:169:ARG:HD3	1.94	0.50
1:A:160:PRO:C	1:A:162:SER:H	2.15	0.50
1:B:351:ARG:NH1	1:B:368:GLU:CD	2.65	0.50
1:A:92:VAL:HG22	1:A:134:LEU:HD11	1.93	0.50
1:B:17:VAL:HG11	1:B:20:LYS:HD3	1.94	0.49
1:B:241:PHE:C	1:B:242:LEU:HD23	2.32	0.49
1:A:283:LEU:C	1:A:283:LEU:HD23	2.31	0.49
1:A:67:MET:HE2	1:A:75:ARG:HB3	1.94	0.49
1:B:288:GLU:HG3	1:B:330:ARG:HD3	1.95	0.49
1:A:247:THR:OG1	1:A:255:GLN:HG2	2.12	0.49
1:A:354:LEU:HB3	1:A:362:CYS:HG	1.74	0.49
1:A:369:PRO:HB2	1:B:334:VAL:HG11	1.86	0.49
1:A:253:GLN:O	1:A:254:VAL:HB	2.13	0.48
1:B:65:LYS:HG3	1:B:67:MET:CE	2.43	0.48
1:B:164:LEU:HD21	1:B:169:ARG:NH2	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ILE:HD11	1:A:172:MET:CE	2.44	0.48
1:A:91:SER:HA	1:A:129:ARG:O	2.14	0.48
1:A:164:LEU:HD11	1:A:193:ASP:CG	2.34	0.48
1:A:138:GLN:O	1:A:142:VAL:HG23	2.12	0.48
1:B:68:ASN:ND2	1:B:68:ASN:N	2.42	0.48
1:A:263:ARG:HG3	1:A:263:ARG:HH11	1.78	0.48
1:B:143:ALA:O	1:B:146:ARG:HB3	2.13	0.48
1:B:351:ARG:NH1	1:B:368:GLU:OE2	2.47	0.48
1:A:35:VAL:HG11	1:A:197:ALA:CB	2.42	0.48
1:A:10:THR:HG22	1:A:21:ASP:CA	2.43	0.47
1:B:87:TYR:O	1:B:90:LEU:HB2	2.14	0.47
1:B:164:LEU:HD11	1:B:193:ASP:HB3	1.94	0.47
1:B:334:VAL:O	1:B:335:VAL:HB	2.14	0.47
1:A:271:GLU:O	1:A:272:SER:HB3	2.14	0.47
1:B:45:LEU:HA	1:B:48:MET:CE	2.44	0.47
1:B:92:VAL:HG22	1:B:134:LEU:HD11	1.96	0.47
1:A:275:VAL:O	1:A:276:GLN:CB	2.61	0.47
1:B:138:GLN:O	1:B:142:VAL:HG23	2.15	0.47
1:B:261:PRO:O	1:B:262:ASN:CG	2.53	0.47
1:B:354:LEU:H	1:B:363:ARG:CD	2.25	0.47
1:B:321:PRO:C	1:B:322:SER:OG	2.47	0.47
1:B:233:PHE:HD2	1:B:234:ILE:HD12	1.79	0.47
1:B:92:VAL:HB	1:B:127:LEU:O	2.15	0.46
1:B:243:PRO:O	1:B:259:PRO:HD3	2.15	0.46
1:A:35:VAL:HG12	1:A:191:THR:HG23	1.97	0.46
1:B:10:THR:HG22	1:B:21:ASP:CA	2.45	0.46
1:A:34:PHE:HD2	1:A:205:VAL:HG13	1.80	0.46
1:A:266:VAL:CG2	1:A:298:VAL:HG21	2.44	0.46
1:B:270:VAL:HG22	1:B:364:ARG:H	1.80	0.46
1:A:252:ASP:HA	1:A:272:SER:HB3	1.97	0.46
1:B:318:ILE:HD11	1:B:343:PHE:HD2	1.81	0.46
1:A:20:LYS:O	1:A:21:ASP:C	2.54	0.46
1:B:356:ARG:NH2	1:B:362:CYS:HB2	2.31	0.46
1:A:255:GLN:HB2	1:A:267:TRP:CD2	2.51	0.46
1:B:134:LEU:CA	1:B:138:GLN:HE21	2.28	0.46
1:A:98:PHE:O	1:A:101:LYS:HB2	2.15	0.46
1:B:266:VAL:HG21	1:B:298:VAL:HG21	1.97	0.46
1:A:67:MET:HE3	1:A:75:ARG:HG2	1.97	0.45
1:B:114:VAL:O	1:B:118:ALA:HB2	2.15	0.45
1:A:35:VAL:CG1	1:A:197:ALA:HB2	2.41	0.45
1:A:86:LEU:HD12	1:A:95:ASN:ND2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:VAL:HB	1:A:127:LEU:O	2.15	0.45
1:A:246:VAL:HG22	1:A:281:MET:HE1	1.94	0.45
1:B:270:VAL:CG2	1:B:363:ARG:HB3	2.45	0.45
1:A:87:TYR:CB	1:A:90:LEU:HD12	2.47	0.45
1:B:159:GLU:HG2	1:B:191:THR:HA	1.98	0.45
1:A:144:ILE:HD11	1:A:172:MET:HE3	1.98	0.45
1:B:98:PHE:O	1:B:101:LYS:HB2	2.17	0.45
1:B:162:SER:CB	1:B:169:ARG:HB2	2.47	0.45
1:B:363:ARG:HH11	1:B:363:ARG:CG	2.30	0.45
1:A:270:VAL:CG1	1:A:362:CYS:O	2.56	0.44
1:B:272:SER:O	1:B:273:ARG:C	2.56	0.44
1:A:134:LEU:CA	1:A:138:GLN:HE21	2.30	0.44
1:A:138:GLN:HB3	1:A:141:ARG:HH21	1.82	0.44
1:B:91:SER:OG	1:B:94:GLU:HG3	2.18	0.44
1:B:144:ILE:HD11	1:B:172:MET:CE	2.48	0.44
1:A:285:ILE:HG23	1:A:285:ILE:O	2.17	0.44
1:B:252:ASP:HA	1:B:272:SER:HB3	2.00	0.44
1:B:263:ARG:HH11	1:B:263:ARG:HG3	1.83	0.44
1:A:33:VAL:HA	1:A:189:TYR:O	2.17	0.44
1:B:163:ASN:O	1:B:169:ARG:HD3	2.16	0.44
1:A:245:LYS:HE3	1:A:278:GLY:HA2	1.99	0.44
1:A:288:GLU:HG2	1:B:312:ASN:HB2	2.00	0.43
1:B:261:PRO:HD3	1:B:322:SER:OG	2.18	0.43
1:A:17:VAL:CG1	1:A:20:LYS:HD3	2.48	0.43
1:A:67:MET:CE	1:A:75:ARG:HG2	2.48	0.43
1:B:92:VAL:CG2	1:B:134:LEU:HD11	2.48	0.43
1:B:109:VAL:HG23	1:B:110:ILE:N	2.33	0.43
1:A:45:LEU:HA	1:A:48:MET:HE2	2.00	0.43
1:A:241:PHE:C	1:A:242:LEU:HD23	2.38	0.43
1:A:115:ASN:O	1:A:119:GLU:HG3	2.19	0.43
1:A:167:ALA:O	1:A:171:GLN:HG2	2.18	0.43
1:B:34:PHE:HD2	1:B:205:VAL:HG13	1.84	0.43
1:B:31:PHE:HB2	1:B:180:HIS:CE1	2.53	0.43
1:B:115:ASN:O	1:B:119:GLU:HG3	2.18	0.43
1:A:271:GLU:CG	1:A:273:ARG:HG3	2.47	0.43
1:B:66:ARG:HE	1:B:66:ARG:HB3	1.63	0.43
1:B:134:LEU:HA	1:B:138:GLN:NE2	2.33	0.43
1:B:240:ASN:O	1:B:284:GLY:HA2	2.19	0.43
1:A:87:TYR:O	1:A:90:LEU:HB2	2.19	0.42
1:A:249:THR:O	1:A:249:THR:HG22	2.20	0.42
1:A:295:ILE:N	1:A:295:ILE:CD1	2.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:ARG:NH1	1:A:53:GLU:OE2	2.52	0.42
1:A:87:TYR:HB2	1:A:90:LEU:HD12	2.01	0.42
1:A:159:GLU:HG2	1:A:190:VAL:O	2.19	0.42
1:A:143:ALA:O	1:A:146:ARG:HB3	2.19	0.42
1:A:300:LEU:HB2	1:A:345:ILE:HD12	2.01	0.42
1:A:369:PRO:HG3	1:B:334:VAL:HG11	1.88	0.42
1:B:233:PHE:HD2	1:B:234:ILE:CD1	2.32	0.42
1:B:242:LEU:HD23	1:B:242:LEU:N	2.34	0.42
1:A:6:LEU:HD22	1:A:24:LEU:HD12	2.00	0.42
1:A:92:VAL:O	1:A:96:MET:HG3	2.20	0.42
1:A:181:LYS:HD2	1:B:339:GLU:OE1	2.20	0.42
1:A:240:ASN:O	1:A:284:GLY:HA2	2.20	0.42
1:A:68:ASN:O	1:A:69:ASP:CB	2.59	0.42
1:A:134:LEU:HD23	1:A:138:GLN:NE2	2.35	0.42
1:B:368:GLU:HA	1:B:368:GLU:OE1	2.20	0.42
1:B:82:GLN:HE22	1:B:160:PRO:HA	1.85	0.41
1:B:164:LEU:HD23	1:B:164:LEU:HA	1.76	0.41
1:A:22:ILE:HD13	1:A:45:LEU:HD11	2.01	0.41
1:A:275:VAL:HG23	1:A:276:GLN:N	2.35	0.41
1:B:17:VAL:CG1	1:B:20:LYS:HD3	2.50	0.41
1:A:92:VAL:HG12	1:A:96:MET:HE1	2.03	0.41
1:B:160:PRO:C	1:B:162:SER:H	2.23	0.41
1:B:362:CYS:O	1:B:363:ARG:HB3	2.20	0.41
1:A:219:LEU:HD12	1:A:219:LEU:HA	1.80	0.41
1:A:251:ILE:O	1:A:252:ASP:HB2	2.20	0.41
1:B:20:LYS:HB3	1:B:211:ARG:NH1	2.36	0.41
1:B:224:TYR:O	1:B:225:PRO:C	2.58	0.41
1:A:20:LYS:HG3	1:A:211:ARG:CG	2.49	0.41
1:A:61:PHE:CE2	1:A:66:ARG:HG2	2.56	0.41
1:A:109:VAL:HG23	1:A:110:ILE:N	2.35	0.41
1:B:161:LEU:HD12	1:B:161:LEU:N	2.35	0.41
1:A:164:LEU:HD23	1:A:169:ARG:NE	2.35	0.41
1:B:245:LYS:HE3	1:B:278:GLY:HA2	2.02	0.41
1:B:300:LEU:HB2	1:B:345:ILE:HD12	2.02	0.41
1:A:108:GLU:O	1:A:112:GLN:HB2	2.21	0.40
1:A:134:LEU:CD2	1:A:138:GLN:NE2	2.84	0.40
1:A:262:ASN:O	1:A:263:ARG:C	2.60	0.40
1:B:275:VAL:HG23	1:B:276:GLN:N	2.36	0.40
1:A:253:GLN:O	1:A:254:VAL:CB	2.68	0.40
1:A:275:VAL:HG23	1:A:276:GLN:H	1.87	0.40
1:B:11:LYS:HB3	1:B:19:SER:HB3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:ILE:HA	1:B:321:PRO:HD3	1.89	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	365/381 (96%)	321 (88%)	27 (7%)	17 (5%)	2 8
1	B	365/381 (96%)	322 (88%)	28 (8%)	15 (4%)	3 11
All	All	730/762 (96%)	643 (88%)	55 (8%)	32 (4%)	2 10

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	64	GLU
1	A	162	SER
1	A	164	LEU
1	A	263	ARG
1	A	296	ALA
1	A	322	SER
1	A	363	ARG
1	B	21	ASP
1	B	64	GLU
1	B	162	SER
1	B	263	ARG
1	B	296	ALA
1	B	322	SER
1	B	364	ARG
1	A	21	ASP
1	A	182	ARG
1	A	273	ARG

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Mol	Chain	Res	Type
1	B	182	ARG
1	B	273	ARG
1	B	369	PRO
1	A	161	LEU
1	B	363	ARG
1	A	163	ASN
1	A	254	VAL
1	A	333	ASP
1	B	112	GLN
1	B	271	GLU
1	B	292	PRO
1	A	112	GLN
1	A	335	VAL
1	B	335	VAL
1	A	334	VAL

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	312/323 (97%)	292 (94%)	20 (6%)	17	45
1	B	312/323 (97%)	286 (92%)	26 (8%)	11	32
All	All	624/646 (97%)	578 (93%)	46 (7%)	13	38

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	TRP
1	A	15	GLU
1	A	31	PHE
1	A	33	VAL
1	A	45	LEU
1	A	64	GLU
1	A	66	ARG
1	A	68	ASN

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Mol	Chain	Res	Type
1	A	102	LEU
1	A	129	ARG
1	A	146	ARG
1	A	159	GLU
1	A	163	ASN
1	A	205	VAL
1	A	219	LEU
1	A	249	THR
1	A	270	VAL
1	A	274	ASP
1	A	358	ASP
1	A	360	THR
1	B	13	TRP
1	B	15	GLU
1	B	31	PHE
1	B	33	VAL
1	B	45	LEU
1	B	64	GLU
1	B	66	ARG
1	B	68	ASN
1	B	102	LEU
1	B	129	ARG
1	B	146	ARG
1	B	147	THR
1	B	153	SER
1	B	159	GLU
1	B	163	ASN
1	B	205	VAL
1	B	219	LEU
1	B	249	THR
1	B	251	ILE
1	B	274	ASP
1	B	322	SER
1	B	350	GLU
1	B	351	ARG
1	B	358	ASP
1	B	360	THR
1	B	363	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	95	ASN
1	A	116	GLN
1	A	138	GLN
1	A	223	HIS
1	A	240	ASN
1	A	264	GLN
1	A	265	GLN
1	A	331	GLN
1	A	332	ASN
1	B	68	ASN
1	B	82	GLN
1	B	95	ASN
1	B	116	GLN
1	B	138	GLN
1	B	240	ASN
1	B	264	GLN
1	B	331	GLN
1	B	332	ASN
1	B	353	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	367/381 (96%)	-0.19	10 (2%)	54	50	12, 37, 72, 88	0
1	B	367/381 (96%)	-0.17	13 (3%)	44	38	13, 38, 71, 83	0
All	All	734/762 (96%)	-0.18	23 (3%)	49	44	12, 38, 72, 88	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	163	ASN	5.3
1	A	362	CYS	3.7
1	B	89	HIS	3.5
1	B	262	ASN	2.7
1	A	165	ASP	2.7
1	B	122	GLN	2.7
1	B	165	ASP	2.6
1	A	89	HIS	2.6
1	A	370	GLY	2.3
1	B	129	ARG	2.3
1	B	324	ARG	2.2
1	A	368	GLU	2.2
1	A	164	LEU	2.2
1	B	182	ARG	2.1
1	B	128	ASP	2.1
1	A	296	ALA	2.1
1	B	162	SER	2.1
1	A	294	ASP	2.1
1	B	125	HIS	2.1
1	B	273	ARG	2.1
1	B	367	LYS	2.1
1	A	15	GLU	2.0
1	A	369	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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.