



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 03:41 PM EDT

PDB ID : 3TUJ / pdb_00003tuj
Title : Inward facing conformations of the MetNI methionine ABC transporter: DM crystal form
Authors : Johnson, E.; Nguyen, P.; Rees, D.C.
Deposited on : 2011-09-16
Resolution : 4.00 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

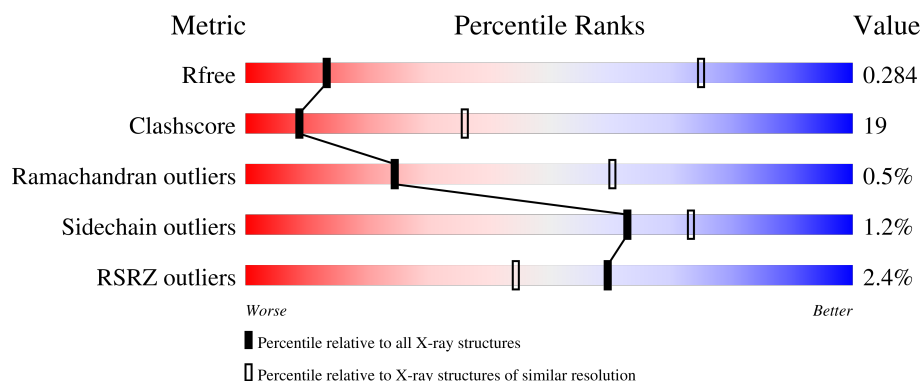
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1082 (4.20-3.80)
Clashscore	190562	1129 (4.20-3.80)
Ramachandran outliers	187476	1064 (4.20-3.80)
Sidechain outliers	187428	1055 (4.20-3.80)
RSRZ outliers	180081	1082 (4.20-3.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	217	0% 63% 35% .
1	B	217	0% 56% 41% ..
2	C	366	2% 66% 27% . 6%
2	D	366	4% 62% 30% . 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8559 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 D-methionine transport system permease protein metI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	Se	0	0	0
			1621	1076	263	272	10			
1	B	215	Total	C	N	O	Se	0	0	0
			1616	1073	262	271	10			

- Molecule 2 is a protein called Methionine import ATP-binding protein MetN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	344	Total	C	N	O	S	Se	0	0
			2661	1670	471	508	5	7		
2	D	344	Total	C	N	O	S	Se	0	0
			2661	1670	471	508	5	7		

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-22	MSE	-	expression tag	UNP P30750
C	-21	GLY	-	expression tag	UNP P30750
C	-20	HIS	-	expression tag	UNP P30750
C	-19	HIS	-	expression tag	UNP P30750
C	-18	HIS	-	expression tag	UNP P30750
C	-17	HIS	-	expression tag	UNP P30750
C	-16	HIS	-	expression tag	UNP P30750
C	-15	HIS	-	expression tag	UNP P30750
C	-14	HIS	-	expression tag	UNP P30750
C	-13	HIS	-	expression tag	UNP P30750
C	-12	HIS	-	expression tag	UNP P30750
C	-11	HIS	-	expression tag	UNP P30750
C	-10	SER	-	expression tag	UNP P30750
C	-9	SER	-	expression tag	UNP P30750
C	-8	GLY	-	expression tag	UNP P30750
C	-7	HIS	-	expression tag	UNP P30750

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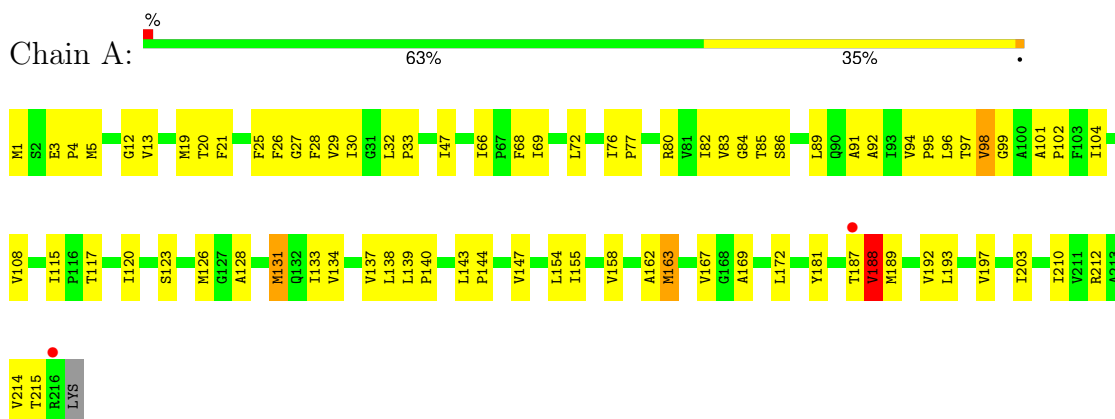
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Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	ILE	-	expression tag	UNP P30750
C	-5	ASP	-	expression tag	UNP P30750
C	-4	ASP	-	expression tag	UNP P30750
C	-3	ASP	-	expression tag	UNP P30750
C	-2	ASP	-	expression tag	UNP P30750
C	-1	LYS	-	expression tag	UNP P30750
C	0	HIS	-	expression tag	UNP P30750
C	166	GLN	GLU	engineered mutation	UNP P30750
D	-22	MSE	-	expression tag	UNP P30750
D	-21	GLY	-	expression tag	UNP P30750
D	-20	HIS	-	expression tag	UNP P30750
D	-19	HIS	-	expression tag	UNP P30750
D	-18	HIS	-	expression tag	UNP P30750
D	-17	HIS	-	expression tag	UNP P30750
D	-16	HIS	-	expression tag	UNP P30750
D	-15	HIS	-	expression tag	UNP P30750
D	-14	HIS	-	expression tag	UNP P30750
D	-13	HIS	-	expression tag	UNP P30750
D	-12	HIS	-	expression tag	UNP P30750
D	-11	HIS	-	expression tag	UNP P30750
D	-10	SER	-	expression tag	UNP P30750
D	-9	SER	-	expression tag	UNP P30750
D	-8	GLY	-	expression tag	UNP P30750
D	-7	HIS	-	expression tag	UNP P30750
D	-6	ILE	-	expression tag	UNP P30750
D	-5	ASP	-	expression tag	UNP P30750
D	-4	ASP	-	expression tag	UNP P30750
D	-3	ASP	-	expression tag	UNP P30750
D	-2	ASP	-	expression tag	UNP P30750
D	-1	LYS	-	expression tag	UNP P30750
D	0	HIS	-	expression tag	UNP P30750
D	166	GLN	GLU	engineered mutation	UNP P30750

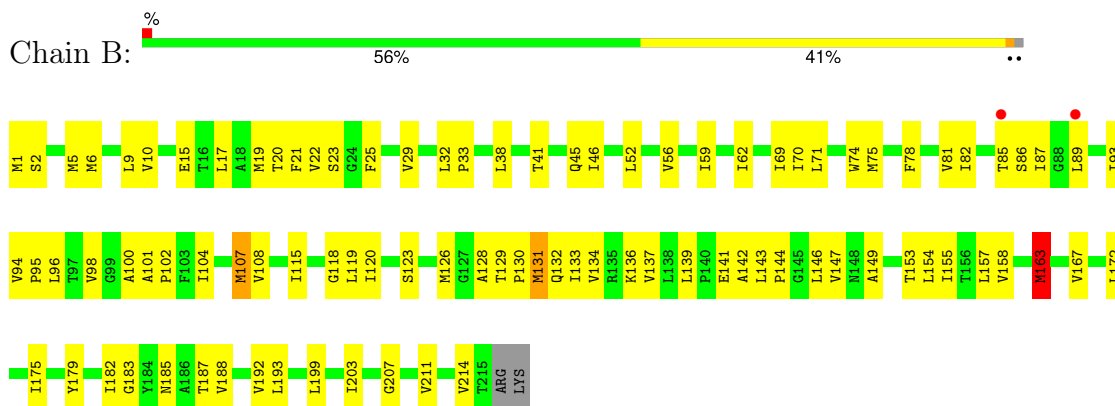
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

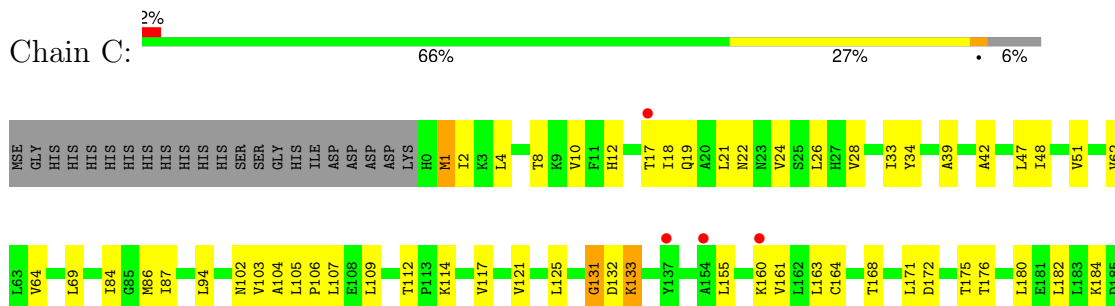
- Molecule 1: D-methionine transport system permease protein metI

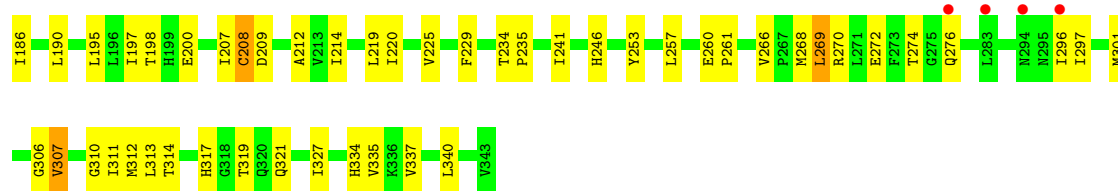


- Molecule 1: D-methionine transport system permease protein metI

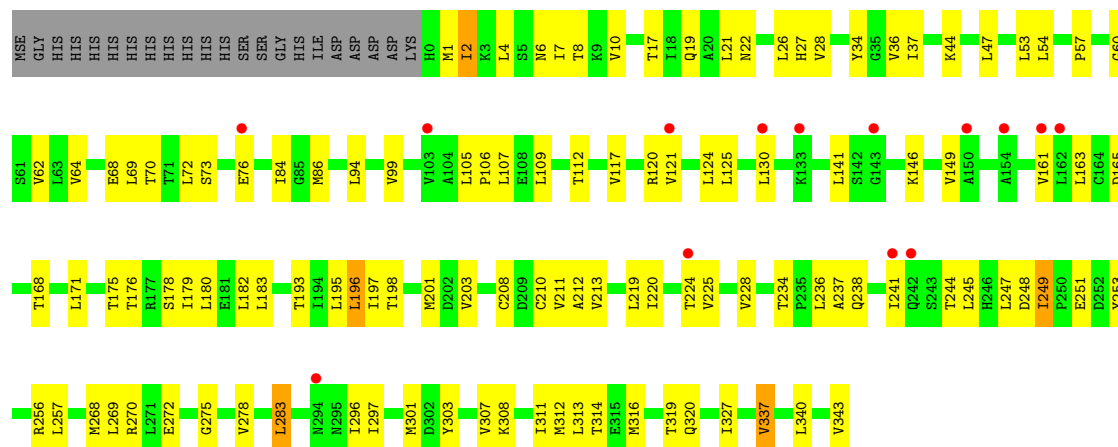


- Molecule 2: Methionine import ATP-binding protein MetN





• Molecule 2: Methionine import ATP-binding protein MetN



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	155.75Å 164.13Å 212.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.85 – 4.00 19.85 – 4.00	Depositor EDS
% Data completeness (in resolution range)	98.0 (19.85-4.00) 97.1 (19.85-4.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 3.82Å)	Xtriage
Refinement program	PHENIX 1.7.1 _743	Depositor
R, R_{free}	0.259 , 0.295 0.248 , 0.284	Depositor DCC
R_{free} test set	1753 reflections (7.33%)	wwPDB-VP
Wilson B-factor (Å ²)	153.5	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.22 , 84.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8559	wwPDB-VP
Average B, all atoms (Å ²)	190.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	1.02	3/1645 (0.2%)	0.97	0/2233
1	B	0.86	1/1640 (0.1%)	0.95	4/2226 (0.2%)
2	C	0.80	2/2694 (0.1%)	0.94	3/3641 (0.1%)
2	D	0.73	2/2694 (0.1%)	0.89	0/3641
All	All	0.84	8/8673 (0.1%)	0.93	7/11741 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	167	VAL	CA-CB	-6.65	1.47	1.54
1	A	163	MSE	SE-CE	6.31	2.14	1.95
2	C	1	MSE	CG-SE	-5.73	1.78	1.95
2	C	1	MSE	SE-CE	-5.56	1.78	1.95
2	D	1	MSE	SE-CE	-5.24	1.79	1.95
1	A	131	MSE	SE-CE	-5.23	1.79	1.95
1	B	131	MSE	SE-CE	-5.01	1.80	1.95
2	D	337	VAL	CA-CB	-5.01	1.48	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	131	GLY	N-CA-C	-7.35	95.76	113.18
2	C	246	HIS	N-CA-C	7.13	120.58	108.67
1	B	163	MSE	N-CA-C	-7.04	104.16	112.89
1	B	136	LYS	N-CA-C	6.84	121.70	112.88
1	B	163	MSE	CG-SE-CE	6.15	112.45	98.92
1	B	107	MSE	CB-CG-SE	-6.12	94.34	112.70
2	C	208	CYS	N-CA-C	5.29	117.95	110.50

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1621	0	1735	84	0
1	B	1616	0	1733	86	0
2	C	2661	0	2714	78	0
2	D	2661	0	2714	104	0
All	All	8559	0	8896	324	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 19.

All (324) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:MSE:SE	1:B:163:MSE:HG3	1.95	1.16
1:B:147:VAL:HG23	1:B:211:VAL:HG22	1.40	1.03
1:A:69:ILE:HG21	1:B:163:MSE:CE	1.90	1.00
1:A:143:LEU:HD22	1:A:214:VAL:HG21	1.47	0.94
1:B:104:ILE:HG22	1:B:153:THR:HG21	1.52	0.92
2:D:107:LEU:HD23	2:D:117:VAL:HG23	1.52	0.91
1:A:69:ILE:HG21	1:B:163:MSE:HE2	1.51	0.91
2:D:37:ILE:HD13	2:D:201:MSE:HE2	1.54	0.89
1:B:104:ILE:CG2	1:B:153:THR:HG21	2.04	0.86
1:A:128:ALA:HB2	2:C:109:LEU:HD21	1.57	0.85
2:D:198:THR:HG21	2:D:203:VAL:HG11	1.59	0.85
2:D:64:VAL:HG11	2:D:84:ILE:HD11	1.59	0.82
1:B:154:LEU:HD23	1:B:203:ILE:HG23	1.63	0.81
1:A:19:MSE:HE2	1:A:96:LEU:HD11	1.64	0.79
2:D:249:ILE:HG22	2:D:313:LEU:HD11	1.64	0.79
1:A:187:THR:O	1:A:188:VAL:HG12	1.82	0.78
1:A:19:MSE:CE	1:A:96:LEU:HD11	2.13	0.78
2:D:256:ARG:HD2	2:D:343:VAL:HG23	1.65	0.77
1:A:163:MSE:HB3	1:B:163:MSE:HE3	1.66	0.76
1:A:162:ALA:HB1	1:B:70:ILE:HD13	1.68	0.76
2:C:253:TYR:CD2	2:C:268:MSE:HE1	2.21	0.76
2:D:241:ILE:O	2:D:244:THR:HG22	1.86	0.76
1:A:133:ILE:HG22	1:A:138:LEU:HD13	1.66	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:ALA:HB2	1:A:169:ALA:HB1	1.69	0.74
1:B:75:MSE:HE1	1:B:98:VAL:HG11	1.70	0.73
1:B:19:MSE:HE2	1:B:96:LEU:HD11	1.68	0.73
2:D:130:LEU:HD13	2:D:149:VAL:HG12	1.71	0.73
2:D:168:THR:HG21	2:D:180:LEU:HD12	1.68	0.73
1:B:17:LEU:HD21	1:B:199:LEU:HD11	1.71	0.72
1:A:20:THR:HG21	1:A:158:VAL:HG22	1.70	0.72
2:D:10:VAL:HG11	2:D:17:THR:HG23	1.70	0.72
2:D:4:LEU:HD13	2:D:7:ILE:HD13	1.72	0.72
2:C:103:VAL:HG21	2:C:125:LEU:HD11	1.72	0.71
2:D:275:GLY:C	2:D:307:VAL:HG21	2.16	0.71
2:C:131:GLY:O	2:C:133:LYS:N	2.25	0.70
1:A:69:ILE:CG2	1:B:163:MSE:CE	2.68	0.69
2:C:229:PHE:CD2	2:C:241:ILE:HD11	2.27	0.69
2:D:180:LEU:HD21	2:D:203:VAL:HG13	1.74	0.69
2:C:297:ILE:HG22	2:C:313:LEU:O	1.92	0.69
1:B:20:THR:HG21	1:B:158:VAL:HG22	1.73	0.69
2:D:171:LEU:HD23	2:D:175:THR:HG22	1.74	0.68
2:D:28:VAL:HG22	2:D:34:TYR:CD2	2.28	0.68
1:B:123:SER:HA	1:B:126:MSE:HE2	1.74	0.68
1:A:137:VAL:HG13	2:C:94:LEU:HD21	1.75	0.67
2:C:10:VAL:HG12	2:C:19:GLN:HA	1.77	0.66
2:D:249:ILE:HG22	2:D:313:LEU:CD1	2.24	0.66
1:B:41:THR:HG21	1:B:52:LEU:HD11	1.77	0.64
1:B:129:THR:HG22	1:B:132:GLN:HG3	1.79	0.64
1:B:147:VAL:CG2	1:B:211:VAL:HG22	2.21	0.64
2:D:168:THR:HG21	2:D:180:LEU:CD1	2.27	0.64
1:A:117:THR:O	1:A:120:ILE:HG22	1.97	0.64
2:D:248:ASP:O	2:D:249:ILE:HG13	1.97	0.64
1:B:5:MSE:HE1	1:B:179:TYR:CZ	2.32	0.64
2:C:12:HIS:HB3	2:C:17:THR:HG22	1.80	0.64
2:D:253:TYR:CE2	2:D:297:ILE:HD13	2.33	0.64
1:A:21:PHE:CE2	1:A:154:LEU:HD11	2.33	0.63
1:A:97:THR:O	1:A:98:VAL:HB	1.99	0.63
2:C:168:THR:HG23	2:C:176:THR:CG2	2.28	0.63
2:C:171:LEU:HB3	2:C:175:THR:HG23	1.81	0.62
2:C:301:MSE:HE2	2:C:312:MSE:HE3	1.81	0.62
1:B:5:MSE:HE1	1:B:179:TYR:OH	1.99	0.62
2:C:24:VAL:HG13	2:C:219:LEU:HB2	1.79	0.62
2:D:107:LEU:HG	2:D:112:THR:HG21	1.81	0.62
1:A:32:LEU:HD12	1:A:108:VAL:HG11	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:10:VAL:CG1	2:D:17:THR:HG23	2.31	0.61
1:A:85:THR:HG23	1:A:91:ALA:HB2	1.83	0.61
2:C:301:MSE:CE	2:C:312:MSE:HE3	2.31	0.61
2:D:272:GLU:HG2	2:D:311:ILE:HG22	1.81	0.61
2:D:36:VAL:HB	2:D:197:ILE:HG22	1.84	0.60
2:C:180:LEU:HD22	2:C:207:ILE:HD13	1.83	0.60
2:C:48:ILE:HD12	2:C:197:ILE:HG12	1.84	0.60
2:C:297:ILE:HD11	2:D:303:TYR:CD1	2.37	0.59
2:D:245:LEU:HD12	2:D:311:ILE:HG23	1.85	0.59
2:D:296:ILE:HA	2:D:314:THR:HG22	1.83	0.59
2:D:249:ILE:HD12	2:D:249:ILE:O	2.03	0.58
1:A:25:PHE:O	1:A:29:VAL:HG23	2.04	0.58
1:A:155:ILE:HG22	1:A:203:ILE:CD1	2.33	0.58
1:A:69:ILE:CG2	1:B:163:MSE:HE3	2.33	0.58
2:C:28:VAL:HG22	2:C:34:TYR:CD2	2.39	0.58
1:A:82:ILE:HG23	1:A:83:VAL:HG23	1.84	0.58
1:A:69:ILE:HG21	1:B:163:MSE:HE3	1.83	0.58
1:A:101:ALA:HB3	1:A:102:PRO:CD	2.34	0.58
2:D:311:ILE:HD12	2:D:311:ILE:C	2.29	0.57
2:C:163:LEU:HD23	2:C:195:LEU:HB2	1.86	0.57
1:B:104:ILE:HG23	1:B:153:THR:HG21	1.85	0.57
2:C:64:VAL:HG11	2:C:84:ILE:HD11	1.87	0.57
2:D:301:MSE:HE2	2:D:312:MSE:CE	2.35	0.57
1:A:19:MSE:HE3	1:A:92:ALA:HB1	1.86	0.56
1:A:69:ILE:CG2	1:B:163:MSE:HE2	2.31	0.56
1:A:172:LEU:HG	1:A:192:VAL:HG23	1.86	0.56
1:B:78:PHE:CE2	1:B:82:ILE:HD11	2.40	0.56
1:B:23:SER:HB2	1:B:157:LEU:HD13	1.86	0.56
1:B:19:MSE:CE	1:B:96:LEU:HD11	2.36	0.56
1:A:20:THR:HG21	1:A:158:VAL:CG2	2.35	0.56
2:C:319:THR:HG22	2:C:321:GLN:H	1.70	0.56
1:B:207:GLY:O	1:B:211:VAL:HG23	2.05	0.56
2:D:168:THR:HB	2:D:176:THR:HG23	1.87	0.56
1:B:182:ILE:HD12	1:B:183:GLY:N	2.21	0.56
1:B:100:ALA:HB2	1:B:157:LEU:HD21	1.88	0.55
2:C:2:ILE:CG2	2:C:161:VAL:HG21	2.37	0.55
2:C:212:ALA:HB1	2:C:219:LEU:HD11	1.88	0.55
2:D:120:ARG:O	2:D:124:LEU:HD13	2.07	0.55
1:B:120:ILE:HD11	1:B:133:ILE:HD12	1.89	0.55
2:C:297:ILE:HD13	2:D:278:VAL:HG21	1.89	0.55
1:A:181:TYR:CE1	1:B:167:VAL:HG13	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:ILE:HA	1:B:137:VAL:HG22	1.89	0.54
1:A:181:TYR:CD1	1:B:167:VAL:HG13	2.42	0.54
1:A:80:ARG:O	1:A:84:GLY:HA2	2.06	0.54
2:C:229:PHE:CG	2:C:241:ILE:HD11	2.42	0.54
2:D:224:THR:O	2:D:228:VAL:HG23	2.07	0.54
1:B:128:ALA:HB2	2:D:109:LEU:HD11	1.90	0.54
2:D:196:LEU:C	2:D:196:LEU:HD22	2.33	0.54
1:A:123:SER:HA	1:A:126:MSE:HE2	1.90	0.53
1:B:22:VAL:HG11	1:B:93:ILE:HG21	1.90	0.53
1:B:78:PHE:CZ	1:B:82:ILE:HD11	2.44	0.53
2:C:260:GLU:HB3	2:C:261:PRO:CD	2.38	0.53
2:D:54:LEU:HD11	2:D:86:MSE:SE	2.59	0.53
1:A:21:PHE:CD2	1:A:154:LEU:HD11	2.43	0.53
1:A:94:VAL:HB	1:A:95:PRO:HD3	1.91	0.52
1:B:172:LEU:HB3	1:B:192:VAL:HG13	1.92	0.52
2:C:296:ILE:HA	2:C:314:THR:HG22	1.92	0.52
1:B:129:THR:OG1	1:B:130:PRO:HD2	2.08	0.52
2:C:297:ILE:HG22	2:C:313:LEU:C	2.35	0.52
2:D:179:ILE:O	2:D:183:LEU:HD23	2.09	0.52
1:A:214:VAL:HG22	1:A:214:VAL:O	2.10	0.52
2:C:260:GLU:HB3	2:C:261:PRO:HD2	1.92	0.52
2:D:105:LEU:N	2:D:106:PRO:HD2	2.24	0.52
2:D:26:LEU:HD23	2:D:27:HIS:N	2.25	0.51
1:B:15:GLU:HA	1:B:89:LEU:HD21	1.91	0.51
1:B:115:ILE:C	1:B:115:ILE:HD12	2.35	0.51
2:C:104:ALA:HB2	2:C:121:VAL:HG21	1.92	0.51
2:C:107:LEU:HD23	2:C:112:THR:HG21	1.93	0.51
2:D:253:TYR:HB3	2:D:268:MSE:HE1	1.92	0.51
1:B:93:ILE:HA	1:B:96:LEU:HD12	1.93	0.51
2:C:220:ILE:HG21	2:C:234:THR:HG21	1.92	0.51
1:A:19:MSE:HE3	1:A:96:LEU:HD11	1.89	0.50
1:A:83:VAL:HG21	1:A:91:ALA:HA	1.93	0.50
2:C:276:GLN:OE1	2:C:335:VAL:HG22	2.11	0.50
1:A:212:ARG:O	1:A:215:THR:HG22	2.11	0.50
1:B:101:ALA:HB3	1:B:102:PRO:CD	2.41	0.50
2:C:209:ASP:O	2:C:225:VAL:HG23	2.12	0.50
1:B:41:THR:CB	1:B:52:LEU:HD11	2.42	0.50
2:D:247:LEU:HD22	2:D:270:ARG:HD2	1.92	0.50
2:D:269:LEU:HD22	2:D:316:MSE:HE3	1.94	0.50
1:B:119:LEU:HD11	2:D:94:LEU:HD23	1.93	0.50
2:D:210:CYS:SG	2:D:224:THR:HG22	2.52	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:LEU:C	1:A:193:LEU:HD23	2.37	0.50
2:D:21:LEU:HD21	2:D:47:LEU:HB2	1.94	0.49
1:A:72:LEU:C	1:A:72:LEU:HD23	2.37	0.49
2:C:257:LEU:HD11	2:C:340:LEU:O	2.12	0.49
2:D:8:THR:HG23	2:D:22:ASN:HA	1.95	0.49
2:D:37:ILE:HG21	2:D:201:MSE:HE1	1.93	0.49
1:A:13:VAL:HG12	1:A:172:LEU:HD21	1.94	0.49
1:A:97:THR:C	1:A:99:GLY:H	2.18	0.49
2:D:84:ILE:HG23	2:D:163:LEU:HD13	1.93	0.49
1:A:26:PHE:O	1:A:30:ILE:HG22	2.12	0.49
1:A:80:ARG:O	1:A:84:GLY:CA	2.61	0.49
1:B:1:MSE:N	1:B:187:THR:HG21	2.27	0.49
1:B:41:THR:CG2	1:B:52:LEU:HD11	2.42	0.49
2:C:87:ILE:HD11	2:C:164:CYS:SG	2.53	0.49
2:C:168:THR:HG23	2:C:176:THR:HG22	1.93	0.49
1:A:143:LEU:C	1:A:143:LEU:HD23	2.37	0.49
1:A:155:ILE:HG22	1:A:203:ILE:HG13	1.95	0.49
2:D:245:LEU:HD22	2:D:245:LEU:N	2.28	0.49
2:D:37:ILE:HD13	2:D:201:MSE:CE	2.34	0.48
2:D:249:ILE:O	2:D:249:ILE:CG1	2.61	0.48
1:B:32:LEU:HD12	1:B:108:VAL:HG11	1.95	0.48
2:C:172:ASP:O	2:C:176:THR:HG23	2.13	0.48
2:D:2:ILE:HD12	2:D:161:VAL:HG11	1.95	0.48
2:D:62:VAL:O	2:D:68:GLU:HA	2.13	0.48
1:B:107:MSE:HE1	1:B:149:ALA:HA	1.96	0.48
2:D:171:LEU:HD23	2:D:175:THR:CG2	2.42	0.48
1:B:1:MSE:HG3	1:B:2:SER:H	1.77	0.48
2:C:26:LEU:CD2	2:C:28:VAL:HG23	2.43	0.48
2:C:105:LEU:N	2:C:106:PRO:CD	2.77	0.48
1:B:9:LEU:HD12	1:B:175:ILE:HD13	1.96	0.48
1:A:27:GLY:C	1:A:104:ILE:HD11	2.39	0.48
1:A:66:ILE:HD13	1:B:193:LEU:HD11	1.95	0.48
2:C:86:MSE:HE2	2:C:163:LEU:HD13	1.96	0.48
1:A:154:LEU:CD2	1:A:203:ILE:HD12	2.44	0.47
1:B:185:ASN:ND2	1:B:188:VAL:HG23	2.28	0.47
2:C:234:THR:CG2	2:C:235:PRO:HD2	2.44	0.47
2:D:10:VAL:HG22	2:D:19:GLN:HG2	1.96	0.47
2:D:301:MSE:HE2	2:D:312:MSE:HE1	1.96	0.47
1:A:143:LEU:CD2	1:A:214:VAL:HG11	2.44	0.47
1:B:143:LEU:N	1:B:144:PRO:HD2	2.29	0.47
2:C:269:LEU:HD13	2:C:269:LEU:N	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:LEU:HB3	1:A:33:PRO:HD3	1.97	0.47
1:A:189:MSE:HE2	1:B:74:TRP:N	2.29	0.47
1:B:78:PHE:O	1:B:81:VAL:HG22	2.15	0.47
2:D:301:MSE:HE2	2:D:312:MSE:HE3	1.97	0.47
1:A:76:ILE:HB	1:A:77:PRO:HD3	1.96	0.47
1:B:32:LEU:N	1:B:33:PRO:CD	2.78	0.47
1:B:46:ILE:CD1	1:B:133:ILE:HD11	2.45	0.47
2:C:1:MSE:HE2	2:C:160:LYS:HB2	1.96	0.47
2:C:172:ASP:CG	2:C:175:THR:HG22	2.39	0.47
1:A:143:LEU:N	1:A:144:PRO:HD2	2.30	0.47
2:D:234:THR:HB	2:D:237:ALA:HB3	1.97	0.47
2:D:212:ALA:HB1	2:D:219:LEU:HD11	1.97	0.46
1:A:12:GLY:C	1:A:172:LEU:HD13	2.41	0.46
2:D:178:SER:O	2:D:182:LEU:HD13	2.16	0.46
1:A:139:LEU:N	1:A:140:PRO:HD2	2.30	0.46
2:D:211:VAL:CG1	2:D:228:VAL:HG21	2.46	0.46
1:B:20:THR:HG21	1:B:158:VAL:CG2	2.43	0.46
2:D:94:LEU:CD1	2:D:105:LEU:HD22	2.46	0.46
1:A:163:MSE:HB3	1:B:163:MSE:CE	2.40	0.46
1:B:46:ILE:HD11	1:B:120:ILE:HD11	1.97	0.46
1:B:118:GLY:H	1:B:120:ILE:HG22	1.80	0.46
1:B:134:VAL:O	1:B:139:LEU:HD23	2.16	0.46
2:C:270:ARG:HB3	2:C:313:LEU:HD13	1.98	0.46
2:C:198:THR:HG22	2:C:200:GLU:H	1.80	0.46
2:C:297:ILE:CD1	2:D:278:VAL:HG21	2.45	0.46
1:A:85:THR:O	1:A:86:SER:OG	2.34	0.46
2:C:4:LEU:HD21	2:C:51:VAL:HG22	1.98	0.46
2:D:73:SER:HB3	2:D:76:GLU:HB2	1.97	0.46
1:B:25:PHE:O	1:B:29:VAL:HG23	2.16	0.45
2:C:33:ILE:HB	2:C:208:CYS:HA	1.97	0.45
2:C:62:VAL:HG12	2:C:69:LEU:HD12	1.97	0.45
2:D:94:LEU:HD12	2:D:105:LEU:HD22	1.97	0.45
1:A:80:ARG:O	1:A:84:GLY:N	2.49	0.45
1:B:94:VAL:HB	1:B:95:PRO:HD3	1.98	0.45
2:C:186:ILE:HG23	2:C:190:LEU:HD12	1.97	0.45
2:C:272:GLU:HG2	2:C:311:ILE:HG22	1.98	0.45
2:D:319:THR:HG22	2:D:320:GLN:N	2.32	0.45
1:A:85:THR:O	1:A:86:SER:CB	2.63	0.45
1:A:155:ILE:HG22	1:A:203:ILE:HD11	1.98	0.45
1:B:41:THR:CG2	1:B:52:LEU:HD21	2.46	0.45
1:B:59:ILE:O	1:B:62:ILE:HG22	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:8:THR:HG23	2:C:22:ASN:HA	1.97	0.45
2:D:37:ILE:CG2	2:D:213:VAL:HG22	2.47	0.45
2:D:208:CYS:O	2:D:225:VAL:HG21	2.17	0.45
2:D:220:ILE:HD12	2:D:234:THR:HG21	1.99	0.45
2:C:62:VAL:CG1	2:C:69:LEU:HD12	2.48	0.44
2:D:99:VAL:HG11	2:D:141:LEU:HD21	1.99	0.44
2:C:103:VAL:HG21	2:C:125:LEU:CD1	2.44	0.44
1:A:85:THR:HG23	1:A:86:SER:N	2.32	0.44
1:A:163:MSE:O	1:B:163:MSE:CE	2.66	0.44
2:C:114:LYS:O	2:C:117:VAL:HG12	2.17	0.44
1:B:155:ILE:C	1:B:155:ILE:HD12	2.43	0.44
2:D:313:LEU:C	2:D:313:LEU:HD23	2.42	0.44
2:D:327:ILE:HG13	2:D:337:VAL:HG11	1.99	0.44
2:C:306:GLY:C	2:C:307:VAL:CG2	2.90	0.44
2:D:2:ILE:HG23	2:D:28:VAL:HB	1.99	0.44
2:D:72:LEU:HD13	2:D:72:LEU:O	2.18	0.44
2:D:171:LEU:N	2:D:171:LEU:HD12	2.32	0.44
1:A:3:GLU:N	1:A:4:PRO:HD2	2.33	0.44
2:D:2:ILE:CG2	2:D:193:THR:HG21	2.48	0.44
2:D:234:THR:HG22	2:D:236:LEU:H	1.83	0.44
2:D:121:VAL:O	2:D:125:LEU:HD13	2.18	0.43
2:C:207:ILE:HG13	2:C:208:CYS:N	2.32	0.43
1:A:187:THR:O	1:A:188:VAL:CG1	2.61	0.43
1:B:85:THR:O	1:B:86:SER:CB	2.67	0.43
1:B:21:PHE:CD2	1:B:154:LEU:HD11	2.53	0.43
2:C:225:VAL:HG13	2:C:229:PHE:CE1	2.53	0.43
2:D:249:ILE:O	2:D:249:ILE:CD1	2.66	0.43
1:B:87:ILE:HD12	1:B:87:ILE:C	2.43	0.43
1:B:129:THR:HG23	1:B:131:MSE:H	1.83	0.43
2:C:18:ILE:N	2:C:18:ILE:HD12	2.34	0.43
2:D:2:ILE:HG21	2:D:193:THR:HG21	2.01	0.43
2:C:39:ALA:HB3	2:C:42:ALA:HB2	2.01	0.43
2:C:155:LEU:HD11	2:C:186:ILE:HD13	2.01	0.43
2:C:301:MSE:HG2	2:C:310:GLY:HA3	2.01	0.43
2:D:34:TYR:CD2	2:D:210:CYS:HB2	2.54	0.43
2:D:7:ILE:HG23	2:D:57:PRO:HB3	2.01	0.43
2:D:307:VAL:HG22	2:D:308:LYS:O	2.19	0.43
2:D:2:ILE:HG23	2:D:2:ILE:O	2.19	0.42
1:B:119:LEU:HD21	2:D:94:LEU:HD23	2.01	0.42
2:D:283:LEU:H	2:D:283:LEU:HD12	1.84	0.42
1:A:154:LEU:HD23	1:A:203:ILE:HD12	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:LEU:HD21	1:B:56:VAL:HG12	2.01	0.42
2:D:53:LEU:HD12	2:D:69:LEU:HB3	2.01	0.42
2:D:238:GLN:O	2:D:241:ILE:HG22	2.19	0.42
1:A:28:PHE:CD1	1:A:104:ILE:HD12	2.54	0.42
1:A:68:PHE:CZ	1:A:96:LEU:HD23	2.54	0.42
1:B:52:LEU:HD13	1:B:56:VAL:HG21	2.02	0.42
1:A:115:ILE:HD13	1:A:115:ILE:HG21	1.83	0.42
2:C:64:VAL:CG2	2:C:69:LEU:HD11	2.49	0.42
2:C:274:THR:HG21	2:C:334:HIS:HB3	2.01	0.42
2:C:327:ILE:HG23	2:C:337:VAL:HG11	2.02	0.42
1:B:85:THR:O	1:B:86:SER:HB3	2.19	0.42
2:D:195:LEU:HD13	2:D:196:LEU:N	2.35	0.42
1:A:97:THR:C	1:A:99:GLY:N	2.79	0.41
1:A:193:LEU:O	1:A:197:VAL:HG23	2.20	0.41
2:C:21:LEU:HD21	2:C:47:LEU:HB2	2.02	0.41
2:D:44:LYS:O	2:D:197:ILE:HD13	2.20	0.41
2:D:165:ASP:OD1	2:D:197:ILE:HD11	2.20	0.41
2:D:270:ARG:HB2	2:D:340:LEU:HD11	2.02	0.41
1:A:126:MSE:SE	2:C:105:LEU:HD22	2.70	0.41
1:A:163:MSE:HG2	1:B:69:ILE:HB	2.01	0.41
1:B:142:ALA:O	1:B:146:LEU:HD13	2.21	0.41
2:D:211:VAL:HG13	2:D:228:VAL:HG21	2.02	0.41
2:D:307:VAL:HG22	2:D:308:LYS:N	2.34	0.41
1:A:1:MSE:HE3	1:A:5:MSE:HE3	2.02	0.41
2:D:2:ILE:HD12	2:D:161:VAL:CG1	2.51	0.41
2:C:266:VAL:HG11	2:C:317:HIS:CG	2.56	0.41
1:A:147:VAL:HG11	1:A:210:ILE:HG23	2.02	0.41
2:D:168:THR:CB	2:D:176:THR:HG23	2.49	0.41
2:D:327:ILE:CG1	2:D:337:VAL:HG11	2.50	0.41
1:A:85:THR:CG2	1:A:91:ALA:HB2	2.51	0.41
2:C:214:ILE:O	2:C:214:ILE:HG23	2.21	0.41
2:D:234:THR:HB	2:D:237:ALA:CB	2.51	0.41
1:A:126:MSE:HE1	2:C:105:LEU:HD22	2.03	0.41
1:B:71:LEU:HD11	1:B:75:MSE:CE	2.51	0.41
2:C:103:VAL:CG2	2:C:125:LEU:HD11	2.46	0.41
2:C:301:MSE:HE1	2:C:312:MSE:HE3	2.02	0.41
2:D:60:GLY:O	2:D:70:THR:HG21	2.21	0.41
2:D:195:LEU:HD13	2:D:195:LEU:C	2.46	0.41
1:A:20:THR:HG21	1:A:158:VAL:HA	2.03	0.41
1:B:193:LEU:C	1:B:193:LEU:HD23	2.46	0.41
2:C:2:ILE:C	2:C:2:ILE:HD12	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:253:TYR:O	2:D:257:LEU:N	2.54	0.40
1:A:69:ILE:HG21	1:B:163:MSE:SE	2.70	0.40
1:A:123:SER:HA	1:A:126:MSE:CE	2.51	0.40
1:B:101:ALA:HB3	1:B:102:PRO:HD3	2.02	0.40
1:B:143:LEU:HG	1:B:214:VAL:HG21	2.02	0.40
2:C:182:LEU:C	2:C:182:LEU:HD13	2.47	0.40
2:D:146:LYS:O	2:D:149:VAL:HG22	2.21	0.40
1:A:131:MSE:HA	1:A:134:VAL:HG12	2.04	0.40
1:B:137:VAL:O	1:B:141:GLU:HB2	2.21	0.40
2:C:48:ILE:HG21	2:C:197:ILE:HD11	2.03	0.40
2:C:184:LYS:HG2	2:D:251:GLU:HG3	2.02	0.40
1:A:32:LEU:N	1:A:33:PRO:CD	2.85	0.40
1:B:6:MSE:O	1:B:10:VAL:HG23	2.21	0.40
2:D:196:LEU:N	2:D:196:LEU:HD13	2.36	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	214/217 (99%)	205 (96%)	6 (3%)	3 (1%)	9	39
1	B	213/217 (98%)	204 (96%)	9 (4%)	0	100	100
2	C	342/366 (93%)	318 (93%)	22 (6%)	2 (1%)	21	57
2	D	342/366 (93%)	324 (95%)	17 (5%)	1 (0%)	36	70
All	All	1111/1166 (95%)	1051 (95%)	54 (5%)	6 (0%)	24	60

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	98	VAL

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Mol	Chain	Res	Type
1	A	188	VAL
2	C	132	ASP
2	C	133	LYS
1	A	47	ILE
2	D	2	ILE

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	170/162 (105%)	168 (99%)	2 (1%)	63	73
1	B	170/162 (105%)	168 (99%)	2 (1%)	63	73
2	C	299/311 (96%)	296 (99%)	3 (1%)	68	76
2	D	299/311 (96%)	295 (99%)	4 (1%)	61	72
All	All	938/946 (99%)	927 (99%)	11 (1%)	63	73

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	LEU
1	A	188	VAL
1	B	45	GLN
1	B	163	MSE
2	C	102	ASN
2	C	269	LEU
2	C	307	VAL
2	D	6	ASN
2	D	196	LEU
2	D	249	ILE
2	D	283	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	GLN
1	B	61	ASN
1	B	185	ASN
2	C	23	ASN
2	C	187	ASN
2	C	294	ASN
2	D	145	GLN
2	D	231	HIS
2	D	242	GLN
2	D	276	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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

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	206/217 (94%)	0.04	2 (0%) 79 61	118, 150, 193, 217	0
1	B	205/217 (94%)	0.03	2 (0%) 79 61	124, 176, 223, 233	0
2	C	337/366 (92%)	0.22	8 (2%) 59 44	136, 176, 239, 290	0
2	D	337/366 (92%)	0.47	14 (4%) 40 31	124, 241, 289, 314	0
All	All	1085/1166 (93%)	0.23	26 (2%) 59 44	118, 178, 277, 314	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	294	ASN	3.5
1	B	85	THR	3.3
2	C	17	THR	3.2
2	D	154	ALA	3.2
2	D	130	LEU	3.1
2	D	242	GLN	3.0
1	A	216	ARG	3.0
2	D	161	VAL	2.8
2	C	294	ASN	2.8
2	D	150	ALA	2.7
2	D	121	VAL	2.7
2	C	137	TYR	2.6
2	C	283	LEU	2.6
2	D	103	VAL	2.5
2	C	154	ALA	2.5
1	B	89	LEU	2.4
2	D	143	GLY	2.4
2	D	241	ILE	2.4
1	A	187	THR	2.3
2	C	276	GLN	2.1
2	D	162	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
2	C	160	LYS	2.1
2	C	296	ILE	2.1
2	D	133	LYS	2.0
2	D	76	GLU	2.0
2	D	224	THR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.