



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:10 AM GMT

PDB ID : 2FZ1
Title : Structure of Empty Head Turnip Yellow Mosaic Virus (ATC) at 100 K
Authors : Larson, S.B.; Lucas, R.W.; McPherson, A.
Deposited on : 2006-02-08
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
<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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

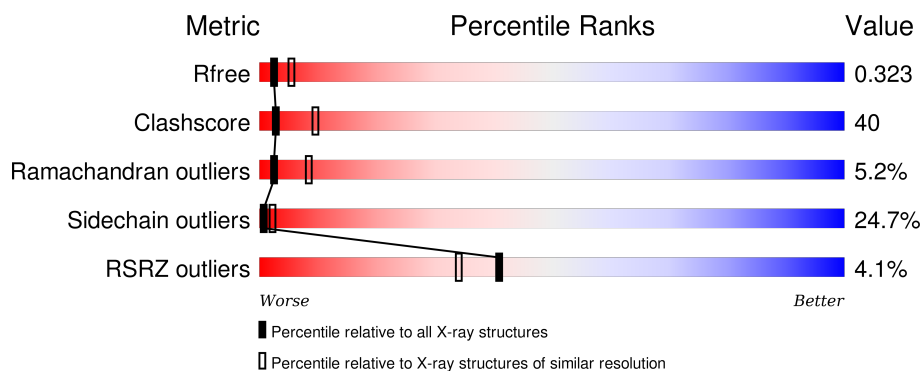
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (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 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	189	 5% 30% 37% 19% • 14%
1	B	189	 3% 40% 38% 21% •
1	C	189	 3% 37% 55% 8% •

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4058 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 Coat protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	163	Total	C	N	O	S	0	0	0
			1224	787	194	236	7			
1	B	189	Total	C	N	O	S	0	0	0
			1417	909	225	275	8			
1	C	189	Total	C	N	O	S	0	0	0
			1417	909	225	275	8			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	99	GLN	ASN	VARIANT	UNP P20125
A	123	GLN	ASN	VARIANT	UNP P20125
A	138	GLN	ASN	VARIANT	UNP P20125
B	99	GLN	ASN	VARIANT	UNP P20125
B	123	GLN	ASN	VARIANT	UNP P20125
B	138	GLN	ASN	VARIANT	UNP P20125
C	99	GLN	ASN	VARIANT	UNP P20125
C	123	GLN	ASN	VARIANT	UNP P20125
C	138	GLN	ASN	VARIANT	UNP P20125

H137	Q138	P139	R140	V141	K142	D143	S144	I145		L148	D149	S150	P151	K152	L153		S156	I157	T158	A159	Q160	P161	T162	A163	P164		S167	T168			I171	T172	V173	S174	G175	T176	L177	S178	M179	H180	S181	P182	L183	I184	T185	D186	T187	S188	T189
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4 Data and refinement statistics

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α , β , γ	511.36Å 511.36Å 304.31Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.64 – 2.90 49.53 – 2.89	Depositor EDS
% Data completeness (in resolution range)	49.2 (48.64-2.90) 55.2 (49.53-2.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.48 (at 2.91Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.309 , 0.315 0.321 , 0.323	Depositor DCC
R_{free} test set	14121 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	49.7	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 322958 reflections	Xtriage
F_o, F_c correlation	0.78	EDS
Total number of atoms	4058	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.83% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.60	0/1255	0.82	0/1724
1	B	0.57	0/1452	0.84	0/1997
1	C	0.29	0/1452	0.60	0/1997
All	All	0.50	0/4159	0.76	0/5718

There are no bond length outliers.

There are no bond angle outliers.

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	1224	0	1250	106	0
1	B	1417	0	1454	125	0
1	C	1417	0	1454	113	0
All	All	4058	0	4158	329	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 40.

All (329) 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:153:LEU:HD13	1:B:155:ILE:HD11	1.30	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:THR:HG23	1:B:83:GLN:H	1.20	1.05
1:A:30:THR:HG23	1:A:180:HIS:HB3	1.34	1.03
1:B:30:THR:HG22	1:B:180:HIS:HB3	1.42	1.01
1:C:184:ILE:HD12	1:C:184:ILE:H	1.23	0.99
1:A:136:MET:HG2	1:B:184:ILE:HD11	1.43	0.98
1:C:97:PRO:HB2	1:C:99:GLN:HE22	1.32	0.94
1:C:81:THR:HG23	1:C:83:GLN:H	1.29	0.94
1:C:134:LEU:HD22	1:C:134:LEU:H	1.38	0.88
1:C:10:GLN:HE22	1:C:94:CYS:HA	1.38	0.87
1:A:118:ILE:HA	1:A:123:GLN:HG2	1.54	0.87
1:A:184:ILE:HD12	1:A:184:ILE:H	1.38	0.87
1:A:102:VAL:HG13	1:A:106:GLN:HG3	1.57	0.86
1:A:59:VAL:O	1:A:63:THR:HG23	1.76	0.84
1:A:88:PRO:HA	1:A:118:ILE:O	1.78	0.84
1:C:5:LYS:HD3	1:C:129:ILE:HD11	1.61	0.83
1:B:153:LEU:CD1	1:B:155:ILE:HD11	2.09	0.83
1:C:56:ILE:HG12	1:C:59:VAL:HG12	1.62	0.82
1:B:71:LEU:HD13	1:B:177:LEU:HD23	1.59	0.82
1:A:77:THR:HB	1:A:129:ILE:HG22	1.62	0.81
1:B:4:ASP:O	1:B:127:PRO:HD2	1.83	0.78
1:A:81:THR:HG23	1:A:83:GLN:H	1.47	0.78
1:B:50:SER:C	1:B:51:LEU:HD12	2.04	0.78
1:B:19:VAL:HG23	1:C:23:VAL:HG12	1.65	0.78
1:A:67:ARG:HH12	1:C:17:ALA:HB2	1.50	0.76
1:B:99:GLN:HE22	1:B:149:ASP:HB3	1.50	0.76
1:B:71:LEU:HD13	1:B:177:LEU:CD2	2.17	0.75
1:B:77:THR:HB	1:B:79:HIS:CE1	2.22	0.74
1:A:126:SER:C	1:A:128:LEU:H	1.91	0.74
1:A:81:THR:HG22	1:A:168:THR:O	1.88	0.73
1:B:63:THR:HA	1:B:179:MET:HE2	1.70	0.73
1:B:115:ILE:HD13	1:B:115:ILE:O	1.89	0.72
1:A:90:THR:HB	1:A:158:THR:HB	1.72	0.72
1:A:125:LEU:C	1:A:127:PRO:HD2	2.09	0.71
1:B:153:LEU:HD13	1:B:155:ILE:CD1	2.16	0.71
1:B:148:LEU:H	1:B:148:LEU:HD12	1.56	0.71
1:A:81:THR:CG2	1:A:83:GLN:HB2	2.21	0.70
1:B:30:THR:HG22	1:B:180:HIS:CB	2.20	0.70
1:B:90:THR:HB	1:B:158:THR:HB	1.73	0.70
1:B:15:THR:HG22	1:B:136:MET:HG3	1.71	0.69
1:C:81:THR:HG21	1:C:167:SER:O	1.91	0.69
1:B:3:ILE:H	1:B:3:ILE:HD12	1.57	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:THR:HG22	1:A:168:THR:C	2.13	0.69
1:B:77:THR:HG23	1:B:129:ILE:HG12	1.74	0.69
1:C:97:PRO:CB	1:C:99:GLN:HE22	2.05	0.69
1:B:95:TRP:CE2	1:B:132:CYS:HB2	2.28	0.69
1:B:74:LEU:HD22	1:B:75:TRP:N	2.08	0.68
1:C:61:THR:HA	1:C:64:THR:HG23	1.76	0.68
1:B:154:LEU:C	1:B:155:ILE:HD12	2.14	0.68
1:A:123:GLN:HA	1:A:123:GLN:HE21	1.58	0.67
1:B:74:LEU:HD22	1:B:75:TRP:H	1.57	0.67
1:B:40:LEU:HD22	1:B:41:PHE:N	2.09	0.67
1:A:118:ILE:HA	1:A:123:GLN:CG	2.25	0.67
1:B:60:SER:O	1:B:64:THR:HG23	1.95	0.66
1:A:87:PHE:CD1	1:A:164:PRO:HB3	2.29	0.66
1:C:90:THR:HB	1:C:158:THR:HB	1.78	0.66
1:B:39:VAL:HG11	1:B:173:VAL:HG23	1.76	0.66
1:A:81:THR:HG21	1:A:167:SER:O	1.96	0.65
1:C:93:VAL:HG11	1:C:130:VAL:HG11	1.78	0.65
1:A:81:THR:HG23	1:A:83:GLN:HB2	1.78	0.65
1:B:32:LYS:HA	1:B:177:LEU:O	1.97	0.64
1:A:103:THR:O	1:A:106:GLN:HG2	1.98	0.64
1:A:39:VAL:HG22	1:A:171:ILE:O	1.98	0.64
1:B:81:THR:HG23	1:B:83:GLN:N	2.04	0.64
1:C:97:PRO:HB2	1:C:99:GLN:NE2	2.07	0.64
1:B:81:THR:HG22	1:B:168:THR:O	1.97	0.64
1:B:103:THR:HG23	1:B:106:GLN:CG	2.27	0.64
1:C:184:ILE:H	1:C:184:ILE:CD1	2.00	0.63
1:C:71:LEU:HD23	1:C:177:LEU:HB3	1.79	0.63
1:A:74:LEU:HG	1:A:75:TRP:N	2.13	0.63
1:B:102:VAL:HA	1:B:106:GLN:HE21	1.62	0.63
1:C:84:ALA:N	1:C:85:PRO:HD2	2.13	0.63
1:B:5:LYS:HD3	1:B:129:ILE:HD11	1.80	0.62
1:A:31:ILE:HD12	1:A:31:ILE:O	1.99	0.62
1:A:126:SER:N	1:A:127:PRO:HD2	2.15	0.62
1:A:98:ALA:HB3	1:A:148:LEU:O	2.00	0.61
1:A:123:GLN:O	1:A:124:THR:HG23	1.99	0.61
1:A:61:THR:HG22	1:A:62:LEU:N	2.14	0.61
1:C:88:PRO:HG3	1:C:120:GLY:HA3	1.82	0.61
1:A:125:LEU:HD13	1:A:127:PRO:HG3	1.82	0.61
1:B:160:GLN:HG3	1:B:161:PRO:HD2	1.83	0.61
1:C:56:ILE:HG12	1:C:59:VAL:CG1	2.31	0.60
1:B:19:VAL:HG11	1:B:140:ARG:CZ	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:56:ILE:CG1	1:C:59:VAL:HG12	2.31	0.60
1:B:99:GLN:NE2	1:B:149:ASP:HB3	2.16	0.60
1:A:126:SER:O	1:A:128:LEU:N	2.34	0.60
1:C:107:ILE:HD12	1:C:156:SER:HB3	1.84	0.60
1:A:30:THR:HG22	1:A:179:MET:O	2.01	0.60
1:C:122:ILE:O	1:C:124:THR:HG23	2.02	0.60
1:C:160:GLN:HG3	1:C:161:PRO:HD2	1.84	0.59
1:A:184:ILE:CD1	1:A:184:ILE:H	2.14	0.59
1:A:125:LEU:HB3	1:A:127:PRO:HD2	1.86	0.58
1:B:81:THR:HG21	1:B:167:SER:O	2.03	0.58
1:A:99:GLN:O	1:A:99:GLN:HG2	2.04	0.57
1:A:154:LEU:HD23	1:A:154:LEU:N	2.19	0.57
1:A:58:SER:O	1:A:62:LEU:HD12	2.05	0.57
1:B:146:GLN:O	1:C:145:ILE:HD13	2.04	0.57
1:A:136:MET:HG2	1:B:184:ILE:CD1	2.28	0.57
1:B:75:TRP:CB	1:B:131:LYS:HA	2.34	0.57
1:B:103:THR:HG23	1:B:106:GLN:CD	2.25	0.57
1:B:89:THR:OG1	1:B:168:THR:HG21	2.05	0.57
1:C:104:PRO:O	1:C:107:ILE:HG13	2.04	0.57
1:C:77:THR:HB	1:C:79:HIS:CE1	2.40	0.57
1:A:153:LEU:HD22	1:A:154:LEU:N	2.19	0.57
1:B:84:ALA:N	1:B:85:PRO:HD2	2.20	0.56
1:A:81:THR:HG23	1:A:83:GLN:N	2.20	0.56
1:B:87:PHE:CG	1:B:164:PRO:HB3	2.41	0.56
1:A:126:SER:C	1:A:128:LEU:N	2.59	0.56
1:B:83:GLN:NE2	1:B:83:GLN:HA	2.20	0.56
1:C:87:PHE:CG	1:C:164:PRO:HB3	2.40	0.56
1:C:77:THR:HG23	1:C:129:ILE:HG12	1.87	0.56
1:B:17:ALA:HB1	1:B:138:GLN:HG2	1.88	0.56
1:C:73:SER:OG	1:C:176:THR:HG23	2.06	0.56
1:C:3:ILE:H	1:C:3:ILE:HD12	1.71	0.55
1:B:31:ILE:H	1:B:31:ILE:HD12	1.71	0.55
1:A:126:SER:N	1:A:127:PRO:CD	2.70	0.55
1:C:171:ILE:N	1:C:171:ILE:HD12	2.21	0.55
1:B:17:ALA:O	1:C:23:VAL:HG13	2.06	0.55
1:C:75:TRP:CE2	1:C:174:SER:HB2	2.42	0.55
1:B:63:THR:HA	1:B:179:MET:CE	2.36	0.55
1:B:20:LEU:HD12	1:B:20:LEU:O	2.06	0.55
1:C:3:ILE:N	1:C:3:ILE:HD12	2.21	0.54
1:C:88:PRO:HA	1:C:118:ILE:O	2.07	0.54
1:C:54:ALA:HA	1:C:141:VAL:HG13	1.87	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:LYS:HA	1:A:177:LEU:O	2.07	0.54
1:A:122:ILE:O	1:A:122:ILE:HG12	2.07	0.54
1:A:95:TRP:CE2	1:A:132:CYS:HB2	2.43	0.54
1:A:89:THR:OG1	1:A:168:THR:HG21	2.08	0.54
1:A:67:ARG:NH1	1:C:17:ALA:HB2	2.19	0.54
1:C:88:PRO:HG3	1:C:120:GLY:CA	2.37	0.54
1:A:49:ALA:HB3	1:A:155:ILE:HB	1.90	0.54
1:B:134:LEU:HD12	1:B:134:LEU:H	1.72	0.54
1:B:40:LEU:HD22	1:B:41:PHE:H	1.72	0.54
1:A:84:ALA:HB3	1:A:85:PRO:HD3	1.88	0.54
1:B:19:VAL:CG2	1:C:23:VAL:HG12	2.36	0.54
1:C:15:THR:HG23	1:C:15:THR:O	2.08	0.54
1:B:81:THR:CG2	1:B:83:GLN:H	2.07	0.53
1:C:184:ILE:HD12	1:C:184:ILE:N	2.07	0.53
1:C:34:PRO:O	1:C:35:PHE:HB3	2.09	0.53
1:A:117:CYS:O	1:A:123:GLN:HG2	2.09	0.53
1:A:40:LEU:HD12	1:A:41:PHE:H	1.72	0.53
1:B:96:VAL:HG13	1:B:100:SER:HB3	1.90	0.53
1:B:96:VAL:CG1	1:B:100:SER:HB3	2.39	0.53
1:B:63:THR:O	1:B:65:PHE:N	2.41	0.53
1:B:103:THR:H	1:B:106:GLN:HG3	1.74	0.53
1:C:74:LEU:HD12	1:C:75:TRP:H	1.74	0.53
1:C:19:VAL:HB	1:C:140:ARG:HD2	1.91	0.53
1:C:59:VAL:HG13	1:C:60:SER:N	2.24	0.52
1:B:18:THR:HB	1:C:20:LEU:HB3	1.91	0.52
1:B:98:ALA:HB3	1:B:148:LEU:O	2.09	0.52
1:A:129:ILE:HD13	1:A:129:ILE:H	1.75	0.52
1:B:77:THR:HA	1:B:128:LEU:O	2.10	0.52
1:B:56:ILE:O	1:B:57:ASP:C	2.48	0.52
1:A:96:VAL:HG12	1:A:152:LYS:HB3	1.91	0.52
1:A:122:ILE:O	1:A:122:ILE:CG1	2.58	0.52
1:C:133:PRO:HA	1:C:135:GLU:OE2	2.10	0.51
1:A:68:HIS:HE1	1:C:16:VAL:O	1.93	0.51
1:C:130:VAL:HG13	1:C:130:VAL:O	2.09	0.51
1:C:7:LEU:CD2	1:C:9:PRO:HD3	2.40	0.51
1:C:31:ILE:O	1:C:31:ILE:HD12	2.10	0.51
1:B:7:LEU:O	1:B:7:LEU:HD12	2.10	0.51
1:A:87:PHE:CG	1:A:164:PRO:HB3	2.45	0.51
1:C:134:LEU:CD2	1:C:134:LEU:H	2.17	0.51
1:A:184:ILE:HD12	1:A:184:ILE:N	2.17	0.51
1:B:20:LEU:HD23	1:B:145:ILE:HD13	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:LYS:HB3	1:C:186:ASP:OD1	2.10	0.50
1:A:95:TRP:CZ3	1:A:153:LEU:HB2	2.46	0.50
1:B:40:LEU:CD2	1:B:41:PHE:H	2.23	0.50
1:B:40:LEU:CD2	1:B:41:PHE:N	2.75	0.50
1:C:10:GLN:NE2	1:C:94:CYS:HA	2.18	0.50
1:B:51:LEU:CD1	1:B:51:LEU:N	2.74	0.50
1:A:33:GLN:HE21	1:A:34:PRO:HD2	1.77	0.50
1:B:180:HIS:HD2	1:B:181:SER:HB2	1.76	0.50
1:B:51:LEU:HD12	1:B:51:LEU:N	2.23	0.50
1:C:161:PRO:HG2	1:C:162:THR:H	1.77	0.50
1:A:58:SER:O	1:A:61:THR:HB	2.12	0.50
1:C:95:TRP:CE2	1:C:132:CYS:HB2	2.47	0.50
1:A:103:THR:HG23	1:A:106:GLN:CD	2.32	0.50
1:C:7:LEU:HD22	1:C:9:PRO:HD3	1.93	0.50
1:B:161:PRO:HG2	1:B:162:THR:H	1.76	0.50
1:A:96:VAL:O	1:A:137:MET:HE1	2.12	0.49
1:A:81:THR:HG21	1:A:83:GLN:HB2	1.92	0.49
1:C:138:GLN:O	1:C:151:PRO:HD3	2.12	0.49
1:B:19:VAL:HG11	1:B:140:ARG:NE	2.26	0.49
1:B:99:GLN:OE1	1:B:99:GLN:N	2.45	0.49
1:B:87:PHE:CD1	1:B:164:PRO:HB3	2.47	0.49
1:A:43:GLY:HA3	1:A:168:THR:HB	1.95	0.49
1:C:75:TRP:CB	1:C:131:LYS:HA	2.42	0.49
1:A:102:VAL:HG12	1:A:103:THR:N	2.26	0.49
1:C:7:LEU:C	1:C:7:LEU:HD22	2.33	0.49
1:B:59:VAL:O	1:B:60:SER:C	2.48	0.49
1:B:18:THR:HG22	1:C:22:THR:HA	1.95	0.49
1:C:177:LEU:HD12	1:C:177:LEU:O	2.13	0.49
1:B:81:THR:CG2	1:B:168:THR:HA	2.43	0.49
1:B:62:LEU:HB3	1:B:179:MET:HE3	1.94	0.49
1:B:102:VAL:HG12	1:B:103:THR:N	2.28	0.49
1:B:3:ILE:N	1:B:3:ILE:HD12	2.27	0.48
1:B:82:LEU:HD23	1:B:82:LEU:N	2.28	0.48
1:A:93:VAL:HG12	1:A:155:ILE:HG12	1.93	0.48
1:A:84:ALA:HB1	1:A:118:ILE:HG22	1.95	0.48
1:B:103:THR:HG23	1:B:106:GLN:HG2	1.95	0.48
1:C:61:THR:HA	1:C:64:THR:CG2	2.43	0.48
1:A:66:TYR:CD1	1:A:66:TYR:N	2.81	0.48
1:C:40:LEU:HD12	1:C:41:PHE:H	1.78	0.48
1:C:74:LEU:HD12	1:C:75:TRP:N	2.28	0.48
1:B:13:THR:HG23	1:B:14:VAL:N	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:81:THR:HG22	1:C:168:THR:O	2.14	0.48
1:C:15:THR:CG2	1:C:136:MET:HB2	2.43	0.48
1:B:19:VAL:C	1:C:20:LEU:HD22	2.33	0.48
1:C:96:VAL:O	1:C:152:LYS:HB3	2.14	0.48
1:B:31:ILE:HD12	1:B:31:ILE:O	2.14	0.47
1:C:183:LEU:C	1:C:183:LEU:HD13	2.35	0.47
1:C:30:THR:HB	1:C:180:HIS:HB3	1.94	0.47
1:B:75:TRP:HB3	1:B:131:LYS:HA	1.95	0.47
1:A:125:LEU:HD22	1:A:127:PRO:CD	2.44	0.47
1:C:68:HIS:HA	1:C:142:LYS:O	2.15	0.47
1:C:81:THR:CG2	1:C:83:GLN:H	2.13	0.47
1:C:81:THR:HG22	1:C:168:THR:HA	1.96	0.47
1:A:81:THR:CG2	1:A:168:THR:HA	2.44	0.47
1:B:81:THR:HG22	1:B:168:THR:C	2.35	0.47
1:B:63:THR:C	1:B:65:PHE:H	2.18	0.47
1:B:155:ILE:N	1:B:155:ILE:HD12	2.29	0.47
1:C:59:VAL:HG13	1:C:60:SER:H	1.80	0.47
1:B:83:GLN:HA	1:B:83:GLN:HE21	1.80	0.46
1:B:115:ILE:HD13	1:B:115:ILE:H	1.80	0.46
1:A:97:PRO:O	1:A:99:GLN:N	2.48	0.46
1:A:125:LEU:CB	1:A:127:PRO:HD2	2.46	0.46
1:C:42:ALA:HB1	1:C:157:ILE:HG21	1.95	0.46
1:C:149:ASP:OD2	1:C:149:ASP:N	2.46	0.46
1:C:43:GLY:HA3	1:C:168:THR:H	1.81	0.46
1:B:138:GLN:O	1:B:150:SER:HA	2.16	0.46
1:B:16:VAL:HG22	1:B:17:ALA:N	2.31	0.46
1:A:98:ALA:HA	1:A:152:LYS:HB2	1.97	0.46
1:A:125:LEU:HD13	1:A:127:PRO:CG	2.45	0.46
1:B:7:LEU:C	1:B:7:LEU:HD12	2.36	0.46
1:B:124:THR:C	1:B:125:LEU:HD23	2.37	0.46
1:C:107:ILE:O	1:C:113:GLY:HA3	2.16	0.45
1:B:75:TRP:HB2	1:B:131:LYS:HA	1.98	0.45
1:C:15:THR:HG22	1:C:136:MET:HB2	1.97	0.45
1:A:102:VAL:CG1	1:A:103:THR:N	2.79	0.45
1:A:32:LYS:NZ	1:A:72:GLU:HG2	2.31	0.45
1:C:123:GLN:HG3	1:C:123:GLN:O	2.15	0.45
1:B:153:LEU:HD22	1:B:154:LEU:N	2.31	0.45
1:C:7:LEU:HB2	1:C:129:ILE:O	2.17	0.45
1:C:56:ILE:O	1:C:57:ASP:C	2.55	0.45
1:B:53:ILE:O	1:B:56:ILE:HG23	2.16	0.45
1:A:123:GLN:HA	1:A:123:GLN:NE2	2.27	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:VAL:O	1:A:60:SER:C	2.55	0.45
1:B:81:THR:HG21	1:B:168:THR:HA	1.98	0.45
1:A:153:LEU:C	1:A:154:LEU:HD23	2.36	0.45
1:A:137:MET:CE	1:A:151:PRO:HA	2.46	0.45
1:B:50:SER:O	1:B:51:LEU:HD12	2.17	0.45
1:C:143:ASP:C	1:C:145:ILE:H	2.19	0.44
1:B:123:GLN:HG3	1:B:124:THR:N	2.31	0.44
1:C:62:LEU:H	1:C:62:LEU:HD12	1.82	0.44
1:C:19:VAL:HG12	1:C:138:GLN:HE22	1.82	0.44
1:C:39:VAL:HB	1:C:171:ILE:O	2.18	0.44
1:C:39:VAL:HG23	1:C:172:THR:HA	2.00	0.44
1:C:114:GLN:O	1:C:115:ILE:HD12	2.17	0.44
1:B:15:THR:HG21	1:C:184:ILE:HD11	1.99	0.44
1:C:60:SER:O	1:C:64:THR:HG23	2.18	0.44
1:C:102:VAL:HG11	1:C:110:THR:OG1	2.17	0.44
1:C:51:LEU:N	1:C:51:LEU:HD23	2.33	0.44
1:B:133:PRO:HB2	1:B:136:MET:HE3	2.00	0.43
1:A:168:THR:O	1:A:169:CYS:HB3	2.18	0.43
1:A:116:PHE:CD1	1:A:116:PHE:N	2.85	0.43
1:A:68:HIS:O	1:A:179:MET:HA	2.18	0.43
1:B:49:ALA:HB1	1:B:51:LEU:HD11	2.00	0.43
1:B:53:ILE:O	1:B:55:ASN:N	2.51	0.43
1:A:109:LYS:HB3	1:B:186:ASP:OD1	2.18	0.43
1:B:81:THR:HG23	1:B:82:LEU:N	2.33	0.43
1:C:99:GLN:NE2	1:C:99:GLN:N	2.66	0.43
1:A:138:GLN:O	1:A:151:PRO:HD3	2.17	0.43
1:A:71:LEU:HD23	1:A:176:THR:O	2.19	0.43
1:B:35:PHE:O	1:B:174:SER:HA	2.18	0.43
1:C:138:GLN:HB3	1:C:149:ASP:C	2.38	0.43
1:A:62:LEU:H	1:A:62:LEU:HD12	1.83	0.43
1:B:79:HIS:HE1	1:B:172:THR:OG1	2.01	0.43
1:C:68:HIS:HB2	1:C:180:HIS:CE1	2.54	0.43
1:C:78:ILE:O	1:C:80:PRO:HD3	2.19	0.43
1:A:138:GLN:HA	1:A:139:PRO:HD3	1.82	0.43
1:A:143:ASP:HB2	1:A:147:TYR:HE1	1.84	0.42
1:A:118:ILE:HD13	1:A:123:GLN:HG3	2.00	0.42
1:A:106:GLN:H	1:A:106:GLN:HG2	1.73	0.42
1:B:149:ASP:OD2	1:B:149:ASP:N	2.45	0.42
1:B:95:TRP:CZ3	1:B:153:LEU:HG	2.54	0.42
1:C:132:CYS:O	1:C:134:LEU:HD22	2.19	0.42
1:B:20:LEU:HD12	1:B:20:LEU:C	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:LEU:H	1:C:51:LEU:HD23	1.83	0.42
1:B:184:ILE:N	1:B:184:ILE:HD12	2.34	0.42
1:B:153:LEU:HD22	1:B:155:ILE:CD1	2.49	0.42
1:C:99:GLN:NE2	1:C:99:GLN:H	2.17	0.42
1:A:33:GLN:NE2	1:A:34:PRO:HD2	2.34	0.42
1:C:51:LEU:HD21	1:C:153:LEU:HB3	2.02	0.42
1:A:59:VAL:HG12	1:A:60:SER:N	2.35	0.42
1:B:56:ILE:HG13	1:B:56:ILE:O	2.20	0.42
1:C:82:LEU:N	1:C:82:LEU:HD23	2.35	0.42
1:B:187:THR:O	1:B:188:SER:C	2.58	0.42
1:C:99:GLN:OE1	1:C:149:ASP:HB3	2.18	0.42
1:A:118:ILE:CA	1:A:123:GLN:HG2	2.36	0.42
1:B:125:LEU:HD23	1:B:125:LEU:N	2.34	0.41
1:A:81:THR:HG21	1:A:168:THR:HA	2.03	0.41
1:A:74:LEU:HD23	1:A:132:CYS:CB	2.51	0.41
1:B:99:GLN:CA	1:B:99:GLN:OE1	2.68	0.41
1:C:183:LEU:O	1:C:183:LEU:HD13	2.20	0.41
1:C:34:PRO:HA	1:C:176:THR:HA	2.02	0.41
1:B:115:ILE:CD1	1:B:115:ILE:O	2.66	0.41
1:B:163:ALA:HB1	1:B:164:PRO:HD2	2.01	0.41
1:C:69:ALA:HA	1:C:178:SER:O	2.21	0.41
1:A:68:HIS:HD2	1:A:143:ASP:OD1	2.04	0.41
1:A:125:LEU:HD22	1:A:127:PRO:HD3	2.01	0.41
1:B:3:ILE:N	1:B:3:ILE:CD1	2.84	0.41
1:A:99:GLN:N	1:A:99:GLN:OE1	2.54	0.41
1:C:87:PHE:CD1	1:C:164:PRO:HB3	2.56	0.41
1:A:56:ILE:O	1:A:57:ASP:C	2.59	0.41
1:A:68:HIS:HA	1:A:142:LYS:O	2.21	0.41
1:C:99:GLN:HE21	1:C:100:SER:N	2.19	0.41
1:A:82:LEU:O	1:A:85:PRO:HD2	2.21	0.41
1:B:19:VAL:HG12	1:B:147:TYR:OH	2.20	0.41
1:C:81:THR:HG22	1:C:168:THR:C	2.41	0.40
1:B:71:LEU:HD13	1:B:177:LEU:HD21	2.00	0.40
1:C:50:SER:HA	1:C:153:LEU:O	2.21	0.40
1:C:19:VAL:CG1	1:C:138:GLN:HE22	2.33	0.40
1:B:82:LEU:CD2	1:B:82:LEU:H	2.35	0.40
1:C:185:THR:OG1	1:C:186:ASP:N	2.54	0.40
1:A:54:ALA:HA	1:A:59:VAL:HG11	2.02	0.40
1:A:91:VAL:HB	1:A:157:ILE:HD12	2.03	0.40
1:A:53:ILE:O	1:A:55:ASN:N	2.54	0.40
1:A:186:ASP:C	1:A:186:ASP:OD2	2.59	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	161/189 (85%)	129 (80%)	23 (14%)	9 (6%)	2	7
1	B	187/189 (99%)	158 (84%)	19 (10%)	10 (5%)	2	8
1	C	187/189 (99%)	164 (88%)	14 (8%)	9 (5%)	3	10
All	All	535/567 (94%)	451 (84%)	56 (10%)	28 (5%)	2	8

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	98	ALA
1	B	64	THR
1	B	188	SER
1	B	54	ALA
1	B	111	TYR
1	C	5	LYS
1	C	58	SER
1	C	59	VAL
1	C	64	THR
1	C	188	SER
1	A	122	ILE
1	A	124	THR
1	B	57	ASP
1	C	57	ASP
1	A	54	ALA
1	A	111	TYR
1	B	4	ASP
1	B	98	ALA
1	A	181	SER
1	A	188	SER
1	B	181	SER

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Mol	Chain	Res	Type
1	C	111	TYR
1	C	113	GLY
1	C	181	SER
1	A	118	ILE
1	B	101	PRO
1	B	113	GLY
1	A	126	SER

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	144/167 (86%)	97 (67%)	47 (33%)	0	1
1	B	167/167 (100%)	118 (71%)	49 (29%)	0	1
1	C	167/167 (100%)	145 (87%)	22 (13%)	5	14
All	All	478/501 (95%)	360 (75%)	118 (25%)	1	2

All (118) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	SER
1	A	30	THR
1	A	44	THR
1	A	45	LYS
1	A	46	ASP
1	A	57	ASP
1	A	60	SER
1	A	62	LEU
1	A	70	SER
1	A	72	GLU
1	A	74	LEU
1	A	77	THR
1	A	82	LEU
1	A	83	GLN
1	A	91	VAL

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Mol	Chain	Res	Type
1	A	96	VAL
1	A	99	GLN
1	A	103	THR
1	A	106	GLN
1	A	108	THR
1	A	110	THR
1	A	123	GLN
1	A	124	THR
1	A	125	LEU
1	A	128	LEU
1	A	129	ILE
1	A	131	LYS
1	A	134	LEU
1	A	136	MET
1	A	137	MET
1	A	143	ASP
1	A	145	ILE
1	A	148	LEU
1	A	150	SER
1	A	153	LEU
1	A	154	LEU
1	A	156	SER
1	A	162	THR
1	A	167	SER
1	A	168	THR
1	A	174	SER
1	A	176	THR
1	A	179	MET
1	A	184	ILE
1	A	187	THR
1	A	188	SER
1	A	189	THR
1	B	1	MET
1	B	3	ILE
1	B	13	THR
1	B	14	VAL
1	B	18	THR
1	B	19	VAL
1	B	23	VAL
1	B	29	PHE
1	B	31	ILE
1	B	39	VAL

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Mol	Chain	Res	Type
1	B	40	LEU
1	B	44	THR
1	B	45	LYS
1	B	46	ASP
1	B	50	SER
1	B	52	THR
1	B	55	ASN
1	B	60	SER
1	B	62	LEU
1	B	71	LEU
1	B	74	LEU
1	B	77	THR
1	B	99	GLN
1	B	103	THR
1	B	106	GLN
1	B	114	GLN
1	B	115	ILE
1	B	122	ILE
1	B	123	GLN
1	B	124	THR
1	B	134	LEU
1	B	136	MET
1	B	137	MET
1	B	138	GLN
1	B	140	ARG
1	B	143	ASP
1	B	145	ILE
1	B	148	LEU
1	B	150	SER
1	B	153	LEU
1	B	154	LEU
1	B	156	SER
1	B	167	SER
1	B	168	THR
1	B	174	SER
1	B	179	MET
1	B	181	SER
1	B	183	LEU
1	B	187	THR
1	C	3	ILE
1	C	7	LEU
1	C	13	THR

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Mol	Chain	Res	Type
1	C	14	VAL
1	C	16	VAL
1	C	30	THR
1	C	44	THR
1	C	45	LYS
1	C	57	ASP
1	C	70	SER
1	C	99	GLN
1	C	102	VAL
1	C	103	THR
1	C	108	THR
1	C	134	LEU
1	C	148	LEU
1	C	156	SER
1	C	176	THR
1	C	177	LEU
1	C	179	MET
1	C	184	ILE
1	C	189	THR

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

Mol	Chain	Res	Type
1	A	33	GLN
1	A	55	ASN
1	A	68	HIS
1	A	114	GLN
1	A	123	GLN
1	A	180	HIS
1	B	36	GLN
1	B	68	HIS
1	B	79	HIS
1	B	83	GLN
1	B	106	GLN
1	B	180	HIS
1	C	10	GLN
1	C	33	GLN
1	C	55	ASN
1	C	99	GLN
1	C	106	GLN
1	C	123	GLN
1	C	138	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 carbohydrates 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	163/189 (86%)	0.26	10 (6%) 25 18	19, 37, 112, 177	0
1	B	189/189 (100%)	0.35	6 (3%) 51 43	20, 38, 89, 153	0
1	C	189/189 (100%)	0.31	6 (3%) 51 43	20, 38, 83, 109	0
All	All	541/567 (95%)	0.31	22 (4%) 41 34	19, 37, 96, 177	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	123	GLN	8.2
1	B	25	GLY	7.4
1	B	24	PRO	6.2
1	A	125	LEU	5.5
1	B	23	VAL	5.1
1	A	122	ILE	5.0
1	C	25	GLY	3.8
1	A	124	THR	3.6
1	A	126	SER	3.4
1	B	26	PRO	3.4
1	C	22	THR	3.1
1	A	27	SER	2.9
1	A	127	PRO	2.8
1	A	120	GLY	2.7
1	A	163	ALA	2.6
1	B	1	MET	2.5
1	B	19	VAL	2.4
1	C	27	SER	2.3
1	C	24	PRO	2.3
1	C	26	PRO	2.2
1	C	29	PHE	2.2
1	A	29	PHE	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.