



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 30, 2026 – 11:18 PM UTC

PDB ID : 2IZW / pdb_00002izw
Title : Crystal structure of Ryegrass Mottle Virus
Authors : Plevka, P.; Tars, K.; Zeltins, A.; Balke, I.; Truve, E.; Liljas, L.
Deposited on : 2006-07-27
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 ⓘ 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
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

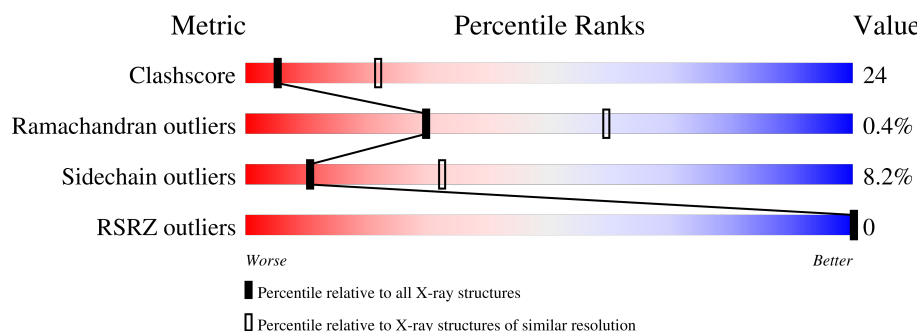
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (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 $\leq 5\%$. 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	234	
1	B	234	
1	C	234	

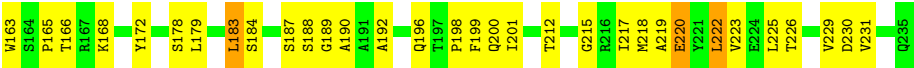
2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4324 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 RYEGRASS MOTTLE VIRUS COAT PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	178	Total	C	N	O	S	0	0	0
			1366	861	235	266	4			
1	B	179	Total	C	N	O	S	0	0	0
			1373	865	236	268	4			
1	C	208	Total	C	N	O	S	0	0	0
			1585	996	276	308	5			



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	277.64Å 298.72Å 392.49Å 90.00° 92.74° 90.00°	Depositor
Resolution (Å)	34.90 – 2.90 34.90 – 2.90	Depositor EDS
% Data completeness (in resolution range)	40.5 (34.90-2.90) 34.4 (34.90-2.90)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.294 , (Not available) 0.322 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	19.4	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 245.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.28$, $\langle L^2 \rangle = 0.12$	Xtriage
Estimated twinning fraction	0.209 for h,-k,-l	Xtriage
F_o, F_c correlation	0.23	EDS
Total number of atoms	4324	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.42 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.7261e-03.*

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 ⓘ

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	0.53	0/1399	1.07	8/1912 (0.4%)
1	B	0.55	0/1406	1.11	14/1922 (0.7%)
1	C	0.57	0/1623	1.13	15/2221 (0.7%)
All	All	0.55	0/4428	1.10	37/6055 (0.6%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	92	VAL	N-CA-C	-9.82	101.50	112.80
1	C	123	ILE	N-CA-C	-8.67	104.69	113.10
1	C	188	SER	N-CA-C	8.64	121.85	111.82
1	C	98	LEU	CA-C-N	8.28	127.58	118.97
1	C	98	LEU	C-N-CA	8.28	127.58	118.97
1	B	173	ASP	CA-C-N	8.23	128.18	119.05
1	B	173	ASP	C-N-CA	8.23	128.18	119.05
1	A	179	LEU	CA-C-N	8.16	127.72	119.24
1	A	179	LEU	C-N-CA	8.16	127.72	119.24
1	B	179	LEU	CA-C-N	7.87	127.42	119.24
1	B	179	LEU	C-N-CA	7.87	127.42	119.24
1	A	123	ILE	N-CA-C	-7.09	106.97	113.71
1	C	60	GLY	N-CA-C	-6.87	100.44	112.53
1	B	58	GLN	N-CA-C	6.74	119.65	110.55
1	C	92	VAL	N-CA-C	-6.73	103.92	113.07
1	A	234	ASN	CA-CB-CG	-6.66	105.94	112.60
1	C	128	THR	N-CA-C	6.66	119.41	110.35
1	B	164	SER	N-CA-C	6.49	118.26	109.24
1	C	61	ASP	N-CA-C	-6.47	104.64	113.30
1	C	63	THR	CA-C-N	6.44	128.16	120.23
1	C	63	THR	C-N-CA	6.44	128.16	120.23
1	B	197	THR	CA-C-N	6.18	125.75	119.19
1	B	197	THR	C-N-CA	6.18	125.75	119.19
1	B	162	ILE	N-CA-C	6.13	117.61	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	98	LEU	CA-C-N	5.88	127.20	119.84
1	A	98	LEU	C-N-CA	5.88	127.20	119.84
1	A	164	SER	N-CA-C	5.79	116.95	109.72
1	B	125	GLY	N-CA-C	5.78	118.69	112.33
1	B	76	SER	N-CA-C	5.74	117.94	110.43
1	B	155	ARG	N-CA-C	5.70	119.08	110.64
1	B	171	SER	N-CA-C	5.68	118.47	109.50
1	C	220	GLU	N-CA-C	-5.63	99.72	108.90
1	C	190	ALA	N-CA-C	-5.55	104.86	111.69
1	C	121	GLU	CA-C-N	5.43	125.19	119.76
1	C	121	GLU	C-N-CA	5.43	125.19	119.76
1	C	183	LEU	N-CA-C	5.28	118.41	109.76
1	A	210	VAL	N-CA-C	5.26	115.48	108.11

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	1366	0	1333	62	0
1	B	1373	0	1340	87	0
1	C	1585	0	1556	70	0
All	All	4324	0	4229	208	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 24.

All (208) 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:154:THR:HG22	1:C:230:ASP:H	1.30	0.95
1:B:91:ASN:HD21	1:B:197:THR:HB	1.33	0.93
1:B:92:VAL:HA	1:B:101:LEU:HD12	1.53	0.90
1:C:89:LEU:HD11	1:C:140:VAL:HG23	1.55	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:ASN:ND2	1:B:197:THR:HB	1.88	0.87
1:C:125:GLY:O	1:C:128:THR:HG22	1.77	0.83
1:A:66:LYS:HB2	1:A:222:LEU:HD22	1.59	0.82
1:A:147:THR:HG23	1:A:150:ARG:HD2	1.64	0.80
1:C:114:ARG:HB2	1:C:222:LEU:HD12	1.61	0.80
1:B:91:ASN:HD21	1:B:197:THR:CB	1.98	0.76
1:C:114:ARG:HA	1:C:114:ARG:HH11	1.50	0.75
1:B:92:VAL:O	1:B:102:SER:HA	1.88	0.73
1:C:76:SER:HB3	1:C:212:THR:H	1.52	0.73
1:B:119:THR:HG23	1:B:218:MET:HB2	1.71	0.73
1:A:124:VAL:HB	1:A:128:THR:HG21	1.70	0.72
1:C:97:GLU:C	1:C:99:PRO:HD3	2.14	0.72
1:B:91:ASN:HB3	1:B:93:ARG:H	1.55	0.72
1:B:216:ARG:HB2	1:B:218:MET:HE3	1.72	0.71
1:A:125:GLY:O	1:A:128:THR:HG23	1.92	0.69
1:A:130:GLY:C	1:A:162:ILE:HG13	2.18	0.68
1:B:145:SER:HB3	1:B:150:ARG:NH1	2.07	0.68
1:B:136:MET:HE1	1:B:172:TYR:HD1	1.58	0.67
1:B:136:MET:HE2	1:B:138:TYR:CZ	2.29	0.67
1:B:92:VAL:HA	1:B:101:LEU:CD1	2.24	0.67
1:B:154:THR:HG22	1:C:230:ASP:N	2.05	0.67
1:A:92:VAL:O	1:A:102:SER:HA	1.95	0.67
1:B:87:THR:HG21	1:B:143:VAL:HG13	1.77	0.66
1:C:92:VAL:O	1:C:102:SER:HA	1.94	0.66
1:A:136:MET:HE2	1:A:138:TYR:CZ	2.31	0.66
1:C:96:THR:HB	1:C:97:GLU:OE1	1.94	0.66
1:C:136:MET:HE2	1:C:138:TYR:CZ	2.31	0.66
1:B:167:ARG:HH11	1:B:167:ARG:HG3	1.62	0.64
1:B:154:THR:HB	1:C:229:VAL:HA	1.79	0.64
1:C:136:MET:HE3	1:C:199:PHE:HB2	1.80	0.64
1:B:66:LYS:HB2	1:B:222:LEU:CD2	2.28	0.63
1:A:216:ARG:HB2	1:A:218:MET:HE3	1.81	0.63
1:B:93:ARG:NH1	1:B:186:VAL:HG13	2.14	0.62
1:C:48:ILE:HD12	1:C:218:MET:HE3	1.81	0.62
1:A:234:ASN:O	1:A:235:GLN:HG3	2.00	0.62
1:A:155:ARG:HG3	1:B:227:ASP:HA	1.82	0.61
1:B:136:MET:HE2	1:B:138:TYR:OH	2.00	0.61
1:C:114:ARG:CB	1:C:222:LEU:HD12	2.30	0.61
1:B:150:ARG:O	1:B:154:THR:HG23	2.01	0.61
1:C:92:VAL:HG23	1:C:93:ARG:HG3	1.81	0.61
1:A:158:THR:HG23	1:A:168:LYS:HG3	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:VAL:HG22	1:A:235:GLN:O	2.02	0.60
1:B:114:ARG:CB	1:B:222:LEU:HD12	2.32	0.60
1:C:145:SER:HB3	1:C:150:ARG:NH1	2.16	0.59
1:C:123:ILE:O	1:C:123:ILE:HG22	2.01	0.59
1:B:234:ASN:O	1:B:235:GLN:HG3	2.02	0.59
1:A:91:ASN:HD22	1:A:93:ARG:H	1.50	0.59
1:A:115:TYR:CE2	1:A:222:LEU:HB2	2.37	0.59
1:C:199:PHE:HE1	1:C:201:ILE:HD11	1.68	0.59
1:B:148:ILE:HG22	1:B:152:MET:HE2	1.85	0.58
1:B:192:ALA:O	1:B:196:GLN:HG3	2.03	0.58
1:B:71:LEU:O	1:B:97:GLU:HB3	2.04	0.58
1:B:94:ASN:ND2	1:B:97:GLU:H	2.02	0.58
1:A:97:GLU:C	1:A:99:PRO:HD3	2.29	0.58
1:B:172:TYR:O	1:B:174:PRO:HD3	2.04	0.58
1:C:123:ILE:HG13	1:C:215:GLY:HA2	1.85	0.57
1:C:41:PRO:HG3	1:C:44:MET:HG3	1.85	0.57
1:C:134:MET:HE1	1:C:217:ILE:HG23	1.87	0.56
1:B:112:ARG:HG2	1:B:181:TRP:CZ3	2.40	0.56
1:B:179:LEU:HB2	1:B:180:PRO:HD2	1.86	0.56
1:C:187:SER:H	1:C:196:GLN:HE22	1.53	0.56
1:A:162:ILE:HD11	1:A:206:GLN:O	2.06	0.56
1:C:158:THR:HG21	1:C:168:LYS:O	2.06	0.55
1:C:192:ALA:O	1:C:196:GLN:HG3	2.06	0.55
1:B:91:ASN:HD21	1:B:197:THR:CG2	2.18	0.55
1:B:167:ARG:HG3	1:B:167:ARG:NH1	2.22	0.55
1:A:91:ASN:ND2	1:A:93:ARG:H	2.05	0.54
1:A:134:MET:O	1:A:157:GLY:HA2	2.07	0.54
1:C:97:GLU:O	1:C:99:PRO:HD3	2.07	0.54
1:C:54:GLN:HB2	1:C:66:LYS:HB3	1.88	0.54
1:C:68:SER:HA	1:C:219:ALA:O	2.07	0.54
1:B:193:GLY:O	1:B:197:THR:HG23	2.07	0.54
1:B:114:ARG:HB2	1:B:222:LEU:HD12	1.88	0.54
1:C:106:SER:HB2	1:C:231:VAL:HG23	1.90	0.54
1:C:30:GLN:HB3	1:C:31:PRO:HD2	1.90	0.53
1:A:163:TRP:H	1:A:163:TRP:CD1	2.26	0.53
1:A:155:ARG:HG2	1:B:228:PRO:HD2	1.91	0.53
1:B:94:ASN:ND2	1:B:97:GLU:HB2	2.24	0.53
1:B:66:LYS:HB2	1:B:222:LEU:HD23	1.91	0.52
1:B:97:GLU:C	1:B:99:PRO:HD3	2.35	0.52
1:B:216:ARG:HB2	1:B:218:MET:CE	2.37	0.52
1:B:71:LEU:HB2	1:B:217:ILE:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:MET:HE3	1:A:199:PHE:HB2	1.91	0.52
1:A:148:ILE:O	1:A:152:MET:HG3	2.09	0.52
1:C:61:ASP:HB3	1:C:226:THR:OG1	2.10	0.52
1:B:195:ILE:O	1:B:195:ILE:HG22	2.08	0.52
1:A:106:SER:HB2	1:A:231:VAL:HG13	1.92	0.52
1:C:56:VAL:O	1:C:64:PRO:HD2	2.09	0.51
1:A:136:MET:HE3	1:A:199:PHE:CD1	2.45	0.51
1:B:229:VAL:HG11	1:B:234:ASN:HB2	1.92	0.51
1:A:179:LEU:HB2	1:A:180:PRO:HD2	1.91	0.51
1:A:67:ASN:CG	1:A:68:SER:H	2.17	0.51
1:B:86:GLY:HA3	1:B:203:TRP:CE2	2.45	0.51
1:C:115:TYR:CE2	1:C:222:LEU:HB2	2.46	0.51
1:C:148:ILE:O	1:C:152:MET:HG3	2.10	0.51
1:A:89:LEU:HD11	1:A:140:VAL:HG13	1.92	0.50
1:A:93:ARG:HH12	1:A:235:GLN:HA	1.75	0.50
1:C:91:ASN:HB2	1:C:94:ASN:HB2	1.93	0.50
1:A:76:SER:HB3	1:A:212:THR:H	1.76	0.50
1:B:229:VAL:HG12	1:B:230:ASP:N	2.27	0.50
1:A:70:SER:HB2	1:A:218:MET:HE1	1.94	0.49
1:A:162:ILE:HD13	1:A:214:LEU:HD23	1.94	0.49
1:C:44:MET:HE2	1:C:165:PRO:HD2	1.93	0.49
1:C:66:LYS:HB2	1:C:222:LEU:HD23	1.93	0.49
1:A:216:ARG:HB2	1:A:218:MET:CE	2.42	0.49
1:C:120:TRP:HB3	1:C:166:THR:HB	1.93	0.49
1:C:108:TYR:CG	1:C:225:LEU:HB3	2.47	0.49
1:A:130:GLY:HA2	1:A:206:GLN:NE2	2.27	0.48
1:A:109:SER:OG	1:A:229:VAL:HG21	2.13	0.48
1:C:119:THR:HG23	1:C:218:MET:HB2	1.94	0.48
1:C:178:SER:O	1:C:179:LEU:HD12	2.12	0.48
1:A:92:VAL:HG23	1:A:93:ARG:HG3	1.94	0.48
1:C:92:VAL:HA	1:C:101:LEU:HG	1.94	0.48
1:C:198:PRO:O	1:C:199:PHE:HB3	2.13	0.48
1:B:62:ILE:HG23	1:B:224:GLU:HB3	1.95	0.48
1:C:91:ASN:HB3	1:C:93:ARG:H	1.79	0.48
1:A:72:VAL:CG1	1:A:73:ARG:N	2.76	0.48
1:C:66:LYS:HB2	1:C:222:LEU:CD2	2.44	0.48
1:C:97:GLU:OE1	1:C:97:GLU:N	2.46	0.48
1:B:62:ILE:CG2	1:B:224:GLU:HB3	2.44	0.48
1:B:134:MET:O	1:B:157:GLY:HA2	2.14	0.48
1:B:154:THR:CB	1:C:229:VAL:HA	2.42	0.48
1:A:71:LEU:HD21	1:A:98:LEU:HD21	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:175:GLU:CD	1:B:175:GLU:H	2.22	0.48
1:A:70:SER:CB	1:A:218:MET:HE1	2.45	0.47
1:C:120:TRP:H	1:C:166:THR:CG2	2.28	0.46
1:A:76:SER:HB2	1:A:214:LEU:HD11	1.97	0.46
1:B:112:ARG:NH1	1:B:224:GLU:OE2	2.45	0.46
1:A:118:PHE:HB3	1:A:134:MET:HE1	1.98	0.46
1:B:92:VAL:HB	1:B:105:GLY:HA3	1.97	0.46
1:A:136:MET:HE3	1:A:199:PHE:CG	2.50	0.46
1:B:66:LYS:HB2	1:B:222:LEU:HD22	1.98	0.46
1:C:89:LEU:HD13	1:C:200:GLN:CG	2.46	0.46
1:A:230:ASP:OD2	1:C:150:ARG:HD3	2.16	0.46
1:B:133:ALA:HA	1:B:158:THR:O	2.16	0.46
1:C:41:PRO:CG	1:C:44:MET:HG3	2.45	0.46
1:B:98:LEU:N	1:B:99:PRO:HD3	2.30	0.45
1:B:139:ASP:HB3	1:B:142:ASP:OD1	2.16	0.45
1:B:229:VAL:CG1	1:B:234:ASN:HB2	2.45	0.45
1:C:110:LYS:HG2	1:C:183:LEU:CD2	2.47	0.45
1:A:109:SER:OG	1:A:229:VAL:CG2	2.64	0.45
1:A:214:LEU:N	1:A:214:LEU:CD1	2.80	0.45
1:A:104:GLN:OE1	1:A:104:GLN:HA	2.17	0.45
1:B:135:ALA:HB1	1:B:154:THR:OG1	2.16	0.45
1:C:87:THR:HA	1:C:201:ILE:O	2.16	0.44
1:A:126:SER:HA	1:A:163:TRP:CG	2.53	0.44
1:A:120:TRP:CD1	1:A:122:PRO:HG3	2.53	0.44
1:B:234:ASN:O	1:B:235:GLN:CG	2.65	0.44
1:B:154:THR:CG2	1:C:230:ASP:H	2.16	0.44
1:C:46:THR:HB	1:C:121:GLU:OE1	2.18	0.44
1:C:120:TRP:H	1:C:166:THR:HG22	1.83	0.44
1:A:101:LEU:HD13	1:A:221:TYR:CE2	2.53	0.44
1:A:136:MET:HE3	1:A:199:PHE:CB	2.47	0.44
1:B:94:ASN:HD21	1:B:97:GLU:CB	2.31	0.44
1:A:98:LEU:N	1:A:99:PRO:HD3	2.31	0.44
1:B:139:ASP:HB3	1:B:142:ASP:CG	2.43	0.44
1:B:179:LEU:HD13	1:B:181:TRP:O	2.18	0.44
1:B:143:VAL:HG12	1:B:143:VAL:O	2.18	0.43
1:A:130:GLY:HA2	1:A:206:GLN:HE22	1.82	0.43
1:B:131:GLU:HG2	1:B:206:GLN:HG3	2.01	0.43
1:A:89:LEU:HD13	1:A:200:GLN:HG2	2.01	0.43
1:B:114:ARG:HB3	1:B:222:LEU:HD12	1.99	0.43
1:B:136:MET:HE2	1:B:138:TYR:CE2	2.54	0.43
1:C:222:LEU:C	1:C:222:LEU:HD13	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:136:MET:HE2	1:C:136:MET:HB3	1.74	0.43
1:B:92:VAL:HB	1:B:105:GLY:CA	2.49	0.43
1:B:94:ASN:HD21	1:B:97:GLU:HB2	1.82	0.42
1:B:70:SER:HA	1:B:218:MET:CE	2.49	0.42
1:B:75:THR:HA	1:B:213:THR:HA	2.00	0.42
1:C:89:LEU:HD13	1:C:200:GLN:HG3	2.01	0.42
1:C:126:SER:HA	1:C:163:TRP:CG	2.53	0.42
1:B:222:LEU:HD13	1:B:222:LEU:C	2.45	0.42
1:B:134:MET:HE3	1:B:217:ILE:HD13	2.01	0.42
1:B:186:VAL:O	1:B:187:SER:C	2.63	0.42
1:C:134:MET:HE2	1:C:217:ILE:HD13	2.00	0.42
1:A:92:VAL:HB	1:A:105:GLY:HA3	2.00	0.42
1:A:196:GLN:HE21	1:A:196:GLN:HB2	1.64	0.42
1:C:124:VAL:CG1	1:C:128:THR:HG21	2.50	0.42
1:C:132:VAL:CG2	1:C:133:ALA:N	2.82	0.42
1:A:172:TYR:O	1:A:174:PRO:HD3	2.20	0.42
1:B:161:PRO:HG2	1:B:164:SER:HB3	2.02	0.41
1:C:44:MET:CE	1:C:165:PRO:HD2	2.49	0.41
1:C:113:VAL:HB	1:C:172:TYR:CE2	2.55	0.41
1:C:132:VAL:HG22	1:C:133:ALA:N	2.34	0.41
1:B:68:SER:HA	1:B:219:ALA:O	2.20	0.41
1:B:94:ASN:HD21	1:B:97:GLU:H	1.68	0.41
1:A:178:SER:OG	1:B:110:LYS:HE2	2.20	0.41
1:B:140:VAL:HG23	1:B:194:ASN:OD1	2.21	0.41
1:B:234:ASN:C	1:B:235:GLN:HG3	2.45	0.41
1:A:158:THR:HG22	1:A:159:TRP:N	2.36	0.41
1:A:120:TRP:HB3	1:A:166:THR:HG23	2.03	0.41
1:A:155:ARG:CG	1:B:227:ASP:HA	2.51	0.41
1:A:71:LEU:CD2	1:A:98:LEU:HD21	2.51	0.41
1:B:117:HIS:HB3	1:B:169:ARG:HD2	2.03	0.41
1:B:134:MET:CE	1:B:217:ILE:HG23	2.50	0.41
1:C:115:TYR:CE2	1:C:222:LEU:CB	3.03	0.41
1:B:115:TYR:CD1	1:B:115:TYR:C	2.99	0.41
1:A:110:LYS:HE3	1:C:138:TYR:CD2	2.56	0.40
1:B:85:SER:HB3	1:B:146:ILE:HB	2.03	0.40
1:B:222:LEU:C	1:B:222:LEU:CD1	2.94	0.40
1:A:136:MET:HE2	1:A:136:MET:HB3	1.80	0.40
1:B:84:VAL:O	1:B:204:ALA:HA	2.21	0.40
1:C:93:ARG:HD2	1:C:184:SER:OG	2.20	0.40
1:C:139:ASP:HB3	1:C:142:ASP:OD2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	176/234 (75%)	158 (90%)	17 (10%)	1 (1%)	21	51
1	B	177/234 (76%)	166 (94%)	11 (6%)	0	100	100
1	C	206/234 (88%)	195 (95%)	10 (5%)	1 (0%)	24	54
All	All	559/702 (80%)	519 (93%)	38 (7%)	2 (0%)	30	59

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	189	GLY
1	A	148	ILE

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	147/194 (76%)	136 (92%)	11 (8%)	12	37
1	B	148/194 (76%)	131 (88%)	17 (12%)	5	18
1	C	171/194 (88%)	161 (94%)	10 (6%)	18	49
All	All	466/582 (80%)	428 (92%)	38 (8%)	10	32

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	87	THR

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Mol	Chain	Res	Type
1	A	89	LEU
1	A	127	ASN
1	A	144	THR
1	A	147	THR
1	A	149	GLU
1	A	169	ARG
1	A	175	GLU
1	A	200	GLN
1	A	216	ARG
1	A	222	LEU
1	B	71	LEU
1	B	75	THR
1	B	101	LEU
1	B	112	ARG
1	B	119	THR
1	B	127	ASN
1	B	143	VAL
1	B	144	THR
1	B	155	ARG
1	B	162	ILE
1	B	175	GLU
1	B	197	THR
1	B	206	GLN
1	B	216	ARG
1	B	222	LEU
1	B	223	VAL
1	B	231	VAL
1	C	32	THR
1	C	63	THR
1	C	71	LEU
1	C	114	ARG
1	C	119	THR
1	C	132	VAL
1	C	155	ARG
1	C	220	GLU
1	C	222	LEU
1	C	223	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	ASN

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Mol	Chain	Res	Type
1	A	206	GLN
1	A	235	GLN
1	B	58	GLN
1	B	91	ASN
1	B	94	ASN
1	B	206	GLN
1	B	235	GLN
1	C	58	GLN
1	C	67	ASN
1	C	91	ASN
1	C	104	GLN
1	C	117	HIS
1	C	196	GLN
1	C	200	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 [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	178/234 (76%)	-1.09	0 100 100	28, 45, 68, 112	0
1	B	179/234 (76%)	-1.09	0 100 100	28, 43, 67, 95	0
1	C	208/234 (88%)	-1.16	0 100 100	21, 43, 71, 108	0
All	All	565/702 (80%)	-1.12	0 100 100	21, 44, 68, 112	0

There are no RSRZ outliers to report.

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.