



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 02:19 PM UTC

PDB ID : 2WWS / pdb_00002wws
Title : Physalis Mottle Virus: Natural Empty Capsid
Authors : Sagurthi, S.R.; Rajaram, V.; Savithri, H.S.; Murthy, M.R.N.
Deposited on : 2009-10-27
Resolution : 3.90 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

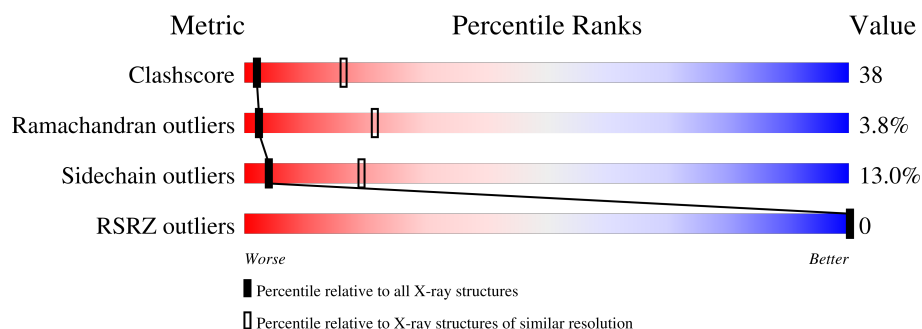
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.90 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	1034 (4.08-3.72)
Ramachandran outliers	187476	1251 (4.10-3.70)
Sidechain outliers	187428	1243 (4.10-3.70)
RSRZ outliers	180081	1269 (4.10-3.70)

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	188	
1	B	188	
1	C	188	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3976 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.

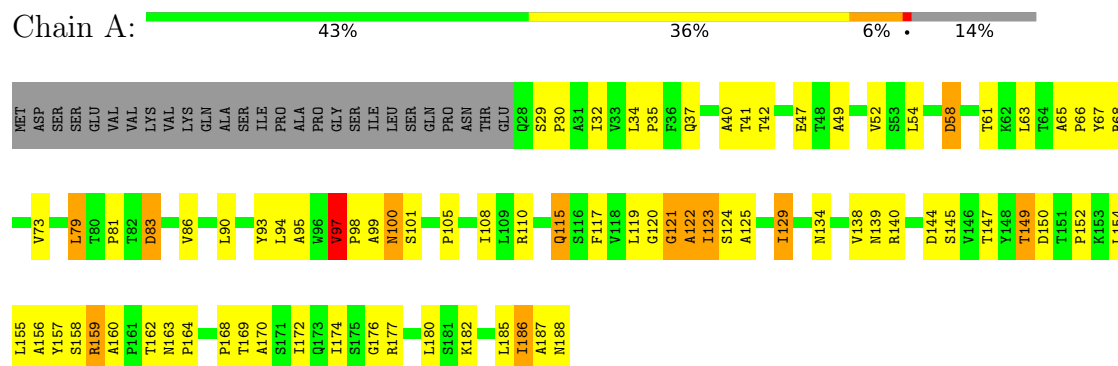
- Molecule 1 is a protein called COAT PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	S	0	0	0
			1210	774	203	230	3			
1	B	185	Total	C	N	O	S	0	0	0
			1383	882	232	266	3			
1	C	185	Total	C	N	O	S	0	0	0
			1383	882	232	266	3			

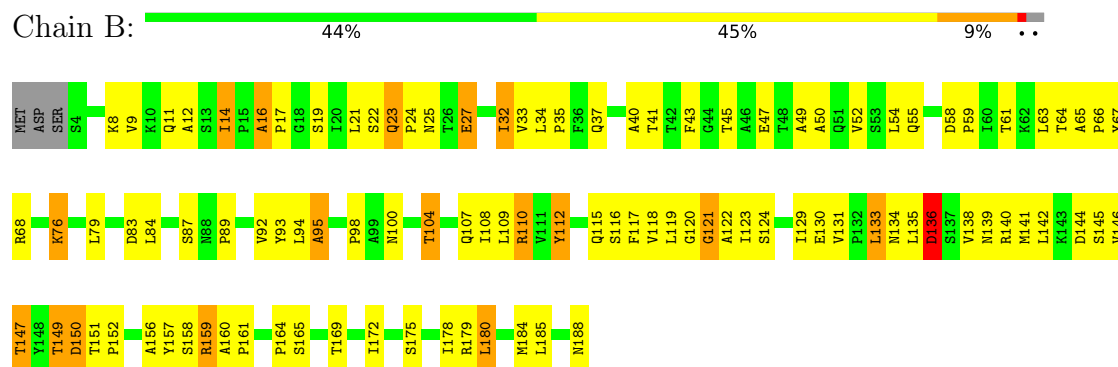
3 Residue-property plots

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.

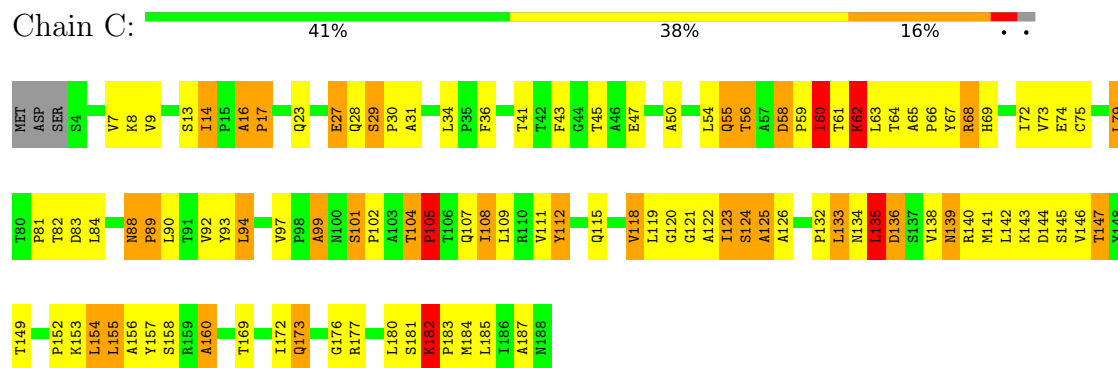
• Molecule 1: COAT PROTEIN



• Molecule 1: COAT PROTEIN



• Molecule 1: COAT PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	325.76Å 325.76Å 738.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.90 20.00 – 3.90	Depositor EDS
% Data completeness (in resolution range)	79.4 (20.00-3.90) 63.9 (20.00-3.90)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 3.94Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.230 , 0.242 0.546 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	59.7	Xtriage
Anisotropy	0.237	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.49 , 212.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.408 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.18	EDS
Total number of atoms	3976	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.64% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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	0.64	0/1236	1.18	9/1695 (0.5%)
1	B	0.66	0/1412	1.06	7/1936 (0.4%)
1	C	0.68	0/1412	1.16	16/1936 (0.8%)
All	All	0.66	0/4060	1.13	32/5567 (0.6%)

There are no bond length outliers.

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	121	GLY	N-CA-C	-10.43	98.84	114.90
1	A	100	ASN	N-CA-C	-10.34	101.16	113.88
1	B	136	ASP	N-CA-C	-9.01	101.60	114.12
1	C	101	SER	CA-C-N	7.92	127.58	119.82
1	C	101	SER	C-N-CA	7.92	127.58	119.82
1	C	29	SER	N-CA-C	7.91	117.23	108.14
1	B	16	ALA	CA-C-N	7.67	129.43	119.84
1	B	16	ALA	C-N-CA	7.67	129.43	119.84
1	C	62	LYS	N-CA-C	-7.48	104.03	113.01
1	B	122	ALA	N-CA-C	-7.48	104.30	113.50
1	C	58	ASP	N-CA-C	-7.28	99.45	110.07
1	C	16	ALA	CA-C-N	7.02	128.62	119.84
1	C	16	ALA	C-N-CA	7.02	128.62	119.84
1	C	182	LYS	CA-C-N	6.77	127.30	119.99
1	C	182	LYS	C-N-CA	6.77	127.30	119.99
1	A	97	VAL	CA-C-N	6.73	126.41	119.82
1	A	97	VAL	C-N-CA	6.73	126.41	119.82
1	C	43	PHE	N-CA-C	6.46	121.77	113.18
1	A	97	VAL	N-CA-C	6.29	114.35	109.19
1	C	136	ASP	N-CA-C	-6.15	106.38	114.31
1	A	182	LYS	CA-C-N	6.03	127.38	119.84
1	A	182	LYS	C-N-CA	6.03	127.38	119.84
1	A	115	GLN	N-CA-C	5.57	117.65	109.24
1	B	108	ILE	N-CA-C	5.50	115.59	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	73	VAL	N-CA-C	-5.45	107.49	111.90
1	C	160	ALA	CA-C-N	5.42	125.36	119.78
1	C	160	ALA	C-N-CA	5.42	125.36	119.78
1	C	88	ASN	CA-C-N	5.18	126.32	119.84
1	C	88	ASN	C-N-CA	5.18	126.32	119.84
1	B	131	VAL	CA-C-N	5.09	125.53	120.14
1	B	131	VAL	C-N-CA	5.09	125.53	120.14
1	C	13	SER	N-CA-C	5.01	115.92	108.60

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	1210	0	1254	81	0
1	B	1383	0	1435	102	0
1	C	1383	0	1435	132	0
All	All	3976	0	4124	305	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 38.

All (305) 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:89:PRO:HA	1:B:120:GLY:HA3	1.17	1.16
1:C:68:ARG:HG2	1:C:182:LYS:HZ2	1.08	1.09
1:C:182:LYS:HE2	1:C:182:LYS:C	1.81	1.05
1:C:139:ASN:HD22	1:C:140:ARG:N	1.58	1.01
1:C:31:ALA:HB2	1:C:181:SER:HB3	1.45	0.96
1:C:146:VAL:HG12	1:C:147:THR:H	1.31	0.94
1:A:138:VAL:HG21	1:A:152:PRO:HG3	1.50	0.93
1:B:11:GLN:HG2	1:B:12:ALA:H	1.33	0.91
1:A:159:ARG:HB3	1:A:159:ARG:HH11	1.34	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:68:ARG:HG2	1:C:182:LYS:NZ	1.90	0.87
1:C:68:ARG:CG	1:C:182:LYS:HZ2	1.87	0.86
1:B:139:ASN:HD22	1:B:140:ARG:N	1.73	0.86
1:B:32:ILE:HD13	1:B:33:VAL:H	1.40	0.83
1:C:59:PRO:C	1:C:61:THR:H	1.85	0.83
1:B:139:ASN:HB2	1:B:150:ASP:HB2	1.60	0.82
1:A:94:LEU:HD21	1:A:172:ILE:HD13	1.61	0.80
1:C:139:ASN:ND2	1:C:141:MET:H	1.81	0.79
1:B:12:ALA:HB1	1:B:112:TYR:CE2	2.18	0.79
1:C:146:VAL:HG12	1:C:147:THR:N	1.97	0.78
1:B:159:ARG:HH11	1:B:159:ARG:HB2	1.47	0.78
1:C:90:LEU:HD21	1:C:119:LEU:HB2	1.65	0.78
1:C:72:ILE:HD11	1:C:75:CYS:HB2	1.67	0.77
1:A:97:VAL:HG12	1:A:99:ALA:H	1.51	0.76
1:A:52:VAL:CG2	1:A:154:LEU:HB3	2.16	0.75
1:A:40:ALA:HB1	1:A:52:VAL:HG11	1.68	0.75
1:B:49:ALA:HB2	1:B:157:TYR:HD2	1.52	0.74
1:B:24:PRO:HA	1:B:27:GLU:CG	2.17	0.74
1:B:43:PHE:CE1	1:B:172:ILE:HD13	2.23	0.73
1:B:32:ILE:HD13	1:B:33:VAL:N	2.04	0.73
1:C:14:ILE:HD13	1:C:14:ILE:H	1.54	0.72
1:B:89:PRO:HA	1:B:120:GLY:CA	2.10	0.72
1:C:108:ILE:O	1:C:111:VAL:HG22	1.89	0.71
1:A:41:THR:HG22	1:A:42:THR:N	2.04	0.71
1:B:180:LEU:HD22	1:B:180:LEU:H	1.56	0.70
1:C:31:ALA:CB	1:C:181:SER:HB3	2.21	0.70
1:A:97:VAL:CG1	1:A:101:SER:HB3	2.21	0.70
1:C:182:LYS:HE2	1:C:182:LYS:O	1.91	0.70
1:C:64:THR:HG22	1:C:180:LEU:HD21	1.73	0.70
1:A:129:ILE:HD12	1:A:129:ILE:N	2.06	0.69
1:C:139:ASN:HD22	1:C:139:ASN:C	1.98	0.69
1:B:139:ASN:HD22	1:B:139:ASN:C	1.99	0.68
1:A:47:GLU:HB2	1:A:160:ALA:HB2	1.77	0.67
1:C:138:VAL:HG21	1:C:152:PRO:HB3	1.75	0.67
1:B:24:PRO:HA	1:B:27:GLU:HG3	1.76	0.67
1:A:81:PRO:HB3	1:A:119:LEU:HD13	1.77	0.67
1:C:139:ASN:HD22	1:C:140:ARG:H	1.40	0.67
1:B:37:GLN:HG3	1:B:175:SER:HB3	1.76	0.66
1:B:159:ARG:HB2	1:B:159:ARG:NH1	2.10	0.66
1:A:95:ALA:HB2	1:A:108:ILE:HG23	1.77	0.66
1:A:186:ILE:HD13	1:A:186:ILE:H	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:LYS:HG2	1:B:130:GLU:HB3	1.77	0.66
1:B:43:PHE:HE1	1:B:172:ILE:HD13	1.61	0.66
1:B:180:LEU:HD22	1:B:180:LEU:N	2.10	0.66
1:A:100:ASN:HD21	1:A:149:THR:HB	1.61	0.66
1:C:54:LEU:HD12	1:C:60:ILE:HD13	1.77	0.66
1:C:82:THR:HG22	1:C:84:LEU:H	1.59	0.66
1:A:97:VAL:HG11	1:A:101:SER:HB3	1.77	0.66
1:A:121:GLY:O	1:A:123:ILE:HG13	1.95	0.65
1:B:49:ALA:HB2	1:B:157:TYR:CD2	2.31	0.65
1:A:83:ASP:O	1:A:86:VAL:HG22	1.97	0.65
1:C:9:VAL:HG12	1:C:115:GLN:HE22	1.62	0.64
1:A:105:PRO:O	1:A:108:ILE:HG12	1.97	0.64
1:C:47:GLU:HB2	1:C:160:ALA:HB2	1.80	0.63
1:C:90:LEU:CD2	1:C:119:LEU:HB2	2.28	0.63
1:B:185:LEU:HD12	1:B:185:LEU:H	1.63	0.63
1:C:60:ILE:HD12	1:C:142:LEU:HD13	1.80	0.63
1:B:139:ASN:ND2	1:B:141:MET:H	1.96	0.63
1:B:150:ASP:OD2	1:C:145:SER:HB3	1.99	0.63
1:B:185:LEU:HD12	1:B:185:LEU:N	2.14	0.63
1:C:59:PRO:O	1:C:61:THR:N	2.32	0.62
1:A:108:ILE:HD11	1:A:155:LEU:HB3	1.82	0.62
1:C:79:LEU:HD22	1:C:119:LEU:HD11	1.81	0.62
1:A:41:THR:CG2	1:A:42:THR:N	2.63	0.62
1:A:98:PRO:HA	1:A:138:VAL:HG12	1.80	0.62
1:A:122:ALA:C	1:A:124:SER:H	2.08	0.62
1:C:68:ARG:HG2	1:C:182:LYS:CD	2.30	0.61
1:A:34:LEU:HD23	1:A:34:LEU:H	1.64	0.61
1:B:104:THR:HG23	1:B:107:GLN:OE1	2.01	0.61
1:C:72:ILE:HD11	1:C:75:CYS:CB	2.30	0.61
1:B:79:LEU:HD22	1:B:119:LEU:HD21	1.83	0.61
1:C:69:HIS:CE1	1:C:141:MET:HE1	2.36	0.60
1:C:123:ILE:HD13	1:C:123:ILE:O	2.01	0.60
1:A:129:ILE:HD12	1:A:129:ILE:H	1.65	0.60
1:A:93:TYR:O	1:A:156:ALA:HA	2.02	0.60
1:B:23:GLN:N	1:B:24:PRO:CD	2.64	0.60
1:B:11:GLN:HG2	1:B:12:ALA:N	2.10	0.60
1:C:139:ASN:ND2	1:C:140:ARG:N	2.41	0.60
1:B:55:GLN:HB2	1:B:151:THR:HG21	1.84	0.60
1:C:34:LEU:HD12	1:C:180:LEU:HD12	1.84	0.60
1:B:138:VAL:HG21	1:B:152:PRO:HB3	1.82	0.59
1:C:16:ALA:N	1:C:17:PRO:HD3	2.18	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:41:THR:HG21	1:C:50:ALA:HB3	1.83	0.59
1:A:54:LEU:HD21	1:A:174:ILE:HG21	1.85	0.59
1:A:97:VAL:HG13	1:A:98:PRO:HD2	1.85	0.58
1:A:159:ARG:HB3	1:A:159:ARG:NH1	2.13	0.58
1:B:117:PHE:CD1	1:B:129:ILE:HD13	2.37	0.58
1:B:120:GLY:O	1:B:121:GLY:C	2.44	0.58
1:C:155:LEU:HD23	1:C:155:LEU:H	1.67	0.58
1:C:9:VAL:CG1	1:C:115:GLN:HE22	2.16	0.58
1:C:81:PRO:HG3	1:C:119:LEU:HD22	1.85	0.58
1:B:135:LEU:HD23	1:B:135:LEU:H	1.67	0.57
1:C:58:ASP:HB2	1:C:59:PRO:HA	1.85	0.57
1:C:59:PRO:C	1:C:61:THR:N	2.53	0.57
1:A:98:PRO:O	1:A:99:ALA:HB3	2.05	0.57
1:B:41:THR:HG21	1:B:50:ALA:HB3	1.86	0.57
1:C:93:TYR:CZ	1:C:109:LEU:HD11	2.40	0.57
1:A:187:ALA:O	1:A:188:ASN:HB2	2.02	0.57
1:B:23:GLN:N	1:B:24:PRO:HD2	2.20	0.57
1:C:139:ASN:ND2	1:C:139:ASN:C	2.63	0.57
1:C:93:TYR:CE2	1:C:109:LEU:HD11	2.40	0.57
1:C:82:THR:HB	1:C:169:THR:O	2.04	0.57
1:B:178:ILE:HG22	1:B:179:ARG:N	2.20	0.57
1:C:120:GLY:H	1:C:125:ALA:HB1	1.68	0.56
1:B:24:PRO:HA	1:B:27:GLU:HG2	1.87	0.56
1:C:146:VAL:CG1	1:C:147:THR:H	2.11	0.56
1:C:93:TYR:O	1:C:156:ALA:HA	2.05	0.56
1:B:67:TYR:CG	1:B:180:LEU:HB3	2.41	0.55
1:B:49:ALA:CB	1:B:157:TYR:HD2	2.17	0.55
1:C:36:PHE:CD2	1:C:60:ILE:HG22	2.41	0.55
1:A:42:THR:OG1	1:A:168:PRO:HB3	2.07	0.55
1:A:52:VAL:HG22	1:A:154:LEU:HB3	1.88	0.55
1:A:86:VAL:HA	1:A:121:GLY:H	1.72	0.55
1:B:123:ILE:HG13	1:B:124:SER:N	2.21	0.55
1:C:118:VAL:HG22	1:C:118:VAL:O	2.05	0.55
1:C:133:LEU:HD13	1:C:135:LEU:CD1	2.36	0.55
1:C:68:ARG:HG2	1:C:182:LYS:HD3	1.88	0.55
1:B:93:TYR:O	1:B:156:ALA:HA	2.06	0.55
1:B:110:ARG:CB	1:B:110:ARG:HH11	2.19	0.55
1:C:155:LEU:HD23	1:C:155:LEU:N	2.20	0.55
1:C:73:VAL:HG13	1:C:74:GLU:N	2.22	0.55
1:C:58:ASP:HB3	1:C:61:THR:OG1	2.06	0.55
1:B:16:ALA:HA	1:C:182:LYS:HG2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:92:VAL:HG12	1:C:93:TYR:N	2.22	0.55
1:A:52:VAL:HG23	1:A:154:LEU:HB3	1.86	0.54
1:C:65:ALA:N	1:C:66:PRO:HD2	2.23	0.54
1:A:120:GLY:C	1:A:122:ALA:N	2.61	0.54
1:B:161:PRO:HG2	1:B:164:PRO:HG3	1.89	0.54
1:C:105:PRO:O	1:C:108:ILE:HG12	2.08	0.53
1:A:159:ARG:HH11	1:A:159:ARG:CB	2.13	0.53
1:A:186:ILE:HD13	1:A:186:ILE:N	2.23	0.53
1:C:68:ARG:HA	1:C:143:LYS:HE2	1.90	0.53
1:C:120:GLY:N	1:C:125:ALA:HB1	2.23	0.53
1:A:97:VAL:HG13	1:A:101:SER:HB3	1.91	0.52
1:B:119:LEU:HD12	1:B:119:LEU:H	1.74	0.52
1:A:41:THR:CG2	1:A:42:THR:H	2.22	0.52
1:C:73:VAL:CG1	1:C:177:ARG:HB2	2.39	0.52
1:C:104:THR:HG23	1:C:107:GLN:OE1	2.10	0.52
1:C:154:LEU:HD22	1:C:155:LEU:N	2.24	0.52
1:A:34:LEU:HD23	1:A:34:LEU:N	2.25	0.52
1:A:58:ASP:O	1:A:61:THR:HB	2.10	0.52
1:C:60:ILE:O	1:C:60:ILE:HG13	2.08	0.52
1:B:144:ASP:OD1	1:B:145:SER:N	2.38	0.51
1:C:67:TYR:CD2	1:C:180:LEU:HB3	2.44	0.51
1:C:107:GLN:O	1:C:108:ILE:C	2.50	0.51
1:C:154:LEU:HD22	1:C:154:LEU:C	2.34	0.51
1:B:159:ARG:HH11	1:B:159:ARG:CB	2.20	0.51
1:C:61:THR:C	1:C:63:LEU:N	2.68	0.51
1:C:58:ASP:HA	1:C:61:THR:OG1	2.11	0.51
1:C:111:VAL:O	1:C:112:TYR:C	2.53	0.51
1:C:176:GLY:C	1:C:177:ARG:HG2	2.36	0.51
1:B:110:ARG:HH12	1:C:187:ALA:HB1	1.75	0.51
1:A:61:THR:C	1:A:63:LEU:N	2.69	0.51
1:A:157:TYR:CG	1:A:158:SER:N	2.79	0.50
1:B:55:GLN:CB	1:B:151:THR:HG21	2.41	0.50
1:A:122:ALA:O	1:A:124:SER:N	2.45	0.50
1:B:40:ALA:HB1	1:B:52:VAL:HG21	1.94	0.50
1:B:146:VAL:HG12	1:B:147:THR:N	2.26	0.50
1:B:123:ILE:HG13	1:B:124:SER:H	1.77	0.49
1:A:108:ILE:CD1	1:A:155:LEU:HB3	2.42	0.49
1:B:19:SER:OG	1:C:23:GLN:NE2	2.44	0.49
1:B:19:SER:HB2	1:C:145:SER:HB2	1.94	0.49
1:B:178:ILE:CG2	1:B:179:ARG:N	2.76	0.49
1:C:14:ILE:HD13	1:C:14:ILE:N	2.23	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:182:LYS:HE2	1:C:183:PRO:N	2.26	0.49
1:B:150:ASP:OD1	1:C:68:ARG:HD2	2.13	0.49
1:B:32:ILE:CD1	1:B:33:VAL:N	2.74	0.49
1:B:98:PRO:HB3	1:B:138:VAL:HG12	1.95	0.49
1:C:111:VAL:O	1:C:111:VAL:HG23	2.13	0.49
1:A:129:ILE:H	1:A:129:ILE:CD1	2.26	0.49
1:B:16:ALA:HB1	1:C:68:ARG:NH2	2.28	0.49
1:C:94:LEU:HD11	1:C:172:ILE:HD13	1.95	0.49
1:B:22:SER:C	1:B:24:PRO:HD2	2.38	0.48
1:C:157:TYR:CG	1:C:158:SER:N	2.81	0.48
1:C:29:SER:HB2	1:C:30:PRO:HD2	1.94	0.48
1:B:66:PRO:HG2	1:B:67:TYR:CE2	2.48	0.48
1:B:110:ARG:NH1	1:C:187:ALA:HB1	2.28	0.48
1:C:68:ARG:O	1:C:68:ARG:HG3	2.13	0.48
1:C:181:SER:O	1:C:182:LYS:O	2.31	0.48
1:B:9:VAL:HG23	1:B:9:VAL:O	2.13	0.48
1:B:34:LEU:HD13	1:B:35:PRO:HD2	1.95	0.48
1:C:133:LEU:HD11	1:C:152:PRO:HB3	1.96	0.48
1:B:100:ASN:HD22	1:B:149:THR:HB	1.79	0.48
1:C:82:THR:HG22	1:C:83:ASP:N	2.28	0.48
1:C:82:THR:CG2	1:C:83:ASP:N	2.76	0.47
1:C:134:ASN:OD1	1:C:136:ASP:HB2	2.13	0.47
1:A:100:ASN:ND2	1:A:149:THR:HB	2.28	0.47
1:A:185:LEU:N	1:A:185:LEU:HD12	2.29	0.47
1:C:61:THR:C	1:C:63:LEU:H	2.21	0.47
1:B:12:ALA:HB1	1:B:112:TYR:CD2	2.49	0.47
1:B:79:LEU:HD22	1:B:119:LEU:CD2	2.45	0.47
1:A:163:ASN:CG	1:A:163:ASN:O	2.58	0.47
1:B:93:TYR:CE2	1:B:116:SER:HB2	2.50	0.47
1:A:144:ASP:CG	1:A:145:SER:H	2.22	0.47
1:B:119:LEU:HD12	1:B:119:LEU:N	2.31	0.46
1:A:110:ARG:HD2	1:B:188:ASN:OXT	2.16	0.46
1:C:185:LEU:HD12	1:C:185:LEU:H	1.80	0.46
1:A:97:VAL:HG13	1:A:98:PRO:CD	2.45	0.46
1:C:144:ASP:CG	1:C:145:SER:H	2.23	0.46
1:A:41:THR:HG22	1:A:42:THR:H	1.79	0.46
1:B:14:ILE:HG23	1:B:14:ILE:O	2.16	0.46
1:B:58:ASP:N	1:B:59:PRO:HD2	2.31	0.46
1:C:182:LYS:O	1:C:182:LYS:CE	2.62	0.46
1:C:72:ILE:HG23	1:C:72:ILE:O	2.16	0.45
1:C:69:HIS:HA	1:C:143:LYS:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:GLY:C	1:A:122:ALA:H	2.21	0.45
1:B:61:THR:C	1:B:63:LEU:N	2.73	0.45
1:B:94:LEU:N	1:B:94:LEU:HD12	2.31	0.45
1:A:176:GLY:C	1:A:177:ARG:HG3	2.42	0.45
1:A:90:LEU:HD23	1:A:90:LEU:N	2.31	0.45
1:A:68:ARG:O	1:A:68:ARG:HG2	2.17	0.45
1:A:67:TYR:CD2	1:A:180:LEU:HD13	2.51	0.45
1:B:135:LEU:H	1:B:135:LEU:CD2	2.30	0.45
1:C:8:LYS:HE2	1:C:132:PRO:HG3	1.99	0.45
1:C:72:ILE:HG22	1:C:140:ARG:O	2.17	0.45
1:B:65:ALA:HB3	1:B:66:PRO:HD3	1.99	0.44
1:B:94:LEU:O	1:B:95:ALA:HB2	2.17	0.44
1:C:97:VAL:HG22	1:C:155:LEU:CD2	2.47	0.44
1:A:81:PRO:HA	1:A:170:ALA:CB	2.47	0.44
1:C:185:LEU:HD12	1:C:185:LEU:N	2.32	0.44
1:B:61:THR:O	1:B:63:LEU:N	2.51	0.44
1:C:16:ALA:N	1:C:17:PRO:CD	2.79	0.44
1:C:123:ILE:CG2	1:C:124:SER:N	2.80	0.44
1:A:79:LEU:CD2	1:A:119:LEU:HD11	2.47	0.44
1:A:122:ALA:C	1:A:124:SER:N	2.76	0.44
1:A:47:GLU:HB2	1:A:160:ALA:CB	2.45	0.44
1:A:117:PHE:CE2	1:A:129:ILE:HG23	2.53	0.43
1:A:187:ALA:O	1:A:188:ASN:CB	2.66	0.43
1:C:92:VAL:CG1	1:C:93:TYR:N	2.81	0.43
1:B:14:ILE:O	1:B:14:ILE:HD13	2.18	0.43
1:B:16:ALA:N	1:B:17:PRO:CD	2.80	0.43
1:A:97:VAL:HA	1:A:98:PRO:HD3	1.87	0.43
1:C:65:ALA:N	1:C:66:PRO:CD	2.82	0.43
1:B:157:TYR:CG	1:B:158:SER:N	2.86	0.43
1:C:27:GLU:CD	1:C:27:GLU:H	2.27	0.43
1:A:37:GLN:HA	1:A:174:ILE:O	2.19	0.43
1:C:14:ILE:N	1:C:14:ILE:CD1	2.82	0.43
1:C:31:ALA:HB1	1:C:180:LEU:O	2.18	0.43
1:C:68:ARG:HG2	1:C:182:LYS:CE	2.49	0.43
1:C:184:MET:HG3	1:C:185:LEU:N	2.33	0.43
1:A:138:VAL:HB	1:A:139:ASN:H	1.61	0.43
1:B:84:LEU:HB3	1:B:169:THR:HG22	2.01	0.43
1:A:98:PRO:O	1:A:150:ASP:HA	2.18	0.43
1:B:120:GLY:O	1:B:121:GLY:O	2.37	0.43
1:C:89:PRO:HB3	1:C:121:GLY:HA3	2.00	0.43
1:B:47:GLU:HB2	1:B:160:ALA:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:92:VAL:O	1:B:93:TYR:HD2	2.02	0.43
1:B:185:LEU:H	1:B:185:LEU:CD1	2.30	0.43
1:A:49:ALA:HB2	1:A:157:TYR:HD1	1.83	0.42
1:B:110:ARG:HH11	1:B:110:ARG:HB2	1.84	0.42
1:C:7:VAL:HG13	1:C:7:VAL:O	2.17	0.42
1:C:59:PRO:HA	1:C:62:LYS:HG3	2.00	0.42
1:C:125:ALA:O	1:C:126:ALA:HB3	2.18	0.42
1:B:45:THR:HA	1:B:165:SER:O	2.19	0.42
1:B:87:SER:C	1:B:89:PRO:HD3	2.45	0.42
1:C:99:ALA:HB3	1:C:149:THR:O	2.18	0.42
1:A:61:THR:C	1:A:63:LEU:H	2.28	0.42
1:B:92:VAL:HG12	1:B:93:TYR:N	2.34	0.42
1:B:129:ILE:HG22	1:B:130:GLU:N	2.35	0.42
1:C:73:VAL:CG1	1:C:74:GLU:N	2.82	0.42
1:A:154:LEU:HD12	1:A:155:LEU:N	2.35	0.42
1:C:82:THR:C	1:C:84:LEU:N	2.78	0.42
1:A:29:SER:OG	1:A:30:PRO:CD	2.68	0.42
1:C:54:LEU:O	1:C:56:THR:N	2.52	0.42
1:C:101:SER:HA	1:C:102:PRO:HD3	1.71	0.42
1:C:133:LEU:HD13	1:C:135:LEU:HD12	2.01	0.42
1:C:97:VAL:O	1:C:153:LYS:HB3	2.20	0.42
1:C:173:GLN:HE21	1:C:173:GLN:HB3	1.65	0.42
1:A:140:ARG:H	1:A:140:ARG:HG3	1.65	0.41
1:A:65:ALA:N	1:A:66:PRO:CD	2.83	0.41
1:B:64:THR:HG22	1:B:180:LEU:HD11	2.03	0.41
1:B:66:PRO:HG2	1:B:67:TYR:CD2	2.55	0.41
1:C:54:LEU:O	1:C:55:GLN:C	2.62	0.41
1:C:14:ILE:H	1:C:14:ILE:CD1	2.30	0.41
1:C:133:LEU:C	1:C:133:LEU:CD2	2.93	0.41
1:C:58:ASP:CB	1:C:61:THR:OG1	2.69	0.41
1:B:76:LYS:HD3	1:B:130:GLU:HG2	2.03	0.41
1:C:74:GLU:HG2	1:C:177:ARG:HG3	2.03	0.41
1:C:157:TYR:CD2	1:C:158:SER:N	2.89	0.41
1:A:67:TYR:CD2	1:A:67:TYR:N	2.89	0.41
1:B:134:ASN:C	1:B:136:ASP:H	2.29	0.41
1:B:150:ASP:OD2	1:C:145:SER:CB	2.68	0.41
1:B:68:ARG:HG2	1:B:184:MET:HB2	2.02	0.40
1:C:133:LEU:C	1:C:133:LEU:HD22	2.45	0.40
1:A:95:ALA:CB	1:A:108:ILE:HG23	2.47	0.40
1:A:154:LEU:CD2	1:A:174:ILE:HD11	2.51	0.40
1:B:133:LEU:HD22	1:B:134:ASN:N	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:ASN:C	1:B:139:ASN:ND2	2.69	0.40
1:A:47:GLU:HB2	1:A:160:ALA:CA	2.52	0.40
1:A:94:LEU:HD21	1:A:172:ILE:CD1	2.42	0.40
1:A:134:ASN:OD1	1:A:134:ASN:C	2.64	0.40
1:B:93:TYR:C	1:B:94:LEU:HD12	2.46	0.40
1:C:146:VAL:CG1	1:C:147:THR:N	2.69	0.40
1:B:83:ASP:OD1	1:B:84:LEU:HD12	2.20	0.40
1:A:97:VAL:CG1	1:A:98:PRO:N	2.81	0.40
1:C:84:LEU:HB3	1:C:169:THR:HG22	2.03	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	159/188 (85%)	138 (87%)	17 (11%)	4 (2%)	4	29
1	B	183/188 (97%)	154 (84%)	26 (14%)	3 (2%)	7	36
1	C	183/188 (97%)	146 (80%)	24 (13%)	13 (7%)	1	13
All	All	525/564 (93%)	438 (83%)	67 (13%)	20 (4%)	2	21

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	ILE
1	A	164	PRO
1	C	55	GLN
1	C	122	ALA
1	C	125	ALA
1	A	122	ALA
1	A	125	ALA
1	B	95	ALA

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Mol	Chain	Res	Type
1	B	121	GLY
1	C	60	ILE
1	C	99	ALA
1	C	135	LEU
1	B	112	TYR
1	C	45	THR
1	C	17	PRO
1	C	108	ILE
1	C	112	TYR
1	C	182	LYS
1	C	105	PRO
1	C	89	PRO

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	135/159 (85%)	121 (90%)	14 (10%)	7	26
1	B	156/159 (98%)	135 (86%)	21 (14%)	4	18
1	C	156/159 (98%)	133 (85%)	23 (15%)	3	16
All	All	447/477 (94%)	389 (87%)	58 (13%)	4	19

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ILE
1	A	35	PRO
1	A	58	ASP
1	A	79	LEU
1	A	83	ASP
1	A	97	VAL
1	A	115	GLN
1	A	129	ILE
1	A	147	THR
1	A	149	THR

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Mol	Chain	Res	Type
1	A	159	ARG
1	A	162	THR
1	A	169	THR
1	A	186	ILE
1	B	14	ILE
1	B	21	LEU
1	B	23	GLN
1	B	25	ASN
1	B	27	GLU
1	B	32	ILE
1	B	54	LEU
1	B	76	LYS
1	B	104	THR
1	B	109	LEU
1	B	110	ARG
1	B	115	GLN
1	B	118	VAL
1	B	133	LEU
1	B	136	ASP
1	B	142	LEU
1	B	147	THR
1	B	149	THR
1	B	150	ASP
1	B	159	ARG
1	B	180	LEU
1	C	14	ILE
1	C	27	GLU
1	C	28	GLN
1	C	56	THR
1	C	60	ILE
1	C	62	LYS
1	C	68	ARG
1	C	79	LEU
1	C	88	ASN
1	C	94	LEU
1	C	104	THR
1	C	105	PRO
1	C	118	VAL
1	C	123	ILE
1	C	124	SER
1	C	133	LEU
1	C	135	LEU

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Mol	Chain	Res	Type
1	C	139	ASN
1	C	147	THR
1	C	154	LEU
1	C	155	LEU
1	C	173	GLN
1	C	182	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	51	GLN
1	A	69	HIS
1	A	100	ASN
1	A	163	ASN
1	B	37	GLN
1	B	69	HIS
1	B	88	ASN
1	B	107	GLN
1	B	115	GLN
1	B	139	ASN
1	B	173	GLN
1	C	23	GLN
1	C	37	GLN
1	C	69	HIS
1	C	115	GLN
1	C	139	ASN
1	C	173	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	161/188 (85%)	-0.78	0 100 100	1, 13, 45, 59	0
1	B	185/188 (98%)	-0.75	0 100 100	1, 13, 46, 74	0
1	C	185/188 (98%)	-0.69	0 100 100	11, 23, 39, 52	0
All	All	531/564 (94%)	-0.74	0 100 100	1, 18, 44, 74	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.