



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 07:14 AM GMT

PDB ID : 1CWP
Title : STRUCTURES OF THE NATIVE AND SWOLLEN FORMS OF COW-
PEA CHLOROTIC MOTTLE VIRUS DETERMINED BY X-RAY CRYSTALLOGRAPHY AND CRYO-ELECTRON MICROSCOPY
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Deposited on : 1995-05-22
Resolution : 3.20 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

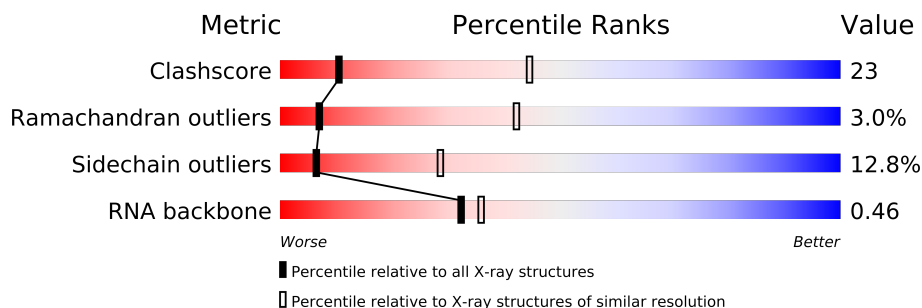
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	21963
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.20 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	79885	1078 (3.26-3.14)
Ramachandran outliers	78287	1059 (3.26-3.14)
Sidechain outliers	78261	1058 (3.26-3.14)
RNA backbone	1838	1002 (3.72-2.68)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	D	4	
1	F	4	
2	E	2	
3	A	190	
3	B	190	
3	C	190	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3784 atoms, of which 0 are hydrogen and 0 are deuterium.

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 RNA chain called RNA (5'-R(*AP*UP*AP*U)-3').

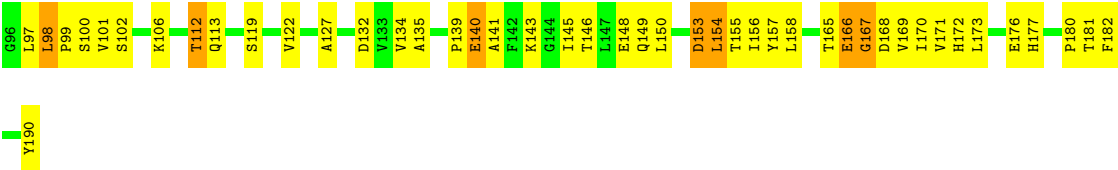
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	4	Total	C	N	O	P	0	0	0
			84	38	14	28	4			
1	F	4	Total	C	N	O	P	0	0	0
			84	38	14	28	4			

- Molecule 2 is a RNA chain called RNA (5'-R(*AP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	2	Total	C	N	O	P	0	0	0
			42	19	7	14	2			

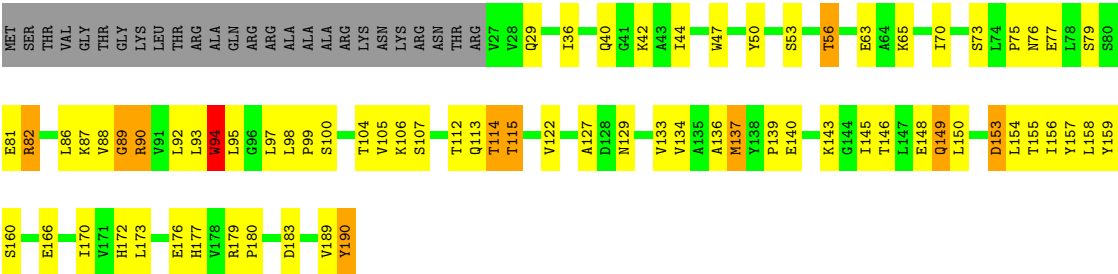
- Molecule 3 is a protein called Coat protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	149	Total	C	N	O	S	0	0	0
			1122	717	181	221	3			
3	B	164	Total	C	N	O	S	0	0	0
			1226	784	198	241	3			
3	C	164	Total	C	N	O	S	0	0	0
			1226	784	198	241	3			



● Molecule 3: Coat protein

Chain C:



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	381.30Å 381.30Å 408.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 3.20	Depositor
% Data completeness (in resolution range)	57.0 (7.00-3.20)	Depositor
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.310 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3784	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

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	D	4.32	4/93 (4.3%)	4.63	10/142 (7.0%)
1	F	1.03	0/93	1.70	1/142 (0.7%)
2	E	1.80	1/46 (2.2%)	1.44	0/69
3	A	0.55	0/1143	0.88	2/1560 (0.1%)
3	B	0.56	1/1249 (0.1%)	0.96	3/1707 (0.2%)
3	C	0.58	1/1249 (0.1%)	0.96	2/1707 (0.1%)
All	All	0.90	7/3873 (0.2%)	1.22	18/5327 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	A	0	1
3	C	0	1
All	All	0	2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	203	A	O3'-P	28.80	1.95	1.61
1	D	202	U	O3'-P	-24.04	1.32	1.61
1	D	203	A	N9-C4	-10.58	1.31	1.37
3	B	94	TRP	CB-CG	-6.57	1.38	1.50
3	C	94	TRP	CB-CG	-6.20	1.39	1.50
2	E	202	U	C5'-C4'	5.66	1.58	1.51
1	D	202	U	C5'-C4'	5.60	1.58	1.51

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	203	A	O3'-P-O5'	-24.17	58.08	104.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	203	A	OP2-P-O3'	24.04	158.08	105.20
1	D	202	U	OP2-P-O3'	-22.53	55.64	105.20
1	D	203	A	OP1-P-O3'	-20.73	59.59	105.20
1	D	203	A	P-O3'-C3'	14.89	137.57	119.70
1	D	202	U	O3'-P-O5'	12.97	128.64	104.00
1	F	201	A	O5'-P-OP1	12.12	125.24	110.70
3	B	167	GLY	N-CA-C	10.03	138.18	113.10
3	C	115	THR	N-CA-C	-8.58	87.83	111.00
1	D	202	U	OP1-P-O3'	7.86	122.49	105.20
1	D	203	A	C8-N9-C4	6.90	108.56	105.80
1	D	203	A	C4-C5-C6	-6.08	113.96	117.00
3	A	140	GLU	N-CA-C	5.71	126.41	111.00
3	A	173	LEU	CA-CB-CG	5.68	128.37	115.30
1	D	203	A	N9-C1'-C2'	-5.42	106.04	112.00
3	B	154	LEU	CA-CB-CG	5.29	127.46	115.30
3	C	114	THR	CA-CB-CG2	-5.24	105.07	112.40
3	B	98	LEU	CA-CB-CG	5.07	126.96	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	139	PRO	Mainchain
3	C	190	TYR	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	84	0	42	12	0
1	F	84	0	41	28	0
2	E	42	0	22	2	0
3	A	1122	0	1133	40	0
3	B	1226	0	1243	45	0
3	C	1226	0	1243	60	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3784	0	3724	174	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 23.

All (174) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:202:U:H2'	1:D:203:A:C8	1.58	1.37
1:F:203:A:O2'	1:F:204:U:O4'	1.54	1.24
1:F:203:A:H2'	1:F:204:U:C6	1.75	1.20
1:F:203:A:O2'	1:F:204:U:H5'	1.55	1.07
1:D:202:U:C2'	1:D:203:A:H8	1.71	1.03
1:F:203:A:H2'	1:F:204:U:H6	1.05	1.03
1:D:203:A:H2'	1:D:204:U:C6	1.94	1.03
1:D:202:U:H2'	1:D:203:A:H8	0.85	1.02
1:F:201:A:O2'	1:F:202:U:H5'	1.61	1.00
3:B:98:LEU:HB2	3:B:99:PRO:HD2	1.44	0.97
1:D:202:U:C2'	1:D:203:A:C8	2.45	0.97
2:E:201:A:H8	3:B:143:LYS:HZ1	1.06	0.96
1:F:201:A:H8	3:C:143:LYS:HZ3	1.05	0.95
3:C:98:LEU:HB2	3:C:99:PRO:HD2	1.50	0.91
3:A:81:GLU:HG3	3:C:145:ILE:HD11	1.53	0.90
1:F:203:A:O2'	1:F:204:U:C5'	2.19	0.89
1:F:203:A:C2'	1:F:204:U:H6	1.86	0.88
1:F:202:U:H2'	1:F:203:A:C8	2.09	0.88
1:F:202:U:C2'	1:F:203:A:O4'	2.21	0.87
1:F:202:U:H2'	1:F:203:A:H8	1.39	0.86
1:F:203:A:C2'	1:F:204:U:C6	2.59	0.85
3:C:94:TRP:CB	3:C:172:HIS:HB2	2.08	0.82
3:C:145:ILE:HG23	3:C:149:GLN:HB3	1.61	0.82
3:C:86:LEU:HD21	3:C:180:PRO:HG3	1.61	0.82
3:B:36:ILE:HG23	3:B:40:GLN:HB2	1.63	0.81
1:F:203:A:O2'	1:F:204:U:C4'	2.28	0.80
3:B:94:TRP:CB	3:B:172:HIS:HB2	2.12	0.80
1:D:203:A:H2'	1:D:204:U:C5	2.16	0.80
3:B:146:THR:HB	3:B:149:GLN:HE21	1.48	0.78
3:C:150:LEU:HD12	3:C:154:LEU:HD12	1.66	0.78
3:A:145:ILE:HG23	3:A:149:GLN:HB2	1.68	0.76
3:B:36:ILE:HG22	3:B:37:ALA:O	1.86	0.76
1:F:201:A:C2'	1:F:202:U:H5'	2.15	0.75
3:B:94:TRP:HB2	3:B:172:HIS:HB2	1.68	0.75
3:A:94:TRP:HB2	3:A:172:HIS:HB2	1.70	0.74

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:A:49:GLY:O	3:A:179:ARG:HB2	1.89	0.72
1:D:203:A:C2'	1:D:204:U:C6	2.73	0.71
3:B:56:THR:HG23	3:B:170:ILE:HG23	1.72	0.70
3:C:89:GLY:O	3:C:90:ARG:HG3	1.91	0.70
3:C:94:TRP:HB3	3:C:172:HIS:HB2	1.75	0.68
3:C:145:ILE:CG2	3:C:149:GLN:HB3	2.23	0.68
3:C:94:TRP:HB2	3:C:172:HIS:HB2	1.75	0.67
3:C:129:ASN:HA	3:C:134:VAL:HG22	1.76	0.67
3:A:93:LEU:HB3	3:A:136:ALA:HB3	1.75	0.66
3:B:150:LEU:HD12	3:B:154:LEU:HD12	1.76	0.66
1:F:201:A:H8	3:C:143:LYS:NZ	1.90	0.65
1:D:203:A:O2'	1:D:204:U:H5'	1.95	0.65
1:F:202:U:H2'	1:F:203:A:O4'	1.95	0.65
3:B:59:CYS:O	3:B:59:CYS:SG	2.55	0.65
3:C:53:SER:OG	3:C:177:HIS:HE1	1.81	0.64
1:D:202:U:O2'	1:D:203:A:O4'	2.15	0.63
3:A:143:LYS:HD3	3:B:82:ARG:NH2	2.14	0.63
3:A:59:CYS:HB2	3:A:169:VAL:O	1.99	0.63
1:F:203:A:C2	1:F:204:U:C4	2.88	0.61
3:B:85:GLN:HA	3:B:146:THR:HG23	1.81	0.61
3:B:146:THR:HG22	3:B:148:GLU:H	1.65	0.61
1:F:202:U:C1'	1:F:203:A:O4'	2.48	0.61
3:B:59:CYS:O	3:B:166:GLU:O	2.19	0.61
3:A:81:GLU:HG2	3:C:154:LEU:HD21	1.82	0.61
3:A:139:PRO:O	3:A:141:ALA:N	2.31	0.61
1:F:203:A:C2'	1:F:204:U:H5'	2.32	0.60
3:C:146:THR:HG22	3:C:148:GLU:H	1.66	0.60
1:F:201:A:H2'	1:F:202:U:H5'	1.83	0.60
3:B:143:LYS:HD2	3:C:82:ARG:NH2	2.17	0.60
3:C:93:LEU:HB3	3:C:136:ALA:HB3	1.82	0.60
3:C:113:GLN:HA	3:C:113:GLN:OE1	2.00	0.59
3:C:56:THR:HG23	3:C:170:ILE:HG23	1.84	0.59
1:D:203:A:O2'	1:D:204:U:C5'	2.50	0.59
3:B:158:LEU:HD11	3:B:169:VAL:HG11	1.85	0.59
3:C:146:THR:HB	3:C:149:GLN:HB2	1.85	0.58
1:F:201:A:C2'	1:F:202:U:C5'	2.82	0.58
3:A:94:TRP:CB	3:A:172:HIS:HB2	2.34	0.57
1:F:201:A:H2'	1:F:202:U:C5'	2.35	0.57
3:C:97:LEU:HD23	3:C:97:LEU:H	1.70	0.57
3:A:61:ALA:HB2	3:A:166:GLU:N	2.20	0.56
3:C:88:VAL:HG12	3:C:89:GLY:N	2.21	0.56
3:B:146:THR:HB	3:B:149:GLN:NE2	2.19	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:B:86:LEU:HD21	3:B:180:PRO:HB3	1.89	0.55
3:C:90:ARG:HD2	3:C:176:GLU:OE2	2.07	0.55
3:A:85:GLN:HA	3:A:146:THR:HG23	1.88	0.55
3:C:47:TRP:HH2	3:C:143:LYS:HE2	1.72	0.55
3:B:165:THR:O	3:B:166:GLU:C	2.45	0.54
3:B:145:ILE:HD11	3:C:81:GLU:HB3	1.90	0.54
3:B:94:TRP:HB3	3:B:172:HIS:HB2	1.89	0.54
3:A:109:VAL:HG22	3:A:156:ILE:HG12	1.89	0.54
2:E:201:A:O2'	2:E:202:U:H5'	2.09	0.53
3:C:36:ILE:HG12	3:C:40:GLN:HB2	1.91	0.53
1:F:202:U:H1'	1:F:203:A:O4'	2.08	0.53
3:A:146:THR:HG22	3:A:148:GLU:H	1.73	0.53
1:D:201:A:O2'	1:D:202:U:H5'	2.09	0.52
3:B:153:ASP:OD1	3:B:153:ASP:N	2.42	0.52
3:A:74:LEU:HD21	3:A:147:LEU:HB3	1.90	0.52
1:D:203:A:C2'	1:D:204:U:H6	2.20	0.52
3:A:53:SER:OG	3:A:177:HIS:HE1	1.92	0.52
3:A:83:ASN:O	3:A:86:LEU:HB2	2.09	0.52
3:B:145:ILE:HG23	3:B:149:GLN:HB3	1.92	0.52
3:C:47:TRP:CH2	3:C:143:LYS:HE2	2.45	0.51
1:F:203:A:C3'	1:F:204:U:H5'	2.36	0.51
3:C:98:LEU:CB	3:C:99:PRO:HD2	2.29	0.51
3:A:143:LYS:HD3	3:B:82:ARG:HH21	1.73	0.51
3:C:77:GLU:H	3:C:77:GLU:CD	2.14	0.51
1:F:201:A:H1'	3:C:140:GLU:OE2	2.11	0.51
3:A:81:GLU:HG3	3:C:145:ILE:CD1	2.34	0.51
3:B:156:ILE:HD11	3:B:173:LEU:HD22	1.93	0.51
3:C:47:TRP:HE1	3:C:90:ARG:NH2	2.09	0.51
3:B:113:GLN:HA	3:B:113:GLN:OE1	2.12	0.50
3:C:104:THR:O	3:C:160:SER:HA	2.11	0.50
3:A:110:THR:HG22	3:A:111:GLU:N	2.26	0.50
3:B:140:GLU:O	3:C:82:ARG:NH1	2.45	0.49
3:A:97:LEU:HB3	3:A:169:VAL:HG22	1.95	0.49
3:A:53:SER:OG	3:A:177:HIS:CE1	2.65	0.49
3:B:112:THR:OG1	3:B:155:THR:HG23	2.13	0.49
3:B:101:VAL:HG12	3:B:102:SER:N	2.28	0.49
3:C:44:ILE:CG2	3:C:90:ARG:HH12	2.26	0.48
3:A:84:LYS:O	3:A:146:THR:HG22	2.13	0.48
3:C:65:LYS:HA	3:C:159:TYR:CE2	2.48	0.48
3:A:94:TRP:CG	3:A:172:HIS:HB2	2.48	0.48
3:C:36:ILE:HD11	3:C:40:GLN:O	2.14	0.48
3:A:87:LYS:O	3:A:177:HIS:HA	2.13	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:95:LEU:HB3	3:C:134:VAL:CG1	2.44	0.47
1:F:203:A:C4	1:F:204:U:C5	3.02	0.47
3:C:95:LEU:HB3	3:C:134:VAL:HG12	1.96	0.47
3:A:64:ALA:HB1	3:A:161:SER:O	2.14	0.47
3:C:95:LEU:HD23	3:C:97:LEU:HD22	1.96	0.47
3:C:36:ILE:CG1	3:C:40:GLN:HB2	2.44	0.47
3:C:44:ILE:HG22	3:C:90:ARG:HH12	1.80	0.47
3:B:36:ILE:CG2	3:B:40:GLN:HB2	2.41	0.47
3:B:84:LYS:HG2	3:B:148:GLU:HB2	1.97	0.47
3:C:88:VAL:HG12	3:C:89:GLY:H	1.80	0.47
3:A:74:LEU:HD12	3:A:151:ALA:HB2	1.97	0.46
3:C:106:LYS:HA	3:C:127:ALA:O	2.16	0.46
3:C:189:VAL:O	3:C:190:TYR:CB	2.63	0.46
3:B:93:LEU:HD11	3:B:171:VAL:HG13	1.96	0.46
3:A:84:LYS:O	3:A:146:THR:CG2	2.64	0.46
3:A:106:LYS:HA	3:A:127:ALA:O	2.16	0.46
3:A:110:THR:CG2	3:A:111:GLU:N	2.79	0.45
3:C:146:THR:HB	3:C:149:GLN:HE21	1.80	0.45
3:C:87:LYS:O	3:C:177:HIS:HA	2.16	0.45
3:A:95:LEU:HA	3:A:95:LEU:HD13	1.87	0.45
3:A:98:LEU:HG	3:A:99:PRO:HD2	1.99	0.45
3:A:60:ALA:O	3:A:61:ALA:HB3	2.17	0.44
3:A:86:LEU:HG	3:A:178:VAL:O	2.17	0.44
3:B:139:PRO:O	3:B:141:ALA:N	2.50	0.44
3:B:95:LEU:HB3	3:B:134:VAL:HG12	1.98	0.44
3:C:137:MET:HG2	3:C:139:PRO:HG3	1.99	0.44
3:B:90:ARG:O	3:B:176:GLU:HB2	2.18	0.44
3:C:98:LEU:HD12	3:C:100:SER:H	1.83	0.44
3:C:107:SER:HA	3:C:157:TYR:O	2.17	0.44
3:A:131:LYS:HB3	3:B:190:TYR:OH	2.17	0.44
3:C:156:ILE:HD11	3:C:173:LEU:HD22	1.98	0.44
3:B:106:LYS:HA	3:B:127:ALA:O	2.18	0.43
3:B:78:LEU:HA	3:B:83:ASN:HB3	2.00	0.43
3:B:93:LEU:O	3:B:135:ALA:HA	2.19	0.43
3:B:82:ARG:HH11	3:B:182:PHE:HD1	1.65	0.43
3:B:95:LEU:HB3	3:B:134:VAL:CG1	2.49	0.43
3:C:153:ASP:N	3:C:153:ASP:OD1	2.51	0.42
3:C:47:TRP:HE1	3:C:90:ARG:CZ	2.32	0.42
3:A:129:ASN:OD1	3:A:129:ASN:N	2.53	0.42
3:A:61:ALA:HB2	3:A:166:GLU:CA	2.49	0.42
3:C:105:VAL:HA	3:C:159:TYR:O	2.20	0.42
3:C:70:ILE:HB	3:C:156:ILE:HB	2.01	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:A:98:LEU:CG	3:A:99:PRO:HD2	2.50	0.42
1:F:201:A:O2'	1:F:202:U:C5'	2.49	0.41
3:C:179:ARG:HA	3:C:180:PRO:HD3	1.92	0.41
3:B:53:SER:OG	3:B:177:HIS:HE1	2.02	0.41
3:C:92:LEU:HD12	3:C:136:ALA:O	2.21	0.41
3:A:88:VAL:HG12	3:A:89:GLY:N	2.36	0.41
3:B:122:VAL:HG12	3:B:122:VAL:O	2.21	0.41
3:B:53:SER:OG	3:B:177:HIS:CE1	2.74	0.40
3:C:42:LYS:HA	3:C:42:LYS:HD3	1.95	0.40
3:C:47:TRP:HB2	3:C:50:TYR:CD1	2.56	0.40
3:A:138:TYR:C	3:A:139:PRO:O	2.59	0.40
3:B:119:SER:HB3	3:B:157:TYR:CD2	2.56	0.40
1:F:202:U:C6	1:F:203:A:C8	3.09	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	
3	A	147/190 (77%)	136 (92%)	8 (5%)	3 (2%)	11	56
3	B	162/190 (85%)	145 (90%)	12 (7%)	5 (3%)	7	41
3	C	162/190 (85%)	139 (86%)	17 (10%)	6 (4%)	5	34
All	All	471/570 (83%)	420 (89%)	37 (8%)	14 (3%)	7	42

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	116	ALA
3	A	140	GLU
3	B	140	GLU
3	C	90	ARG
3	A	139	PRO
3	B	40	GLN

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Mol	Chain	Res	Type
3	C	166	GLU
3	B	112	THR
3	B	166	GLU
3	C	112	THR
3	C	79	SER
3	C	89	GLY
3	B	167	GLY
3	C	75	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	
3	A	122/154 (79%)	106 (87%)	16 (13%)	6	28
3	B	134/154 (87%)	117 (87%)	17 (13%)	6	29
3	C	134/154 (87%)	117 (87%)	17 (13%)	6	29
All	All	390/462 (84%)	340 (87%)	50 (13%)	6	28

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	68	SER
3	A	74	LEU
3	A	77	GLU
3	A	86	LEU
3	A	95	LEU
3	A	97	LEU
3	A	98	LEU
3	A	100	SER
3	A	113	GLN
3	A	115	THR
3	A	128	ASP
3	A	137	MET
3	A	154	LEU
3	A	158	LEU
3	A	173	LEU

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Mol	Chain	Res	Type
3	A	189	VAL
3	B	29	GLN
3	B	48	THR
3	B	56	THR
3	B	59	CYS
3	B	63	GLU
3	B	77	GLU
3	B	79	SER
3	B	85	GLN
3	B	86	LEU
3	B	94	TRP
3	B	95	LEU
3	B	97	LEU
3	B	100	SER
3	B	132	ASP
3	B	153	ASP
3	B	168	ASP
3	B	181	THR
3	C	29	GLN
3	C	56	THR
3	C	63	GLU
3	C	73	SER
3	C	76	ASN
3	C	82	ARG
3	C	94	TRP
3	C	114	THR
3	C	115	THR
3	C	122	VAL
3	C	133	VAL
3	C	137	MET
3	C	149	GLN
3	C	153	ASP
3	C	155	THR
3	C	158	LEU
3	C	183	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	83	ASN
3	A	113	GLN
3	A	177	HIS

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Mol	Chain	Res	Type
3	B	83	ASN
3	B	149	GLN
3	B	177	HIS
3	C	177	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	D	3/4 (75%)	2 (66%)	2 (66%)
1	F	3/4 (75%)	2 (66%)	1 (33%)
2	E	1/2 (50%)	0	0
All	All	7/10 (70%)	4 (57%)	3 (42%)

All (4) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	D	203	A
1	D	204	U
1	F	203	A
1	F	204	U

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	D	202	U
1	D	203	A
1	F	202	U

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 carbohydrates 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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.