



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:52 PM 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 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

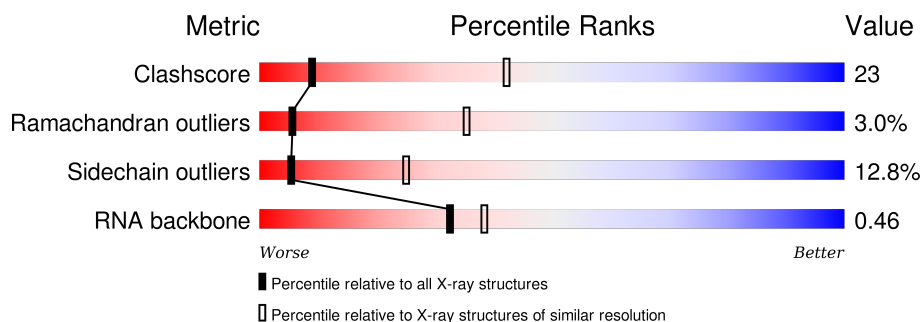
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RNA backbone	2183	1079 (3.70-2.70)

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 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.

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 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 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			

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

- Molecule 1: RNA (5'-R(*AP*UP*AP*U)-3')

Chain D: 

A201
U202
A203
U204

- Molecule 1: RNA (5'-R(*AP*UP*AP*U)-3')

Chain F: 

A201
U202
A203
U204

- Molecule 2: RNA (5'-R(*AP*U)-3')

Chain E: 

A201
U202

- Molecule 3: Coat protein

Chain A: 

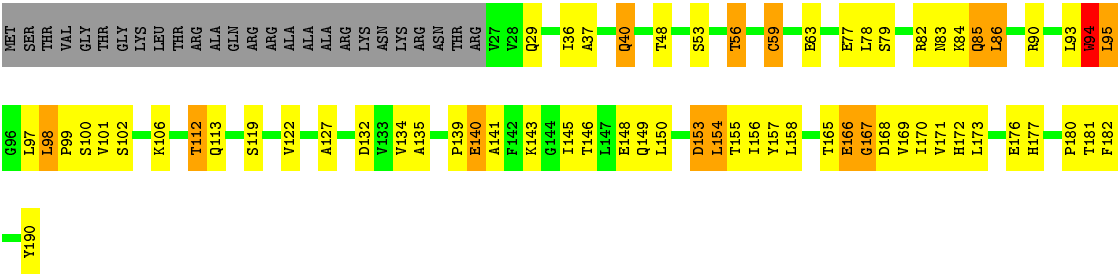
MET SER THR VAL GLY THR GLY LYS LEU THR ARG ALA GIN ARG ARG ALA ALA ARG LYS ASN ASN LYS ARG E111 T110 T112 ASN THR ARG T113 T115 VAL A116 A127 PRQ D128 VAL ILE I129 VAL S130 K131 PRQ I136 I137 M137 Y138 SER F139 E140 A141 GLY K142 G49 S53 C59 A60 A61 A64 S68 L74 E77 E81

H82 N83 K84 Q85 L86 K87 V88 Q89 L93 THR W94 ARG L95 G96 L97 L98 P99 S100 K106 V109 T110 E111 T112 Q113 T114 T115 A116 A127 D128 M129 S130 K131 A136 M137 Y138 F139 E140 A141 F142 K143 G144 I145 T146 L147 E148 Q149 L150 A151 L154 T155 I156 Y157 L158 S161

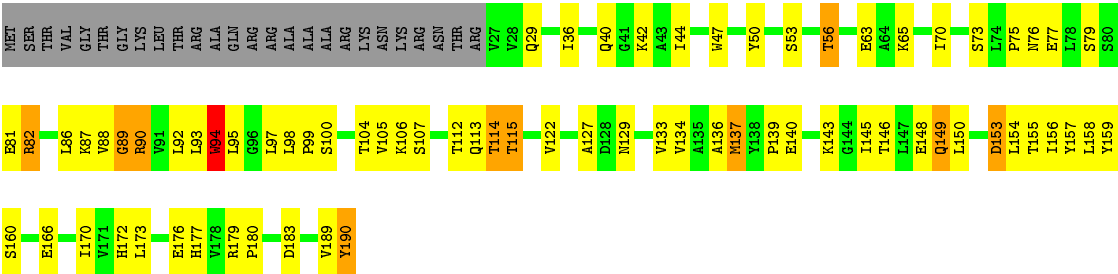
E166 V169 H172 L173 H177 V178 R179 V189 Y190

- Molecule 3: Coat protein

Chain B: 



● Molecule 3: Coat protein



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 [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	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 Too-close contacts [i](#)

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	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
All	All	3784	0	3724	174	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 23.

All (174) 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: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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
3:B:86:LEU:HD21	3:B:180:PRO:HB3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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	Interatomic distance (Å)	Clash overlap (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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:70:ILE:HB	3:C:156:ILE:HB	2.01	0.42
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 [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	
3	A	147/190 (77%)	136 (92%)	8 (5%)	3 (2%)	9	48
3	B	162/190 (85%)	145 (90%)	12 (7%)	5 (3%)	5	34
3	C	162/190 (85%)	139 (86%)	17 (10%)	6 (4%)	4	29
All	All	471/570 (83%)	420 (89%)	37 (8%)	14 (3%)	5	35

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	116	ALA
3	A	140	GLU
3	B	140	GLU

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Mol	Chain	Res	Type
3	C	90	ARG
3	A	139	PRO
3	B	40	GLN
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%)	5	24
3	B	134/154 (87%)	117 (87%)	17 (13%)	5	25
3	C	134/154 (87%)	117 (87%)	17 (13%)	5	25
All	All	390/462 (84%)	340 (87%)	50 (13%)	5	25

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

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Mol	Chain	Res	Type
3	A	137	MET
3	A	154	LEU
3	A	158	LEU
3	A	173	LEU
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
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.