



wwPDB EM Map/Model Validation Report

Apr 10, 2016 – 01:39 PM BST

PDB ID : 2OM3
EMDB ID: : EMD-1316
Title : High-resolution cryo-EM structure of Tobacco Mosaic Virus
Authors : Sachse, C.
Deposited on : 2007-01-20
Resolution : 4.40 Å(reported)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>

MolProbity : 4.02b-467
Mogul : unknown
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) : trunk27241

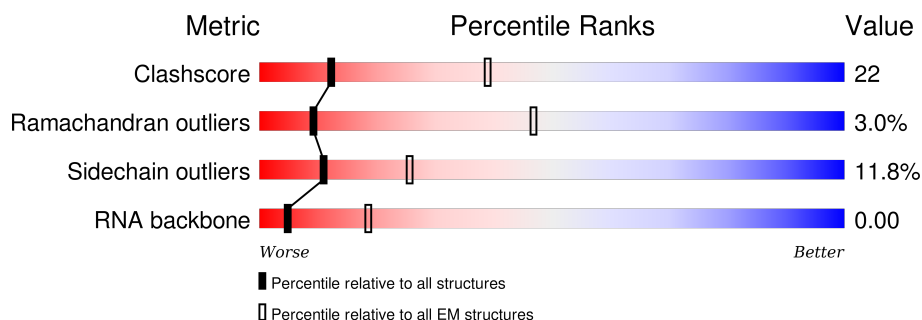
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.40 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)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686
RNA backbone	3027	244

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	1-R	3	33% 67%
1	2-R	3	100%
1	3-R	3	100%
1	4-R	3	33% 67%
1	5-R	3	33% 67%
2	1-A	158	56% 34% 8% ..
2	2-A	158	54% 42% ..
2	3-A	158	59% 35% ..

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Mol	Chain	Length	Quality of chain
2	4-A	158	<div><div></div><div>60%</div><div>35%</div><div>...</div></div>
2	5-A	158	<div><div></div><div>53%</div><div>42%</div><div>..</div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6395 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Tobacco Mosaic Virus RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	1-R	3	Total	C	N	O	P	3	0
			67	30	15	19	3		
1	2-R	3	Total	C	N	O	P	3	0
			67	30	15	19	3		
1	3-R	3	Total	C	N	O	P	3	0
			67	30	15	19	3		
1	4-R	3	Total	C	N	O	P	3	0
			67	30	15	19	3		
1	5-R	3	Total	C	N	O	P	3	0
			67	30	15	19	3		

- Molecule 2 is a protein called Coat protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	1-A	154	Total	C	N	O	S	154	0
			1212	762	211	238	1		
2	2-A	154	Total	C	N	O	S	154	0
			1212	762	211	238	1		
2	3-A	154	Total	C	N	O	S	154	0
			1212	762	211	238	1		
2	4-A	154	Total	C	N	O	S	154	0
			1212	762	211	238	1		
2	5-A	154	Total	C	N	O	S	154	0
			1212	762	211	238	1		

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. 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. 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: Tobacco Mosaic Virus RNA

Chain 1-R: 

G4
A5
A6

- Molecule 1: Tobacco Mosaic Virus RNA

Chain 2-R: 

G4
A5
A6

- Molecule 1: Tobacco Mosaic Virus RNA

Chain 3-R: 

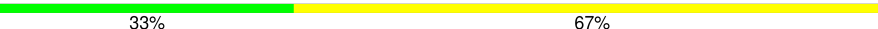
G4
A5
A6

- Molecule 1: Tobacco Mosaic Virus RNA

Chain 4-R: 

G4
A5
A6

- Molecule 1: Tobacco Mosaic Virus RNA

Chain 5-R: 

G4
A5
A6

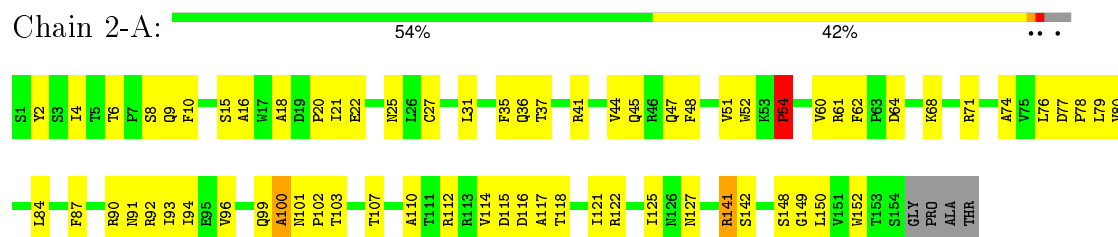
- Molecule 2: Coat protein

Chain 1-A: 

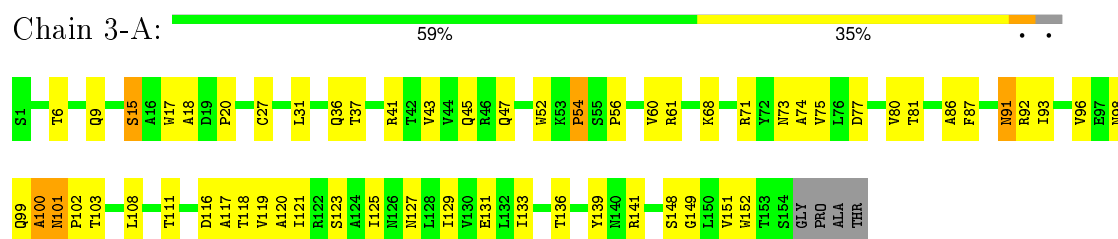
S1 Y2 Q9 F10 L13 S14 S15 A16 W17 A18 D19 F20 I21 N25 L26 C27 L31 F35 Q36 R41 Q47 F48 V51 M52 R53 P54 S55 W56 Q57 V60 R61 F62 R71 Y72 N73 A74 V75 L76 V80 L83 F87 N91 R92 I93 I94 E95

Y96 E97 N98 Q99 A100 M101 P102 T103 T104 L108 D109 A110 D116 A117 T118 V119 A120 I121 I125 L128 I129 L132 I133 R141 S146 G149 L150 V151 W152 T153 S154 GLY PRO ALA THR

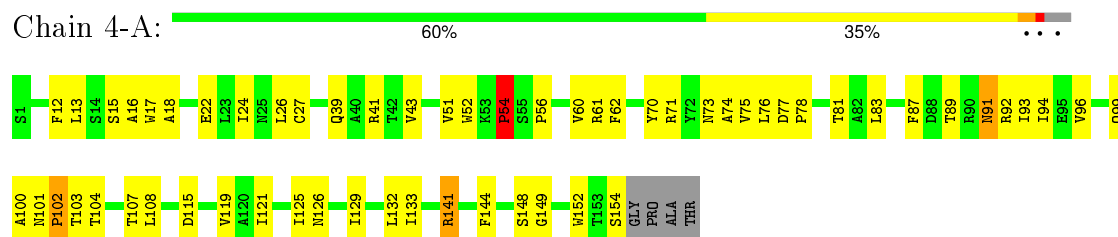
- Molecule 2: Coat protein



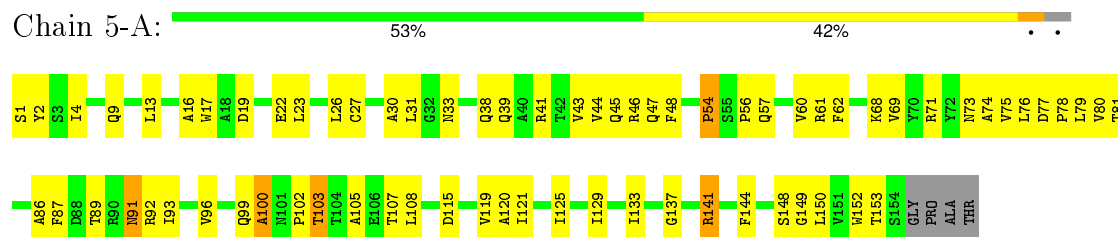
- Molecule 2: Coat protein



- Molecule 2: Coat protein



- Molecule 2: Coat protein



4 Experimental information

Property	Value	Source
Reconstruction method	HELICAL	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	Phase correction and amplitude weighting adapted from Grigorieff 1998(after determination of CTF and specimen tilt)	Depositor
Microscope	Technai F30	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1500	Depositor
Minimum defocus (nm)	1.5	Depositor
Maximum defocus (nm)	4.0	Depositor
Magnification	59000	Depositor
Image detector	Kodak ISO153 film	Depositor

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	$\# Z > 2$	RMSZ	$\# Z > 2$
1	1-R	0.27	0/75	0.78	0/115
2	1-A	0.34	0/1236	0.55	1/1689 (0.1%)
All	All	0.34	0/1311	0.57	1/1804 (0.1%)

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
2	1-A	0	3
2	2-A	0	1
2	3-A	0	2
2	4-A	0	2
2	5-A	0	2
All	All	0	10

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	1-A	14[A]	SER	N-CA-C	5.18	124.98	111.00

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	1-A	13[A]	LEU	Mainchain
2	1-A	54[A]	PRO	Peptide
2	1-A	91[A]	ASN	Peptide
2	2-A	54[B]	PRO	Peptide
2	3-A	54[C]	PRO	Peptide

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Mol	Chain	Res	Type	Group
2	3-A	91[C]	ASN	Peptide
2	4-A	54[D]	PRO	Peptide
2	4-A	91[D]	ASN	Peptide
2	5-A	54[E]	PRO	Peptide
2	5-A	91[E]	ASN	Peptide

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	1-R	67	0	34	13	0
1	2-R	67	0	34	5	0
1	3-R	67	0	34	11	0
1	4-R	67	0	34	4	0
1	5-R	67	0	34	5	0
2	1-A	1212	0	1191	47	0
2	2-A	1212	0	1191	56	0
2	3-A	1212	0	1191	49	0
2	4-A	1212	0	1191	47	0
2	5-A	1212	0	1191	66	0
All	All	6395	0	6125	279	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 22.

All (279) 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)
2:A:54[E]:PRO:HB2	2:A:56[E]:PRO:HD2	1.43	1.00
1:R:4[E]:G:H2'	2:A:119[E]:VAL:HG21	1.40	0.99
2:A:15[D]:SER:HA	2:A:54[D]:PRO:HG3	1.49	0.92
2:A:27[D]:CYS:HB3	2:A:129[D]:ILE:HD11	1.57	0.85
2:A:27[C]:CYS:HB3	2:A:129[C]:ILE:HD11	1.59	0.84
2:A:91[A]:ASN:HB2	2:A:92[A]:ARG:HA	1.64	0.80
2:A:91[B]:ASN:HB3	2:A:94[B]:ILE:HD12	1.64	0.80
2:A:91[C]:ASN:HB2	2:A:92[C]:ARG:HA	1.64	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:5[E]:A:H5''	2:A:119[E]:VAL:CG1	2.14	0.78
2:A:54[C]:PRO:HB2	2:A:56[C]:PRO:HD2	1.64	0.78
2:A:9[A]:GLN:HE22	2:A:150[A]:LEU:HG	1.52	0.75
1:R:6[C]:A:N3	1:R:6[C]:A:H3'	2.02	0.74
2:A:60[D]:VAL:HG12	2:A:61[D]:ARG:H	1.52	0.74
2:A:36[C]:GLN:HG2	2:A:118[C]:THR:OG1	1.88	0.73
2:A:60[E]:VAL:HG12	2:A:61[E]:ARG:H	1.53	0.72
2:A:31[E]:LEU:HD23	2:A:125[E]:ILE:HD12	1.69	0.72
2:A:60[B]:VAL:HG12	2:A:61[B]:ARG:H	1.54	0.72
2:A:60[C]:VAL:HG12	2:A:61[C]:ARG:H	1.53	0.72
2:A:54[D]:PRO:HB2	2:A:56[D]:PRO:HD2	1.72	0.71
2:A:148[D]:SER:OG	2:A:149[D]:GLY:HA2	1.91	0.71
2:A:60[A]:VAL:HG12	2:A:61[A]:ARG:H	1.54	0.71
2:A:91[E]:ASN:HB2	2:A:92[E]:ARG:HA	1.75	0.69
2:A:91[B]:ASN:HB2	2:A:92[B]:ARG:HA	1.74	0.69
2:A:93[D]:ILE:O	2:A:96[D]:VAL:HG12	1.92	0.69
2:A:61[E]:ARG:HG2	2:A:141[E]:ARG:HH12	1.58	0.68
2:A:125[E]:ILE:O	2:A:129[E]:ILE:HB	1.94	0.68
2:A:16[B]:ALA:HB1	2:A:71[B]:ARG:HB3	1.74	0.68
1:R:6[A]:A:H3'	1:R:6[A]:A:N3	2.10	0.67
2:A:54[A]:PRO:HB2	2:A:56[A]:PRO:CD	2.25	0.67
1:R:4[A]:G:H2'	1:R:5[A]:A:H5''	1.76	0.67
2:A:17[C]:TRP:HE1	2:A:56[C]:PRO:HB3	1.58	0.67
2:A:45[B]:GLN:HE21	2:A:87[B]:PHE:HB3	1.60	0.67
2:A:18[D]:ALA:HB2	2:A:52[D]:TRP:HD1	1.60	0.66
2:A:93[A]:ILE:O	2:A:96[A]:VAL:HG12	1.96	0.66
2:A:41[D]:ARG:HH12	2:A:89[D]:THR:HB	1.60	0.66
2:A:129[C]:ILE:O	2:A:133[C]:ILE:HG12	1.96	0.66
2:A:129[D]:ILE:O	2:A:133[D]:ILE:HG12	1.97	0.64
2:A:148[E]:SER:OG	2:A:149[E]:GLY:HA2	1.98	0.63
2:A:54[D]:PRO:HB2	2:A:56[D]:PRO:CD	2.28	0.62
1:R:5[E]:A:H5''	2:A:119[E]:VAL:HG11	1.79	0.62
1:R:4[D]:G:H2'	2:A:119[D]:VAL:HG21	1.81	0.62
2:A:62[B]:PHE:HZ	2:A:68[B]:LYS:HB2	1.63	0.62
2:A:18[D]:ALA:HB2	2:A:52[D]:TRP:CD1	2.35	0.62
2:A:61[E]:ARG:HD3	2:A:152[E]:TRP:CE3	2.35	0.62
2:A:148[A]:SER:OG	2:A:149[A]:GLY:HA2	2.00	0.62
1:R:6[A]:A:C8	2:A:116[A]:ASP:HB2	2.35	0.62
2:A:93[E]:ILE:O	2:A:96[E]:VAL:HG12	2.01	0.61
2:A:52[C]:TRP:CH2	2:A:80[C]:VAL:HG11	2.35	0.61
2:A:87[D]:PHE:HA	2:A:121[D]:ILE:HG21	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:148[C]:SER:OG	2:A:149[C]:GLY:HA2	2.01	0.61
2:A:41[D]:ARG:NH1	2:A:89[D]:THR:HB	2.15	0.60
2:A:83[A]:LEU:HD21	2:A:125[A]:ILE:HG13	1.83	0.60
2:A:31[B]:LEU:HD23	2:A:125[B]:ILE:HD12	1.81	0.60
2:A:77[E]:ASP:HB3	2:A:78[E]:PRO:HD3	1.81	0.60
2:A:76[E]:LEU:HA	2:A:79[E]:LEU:HD12	1.84	0.60
2:A:61[C]:ARG:HD2	2:A:141[C]:ARG:HH22	1.67	0.60
2:A:107[B]:THR:HA	2:A:110[B]:ALA:HB3	1.83	0.60
2:A:90[B]:ARG:HG3	2:A:91[B]:ASN:H	1.67	0.59
2:A:83[D]:LEU:HD21	2:A:125[D]:ILE:HG13	1.83	0.59
2:A:87[E]:PHE:HA	2:A:121[E]:ILE:HG21	1.83	0.59
2:A:45[C]:GLN:HE21	2:A:87[C]:PHE:HB3	1.68	0.59
2:A:54[A]:PRO:HB2	2:A:56[A]:PRO:HD3	1.84	0.59
2:A:22[D]:GLU:O	2:A:26[D]:LEU:HG	2.03	0.59
2:A:47[A]:GLN:O	2:A:51[A]:VAL:HG13	2.02	0.58
2:A:26[A]:LEU:HD12	2:A:51[A]:VAL:HG11	1.85	0.58
2:A:37[C]:THR:O	2:A:41[C]:ARG:HB2	2.04	0.58
2:A:9[E]:GLN:NE2	2:A:150[E]:LEU:HG	2.19	0.58
2:A:22[E]:GLU:O	2:A:26[E]:LEU:HG	2.03	0.57
2:A:16[A]:ALA:HB1	2:A:71[A]:ARG:HB3	1.86	0.57
2:A:22[B]:GLU:HA	2:A:25[B]:ASN:HB2	1.86	0.57
2:A:126[D]:ASN:HA	2:A:129[D]:ILE:HD13	1.86	0.57
2:A:18[C]:ALA:HB2	2:A:52[C]:TRP:HD1	1.69	0.56
2:A:36[B]:GLN:HG3	2:A:115[B]:ASP:HA	1.86	0.56
2:A:94[A]:ILE:HD11	2:A:110[A]:ALA:HB1	1.87	0.56
2:A:35[A]:PHE:HB2	2:A:118[A]:THR:HG23	1.86	0.56
2:A:93[C]:ILE:O	2:A:96[C]:VAL:HG12	2.06	0.56
2:A:87[E]:PHE:HA	2:A:121[E]:ILE:CG2	2.35	0.56
2:A:129[E]:ILE:O	2:A:133[E]:ILE:HG12	2.06	0.55
2:A:107[D]:THR:HG22	2:A:107[D]:THR:O	2.06	0.55
2:A:118[B]:THR:O	2:A:122[B]:ARG:HB2	2.06	0.55
2:A:77[B]:ASP:HB3	2:A:78[B]:PRO:HD3	1.87	0.55
2:A:35[B]:PHE:HB3	2:A:121[B]:ILE:HD11	1.87	0.55
2:A:15[B]:SER:HA	2:A:54[B]:PRO:HB3	1.88	0.55
2:A:91[C]:ASN:HB2	2:A:92[C]:ARG:CA	2.36	0.55
2:A:20[B]:PRO:HD3	2:A:68[B]:LYS:HA	1.88	0.54
2:A:21[B]:ILE:O	2:A:25[B]:ASN:HB2	2.08	0.54
2:A:48[B]:PHE:O	2:A:51[B]:VAL:HG22	2.08	0.54
2:A:99[E]:GLN:H	2:A:100[E]:ALA:HA	1.72	0.54
1:R:6[C]:A:C8	2:A:116[C]:ASP:HB3	2.42	0.54
2:A:37[B]:THR:O	2:A:41[B]:ARG:HB2	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:128[A]:LEU:O	2:A:132[A]:LEU:HB2	2.08	0.54
2:A:100[C]:ALA:O	2:A:101[C]:ASN:HB2	2.07	0.54
2:A:93[B]:ILE:O	2:A:96[B]:VAL:HG12	2.07	0.54
2:A:39[E]:GLN:O	2:A:43[E]:VAL:HG23	2.07	0.53
2:A:54[A]:PRO:HB2	2:A:56[A]:PRO:HD2	1.90	0.53
2:A:129[A]:ILE:O	2:A:133[A]:ILE:HG12	2.07	0.53
2:A:125[C]:ILE:O	2:A:129[C]:ILE:HD13	2.08	0.53
1:R:5[C]:A:OP1	2:A:119[C]:VAL:HB	2.08	0.53
2:A:104[D]:THR:O	2:A:108[D]:LEU:HB2	2.08	0.53
1:R:6[A]:A:H8	2:A:116[A]:ASP:HB2	1.74	0.53
2:A:115[E]:ASP:O	2:A:119[E]:VAL:HG23	2.08	0.53
1:R:4[D]:G:C2	2:A:119[D]:VAL:HG22	2.43	0.53
2:A:54[C]:PRO:HB2	2:A:56[C]:PRO:CD	2.35	0.53
2:A:43[C]:VAL:O	2:A:47[C]:GLN:HB2	2.09	0.53
1:R:5[A]:A:O2'	2:A:120[A]:ALA:HA	2.07	0.53
2:A:4[E]:ILE:HD12	2:A:4[E]:ILE:H	1.74	0.53
2:A:87[D]:PHE:HA	2:A:121[D]:ILE:CG2	2.39	0.52
2:A:87[A]:PHE:HA	2:A:121[A]:ILE:HG21	1.90	0.52
2:A:117[B]:ALA:O	2:A:121[B]:ILE:HG12	2.08	0.52
1:R:4[A]:G:H8	1:R:4[A]:G:H3'	1.74	0.52
2:A:93[A]:ILE:HD12	2:A:94[A]:ILE:N	2.25	0.52
1:R:5[A]:A:H4'	2:A:119[A]:VAL:CG1	2.39	0.52
2:A:125[A]:ILE:O	2:A:129[A]:ILE:HD13	2.10	0.52
1:R:5[B]:A:O2'	1:R:6[B]:A:P	2.68	0.52
2:A:79[B]:LEU:HD21	2:A:127[B]:ASN:OD1	2.10	0.52
2:A:27[E]:CYS:SG	2:A:48[E]:PHE:CZ	3.03	0.52
2:A:102[E]:PRO:HB3	2:A:107[E]:THR:OG1	2.10	0.51
2:A:27[A]:CYS:HB3	2:A:129[A]:ILE:HD11	1.91	0.51
2:A:87[C]:PHE:HA	2:A:121[C]:ILE:HG21	1.92	0.51
2:A:20[C]:PRO:HD3	2:A:68[C]:LYS:HA	1.91	0.51
2:A:17[C]:TRP:HD1	2:A:56[C]:PRO:HD3	1.76	0.51
2:A:2[E]:TYR:HB3	2:A:150[E]:LEU:HD22	1.93	0.51
2:A:2[E]:TYR:OH	2:A:61[E]:ARG:HG3	2.11	0.50
2:A:36[B]:GLN:HE21	2:A:114[B]:VAL:HG23	1.75	0.50
2:A:4[B]:ILE:HG23	2:A:9[B]:GLN:HB3	1.93	0.50
2:A:39[D]:GLN:O	2:A:43[D]:VAL:HG23	2.11	0.50
2:A:9[E]:GLN:O	2:A:13[E]:LEU:HD23	2.12	0.50
2:A:16[D]:ALA:O	2:A:71[D]:ARG:HB2	2.11	0.50
2:A:48[B]:PHE:HB2	2:A:84[B]:LEU:HD21	1.93	0.50
2:A:18[B]:ALA:HB2	2:A:52[B]:TRP:HD1	1.76	0.50
1:R:5[C]:A:O2'	1:R:6[C]:A:P	2.69	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:69[E]:VAL:HG23	2:A:137[E]:GLY:O	2.11	0.50
2:A:73[C]:ASN:ND2	2:A:139[C]:TYR:HE1	2.10	0.50
2:A:99[B]:GLN:N	2:A:100[B]:ALA:HA	2.27	0.49
1:R:6[C]:A:N7	2:A:116[C]:ASP:HB3	2.27	0.49
2:A:99[D]:GLN:H	2:A:100[D]:ALA:HA	1.76	0.49
1:R:4[A]:G:C8	1:R:4[A]:G:H3'	2.47	0.49
2:A:13[D]:LEU:HD13	2:A:56[D]:PRO:O	2.12	0.49
1:R:5[B]:A:HO2'	1:R:6[B]:A:P	2.36	0.49
1:R:4[E]:G:N3	2:A:119[E]:VAL:HG22	2.28	0.49
1:R:4[A]:G:C4	2:A:119[A]:VAL:HG21	2.48	0.49
2:A:93[B]:ILE:HD12	2:A:94[B]:ILE:N	2.28	0.49
2:A:90[B]:ARG:HG3	2:A:91[B]:ASN:N	2.28	0.48
2:A:77[E]:ASP:O	2:A:81[E]:THR:HG22	2.13	0.48
2:A:98[A]:ASN:C	2:A:100[A]:ALA:HB2	2.34	0.48
2:A:16[E]:ALA:O	2:A:71[E]:ARG:HB3	2.13	0.48
2:A:99[D]:GLN:N	2:A:100[D]:ALA:HA	2.29	0.48
2:A:76[A]:LEU:HD21	2:A:132[A]:LEU:HG	1.95	0.48
2:A:15[D]:SER:HA	2:A:54[D]:PRO:CG	2.33	0.48
2:A:94[B]:ILE:HD13	2:A:110[B]:ALA:HB1	1.95	0.48
2:A:91[D]:ASN:HB2	2:A:92[D]:ARG:HA	1.95	0.48
2:A:91[E]:ASN:CB	2:A:92[E]:ARG:HA	2.44	0.47
2:A:35[B]:PHE:H	2:A:122[B]:ARG:HH12	1.62	0.47
2:A:19[E]:ASP:HB3	2:A:22[E]:GLU:HB3	1.97	0.47
2:A:73[D]:ASN:O	2:A:75[D]:VAL:N	2.47	0.47
2:A:61[E]:ARG:HB2	2:A:152[E]:TRP:CZ3	2.49	0.47
2:A:148[B]:SER:OG	2:A:149[B]:GLY:HA2	2.15	0.47
2:A:45[E]:GLN:OE1	2:A:46[E]:ARG:HG2	2.15	0.46
2:A:23[E]:LEU:O	2:A:27[E]:CYS:SG	2.74	0.46
2:A:62[D]:PHE:HE2	2:A:141[D]:ARG:N	2.13	0.46
2:A:36[A]:GLN:HG3	2:A:118[A]:THR:OG1	2.15	0.46
2:A:99[A]:GLN:N	2:A:100[A]:ALA:HB2	2.30	0.46
2:A:18[A]:ALA:HB2	2:A:52[A]:TRP:HA	1.97	0.46
2:A:45[B]:GLN:NE2	2:A:87[B]:PHE:HB3	2.28	0.46
2:A:152[D]:TRP:HE1	2:A:154[D]:SER:HB2	1.80	0.46
2:A:77[C]:ASP:O	2:A:81[C]:THR:HG22	2.15	0.46
2:A:76[B]:LEU:O	2:A:80[B]:VAL:HG23	2.15	0.46
2:A:24[D]:ILE:HG12	2:A:132[D]:LEU:HD13	1.97	0.46
2:A:86[C]:ALA:HB1	2:A:120[C]:ALA:HB1	1.97	0.46
2:A:61[B]:ARG:HG3	2:A:141[B]:ARG:HH22	1.80	0.46
1:R:4[A]:G:C3'	1:R:4[A]:G:C8	2.99	0.46
2:A:98[C]:ASN:C	2:A:100[C]:ALA:HB2	2.37	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:17[A]:TRP:CD1	2:A:56[A]:PRO:HG3	2.51	0.46
2:A:123[C]:SER:O	2:A:127[C]:ASN:HB2	2.16	0.46
2:A:2[E]:TYR:HD2	2:A:57[E]:GLN:O	2.00	0.45
1:R:6[B]:A:N3	1:R:6[B]:A:H5''	2.31	0.45
2:A:77[D]:ASP:O	2:A:81[D]:THR:HG22	2.15	0.45
2:A:17[E]:TRP:NE1	2:A:56[E]:PRO:HG3	2.31	0.45
2:A:115[B]:ASP:O	2:A:118[B]:THR:HB	2.16	0.45
2:A:99[E]:GLN:N	2:A:100[E]:ALA:HA	2.31	0.45
2:A:31[A]:LEU:O	2:A:31[A]:LEU:HD13	2.16	0.45
2:A:54[E]:PRO:HB2	2:A:56[E]:PRO:CD	2.31	0.45
2:A:76[E]:LEU:O	2:A:80[E]:VAL:HG23	2.17	0.45
2:A:141[B]:ARG:HG3	2:A:142[B]:SER:N	2.30	0.45
2:A:73[E]:ASN:O	2:A:76[E]:LEU:N	2.50	0.45
2:A:62[E]:PHE:CZ	2:A:68[E]:LYS:HB2	2.52	0.45
1:R:6[C]:A:N3	1:R:6[C]:A:C3'	2.78	0.45
1:R:4[D]:G:N2	2:A:119[D]:VAL:HG22	2.32	0.45
2:A:104[A]:THR:O	2:A:108[A]:LEU:HB2	2.17	0.45
2:A:131[C]:GLU:OE1	2:A:136[C]:THR:HG21	2.16	0.44
2:A:99[C]:GLN:N	2:A:100[C]:ALA:HA	2.32	0.44
2:A:90[B]:ARG:HG2	2:A:92[B]:ARG:HB3	1.99	0.44
2:A:93[B]:ILE:HD12	2:A:94[B]:ILE:H	1.82	0.44
2:A:64[B]:ASP:HA	2:A:141[B]:ARG:HB2	1.99	0.44
2:A:44[E]:VAL:HA	2:A:47[E]:GLN:HB3	2.00	0.44
2:A:83[A]:LEU:HD21	2:A:125[A]:ILE:CG1	2.47	0.44
2:A:117[A]:ALA:O	2:A:121[A]:ILE:HG12	2.17	0.44
2:A:73[A]:ASN:O	2:A:75[A]:VAL:N	2.50	0.44
2:A:73[C]:ASN:O	2:A:75[C]:VAL:N	2.50	0.44
2:A:31[C]:LEU:O	2:A:31[C]:LEU:HD13	2.16	0.44
2:A:93[D]:ILE:HD12	2:A:94[D]:ILE:N	2.33	0.44
2:A:117[C]:ALA:O	2:A:121[C]:ILE:HG12	2.18	0.44
2:A:6[C]:THR:HB	2:A:9[C]:GLN:HG3	1.99	0.44
1:R:4[E]:G:C2	2:A:119[E]:VAL:HG22	2.52	0.44
2:A:27[E]:CYS:HG	2:A:48[E]:PHE:HZ	1.59	0.44
2:A:77[D]:ASP:HB3	2:A:78[D]:PRO:HD3	1.99	0.44
2:A:91[B]:ASN:CB	2:A:92[B]:ARG:HA	2.41	0.44
2:A:2[B]:TYR:HB3	2:A:150[B]:LEU:HD22	1.99	0.44
2:A:115[D]:ASP:O	2:A:119[D]:VAL:HG23	2.18	0.44
2:A:62[B]:PHE:O	2:A:141[B]:ARG:NH1	2.51	0.44
2:A:9[B]:GLN:HG3	2:A:148[B]:SER:HB2	2.00	0.44
2:A:75[E]:VAL:HG13	2:A:76[E]:LEU:N	2.33	0.43
1:R:4[A]:G:C8	2:A:119[A]:VAL:HG11	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:75[D]:VAL:HG13	2:A:76[D]:LEU:HD12	2.00	0.43
2:A:150[E]:LEU:HD12	2:A:150[E]:LEU:H	1.84	0.43
2:A:83[D]:LEU:HD21	2:A:125[D]:ILE:CG1	2.48	0.43
2:A:9[A]:GLN:NE2	2:A:150[A]:LEU:HG	2.29	0.43
2:A:27[E]:CYS:SG	2:A:48[E]:PHE:HZ	2.41	0.43
2:A:27[E]:CYS:SG	2:A:48[E]:PHE:CE2	3.09	0.43
2:A:12[D]:PHE:HD2	2:A:144[D]:PHE:HE1	1.67	0.43
1:R:4[C]:G:H4'	1:R:4[C]:G:OP2	2.18	0.43
2:A:60[B]:VAL:HG12	2:A:61[B]:ARG:N	2.28	0.43
2:A:1[E]:SER:HB2	2:A:153[E]:THR:O	2.19	0.43
2:A:86[E]:ALA:HB1	2:A:120[E]:ALA:HB1	1.99	0.43
2:A:60[C]:VAL:HG12	2:A:61[C]:ARG:N	2.27	0.43
2:A:105[E]:ALA:O	2:A:108[E]:LEU:HB3	2.18	0.43
2:A:44[B]:VAL:O	2:A:47[B]:GLN:HB3	2.19	0.43
2:A:13[E]:LEU:HD13	2:A:57[E]:GLN:HA	2.00	0.42
2:A:2[E]:TYR:CD2	2:A:57[E]:GLN:O	2.71	0.42
2:A:26[D]:LEU:HD12	2:A:51[D]:VAL:CG1	2.49	0.42
2:A:17[C]:TRP:CD1	2:A:56[C]:PRO:HD3	2.53	0.42
2:A:52[C]:TRP:CZ3	2:A:71[C]:ARG:HG3	2.54	0.42
2:A:21[A]:ILE:O	2:A:25[A]:ASN:HB2	2.19	0.42
2:A:6[B]:THR:HG22	2:A:8[B]:SER:H	1.83	0.42
2:A:62[E]:PHE:HB2	2:A:144[E]:PHE:CD2	2.55	0.42
2:A:30[E]:ALA:O	2:A:33[E]:ASN:HB2	2.19	0.42
2:A:99[D]:GLN:N	2:A:100[D]:ALA:CA	2.83	0.42
1:R:4[C]:G:C8	2:A:119[C]:VAL:HG21	2.55	0.42
2:A:60[E]:VAL:HG12	2:A:61[E]:ARG:N	2.28	0.42
2:A:61[C]:ARG:HB2	2:A:152[C]:TRP:CZ3	2.55	0.42
1:R:4[A]:G:N9	2:A:119[A]:VAL:HG21	2.34	0.42
2:A:151[C]:VAL:HG12	2:A:152[C]:TRP:N	2.35	0.42
2:A:101[D]:ASN:HA	2:A:102[D]:PRO:HD3	1.89	0.42
2:A:2[A]:TYR:CE1	2:A:152[A]:TRP:HE3	2.38	0.42
1:R:5[C]:A:O2'	1:R:6[C]:A:OP1	2.32	0.42
2:A:87[B]:PHE:CE2	2:A:121[B]:ILE:HD12	2.55	0.42
2:A:107[E]:THR:HG22	2:A:107[E]:THR:O	2.19	0.42
2:A:102[E]:PRO:O	2:A:103[E]:THR:HB	2.20	0.41
2:A:91[A]:ASN:HB2	2:A:92[A]:ARG:CA	2.42	0.41
2:A:99[B]:GLN:N	2:A:100[B]:ALA:CA	2.83	0.41
1:R:4[D]:G:H2'	1:R:5[D]:A:OP1	2.20	0.41
2:A:16[D]:ALA:HB1	2:A:71[D]:ARG:CB	2.50	0.41
2:A:73[E]:ASN:O	2:A:75[E]:VAL:N	2.54	0.41
1:R:4[C]:G:H2'	2:A:119[C]:VAL:HG21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:108[C]:LEU:HA	2:A:111[C]:THR:HG22	2.03	0.41
1:R:6[B]:A:C8	2:A:116[B]:ASP:CB	3.04	0.41
1:R:4[C]:G:N9	2:A:119[C]:VAL:HG21	2.36	0.41
2:A:99[C]:GLN:N	2:A:100[C]:ALA:CA	2.84	0.41
2:A:41[E]:ARG:HD3	2:A:87[E]:PHE:O	2.21	0.41
2:A:102[C]:PRO:O	2:A:103[C]:THR:HB	2.20	0.41
1:R:4[B]:G:H1'	2:A:112[B]:ARG:NH2	2.36	0.41
2:A:99[A]:GLN:HA	2:A:99[A]:GLN:HE21	1.85	0.41
2:A:102[A]:PRO:O	2:A:103[A]:THR:HB	2.21	0.41
2:A:27[B]:CYS:SG	2:A:48[B]:PHE:CZ	3.09	0.41
2:A:76[E]:LEU:HG	2:A:79[E]:LEU:HD12	2.02	0.40
2:A:102[D]:PRO:O	2:A:103[D]:THR:HB	2.21	0.40
2:A:102[B]:PRO:O	2:A:103[B]:THR:HB	2.21	0.40
2:A:62[B]:PHE:CZ	2:A:68[B]:LYS:HB2	2.48	0.40
2:A:99[E]:GLN:N	2:A:100[E]:ALA:CA	2.85	0.40
2:A:15[C]:SER:OG	2:A:54[C]:PRO:HD3	2.20	0.40
2:A:61[B]:ARG:HD3	2:A:152[B]:TRP:CE3	2.56	0.40
2:A:38[E]:GLN:NE2	2:A:89[E]:THR:O	2.55	0.40
1:R:4[A]:G:C8	2:A:119[A]:VAL:HG21	2.57	0.40
2:A:76[A]:LEU:O	2:A:80[A]:VAL:HG23	2.21	0.40
2:A:18[C]:ALA:HB2	2:A:52[C]:TRP:CD1	2.54	0.40
2:A:4[B]:ILE:HG21	2:A:10[B]:PHE:HD2	1.86	0.40
2:A:60[D]:VAL:HG12	2:A:61[D]:ARG:N	2.27	0.40
2:A:17[D]:TRP:CZ3	2:A:70[D]:TYR:HB2	2.57	0.40
2:A:133[A]:ILE:O	2:A:133[A]:ILE:HG22	2.21	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 PDB entries followed by that with respect to all EM entries.

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
2	1-A	152/158 (96%)	129 (85%)	17 (11%)	6 (4%)	4 37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	2-A	152/158 (96%)	132 (87%)	15 (10%)	5 (3%)	5	43
2	3-A	152/158 (96%)	135 (89%)	13 (9%)	4 (3%)	7	47
2	4-A	152/158 (96%)	130 (86%)	18 (12%)	4 (3%)	7	47
2	5-A	152/158 (96%)	129 (85%)	19 (12%)	4 (3%)	7	47
All	All	760/790 (96%)	655 (86%)	82 (11%)	23 (3%)	9	44

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	1-A	14[A]	SER
2	3-A	74[C]	ALA
2	4-A	74[D]	ALA
2	4-A	141[D]	ARG
2	1-A	74[A]	ALA
2	1-A	141[A]	ARG
2	2-A	54[B]	PRO
2	2-A	74[B]	ALA
2	2-A	141[B]	ARG
2	3-A	15[C]	SER
2	4-A	102[D]	PRO
2	5-A	74[E]	ALA
2	5-A	141[E]	ARG
2	1-A	103[A]	THR
2	3-A	101[C]	ASN
2	1-A	100[A]	ALA
2	1-A	101[A]	ASN
2	2-A	101[B]	ASN
2	3-A	100[C]	ALA
2	5-A	103[E]	THR
2	2-A	100[B]	ALA
2	5-A	100[E]	ALA
2	4-A	54[D]	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 PDB entries followed by that with respect to all EM entries.

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
2	1-A	136/138 (99%)	120 (88%)	16 (12%)	6 34

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	1-A	2[A]	TYR
2	1-A	10[A]	PHE
2	1-A	14[A]	SER
2	1-A	19[A]	ASP
2	1-A	41[A]	ARG
2	1-A	48[A]	PHE
2	1-A	57[A]	GLN
2	1-A	62[A]	PHE
2	1-A	87[A]	PHE
2	1-A	92[A]	ARG
2	1-A	95[A]	GLU
2	1-A	96[A]	VAL
2	1-A	99[A]	GLN
2	1-A	132[A]	LEU
2	1-A	141[A]	ARG
2	1-A	150[A]	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
2	1-A	9[A]	GLN
2	1-A	45[A]	GLN
2	1-A	99[A]	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	1-R	2/3 (66%)	2 (100%)	0
1	2-R	0/3	-	-
1	3-R	0/3	-	-
1	4-R	0/3	-	-
1	5-R	0/3	-	-
All	All	2/15 (13%)	2 (100%)	0

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	1-R	5[A]	A
1	1-R	6[A]	A

There are no RNA pucker outliers to report.

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.