



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 02:50 am GMT

PDB ID : 1SMV
Title : PRIMARY STRUCTURE OF SESBANIA MOSAIC VIRUS COAT PROTEIN: ITS IMPLICATIONS TO THE ASSEMBLY AND ARCHITECTURE OF THE VIRUS
Authors : Bhuvaneshwari, M.; Murthy, M.R.N.
Deposited on : 1995-06-16
Resolution : 3.00 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

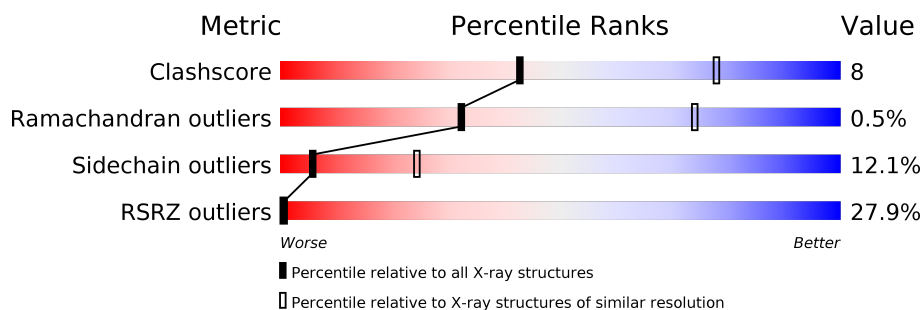
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (3.00-3.00)
RSRZ outliers	101464	1716 (3.00-3.00)

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.

Mol	Chain	Length	Quality of chain
1	A	266	
1	B	266	
1	C	266	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4516 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 SESBANIA MOSAIC VIRUS COAT PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	0	0	0
			1437	911	234	284	8			
1	B	196	Total	C	N	O	S	0	0	0
			1452	924	234	287	7			
1	C	222	Total	C	N	O	S	0	0	0
			1623	1028	269	316	10			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ILE	LYS	CONFLICT	UNP Q9EB06
A	36	PRO	GLN	CONFLICT	UNP Q9EB06
A	38	ILE	THR	CONFLICT	UNP Q9EB06
A	66	ALA	GLY	CONFLICT	UNP Q9EB06
A	?	-	THR	DELETION	UNP Q9EB06
A	72	CYS	SER	CONFLICT	UNP Q9EB06
A	75	THR	SER	CONFLICT	UNP Q9EB06
A	105	ASP	ALA	CONFLICT	UNP Q9EB06
A	149	LYS	GLN	CONFLICT	UNP Q9EB06
A	173	ASN	GLY	CONFLICT	UNP Q9EB06
A	174	SER	THR	CONFLICT	UNP Q9EB06
A	191	GLU	LYS	CONFLICT	UNP Q9EB06
A	218	ASP	PRO	CONFLICT	UNP Q9EB06
A	244	ASP	CYS	CONFLICT	UNP Q9EB06
B	0	ILE	LYS	CONFLICT	UNP Q9EB06
B	36	PRO	GLN	CONFLICT	UNP Q9EB06
B	38	ILE	THR	CONFLICT	UNP Q9EB06
B	66	ALA	GLY	CONFLICT	UNP Q9EB06
B	?	-	THR	DELETION	UNP Q9EB06
B	72	CYS	SER	CONFLICT	UNP Q9EB06
B	75	THR	SER	CONFLICT	UNP Q9EB06
B	105	ASP	ALA	CONFLICT	UNP Q9EB06
B	149	LYS	GLN	CONFLICT	UNP Q9EB06

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Chain	Residue	Modelled	Actual	Comment	Reference
B	173	ASN	GLY	CONFLICT	UNP Q9EB06
B	174	SER	THR	CONFLICT	UNP Q9EB06
B	191	GLU	LYS	CONFLICT	UNP Q9EB06
B	218	ASP	PRO	CONFLICT	UNP Q9EB06
B	244	ASP	CYS	CONFLICT	UNP Q9EB06
C	0	ILE	LYS	CONFLICT	UNP Q9EB06
C	36	PRO	GLN	CONFLICT	UNP Q9EB06
C	38	ILE	THR	CONFLICT	UNP Q9EB06
C	66	ALA	GLY	CONFLICT	UNP Q9EB06
C	?	-	THR	DELETION	UNP Q9EB06
C	72	CYS	SER	CONFLICT	UNP Q9EB06
C	75	THR	SER	CONFLICT	UNP Q9EB06
C	105	ASP	ALA	CONFLICT	UNP Q9EB06
C	149	LYS	GLN	CONFLICT	UNP Q9EB06
C	173	ASN	GLY	CONFLICT	UNP Q9EB06
C	174	SER	THR	CONFLICT	UNP Q9EB06
C	191	GLU	LYS	CONFLICT	UNP Q9EB06
C	218	ASP	PRO	CONFLICT	UNP Q9EB06
C	244	ASP	CYS	CONFLICT	UNP Q9EB06

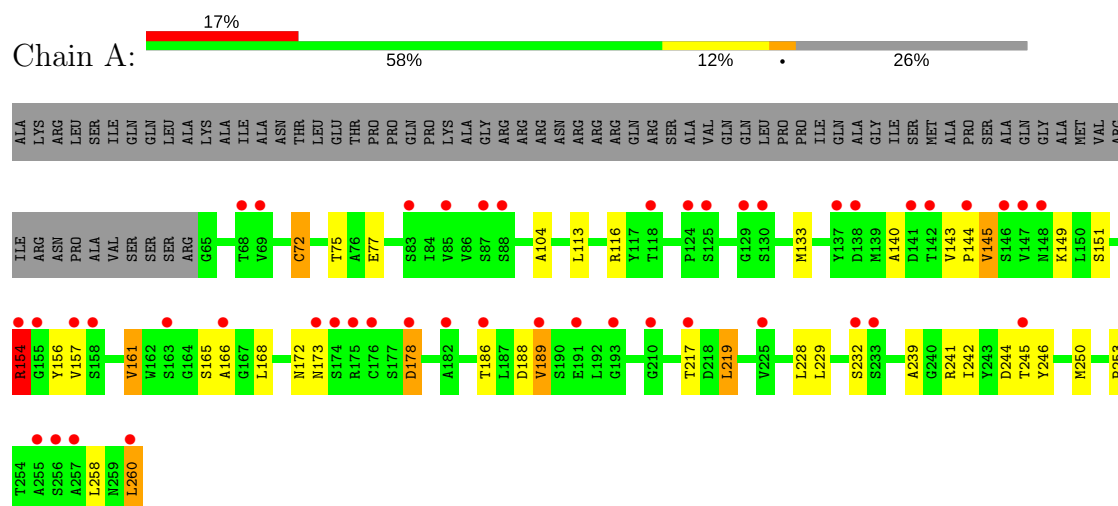
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Ca 3 3	0	0
2	C	1	Total Ca 1 1	0	0

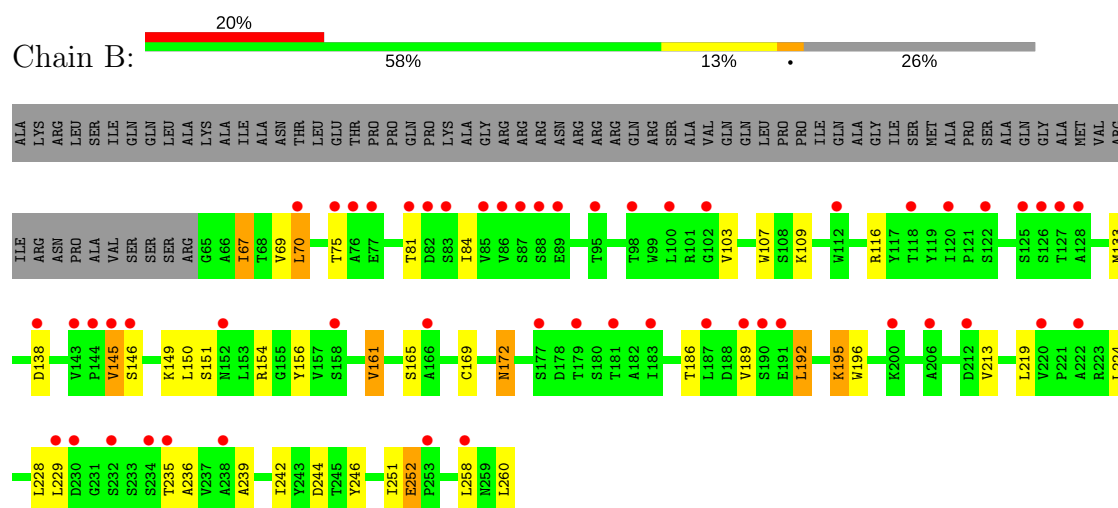
3 Residue-property plots

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

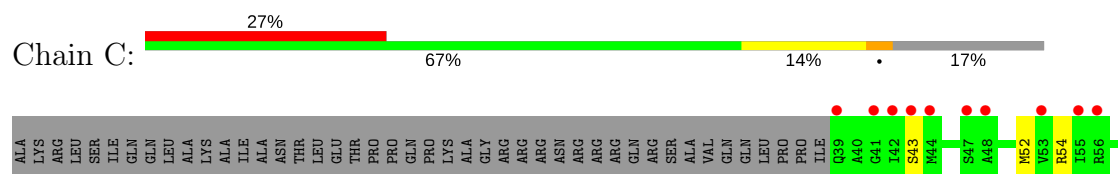
• Molecule 1: SESBANIA MOSAIC VIRUS COAT PROTEIN

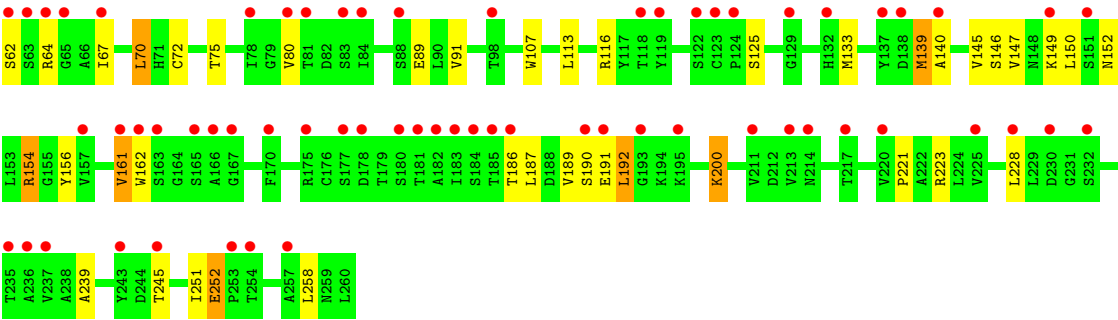


• Molecule 1: SESBANIA MOSAIC VIRUS COAT PROTEIN



• Molecule 1: SESBANIA MOSAIC VIRUS COAT PROTEIN





4 Data and refinement statistics

Property	Value	Source
Space group	R 3	Depositor
Cell constants a, b, c, α , β , γ	291.46Å 291.46Å 291.46Å 61.95° 61.95° 61.95°	Depositor
Resolution (Å)	10.00 – 3.00 51.40 – 2.90	Depositor EDS
% Data completeness (in resolution range)	70.0 (10.00-3.00) 61.1 (51.40-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.227 , (Not available) (Not available) , (Not available)	Depositor DCC
R_{free} test set	NotAvailable	DCC
Wilson B-factor (Å ²)	(Not available)	Xtriage
Anisotropy	(Not available)	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 70.4	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.15	EDS
Total number of atoms	4516	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *(Not available)*

¹Theoretical 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

Bond lengths and bond angles in the following residue types are not validated in this section: CA

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	0.62	0/1468	0.84	2/2012 (0.1%)
1	B	0.59	0/1483	0.83	1/2033 (0.0%)
1	C	0.60	0/1656	0.81	2/2268 (0.1%)
All	All	0.60	0/4607	0.82	5/6313 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	70	LEU	CA-CB-CG	-7.71	97.56	115.30
1	C	70	LEU	CA-CB-CG	-7.02	99.15	115.30
1	A	154	ARG	NE-CZ-NH2	-5.64	117.48	120.30
1	C	70	LEU	CB-CG-CD2	-5.58	101.52	111.00
1	A	219	LEU	CA-CB-CG	5.07	126.96	115.30

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	1437	0	1369	24	0
1	B	1452	0	1416	26	0
1	C	1623	0	1586	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	3	0	0	0	0
2	C	1	0	0	0	0
All	All	4516	0	4371	69	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 8.

All (69) 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:A:145:VAL:HG23	1:A:149:LYS:HE3	1.42	1.00
1:A:72:CYS:SG	1:A:245:THR:HG23	2.08	0.93
1:A:149:LYS:HD3	1:B:258:LEU:HD21	1.47	0.92
1:B:116:ARG:HB2	1:B:186:THR:HG22	1.52	0.91
1:C:72:CYS:SG	1:C:245:THR:HG23	2.16	0.85
1:C:145:VAL:H	1:C:149:LYS:HE2	1.44	0.83
1:C:145:VAL:HG22	1:C:149:LYS:HE2	1.64	0.78
1:A:104:ALA:HA	1:A:250:MET:CE	2.21	0.71
1:A:145:VAL:H	1:A:149:LYS:HD2	1.56	0.70
1:A:75:THR:HG22	1:A:242:ILE:O	1.92	0.70
1:C:67:ILE:HD13	1:C:107:TRP:HE1	1.57	0.70
1:A:104:ALA:HA	1:A:250:MET:HE2	1.76	0.66
1:B:172:ASN:HD22	1:B:172:ASN:H	1.44	0.66
1:C:62:SER:HB3	1:C:67:ILE:HG13	1.80	0.64
1:B:189:VAL:HA	1:B:192:LEU:HD22	1.80	0.62
1:C:161:VAL:HG22	1:C:239:ALA:HB1	1.81	0.62
1:B:75:THR:HG22	1:B:242:ILE:O	1.99	0.62
1:C:133:MET:O	1:C:156:TYR:HA	2.00	0.60
1:A:161:VAL:HG22	1:A:239:ALA:HB1	1.85	0.58
1:C:116:ARG:HB2	1:C:186:THR:HG22	1.86	0.57
1:A:260:LEU:HD23	1:C:140:ALA:O	2.05	0.57
1:A:77:GLU:HG2	1:A:241:ARG:HG2	1.87	0.57
1:C:251:ILE:HG13	1:C:252:GLU:HG2	1.87	0.56
1:A:72:CYS:SG	1:A:245:THR:CG2	2.90	0.55
1:A:151:SER:HA	1:A:156:TYR:CD1	2.41	0.54
1:C:139:MET:HE1	1:C:223:ARG:HG2	1.90	0.54
1:C:67:ILE:HD13	1:C:107:TRP:NE1	2.21	0.53
1:A:244:ASP:OD1	1:A:246:TYR:HD2	1.90	0.53
1:B:251:ILE:HG13	1:B:252:GLU:HG2	1.90	0.53
1:A:258:LEU:HD21	1:C:149:LYS:NZ	2.23	0.53
1:B:229:LEU:N	1:B:229:LEU:HD23	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:154:ARG:HH11	1:C:154:ARG:HG2	1.74	0.51
1:C:145:VAL:H	1:C:149:LYS:CE	2.20	0.51
1:B:109:LYS:HE2	1:B:252:GLU:OE2	2.11	0.51
1:B:172:ASN:HD22	1:B:172:ASN:N	2.06	0.50
1:A:144:PRO:HA	1:A:149:LYS:HD2	1.93	0.50
1:C:189:VAL:HA	1:C:192:LEU:HD22	1.93	0.50
1:C:139:MET:HA	1:C:139:MET:CE	2.42	0.50
1:A:133:MET:O	1:A:156:TYR:HA	2.13	0.48
1:B:145:VAL:HG12	1:B:149:LYS:CE	2.44	0.48
1:C:113:LEU:HA	1:C:113:LEU:HD23	1.67	0.47
1:B:133:MET:O	1:B:156:TYR:HA	2.14	0.47
1:A:140:ALA:O	1:B:260:LEU:HD12	2.16	0.46
1:C:125:SER:HA	1:C:162:TRP:CG	2.50	0.45
1:A:104:ALA:HA	1:A:250:MET:HE1	1.98	0.45
1:B:138:ASP:OD2	1:C:200:LYS:HE3	2.17	0.45
1:A:154:ARG:CZ	1:A:188:ASP:HB3	2.47	0.45
1:B:67:ILE:HG12	1:B:107:TRP:CZ2	2.52	0.44
1:C:145:VAL:HG22	1:C:149:LYS:CE	2.42	0.44
1:A:166:ALA:HB1	1:A:178:ASP:O	2.16	0.44
1:C:146:SER:OG	1:C:149:LYS:HD3	2.17	0.44
1:B:116:ARG:HB2	1:B:186:THR:CG2	2.37	0.44
1:B:195:LYS:HD2	1:B:196:TRP:CE2	2.53	0.44
1:B:169:CYS:HA	1:B:172:ASN:HD21	1.82	0.43
1:A:258:LEU:HD21	1:C:149:LYS:HZ1	1.83	0.43
1:B:81:THR:HG22	1:B:236:ALA:HA	2.01	0.43
1:B:145:VAL:H	1:B:149:LYS:HD2	1.84	0.42
1:B:84:ILE:HA	1:B:229:LEU:HB3	2.01	0.42
1:A:77:GLU:OE2	1:A:241:ARG:NH1	2.53	0.42
1:B:172:ASN:ND2	1:B:172:ASN:H	2.13	0.42
1:B:244:ASP:OD2	1:B:246:TYR:HD2	2.03	0.42
1:B:161:VAL:HB	1:B:239:ALA:HB1	2.01	0.42
1:A:172:ASN:O	1:A:172:ASN:CG	2.58	0.42
1:C:192:LEU:HG	1:C:221:PRO:HB2	2.03	0.41
1:B:151:SER:HA	1:B:156:TYR:CD1	2.55	0.41
1:B:146:SER:H	1:B:149:LYS:HD2	1.86	0.40
1:A:253:PRO:HB2	1:C:152:ASN:O	2.21	0.40
1:B:251:ILE:C	1:B:252:GLU:HG2	2.40	0.40
1:C:139:MET:HA	1:C:139:MET:HE3	2.02	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	194/266 (73%)	180 (93%)	12 (6%)	2 (1%)	18	59
1	B	194/266 (73%)	184 (95%)	9 (5%)	1 (0%)	32	74
1	C	220/266 (83%)	211 (96%)	9 (4%)	0	100	100
All	All	608/798 (76%)	575 (95%)	30 (5%)	3 (0%)	32	74

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	173	ASN
1	B	161	VAL
1	A	189	VAL

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	150/220 (68%)	131 (87%)	19 (13%)	5	22
1	B	157/220 (71%)	140 (89%)	17 (11%)	7	29
1	C	172/220 (78%)	150 (87%)	22 (13%)	5	22
All	All	479/660 (73%)	421 (88%)	58 (12%)	6	24

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

Mol	Chain	Res	Type
1	A	72	CYS
1	A	113	LEU
1	A	116	ARG
1	A	143	VAL
1	A	145	VAL
1	A	154	ARG
1	A	157	VAL
1	A	161	VAL
1	A	165	SER
1	A	168	LEU
1	A	178	ASP
1	A	186	THR
1	A	189	VAL
1	A	217	THR
1	A	219	LEU
1	A	228	LEU
1	A	229	LEU
1	A	232	SER
1	A	260	LEU
1	B	67	ILE
1	B	69	VAL
1	B	70	LEU
1	B	103	VAL
1	B	145	VAL
1	B	150	LEU
1	B	154	ARG
1	B	165	SER
1	B	172	ASN
1	B	192	LEU
1	B	195	LYS
1	B	213	VAL
1	B	219	LEU
1	B	224	LEU
1	B	228	LEU
1	B	235	THR
1	B	252	GLU
1	C	43	SER
1	C	52	MET
1	C	54	ARG
1	C	64	ARG
1	C	70	LEU
1	C	75	THR
1	C	80	VAL

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Mol	Chain	Res	Type
1	C	89	GLU
1	C	91	VAL
1	C	139	MET
1	C	147	VAL
1	C	150	LEU
1	C	154	ARG
1	C	161	VAL
1	C	187	LEU
1	C	190	SER
1	C	191	GLU
1	C	192	LEU
1	C	200	LYS
1	C	228	LEU
1	C	252	GLU
1	C	258	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	172	ASN
1	C	49	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

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	196/266 (73%)	1.28	45 (22%) 1 0	2, 9, 32, 74	0
1	B	196/266 (73%)	1.45	53 (27%) 1 0	3, 10, 33, 58	0
1	C	222/266 (83%)	1.60	73 (32%) 0 0	3, 10, 34, 75	0
All	All	614/798 (76%)	1.45	171 (27%) 1 0	2, 9, 34, 75	0

All (171) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	39	GLN	11.5
1	B	82	ASP	10.0
1	B	145	VAL	9.3
1	A	88	SER	9.2
1	A	176	CYS	8.6
1	C	62	SER	6.9
1	B	235	THR	6.6
1	B	127	THR	6.0
1	C	177	SER	5.8
1	A	118	THR	5.3
1	A	148	ASN	5.3
1	B	98	THR	5.3
1	C	123	CYS	5.1
1	B	253	PRO	4.9
1	A	193	GLY	4.8
1	A	245	THR	4.7
1	C	183	ILE	4.7
1	B	128	ALA	4.7
1	C	167	GLY	4.7
1	C	163	SER	4.7
1	C	175	ARG	4.6
1	C	181	THR	4.5
1	A	155	GLY	4.5

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Mol	Chain	Res	Type	RSRZ
1	B	83	SER	4.5
1	C	48	ALA	4.3
1	A	154	ARG	4.3
1	A	217	THR	4.2
1	A	144	PRO	4.2
1	B	144	PRO	4.2
1	B	234	SER	4.2
1	C	124	PRO	4.2
1	C	122	SER	4.2
1	B	138	ASP	4.1
1	A	210	GLY	4.1
1	C	195	LYS	4.0
1	B	177	SER	4.0
1	C	63	SER	3.9
1	C	230	ASP	3.9
1	C	186	THR	3.8
1	C	65	GLY	3.8
1	C	161	VAL	3.8
1	B	77	GLU	3.8
1	A	138	ASP	3.7
1	C	237	VAL	3.7
1	C	257	ALA	3.6
1	C	182	ALA	3.6
1	C	140	ALA	3.6
1	B	81	THR	3.5
1	C	56	ARG	3.5
1	B	181	THR	3.5
1	B	146	SER	3.4
1	B	158	SER	3.4
1	C	232	SER	3.3
1	C	132	HIS	3.3
1	C	254	THR	3.3
1	A	257	ALA	3.3
1	C	44	MET	3.2
1	B	189	VAL	3.2
1	C	43	SER	3.2
1	A	174	SER	3.2
1	A	166	ALA	3.2
1	C	166	ALA	3.2
1	B	229	LEU	3.2
1	B	179	THR	3.2
1	A	178	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	143	VAL	3.1
1	C	151	SER	3.1
1	C	213	VAL	3.1
1	B	220	VAL	3.1
1	C	118	THR	3.1
1	C	67	ILE	3.1
1	A	141	ASP	3.1
1	B	87	SER	3.1
1	C	191	GLU	3.0
1	A	163	SER	3.0
1	A	225	VAL	3.0
1	C	178	ASP	3.0
1	C	149	LYS	3.0
1	A	69	VAL	2.9
1	C	53	VAL	2.9
1	C	55	ILE	2.9
1	A	142	THR	2.9
1	C	47	SER	2.9
1	C	228	LEU	2.9
1	A	189	VAL	2.9
1	A	146	SER	2.9
1	C	245	THR	2.9
1	A	124	PRO	2.9
1	B	187	LEU	2.8
1	A	175	ARG	2.8
1	C	98	THR	2.8
1	B	126	SER	2.8
1	B	120	ILE	2.8
1	C	193	GLY	2.7
1	A	186	THR	2.7
1	C	162	TRP	2.7
1	A	85	VAL	2.7
1	B	75	THR	2.7
1	B	212	ASP	2.7
1	C	217	THR	2.6
1	A	87	SER	2.6
1	A	256	SER	2.6
1	C	165	SER	2.6
1	B	122	SER	2.6
1	C	243	TYR	2.6
1	C	42	ILE	2.6
1	C	157	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	129	GLY	2.5
1	B	230	ASP	2.5
1	B	191	GLU	2.5
1	C	41	GLY	2.5
1	C	170	PHE	2.5
1	C	184	SER	2.5
1	B	89	GLU	2.5
1	A	182	ALA	2.5
1	C	236	ALA	2.4
1	C	185	THR	2.4
1	B	86	VAL	2.4
1	A	232	SER	2.4
1	B	85	VAL	2.4
1	A	157	VAL	2.4
1	B	118	THR	2.4
1	B	125	SER	2.4
1	C	81	THR	2.4
1	A	147	VAL	2.3
1	C	138	ASP	2.3
1	C	84	ILE	2.3
1	C	220	VAL	2.3
1	B	88	SER	2.3
1	A	125	SER	2.3
1	A	260	LEU	2.3
1	B	222	ALA	2.3
1	C	64	ARG	2.3
1	C	180	SER	2.3
1	C	119	TYR	2.3
1	B	95	THR	2.3
1	C	78	ILE	2.2
1	A	83	SER	2.2
1	B	166	ALA	2.2
1	A	137	TYR	2.2
1	A	173	ASN	2.2
1	B	238	ALA	2.2
1	B	112	TRP	2.2
1	B	183	ILE	2.2
1	B	70	LEU	2.2
1	C	80	VAL	2.2
1	B	200	LYS	2.2
1	B	232	SER	2.2
1	A	255	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	102	GLY	2.1
1	A	233	SER	2.1
1	C	88	SER	2.1
1	A	129	GLY	2.1
1	B	100	LEU	2.1
1	B	152	ASN	2.1
1	B	190	SER	2.1
1	B	258	LEU	2.1
1	C	214	ASN	2.1
1	C	211	VAL	2.1
1	A	191	GLU	2.0
1	B	76	ALA	2.0
1	C	137	TYR	2.0
1	C	225	VAL	2.0
1	A	130	SER	2.0
1	C	83	SER	2.0
1	B	206	ALA	2.0
1	A	68	THR	2.0
1	C	253	PRO	2.0
1	A	158	SER	2.0
1	C	190	SER	2.0
1	C	235	THR	2.0

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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	C	302	1/1	0.73	0.17	-1.63	11,11,11,11	0
2	CA	A	300	1/1	0.90	0.18	-1.65	10,10,10,10	0
2	CA	A	304	1/1	0.88	0.18	-1.92	2,2,2,2	0
2	CA	A	303	1/1	0.94	0.19	-2.00	10,10,10,10	0

6.5 Other polymers [i](#)

There are no such residues in this entry.