



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 03:05 am GMT

PDB ID : 1X35
Title : Recombinant T=3 capsid of a site specific mutant of SeMV CP
Authors : Sangita, V.; Lokesh, G.L.; Satheshkumar, P.S.; Saravanan, V.; Vijay, C.S.;
Savithri, H.S.; Murthy, M.R.
Deposited on : 2005-04-29
Resolution : 4.10 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

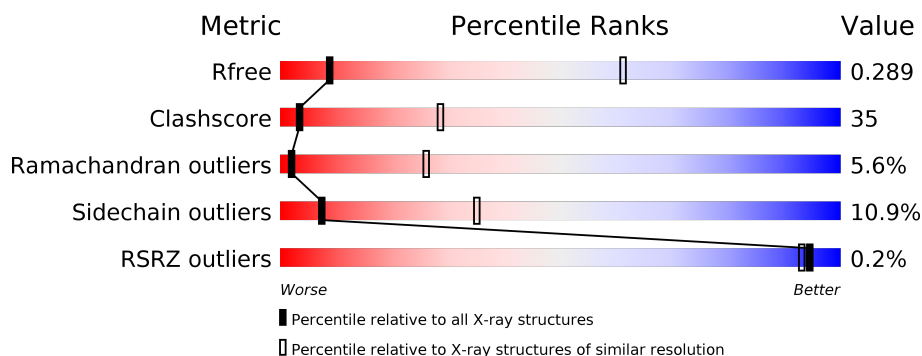
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.10 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))
R_{free}	100719	1153 (4.60-3.60)
Clashscore	112137	1002 (4.54-3.66)
Ramachandran outliers	110173	1000 (4.58-3.62)
Sidechain outliers	110143	1191 (4.60-3.60)
RSRZ outliers	101464	1165 (4.60-3.60)

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	268	
1	B	268	
1	C	268	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4442 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	196	Total	C	N	O	S	0	0	0
			1420	902	232	278	8			
1	B	197	Total	C	N	O	S	0	0	0
			1402	883	229	282	8			
1	C	225	Total	C	N	O	S	0	0	0
			1617	1020	269	318	10			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	53	ALA	PRO	ENGINEERED	UNP Q9EB06
B	53	ALA	PRO	ENGINEERED	UNP Q9EB06
C	53	ALA	PRO	ENGINEERED	UNP Q9EB06

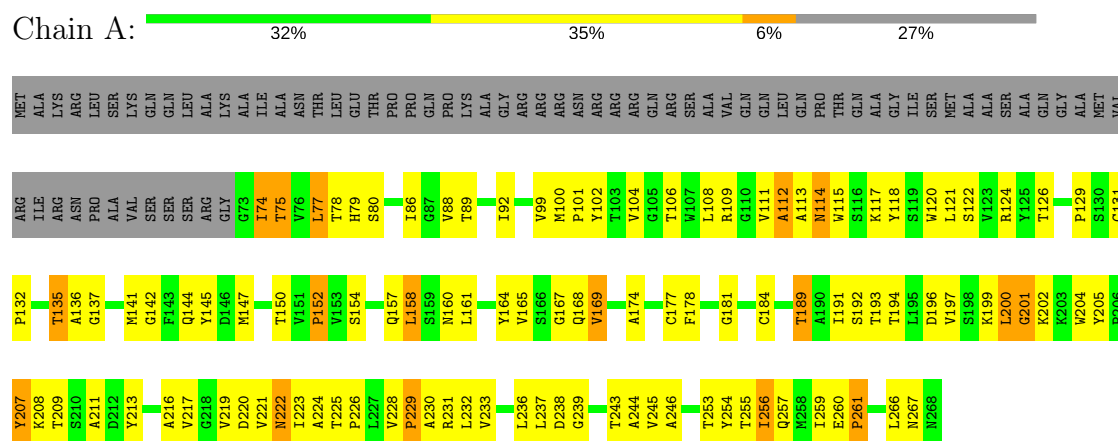
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

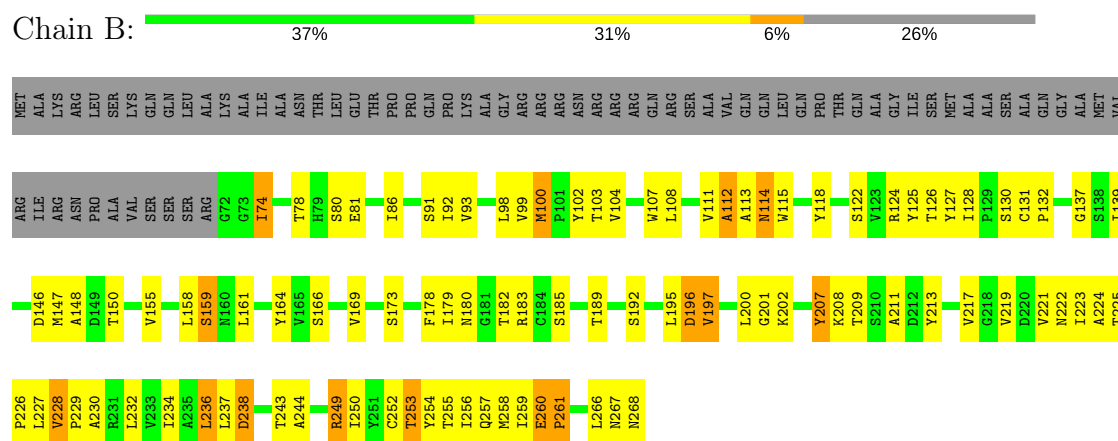
3 Residue-property plots

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

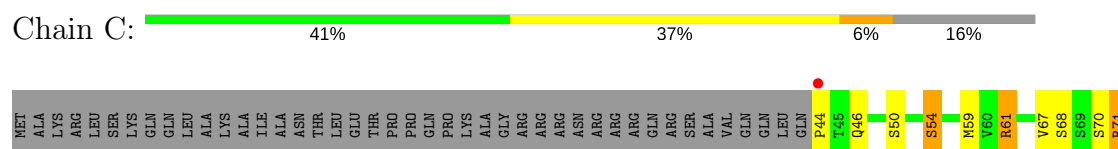
• Molecule 1: Coat protein



• Molecule 1: Coat protein



• Molecule 1: Coat protein





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	471.50Å 330.10Å 351.47Å 90.00° 131.05° 90.00°	Depositor
Resolution (Å)	20.00 – 4.10 25.07 – 4.06	Depositor EDS
% Data completeness (in resolution range)	94.8 (20.00-4.10) 88.9 (25.07-4.06)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.25 (at 4.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.268 , 0.270 0.287 , 0.289	Depositor DCC
R_{free} test set	14986 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	60.1	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , -1.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.23$, $\langle L^2 \rangle = 0.08$	Xtriage
Estimated twinning fraction	0.297 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.76	EDS
Total number of atoms	4442	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.77% 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

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	0.42	0/1452	0.70	0/1994
1	B	0.42	0/1433	0.68	0/1969
1	C	0.41	0/1650	0.74	3/2261 (0.1%)
All	All	0.42	0/4535	0.71	3/6224 (0.0%)

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
1	C	1	0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	54	SER	N-CA-CB	5.94	119.40	110.50
1	C	242	SER	CB-CA-C	5.11	119.81	110.10
1	C	44	PRO	N-CA-CB	5.01	109.31	103.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	C	242	SER	CA

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	1420	0	1348	117	0
1	B	1402	0	1312	106	0
1	C	1617	0	1558	104	0
2	A	2	0	0	0	0
2	C	1	0	0	0	0
All	All	4442	0	4218	304	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 35.

All (304) 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:118:TYR:HB2	1:A:256:ILE:HD11	1.39	1.01
1:C:176:LEU:HD11	1:C:180:ASN:ND2	1.76	0.98
1:B:126:THR:HG23	1:B:192:SER:HB2	1.43	0.98
1:C:109:ARG:HG2	1:C:110:GLY:N	1.79	0.95
1:A:207:TYR:HB3	1:A:267:ASN:HD21	1.34	0.92
1:A:111:VAL:HG13	1:A:112:ALA:H	1.37	0.90
1:A:74:ILE:HG23	1:A:259:ILE:HG23	1.59	0.85
1:C:86:ILE:HG22	1:C:247:ALA:HB3	1.60	0.81
1:A:126:THR:HG23	1:A:192:SER:HB2	1.61	0.80
1:A:217:VAL:HG13	1:A:221:VAL:HA	1.64	0.80
1:A:150:THR:O	1:B:266:LEU:HD22	1.81	0.79
1:A:202:LYS:HZ3	1:B:202:LYS:HZ2	1.28	0.79
1:C:162:ARG:HD3	1:C:163:GLY:N	1.99	0.78
1:C:200:LEU:HD23	1:C:229:PRO:HB2	1.65	0.78
1:C:208:LYS:HA	1:C:208:LYS:HE2	1.64	0.78
1:A:256:ILE:HG12	1:A:257:GLN:N	1.98	0.78
1:A:266:LEU:HD22	1:C:150:THR:O	1.83	0.77
1:A:202:LYS:NZ	1:B:202:LYS:NZ	2.32	0.76
1:B:111:VAL:HG23	1:B:112:ALA:N	1.99	0.76
1:B:249:ARG:HG2	1:B:249:ARG:HH11	1.47	0.76
1:C:99:VAL:HG12	1:C:230:ALA:O	1.86	0.76
1:A:202:LYS:NZ	1:B:202:LYS:HZ2	1.83	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:VAL:HG23	1:B:112:ALA:H	1.54	0.73
1:C:200:LEU:N	1:C:200:LEU:HD12	2.02	0.73
1:A:114:ASN:N	1:A:114:ASN:HD22	1.87	0.73
1:A:102:TYR:CD2	1:A:209:THR:HA	2.25	0.72
1:B:114:ASN:N	1:B:114:ASN:HD22	1.85	0.72
1:A:141:MET:HB3	1:A:232:LEU:HD11	1.72	0.72
1:B:197:VAL:HA	1:B:200:LEU:HD13	1.71	0.71
1:A:79:HIS:CD2	1:A:80:SER:H	2.09	0.71
1:B:80:SER:HB2	1:B:253:THR:HG23	1.72	0.71
1:C:102:TYR:CD2	1:C:209:THR:HA	2.25	0.70
1:C:177:CYS:HG	1:C:184:CYS:HG	1.39	0.70
1:B:159:SER:HA	1:B:164:TYR:CD1	2.26	0.70
1:C:197:VAL:HA	1:C:200:LEU:HD11	1.73	0.70
1:A:200:LEU:H	1:A:200:LEU:HD12	1.56	0.70
1:A:202:LYS:HZ3	1:B:202:LYS:NZ	1.88	0.70
1:A:205:TYR:CE2	1:A:229:PRO:HB3	2.27	0.70
1:A:74:ILE:HG23	1:A:259:ILE:CG2	2.22	0.70
1:A:100:MET:HE1	1:A:208:LYS:H	1.58	0.69
1:A:150:THR:HB	1:B:266:LEU:CD2	2.22	0.69
1:C:139:ILE:HA	1:C:236:LEU:HD23	1.73	0.69
1:C:220:ASP:OD2	1:C:222:ASN:HB2	1.93	0.68
1:C:113:ALA:HA	1:C:207:TYR:CE2	2.29	0.67
1:B:217:VAL:HG13	1:B:221:VAL:HA	1.75	0.66
1:C:224:ALA:HA	1:C:227:LEU:HD12	1.76	0.66
1:C:200:LEU:HD23	1:C:229:PRO:CB	2.26	0.66
1:A:229:PRO:HG2	1:A:230:ALA:H	1.61	0.65
1:A:132:PRO:HD2	1:A:135:THR:OG1	1.95	0.65
1:C:125:TYR:CD2	1:C:232:LEU:HD22	2.31	0.65
1:A:111:VAL:HG13	1:A:112:ALA:N	2.09	0.65
1:A:213:TYR:CE1	1:A:217:VAL:HG21	2.32	0.64
1:C:213:TYR:CE2	1:C:217:VAL:HG21	2.31	0.64
1:C:208:LYS:HD3	1:C:227:LEU:HD22	1.80	0.64
1:B:113:ALA:HA	1:B:207:TYR:CE2	2.34	0.63
1:A:137:GLY:C	1:A:169:VAL:HG23	2.21	0.62
1:B:243:THR:HG22	1:B:244:ALA:N	2.15	0.61
1:A:99:VAL:HG23	1:A:231:ARG:HA	1.82	0.61
1:C:118:TYR:HB2	1:C:256:ILE:HD11	1.82	0.61
1:B:243:THR:HG22	1:B:244:ALA:H	1.66	0.61
1:C:77:LEU:HD22	1:C:78:THR:N	2.16	0.61
1:C:217:VAL:HG13	1:C:221:VAL:HA	1.84	0.60
1:A:117:LYS:HD2	1:A:260:GLU:OE1	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:109:ARG:HG2	1:C:110:GLY:H	1.64	0.60
1:C:80:SER:OG	1:C:253:THR:HG23	2.00	0.60
1:B:207:TYR:HB3	1:B:267:ASN:HD21	1.67	0.59
1:B:249:ARG:HG2	1:B:249:ARG:NH1	2.14	0.59
1:A:108:LEU:HA	1:A:111:VAL:CG1	2.33	0.59
1:B:113:ALA:N	1:B:207:TYR:HE2	2.01	0.58
1:C:137:GLY:C	1:C:169:VAL:HG23	2.22	0.58
1:A:78:THR:HG23	1:A:255:THR:OG1	2.03	0.58
1:A:174:ALA:HB3	1:A:189:THR:HG23	1.85	0.58
1:A:74:ILE:HG22	1:A:75:THR:N	2.18	0.58
1:C:228:VAL:HG23	1:C:229:PRO:HD2	1.84	0.58
1:A:99:VAL:HG23	1:A:230:ALA:O	2.04	0.58
1:C:126:THR:HG23	1:C:192:SER:HB2	1.85	0.58
1:C:167:GLY:HA3	1:C:191:ILE:HD11	1.84	0.58
1:B:180:ASN:HD21	1:B:182:THR:HB	1.69	0.58
1:A:150:THR:HB	1:B:266:LEU:HD23	1.85	0.58
1:C:198:SER:HB2	1:C:199:LYS:HD2	1.84	0.58
1:C:259:ILE:HG13	1:C:260:GLU:HG3	1.86	0.58
1:B:256:ILE:HG12	1:B:257:GLN:N	2.18	0.57
1:A:117:LYS:HE3	1:C:145:TYR:HB3	1.86	0.57
1:B:114:ASN:N	1:B:114:ASN:ND2	2.51	0.57
1:A:200:LEU:N	1:A:200:LEU:HD12	2.20	0.57
1:A:237:LEU:N	1:A:237:LEU:HD23	2.20	0.57
1:A:80:SER:OG	1:A:253:THR:HG23	2.04	0.57
1:B:259:ILE:C	1:B:260:GLU:HG2	2.25	0.57
1:B:126:THR:HG23	1:B:192:SER:CB	2.27	0.57
1:B:127:TYR:O	1:B:128:ILE:HD13	2.04	0.57
1:B:78:THR:HG23	1:B:255:THR:OG1	2.04	0.56
1:A:131:CYS:HB2	1:A:132:PRO:CD	2.35	0.56
1:B:124:ARG:HD3	1:B:178:PHE:CD2	2.40	0.56
1:B:108:LEU:HB2	1:B:254:TYR:OH	2.05	0.56
1:B:100:MET:O	1:B:104:VAL:HG23	2.04	0.56
1:A:202:LYS:HZ1	1:B:202:LYS:NZ	2.04	0.56
1:B:111:VAL:CG2	1:B:112:ALA:H	2.17	0.56
1:C:121:LEU:HD12	1:C:257:GLN:HG3	1.87	0.56
1:B:128:ILE:HG13	1:B:179:ILE:HD11	1.88	0.55
1:C:220:ASP:HB3	1:C:223:ILE:CD1	2.35	0.55
1:C:199:LYS:C	1:C:200:LEU:HD12	2.27	0.55
1:A:126:THR:HG23	1:A:192:SER:CB	2.35	0.55
1:A:254:TYR:N	1:A:254:TYR:CD2	2.74	0.55
1:C:124:ARG:HD3	1:C:178:PHE:HD1	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:GLY:HA3	1:A:191:ILE:HD11	1.88	0.54
1:B:146:ASP:O	1:B:148:ALA:N	2.41	0.54
1:A:77:LEU:HD23	1:A:77:LEU:O	2.08	0.54
1:B:74:ILE:HG22	1:B:259:ILE:CG2	2.38	0.54
1:A:219:VAL:CG1	1:B:219:VAL:HG11	2.37	0.53
1:C:140:HIS:HB2	1:C:235:ALA:HB3	1.90	0.53
1:A:137:GLY:O	1:A:169:VAL:HG23	2.09	0.53
1:A:102:TYR:HA	1:A:109:ARG:HG3	1.90	0.53
1:C:200:LEU:N	1:C:200:LEU:CD1	2.71	0.53
1:A:225:THR:N	1:A:226:PRO:CD	2.72	0.53
1:B:180:ASN:ND2	1:B:182:THR:HB	2.24	0.53
1:A:154:SER:OG	1:A:157:GLN:HG2	2.09	0.53
1:B:225:THR:N	1:B:226:PRO:HD2	2.23	0.53
1:A:266:LEU:HD23	1:C:150:THR:HB	1.89	0.53
1:C:101:PRO:HD3	1:C:118:TYR:CE2	2.44	0.52
1:A:158:LEU:HD21	1:A:233:VAL:HG13	1.91	0.52
1:B:102:TYR:HB2	1:B:207:TYR:HE1	1.75	0.52
1:C:256:ILE:HG12	1:C:257:GLN:N	2.23	0.52
1:C:130:SER:HB3	1:C:249:ARG:NH1	2.23	0.52
1:B:207:TYR:HD1	1:B:208:LYS:N	2.07	0.52
1:B:209:THR:HG23	1:B:268:ASN:HA	1.91	0.52
1:A:225:THR:N	1:A:226:PRO:HD2	2.25	0.52
1:B:137:GLY:HA2	1:B:238:ASP:O	2.09	0.52
1:C:139:ILE:HG13	1:C:236:LEU:HD21	1.92	0.52
1:C:102:TYR:CE1	1:C:210:SER:HB2	2.44	0.51
1:B:200:LEU:HD23	1:B:229:PRO:CB	2.40	0.51
1:A:208:LYS:NZ	1:C:148:ALA:HB3	2.24	0.51
1:A:164:TYR:CG	1:A:165:VAL:N	2.78	0.51
1:A:224:ALA:C	1:A:226:PRO:HD2	2.31	0.51
1:A:114:ASN:N	1:A:114:ASN:ND2	2.59	0.51
1:A:160:ASN:HD22	1:A:160:ASN:N	2.08	0.51
1:C:124:ARG:HD3	1:C:178:PHE:CD1	2.45	0.51
1:C:133:SER:HA	1:C:170:TRP:CG	2.45	0.51
1:C:223:ILE:O	1:C:226:PRO:HD2	2.10	0.51
1:B:232:LEU:HD22	1:B:234:ILE:HG13	1.92	0.51
1:B:223:ILE:O	1:B:223:ILE:HG22	2.11	0.51
1:A:152:PRO:HA	1:B:266:LEU:HD11	1.92	0.50
1:B:213:TYR:CE2	1:B:217:VAL:HG21	2.45	0.50
1:A:100:MET:CE	1:A:208:LYS:H	2.24	0.50
1:B:113:ALA:CA	1:B:207:TYR:CE2	2.95	0.50
1:C:197:VAL:O	1:C:200:LEU:HD13	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:67:VAL:HG22	1:C:77:LEU:HD23	1.93	0.50
1:B:122:SER:O	1:B:254:TYR:HA	2.11	0.50
1:B:227:LEU:N	1:B:227:LEU:HD12	2.27	0.50
1:C:223:ILE:H	1:C:223:ILE:HD12	1.77	0.50
1:C:137:GLY:HA2	1:C:238:ASP:O	2.12	0.50
1:C:86:ILE:HG23	1:C:236:LEU:HD11	1.94	0.50
1:A:120:TRP:O	1:A:197:VAL:HG11	2.11	0.50
1:A:144:GLN:O	1:A:230:ALA:HB1	2.12	0.50
1:B:108:LEU:HA	1:B:111:VAL:HG22	1.94	0.50
1:A:142:GLY:O	1:A:232:LEU:HD12	2.11	0.49
1:B:207:TYR:CD1	1:B:207:TYR:C	2.84	0.49
1:A:145:TYR:N	1:A:145:TYR:CD2	2.79	0.49
1:A:154:SER:H	1:A:157:GLN:HG3	1.77	0.49
1:B:237:LEU:HD23	1:B:237:LEU:H	1.77	0.49
1:C:86:ILE:HG23	1:C:236:LEU:CD1	2.42	0.49
1:B:111:VAL:CG2	1:B:112:ALA:N	2.67	0.49
1:B:232:LEU:HD22	1:B:234:ILE:CG1	2.42	0.49
1:A:207:TYR:HB3	1:A:267:ASN:ND2	2.16	0.49
1:B:225:THR:N	1:B:226:PRO:CD	2.75	0.49
1:A:124:ARG:CZ	1:A:178:PHE:CE1	2.96	0.49
1:B:207:TYR:C	1:B:207:TYR:HD1	2.16	0.49
1:C:158:LEU:O	1:C:158:LEU:HD23	2.13	0.49
1:B:226:PRO:HG2	1:B:227:LEU:HD12	1.94	0.48
1:C:144:GLN:O	1:C:230:ALA:HB1	2.13	0.48
1:A:100:MET:O	1:A:104:VAL:HG23	2.14	0.48
1:A:100:MET:HE1	1:A:208:LYS:N	2.25	0.48
1:A:122:SER:O	1:A:254:TYR:HA	2.13	0.48
1:B:226:PRO:HG2	1:B:227:LEU:CD1	2.43	0.48
1:C:162:ARG:CD	1:C:163:GLY:N	2.73	0.48
1:C:147:MET:SD	1:C:147:MET:N	2.86	0.48
1:C:207:TYR:C	1:C:207:TYR:CD1	2.87	0.48
1:C:131:CYS:HB2	1:C:132:PRO:CD	2.44	0.48
1:C:228:VAL:CG2	1:C:229:PRO:HD2	2.43	0.48
1:A:216:ALA:O	1:A:223:ILE:HD12	2.14	0.48
1:A:220:ASP:OD2	1:A:222:ASN:HB2	2.14	0.48
1:A:266:LEU:CD2	1:C:150:THR:HB	2.44	0.48
1:C:197:VAL:O	1:C:200:LEU:CD1	2.62	0.48
1:C:61:ARG:N	1:C:180:ASN:OD1	2.34	0.47
1:A:121:LEU:HA	1:A:121:LEU:HD12	1.67	0.47
1:B:102:TYR:HB2	1:B:207:TYR:CE1	2.49	0.47
1:C:207:TYR:HD1	1:C:207:TYR:C	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:PRO:HD3	1:A:118:TYR:CE2	2.49	0.47
1:B:158:LEU:HD23	1:B:158:LEU:C	2.35	0.47
1:A:228:VAL:HB	1:A:229:PRO:HD2	1.95	0.47
1:B:131:CYS:HB2	1:B:132:PRO:CD	2.45	0.47
1:B:99:VAL:HG11	1:B:195:LEU:HD22	1.96	0.47
1:A:260:GLU:HB3	1:C:145:TYR:CE1	2.50	0.47
1:B:159:SER:HA	1:B:164:TYR:CG	2.50	0.47
1:C:124:ARG:NH1	1:C:178:PHE:CE1	2.83	0.47
1:A:169:VAL:HG12	1:A:169:VAL:O	2.13	0.46
1:B:209:THR:CG2	1:B:268:ASN:HA	2.45	0.46
1:C:109:ARG:CG	1:C:110:GLY:N	2.62	0.46
1:C:179:ILE:HD13	1:C:251:TYR:HB2	1.97	0.46
1:C:197:VAL:CA	1:C:200:LEU:HD11	2.44	0.46
1:A:111:VAL:CG1	1:A:112:ALA:H	2.19	0.46
1:A:150:THR:HB	1:B:266:LEU:HD22	1.96	0.46
1:A:88:VAL:O	1:A:245:VAL:HG23	2.16	0.46
1:B:221:VAL:HG23	1:B:222:ASN:N	2.31	0.46
1:A:111:VAL:O	1:A:113:ALA:N	2.49	0.46
1:A:89:THR:O	1:A:239:GLY:HA3	2.15	0.46
1:B:158:LEU:HD23	1:B:158:LEU:O	2.15	0.46
1:C:133:SER:HA	1:C:170:TRP:CD2	2.51	0.46
1:A:202:LYS:HZ1	1:B:202:LYS:HZ1	1.63	0.46
1:A:77:LEU:HD21	1:A:256:ILE:CG2	2.46	0.46
1:A:229:PRO:HG2	1:A:230:ALA:N	2.28	0.46
1:A:207:TYR:C	1:A:207:TYR:CD1	2.89	0.46
1:B:125:TYR:CD1	1:B:232:LEU:HD12	2.52	0.46
1:C:156:ASN:O	1:C:159:SER:N	2.49	0.46
1:C:147:MET:CE	1:C:231:ARG:HD3	2.45	0.46
1:B:223:ILE:O	1:B:227:LEU:HD13	2.17	0.45
1:B:236:LEU:HA	1:B:236:LEU:HD22	1.67	0.45
1:A:256:ILE:CG1	1:A:257:GLN:N	2.76	0.45
1:A:99:VAL:O	1:A:120:TRP:NE1	2.49	0.45
1:B:200:LEU:HD23	1:B:229:PRO:HB3	1.98	0.45
1:A:147:MET:SD	1:A:231:ARG:HD3	2.56	0.45
1:A:200:LEU:O	1:A:201:GLY:C	2.55	0.45
1:B:113:ALA:HA	1:B:207:TYR:CD2	2.52	0.45
1:C:120:TRP:HA	1:C:120:TRP:CE3	2.52	0.45
1:A:191:ILE:O	1:A:191:ILE:HG22	2.17	0.45
1:A:259:ILE:HB	1:A:260:GLU:HG3	1.99	0.45
1:C:147:MET:HE2	1:C:231:ARG:HB3	1.98	0.45
1:A:108:LEU:HA	1:A:111:VAL:HG12	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:LEU:HD23	1:B:229:PRO:HB2	1.98	0.44
1:C:77:LEU:HD22	1:C:78:THR:H	1.81	0.44
1:B:224:ALA:C	1:B:226:PRO:HD2	2.37	0.44
1:B:114:ASN:O	1:B:261:PRO:HA	2.17	0.44
1:C:117:LYS:HD2	1:C:260:GLU:OE2	2.17	0.44
1:C:90:ASP:OD1	1:C:90:ASP:O	2.35	0.44
1:A:237:LEU:H	1:A:237:LEU:HD23	1.81	0.44
1:C:107:TRP:CZ3	1:C:111:VAL:HG21	2.53	0.44
1:C:113:ALA:CA	1:C:207:TYR:CE2	2.98	0.44
1:B:80:SER:HA	1:B:252:CYS:O	2.18	0.44
1:B:98:LEU:HD13	1:B:228:VAL:HG11	2.00	0.44
1:A:208:LYS:NZ	1:C:148:ALA:CB	2.81	0.44
1:B:113:ALA:N	1:B:207:TYR:CE2	2.84	0.44
1:B:118:TYR:HB2	1:B:256:ILE:HD11	1.99	0.44
1:C:145:TYR:N	1:C:145:TYR:CD1	2.85	0.44
1:C:169:VAL:HG13	1:C:247:ALA:HB1	2.00	0.44
1:A:205:TYR:OH	1:A:229:PRO:HA	2.18	0.43
1:B:197:VAL:HA	1:B:200:LEU:CD1	2.43	0.43
1:C:174:ALA:HB3	1:C:190:ALA:HB2	2.00	0.43
1:C:259:ILE:HG13	1:C:260:GLU:CG	2.47	0.43
1:B:115:TRP:HA	1:B:261:PRO:HA	1.99	0.43
1:C:154:SER:OG	1:C:157:GLN:HG3	2.18	0.43
1:C:207:TYR:HB3	1:C:267:ASN:HD21	1.84	0.43
1:A:92:ILE:HG12	1:A:237:LEU:HD13	1.99	0.43
1:B:197:VAL:CA	1:B:200:LEU:HD13	2.46	0.43
1:B:229:PRO:HG2	1:B:230:ALA:H	1.83	0.43
1:C:237:LEU:HG	1:C:238:ASP:N	2.33	0.43
1:A:254:TYR:H	1:A:254:TYR:HD2	1.62	0.43
1:A:118:TYR:CB	1:A:256:ILE:HD11	2.28	0.43
1:A:204:TRP:CZ3	1:A:259:ILE:HG12	2.53	0.43
1:C:208:LYS:HD3	1:C:227:LEU:CD2	2.46	0.43
1:A:86:ILE:HG23	1:A:236:LEU:HD11	2.01	0.43
1:B:86:ILE:HD11	1:B:250:ILE:HG13	2.01	0.43
1:C:100:MET:HG2	1:C:103:THR:OG1	2.19	0.43
1:A:100:MET:HE1	1:A:208:LYS:O	2.19	0.42
1:C:113:ALA:N	1:C:207:TYR:HE2	2.16	0.42
1:A:120:TRP:CE3	1:A:120:TRP:HA	2.53	0.42
1:A:197:VAL:HA	1:A:200:LEU:HD11	2.00	0.42
1:C:229:PRO:HG2	1:C:230:ALA:H	1.84	0.42
1:A:106:THR:O	1:A:109:ARG:HB3	2.18	0.42
1:A:99:VAL:O	1:A:99:VAL:HG12	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:67:VAL:HA	1:C:76:VAL:O	2.18	0.42
1:A:137:GLY:HA2	1:A:238:ASP:O	2.20	0.42
1:B:115:TRP:CG	1:B:258:MET:HB3	2.54	0.42
1:B:155:VAL:O	1:B:158:LEU:HB3	2.19	0.42
1:C:70:SER:O	1:C:71:ARG:C	2.57	0.42
1:C:99:VAL:HG12	1:C:230:ALA:C	2.39	0.42
1:A:177:CYS:HG	1:A:184:CYS:CB	2.33	0.42
1:C:223:ILE:O	1:C:223:ILE:HG22	2.20	0.42
1:B:146:ASP:OD2	1:C:208:LYS:HE3	2.19	0.42
1:B:217:VAL:HG12	1:B:217:VAL:O	2.20	0.42
1:C:145:TYR:CD2	1:C:230:ALA:HB2	2.55	0.42
1:A:115:TRP:HA	1:A:261:PRO:HA	2.02	0.41
1:C:142:GLY:O	1:C:232:LEU:HD12	2.21	0.41
1:B:128:ILE:HG13	1:B:179:ILE:CD1	2.49	0.41
1:B:178:PHE:O	1:B:253:THR:HG21	2.19	0.41
1:B:243:THR:CG2	1:B:244:ALA:N	2.83	0.41
1:C:86:ILE:CG2	1:C:87:GLY:N	2.83	0.41
1:B:81:GLU:OE2	1:B:107:TRP:HB3	2.19	0.41
1:B:91:SER:O	1:B:93:VAL:HG23	2.20	0.41
1:A:237:LEU:CD2	1:A:237:LEU:N	2.83	0.41
1:B:146:ASP:C	1:B:148:ALA:N	2.74	0.41
1:A:113:ALA:HA	1:A:207:TYR:CE2	2.56	0.41
1:A:117:LYS:NZ	1:C:200:LEU:HA	2.35	0.41
1:A:88:VAL:HG22	1:A:89:THR:N	2.35	0.41
1:B:182:THR:OG1	1:B:183:ARG:N	2.53	0.41
1:B:139:ILE:O	1:B:166:SER:HA	2.21	0.41
1:C:141:MET:O	1:C:164:TYR:HA	2.21	0.41
1:A:207:TYR:O	1:A:208:LYS:HD2	2.21	0.41
1:B:259:ILE:HG13	1:B:260:GLU:HG2	2.02	0.41
1:C:122:SER:O	1:C:254:TYR:HA	2.21	0.41
1:A:202:LYS:NZ	1:B:202:LYS:HZ1	2.18	0.40
1:B:180:ASN:OD1	1:B:182:THR:HG22	2.21	0.40
1:B:124:ARG:O	1:B:124:ARG:HG2	2.21	0.40
1:B:202:LYS:HZ2	1:C:202:LYS:HE2	1.85	0.40
1:A:243:THR:O	1:A:243:THR:HG23	2.21	0.40
1:B:196:ASP:N	1:B:196:ASP:OD1	2.55	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	194/268 (72%)	152 (78%)	26 (13%)	16 (8%)	1	17
1	B	195/268 (73%)	162 (83%)	20 (10%)	13 (7%)	1	22
1	C	223/268 (83%)	189 (85%)	29 (13%)	5 (2%)	8	46
All	All	612/804 (76%)	503 (82%)	75 (12%)	34 (6%)	2	26

All (34) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	75	THR
1	A	244	ALA
1	B	74	ILE
1	B	92	ILE
1	B	238	ASP
1	A	112	ALA
1	A	136	ALA
1	A	169	VAL
1	A	181	GLY
1	A	201	GLY
1	A	246	ALA
1	B	112	ALA
1	B	130	SER
1	B	147	MET
1	B	161	LEU
1	B	169	VAL
1	C	46	GLN
1	C	211	ALA
1	B	201	GLY
1	B	211	ALA
1	A	161	LEU
1	A	211	ALA
1	A	222	ASN
1	A	261	PRO

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Mol	Chain	Res	Type
1	B	185	SER
1	C	71	ARG
1	A	229	PRO
1	B	261	PRO
1	B	159	SER
1	A	152	PRO
1	A	74	ILE
1	A	129	PRO
1	C	104	VAL
1	C	101	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	146/219 (67%)	133 (91%)	13 (9%)	11	44
1	B	145/219 (66%)	131 (90%)	14 (10%)	9	39
1	C	168/219 (77%)	145 (86%)	23 (14%)	4	26
All	All	459/657 (70%)	409 (89%)	50 (11%)	7	35

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

Mol	Chain	Res	Type
1	A	77	LEU
1	A	114	ASN
1	A	135	THR
1	A	158	LEU
1	A	168	GLN
1	A	189	THR
1	A	193	THR
1	A	194	THR
1	A	196	ASP
1	A	199	LYS
1	A	200	LEU
1	A	207	TYR

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Mol	Chain	Res	Type
1	A	256	ILE
1	B	100	MET
1	B	103	THR
1	B	114	ASN
1	B	150	THR
1	B	173	SER
1	B	189	THR
1	B	196	ASP
1	B	197	VAL
1	B	207	TYR
1	B	228	VAL
1	B	236	LEU
1	B	249	ARG
1	B	253	THR
1	B	260	GLU
1	C	50	SER
1	C	54	SER
1	C	59	MET
1	C	61	ARG
1	C	68	SER
1	C	77	LEU
1	C	80	SER
1	C	103	THR
1	C	109	ARG
1	C	121	LEU
1	C	147	MET
1	C	162	ARG
1	C	165	VAL
1	C	168	GLN
1	C	189	THR
1	C	195	LEU
1	C	199	LYS
1	C	200	LEU
1	C	207	TYR
1	C	208	LYS
1	C	210	SER
1	C	260	GLU
1	C	264	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	HIS
1	A	114	ASN
1	A	160	ASN
1	A	168	GLN
1	B	114	ASN
1	B	160	ASN
1	C	114	ASN
1	C	168	GLN
1	C	222	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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](#)

Of 3 ligands modelled in this entry, 3 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 [i](#)

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	196/268 (73%)	-0.47	0 100 100	9, 37, 83, 177	0
1	B	197/268 (73%)	-0.40	0 100 100	6, 39, 112, 172	0
1	C	225/268 (83%)	-0.40	1 (0%) 92 88	11, 42, 98, 145	0
All	All	618/804 (76%)	-0.42	1 (0%) 94 93	6, 39, 99, 177	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	44	PRO	4.0

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(Å ²)	Q<0.9
2	CA	A	269	1/1	0.96	0.06	-2.59	39,39,39,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	C	270	1/1	0.98	0.06	-2.98	39,39,39,39	0
2	CA	A	271	1/1	0.98	0.06	-3.10	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.