



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 12:46 pm BST

PDB ID : 2WLP  
Title : Sesbania mosaic virus capsid protein dimer mutant (rCP-DEL-N65-W170K)  
Authors : Anju, P.; Subashchandrabose, C.; Satheshkumar, P.S.; Savithri, H.S.; Murthy, M.R.N.  
Deposited on : 2009-06-24  
Resolution : 2.65 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

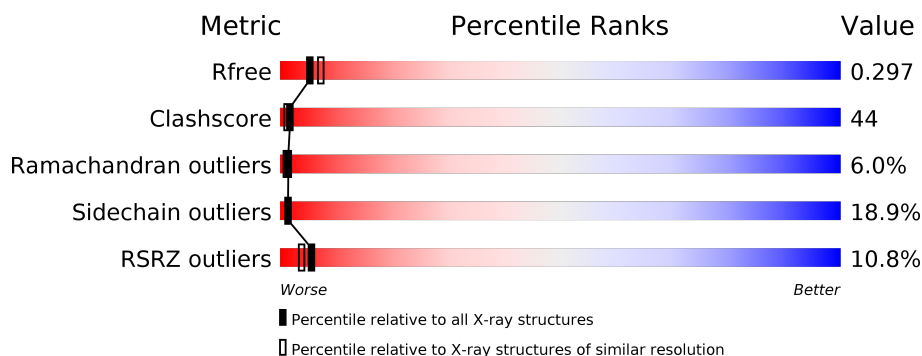
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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))
$R_{free}$	130704	1332 (2.68-2.64)
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	203	
1	B	203	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2821 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	187	Total	C	N	O	S	0	0	0
			1380	880	224	270	6			
1	B	191	Total	C	N	O	S	0	1	0
			1394	890	222	274	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	170	LYS	TRP	engineered mutation	UNP Q9EB06
B	170	LYS	TRP	engineered mutation	UNP Q9EB06

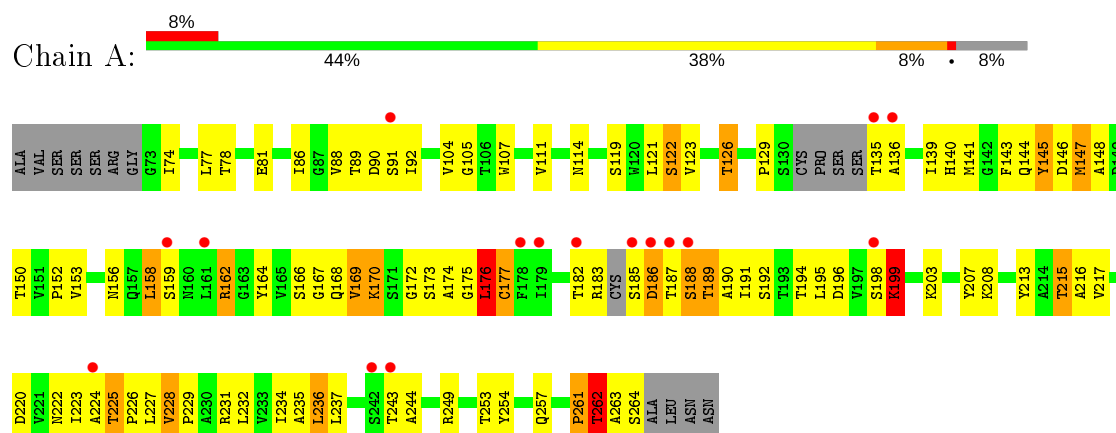
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	26	Total	O	0	0
			26	26		
2	B	21	Total	O	0	0
			21	21		

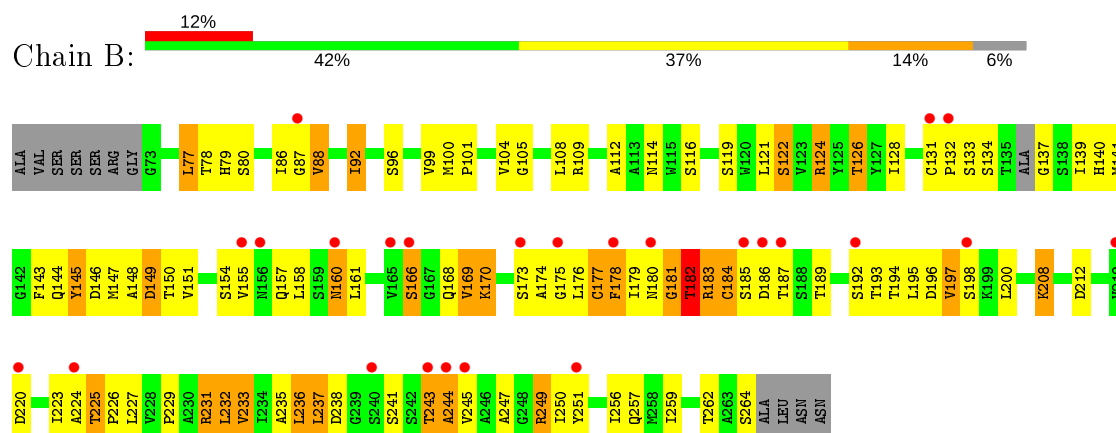
### 3 Residue-property plots [i](#)

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



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	35.88Å 41.95Å 68.65Å 81.07° 75.00° 84.25°	Depositor
Resolution (Å)	65.65 – 2.65 37.34 – 2.65	Depositor EDS
% Data completeness (in resolution range)	93.2 (65.65-2.65) 93.3 (37.34-2.65)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.31 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.4.0069	Depositor
R, $R_{free}$	0.221 , 0.294 0.224 , 0.297	Depositor DCC
$R_{free}$ test set	493 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.1	Xtriage
Anisotropy	0.111	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 72.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	2821	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.49% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

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	0/1407	0.73	0/1924
1	B	0.61	0/1426	0.74	0/1955
All	All	0.62	0/2833	0.73	0/3879

There are no bond length outliers.

There are no bond angle outliers.

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	1380	0	1363	122	0
1	B	1394	0	1356	118	0
2	A	26	0	0	1	0
2	B	21	0	0	1	0
All	All	2821	0	2719	240	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 44.

All (240) 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:183:ARG:O	1:A:185:SER:CB	1.77	1.33

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:GLN:C	1:B:145:TYR:HD1	1.40	1.22
1:A:162:ARG:CD	1:A:162:ARG:H	1.43	1.22
1:A:196:ASP:CB	1:A:199:LYS:HG3	1.71	1.19
1:B:179:ILE:HD11	1:B:251:TYR:HB2	1.16	1.09
1:A:198:SER:C	1:A:199:LYS:HG2	1.75	1.06
1:A:162:ARG:H	1:A:162:ARG:HD3	1.16	1.05
1:B:144:GLN:C	1:B:145:TYR:CD1	2.28	1.05
1:B:179:ILE:HD11	1:B:251:TYR:CB	1.89	1.02
1:B:179:ILE:CD1	1:B:251:TYR:CB	2.38	1.01
1:A:162:ARG:CD	1:A:162:ARG:N	2.25	0.98
1:B:179:ILE:CD1	1:B:251:TYR:HB2	1.95	0.95
1:A:162:ARG:H	1:A:162:ARG:HD2	1.29	0.94
1:A:162:ARG:HH11	1:A:162:ARG:CG	1.82	0.93
1:B:131:CYS:HB2	1:B:132:PRO:HD2	1.51	0.92
1:B:86:ILE:HG23	1:B:236:LEU:HD11	1.50	0.91
1:A:162:ARG:HH11	1:A:162:ARG:HG2	1.35	0.90
1:A:237:LEU:H	1:A:237:LEU:HD23	1.36	0.90
1:A:215:THR:HG22	1:A:216:ALA:N	1.86	0.90
1:B:151:VAL:HG23	2:B:2011:HOH:O	1.70	0.90
1:A:129:PRO:HD3	1:A:172:GLY:HA3	1.56	0.88
1:A:169:VAL:HG13	1:A:169:VAL:O	1.71	0.87
1:B:145:TYR:N	1:B:145:TYR:HD1	1.71	0.87
1:A:121:LEU:HD21	1:A:257:GLN:HE22	1.37	0.87
1:A:162:ARG:N	1:A:162:ARG:HD3	1.88	0.84
1:B:224:ALA:C	1:B:226:PRO:HD2	1.98	0.83
1:A:121:LEU:HD21	1:A:257:GLN:NE2	1.95	0.81
1:B:179:ILE:HD13	1:B:251:TYR:CB	2.10	0.81
1:A:135:THR:CG2	1:A:136:ALA:N	2.42	0.81
1:A:198:SER:O	1:A:199:LYS:HD3	1.79	0.81
1:A:162:ARG:NH1	1:A:162:ARG:HG2	1.93	0.81
1:B:208:LYS:HE2	1:B:212[A]:ASP:OD2	1.82	0.80
1:B:179:ILE:HD13	1:B:251:TYR:CG	2.17	0.79
1:A:225:THR:N	1:A:226:PRO:HD2	1.97	0.79
1:A:198:SER:O	1:A:199:LYS:CG	2.30	0.79
1:A:168:GLN:HG2	1:A:170:LYS:HE3	1.66	0.78
1:A:198:SER:C	1:A:199:LYS:CG	2.49	0.78
1:B:179:ILE:CD1	1:B:251:TYR:CG	2.66	0.78
1:B:92:ILE:HB	1:B:237:LEU:HB3	1.66	0.77
1:B:77:LEU:C	1:B:77:LEU:HD23	2.05	0.77
1:B:225:THR:N	1:B:226:PRO:HD2	2.00	0.77
1:A:77:LEU:HD23	1:A:78:THR:N	2.01	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:ASP:HB3	1:A:223:ILE:HG13	1.67	0.75
1:B:86:ILE:CG2	1:B:236:LEU:HD11	2.17	0.75
1:A:198:SER:O	1:A:199:LYS:CD	2.35	0.74
1:B:139:ILE:HG13	1:B:235:ALA:O	1.87	0.74
1:B:144:GLN:O	1:B:145:TYR:CD1	2.41	0.73
1:A:135:THR:HG23	1:A:136:ALA:H	1.53	0.73
1:A:237:LEU:N	1:A:237:LEU:HD23	2.02	0.73
1:B:184:CYS:O	1:B:185:SER:CB	2.37	0.72
1:B:128:ILE:HD12	1:B:176:LEU:HG	1.72	0.72
1:A:183:ARG:O	1:A:185:SER:CA	2.38	0.71
1:B:131:CYS:HB2	1:B:132:PRO:CD	2.20	0.70
1:A:135:THR:N	1:A:170:LYS:HB3	2.07	0.70
1:B:200:LEU:HD22	1:B:229:PRO:O	1.91	0.69
1:B:144:GLN:OE1	1:B:231:ARG:NH1	2.25	0.69
1:B:225:THR:N	1:B:226:PRO:CD	2.53	0.69
1:A:121:LEU:CD2	1:A:257:GLN:NE2	2.55	0.69
1:A:169:VAL:CG1	1:A:169:VAL:O	2.41	0.69
1:A:198:SER:O	1:A:199:LYS:HG2	1.92	0.69
1:B:86:ILE:HG23	1:B:236:LEU:CD1	2.22	0.68
1:A:225:THR:N	1:A:226:PRO:CD	2.56	0.68
1:B:126:THR:HB	1:B:192:SER:HB2	1.74	0.68
1:B:132:PRO:C	1:B:134:SER:H	1.95	0.68
1:A:121:LEU:CG	1:A:257:GLN:NE2	2.56	0.68
1:B:145:TYR:CD1	1:B:145:TYR:N	2.45	0.68
1:A:148:ALA:HB2	1:A:222:ASN:ND2	2.09	0.67
1:A:121:LEU:HG	1:A:257:GLN:NE2	2.09	0.67
1:B:146:ASP:O	1:B:148:ALA:N	2.27	0.67
1:B:100:MET:O	1:B:104:VAL:HG13	1.94	0.67
1:B:104:VAL:HG23	1:B:105:GLY:N	2.08	0.67
1:B:180:ASN:O	1:B:182:THR:N	2.27	0.67
1:B:140:HIS:HD1	1:B:166:SER:HG	1.40	0.67
1:A:183:ARG:O	1:A:185:SER:N	2.27	0.67
1:B:146:ASP:C	1:B:148:ALA:H	1.97	0.67
1:B:176:LEU:O	1:B:177:CYS:O	2.13	0.67
1:A:135:THR:HG22	1:A:136:ALA:N	2.10	0.66
1:B:220:ASP:HB3	1:B:223:ILE:HD12	1.75	0.66
1:B:96:SER:HB2	1:B:231:ARG:HD3	1.78	0.66
1:A:135:THR:N	1:A:170:LYS:CB	2.59	0.66
1:B:223:ILE:CG2	1:B:223:ILE:O	2.44	0.65
1:A:175:GLY:O	1:A:177:CYS:N	2.29	0.65
1:B:146:ASP:C	1:B:148:ALA:N	2.49	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:SER:HB3	1:B:157:GLN:HG3	1.79	0.65
1:A:176:LEU:O	1:A:177:CYS:C	2.35	0.65
1:B:124:ARG:HB2	1:B:194:THR:HG22	1.79	0.64
1:A:77:LEU:C	1:A:77:LEU:HD23	2.18	0.64
1:B:132:PRO:O	1:B:134:SER:N	2.31	0.64
1:A:121:LEU:HG	1:A:257:GLN:HE21	1.62	0.64
1:A:187:THR:O	1:A:189:THR:N	2.31	0.64
1:A:196:ASP:O	1:A:198:SER:N	2.29	0.63
1:B:224:ALA:C	1:B:226:PRO:CD	2.67	0.63
1:B:169:VAL:O	1:B:169:VAL:HG13	1.98	0.62
1:A:232:LEU:HD23	1:A:234:ILE:HD11	1.80	0.62
1:B:245:VAL:O	1:B:245:VAL:HG12	1.99	0.62
1:A:144:GLN:OE1	1:A:231:ARG:NH1	2.28	0.62
1:B:249:ARG:HD2	1:B:251:TYR:OH	2.01	0.61
1:B:146:ASP:O	1:B:149:ASP:N	2.34	0.60
1:A:158:LEU:O	1:A:158:LEU:HD23	2.01	0.60
1:B:180:ASN:O	1:B:181:GLY:C	2.40	0.60
1:A:140:HIS:HB2	1:A:235:ALA:HB3	1.83	0.60
1:B:126:THR:HB	1:B:192:SER:CB	2.31	0.60
1:A:177:CYS:HG	1:A:185:SER:N	2.00	0.59
1:B:77:LEU:HD22	1:B:256:ILE:HG22	1.85	0.59
1:B:197:VAL:HG12	1:B:198:SER:N	2.18	0.58
1:A:158:LEU:CD2	1:A:158:LEU:O	2.52	0.58
1:A:86:ILE:HG23	1:A:236:LEU:HD11	1.85	0.58
1:A:263:ALA:O	1:A:264:SER:C	2.41	0.58
1:B:104:VAL:CG2	1:B:105:GLY:N	2.66	0.58
1:B:179:ILE:CD1	1:B:251:TYR:HB3	2.33	0.57
1:B:237:LEU:HD23	1:B:237:LEU:H	1.68	0.57
1:B:88:VAL:HG13	1:B:236:LEU:HD23	1.87	0.57
1:A:126:THR:HB	1:A:192:SER:HB2	1.86	0.57
1:A:208:LYS:HE2	1:A:208:LYS:HA	1.87	0.57
1:B:179:ILE:HD13	1:B:251:TYR:HB3	1.85	0.57
1:A:173:SER:C	1:A:175:GLY:H	2.08	0.56
1:B:177:CYS:O	1:B:180:ASN:N	2.37	0.56
1:A:173:SER:O	1:A:175:GLY:N	2.37	0.56
1:A:223:ILE:O	1:A:226:PRO:HG2	2.05	0.56
1:A:167:GLY:HA3	1:A:191:ILE:HD12	1.88	0.56
1:A:162:ARG:HD2	1:A:162:ARG:N	2.04	0.56
1:A:146:ASP:C	1:A:148:ALA:N	2.59	0.55
1:A:121:LEU:HD11	1:A:257:GLN:CD	2.27	0.55
1:B:87:GLY:O	1:B:236:LEU:HD22	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:ASP:O	1:A:148:ALA:N	2.40	0.55
1:A:78:THR:HA	1:A:254:TYR:O	2.06	0.55
1:B:177:CYS:O	1:B:178:PHE:C	2.45	0.55
1:B:157:GLN:O	1:B:160:ASN:HB2	2.07	0.54
1:B:143:PHE:HE1	1:B:194:THR:O	1.90	0.54
1:A:186:ASP:N	1:A:186:ASP:OD1	2.36	0.54
1:A:111:VAL:O	1:A:114:ASN:ND2	2.41	0.54
1:B:224:ALA:O	1:B:226:PRO:N	2.41	0.54
1:B:132:PRO:C	1:B:134:SER:N	2.60	0.54
1:B:223:ILE:HG23	1:B:223:ILE:O	2.05	0.54
1:A:146:ASP:C	1:A:148:ALA:H	2.11	0.54
1:A:135:THR:N	1:A:170:LYS:HD2	2.22	0.53
1:A:148:ALA:HB2	1:A:222:ASN:HD21	1.72	0.53
1:A:129:PRO:CD	1:A:172:GLY:HA3	2.34	0.53
1:A:237:LEU:CD2	1:A:237:LEU:N	2.70	0.53
1:A:135:THR:CG2	1:A:136:ALA:H	2.10	0.53
1:B:77:LEU:C	1:B:77:LEU:CD2	2.76	0.53
1:B:224:ALA:O	1:B:227:LEU:N	2.35	0.53
1:A:224:ALA:C	1:A:226:PRO:HD2	2.29	0.52
1:B:96:SER:HA	1:B:232:LEU:O	2.10	0.52
1:A:196:ASP:CB	1:A:199:LYS:CG	2.66	0.52
1:A:119:SER:HB2	1:A:203:LYS:O	2.10	0.51
1:A:175:GLY:O	1:A:176:LEU:C	2.48	0.51
1:A:135:THR:O	1:A:170:LYS:HD3	2.11	0.51
1:B:173:SER:C	1:B:175:GLY:H	2.14	0.51
1:A:156:ASN:O	1:A:159:SER:N	2.42	0.50
1:A:187:THR:O	1:A:190:ALA:N	2.35	0.50
1:A:77:LEU:C	1:A:77:LEU:CD2	2.80	0.50
1:A:123:VAL:O	1:A:194:THR:HG23	2.10	0.50
1:A:253:THR:HG22	2:A:2025:HOH:O	2.11	0.50
1:B:169:VAL:O	1:B:169:VAL:CG1	2.58	0.50
1:B:243:THR:O	1:B:244:ALA:C	2.49	0.50
1:B:124:ARG:O	1:B:124:ARG:HG2	2.12	0.50
1:B:220:ASP:CB	1:B:223:ILE:HD12	2.41	0.50
1:B:237:LEU:CD2	1:B:237:LEU:H	2.25	0.50
1:B:139:ILE:HG12	1:B:140:HIS:N	2.27	0.50
1:B:77:LEU:HD23	1:B:78:THR:N	2.27	0.50
1:A:92:ILE:HG23	1:A:92:ILE:O	2.12	0.49
1:A:126:THR:HB	1:A:192:SER:CB	2.42	0.49
1:A:145:TYR:N	1:A:145:TYR:CD1	2.80	0.48
1:A:232:LEU:CD2	1:A:234:ILE:HD11	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:THR:O	1:B:187:THR:CG2	2.61	0.48
1:A:145:TYR:N	1:A:145:TYR:HD1	2.11	0.48
1:B:197:VAL:HA	1:B:200:LEU:HG	1.95	0.48
1:B:169:VAL:HG13	1:B:247:ALA:HB1	1.96	0.47
1:A:194:THR:HG22	1:A:195:LEU:O	2.14	0.47
1:B:161:LEU:CD1	1:B:233:VAL:HG13	2.44	0.47
1:B:148:ALA:O	1:B:149:ASP:C	2.52	0.47
1:B:177:CYS:O	1:B:179:ILE:N	2.47	0.47
1:B:187:THR:O	1:B:187:THR:HG22	2.14	0.47
1:B:137:GLY:HA2	1:B:238:ASP:O	2.15	0.47
1:B:176:LEU:O	1:B:177:CYS:C	2.52	0.47
1:A:207:TYR:O	1:A:208:LYS:HE2	2.15	0.46
1:A:144:GLN:C	1:A:145:TYR:HD1	2.19	0.46
1:A:213:TYR:CD1	1:A:228:VAL:HG11	2.50	0.46
1:B:161:LEU:HD23	1:B:161:LEU:HA	1.71	0.46
1:B:237:LEU:CD2	1:B:237:LEU:N	2.79	0.46
1:A:147:MET:HG2	1:A:225:THR:HA	1.98	0.45
1:A:121:LEU:O	1:A:122:SER:HB2	2.16	0.45
1:A:261:PRO:C	1:A:262:THR:HG22	2.36	0.45
1:B:173:SER:C	1:B:175:GLY:N	2.70	0.45
1:A:173:SER:C	1:A:175:GLY:N	2.69	0.45
1:B:224:ALA:O	1:B:225:THR:C	2.54	0.45
1:A:104:VAL:HG23	1:A:105:GLY:N	2.32	0.45
1:B:121:LEU:HD23	1:B:121:LEU:HA	1.71	0.45
1:A:187:THR:O	1:A:188:SER:C	2.55	0.45
1:A:90:ASP:OD1	1:A:91:SER:N	2.50	0.45
1:A:88:VAL:CG1	1:A:236:LEU:HD22	2.47	0.45
1:A:121:LEU:O	1:A:122:SER:CB	2.65	0.44
1:B:196:ASP:O	1:B:198:SER:N	2.45	0.44
1:B:88:VAL:HG13	1:B:236:LEU:CD2	2.46	0.44
1:B:224:ALA:C	1:B:226:PRO:N	2.70	0.44
1:A:162:ARG:HH11	1:A:162:ARG:HG3	1.75	0.44
1:A:175:GLY:C	1:A:177:CYS:N	2.69	0.44
1:A:81:GLU:OE1	1:A:107:TRP:N	2.33	0.44
1:A:77:LEU:HD11	1:A:107:TRP:CZ3	2.53	0.43
1:A:215:THR:HG22	1:A:216:ALA:H	1.79	0.43
1:A:216:ALA:O	1:A:217:VAL:C	2.56	0.43
1:B:121:LEU:HG	1:B:257:GLN:HG3	1.99	0.43
1:B:208:LYS:CE	1:B:212[A]:ASP:OD2	2.62	0.43
1:B:92:ILE:HB	1:B:237:LEU:CB	2.42	0.43
1:A:88:VAL:HG13	1:A:236:LEU:HD22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:MET:O	1:A:164:TYR:HA	2.19	0.43
1:B:140:HIS:CE1	1:B:166:SER:HG	2.35	0.43
1:B:86:ILE:HD11	1:B:250:ILE:HG13	1.99	0.43
1:A:176:LEU:O	1:A:177:CYS:O	2.37	0.43
1:B:168:GLN:O	1:B:170:LYS:N	2.52	0.43
1:B:173:SER:O	1:B:175:GLY:N	2.52	0.43
1:B:99:VAL:O	1:B:99:VAL:HG12	2.19	0.43
1:A:89:THR:HG22	1:A:244:ALA:HA	2.01	0.42
1:B:141:MET:CE	1:B:193:THR:HG22	2.49	0.42
1:A:172:GLY:O	1:A:173:SER:C	2.57	0.42
1:B:148:ALA:O	1:B:149:ASP:O	2.38	0.42
1:B:179:ILE:HD13	1:B:251:TYR:CD2	2.52	0.42
1:B:237:LEU:O	1:B:238:ASP:HB2	2.20	0.42
1:B:131:CYS:CB	1:B:132:PRO:CD	2.92	0.42
1:A:168:GLN:HB2	1:A:168:GLN:HE21	1.65	0.42
1:B:182:THR:O	1:B:183:ARG:O	2.38	0.42
1:A:143:PHE:HB2	1:A:145:TYR:HE1	1.85	0.42
1:A:220:ASP:O	1:A:223:ILE:HB	2.19	0.42
1:A:228:VAL:HA	1:A:229:PRO:HD2	1.83	0.42
1:B:184:CYS:C	1:B:186:ASP:H	2.21	0.42
1:B:108:LEU:O	1:B:109:ARG:C	2.59	0.41
1:A:156:ASN:O	1:A:159:SER:HB3	2.20	0.41
1:A:147:MET:HB2	1:A:225:THR:HG22	2.02	0.41
1:A:81:GLU:OE1	1:A:107:TRP:CB	2.68	0.41
1:B:77:LEU:HD22	1:B:256:ILE:CG2	2.50	0.41
1:B:79:HIS:CG	1:B:80:SER:H	2.39	0.41
1:B:92:ILE:HD12	1:B:237:LEU:HD13	2.02	0.41
1:A:146:ASP:OD1	1:A:225:THR:HG21	2.21	0.41
1:A:81:GLU:OE1	1:A:107:TRP:HB3	2.21	0.41
1:B:79:HIS:CG	1:B:80:SER:N	2.89	0.40
1:B:112:ALA:C	1:B:114:ASN:H	2.23	0.40
1:B:143:PHE:CE1	1:B:195:LEU:HA	2.56	0.40
1:B:119:SER:HB3	1:B:257:GLN:HE21	1.86	0.40
1:A:254:TYR:C	1:A:254:TYR:CD1	2.94	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	181/203 (89%)	150 (83%)	21 (12%)	10 (6%)	2	1
1	B	188/203 (93%)	160 (85%)	16 (8%)	12 (6%)	1	1
All	All	369/406 (91%)	310 (84%)	37 (10%)	22 (6%)	1	1

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	176	LEU
1	A	199	LYS
1	B	177	CYS
1	B	182	THR
1	B	183	ARG
1	A	122	SER
1	A	188	SER
1	B	133	SER
1	B	147	MET
1	B	149	ASP
1	B	178	PHE
1	B	181	GLY
1	B	244	ALA
1	A	147	MET
1	A	152	PRO
1	A	174	ALA
1	A	262	THR
1	B	169	VAL
1	B	174	ALA
1	A	177	CYS
1	B	122	SER
1	A	261	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	149/168 (89%)	125 (84%)	24 (16%)	2	2
1	B	149/168 (89%)	117 (78%)	32 (22%)	1	1
All	All	298/336 (89%)	242 (81%)	56 (19%)	1	1

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	ILE
1	A	126	THR
1	A	139	ILE
1	A	145	TYR
1	A	150	THR
1	A	153	VAL
1	A	158	LEU
1	A	162	ARG
1	A	166	SER
1	A	169	VAL
1	A	170	LYS
1	A	176	LEU
1	A	182	THR
1	A	186	ASP
1	A	189	THR
1	A	199	LYS
1	A	215	THR
1	A	225	THR
1	A	227	LEU
1	A	228	VAL
1	A	236	LEU
1	A	243	THR
1	A	249	ARG
1	A	262	THR
1	B	77	LEU
1	B	88	VAL
1	B	92	ILE

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Mol	Chain	Res	Type
1	B	101	PRO
1	B	116	SER
1	B	122	SER
1	B	124	ARG
1	B	126	THR
1	B	145	TYR
1	B	150	THR
1	B	155	VAL
1	B	158	LEU
1	B	160	ASN
1	B	166	SER
1	B	170	LYS
1	B	182	THR
1	B	184	CYS
1	B	189	THR
1	B	197	VAL
1	B	208	LYS
1	B	225	THR
1	B	231	ARG
1	B	232	LEU
1	B	233	VAL
1	B	236	LEU
1	B	237	LEU
1	B	241	SER
1	B	243	THR
1	B	249	ARG
1	B	259	ILE
1	B	262	THR
1	B	264	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	156	ASN
1	A	160	ASN
1	A	168	GLN
1	A	222	ASN
1	A	257	GLN
1	B	257	GLN

### 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](#)

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.



## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	187/203 (92%)	0.77	16 (8%) 10 8	50, 57, 78, 87	0
1	B	191/203 (94%)	0.91	25 (13%) 3 2	49, 61, 78, 84	0
All	All	378/406 (93%)	0.84	41 (10%) 5 4	49, 59, 78, 87	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	185	SER	6.2
1	B	87	GLY	4.9
1	B	186	ASP	4.4
1	A	198	SER	4.2
1	A	242	SER	4.2
1	A	136	ALA	4.1
1	B	175	GLY	4.1
1	B	173	SER	3.9
1	B	160	ASN	3.6
1	B	185	SER	3.6
1	B	245	VAL	3.4
1	B	243	THR	3.4
1	A	135	THR	3.3
1	B	166	SER	3.3
1	B	244	ALA	3.3
1	A	187	THR	3.1
1	B	155	VAL	3.1
1	A	243	THR	3.0
1	B	156	ASN	3.0
1	B	251	TYR	3.0
1	A	159	SER	3.0
1	A	161	LEU	2.9
1	A	182	THR	2.9
1	A	188	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	240	SER	2.7
1	A	178	PHE	2.6
1	B	187	THR	2.5
1	A	224	ALA	2.5
1	B	220	ASP	2.5
1	B	192	SER	2.4
1	B	198	SER	2.4
1	B	180	ASN	2.4
1	B	132	PRO	2.3
1	A	91	SER	2.3
1	B	224	ALA	2.1
1	A	186	ASP	2.1
1	B	131	CYS	2.1
1	B	219	VAL	2.1
1	A	179	ILE	2.1
1	B	178	PHE	2.1
1	B	165	VAL	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.