



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 22, 2026 – 11:42 AM EDT

PDB ID : 2VQ0 / pdb_00002vq0
Title : Capsid structure of Sesbania mosaic virus coat protein deletion mutant rCP(delta 48 to 59)
Authors : Anju, P.; Subashchandrabose, C.; Satheshkumar, P.S.; Savithri, H.S.; Murthy, M.R.N.
Deposited on : 2008-03-10
Resolution : 3.60 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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

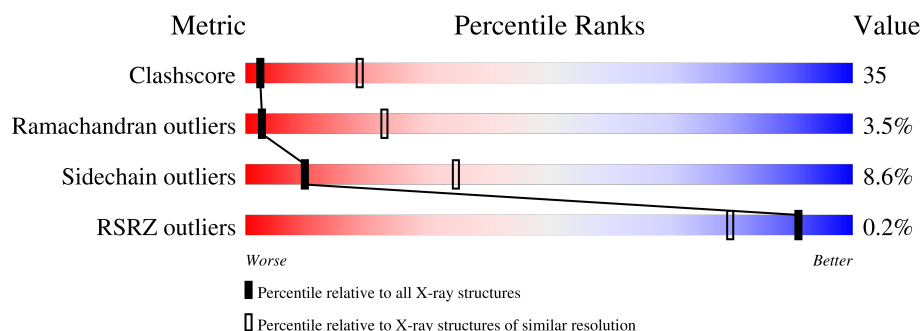
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	256	
1	B	256	
1	C	256	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4454 atoms, of which 0 are hydrogens and 0 are deuteriums.

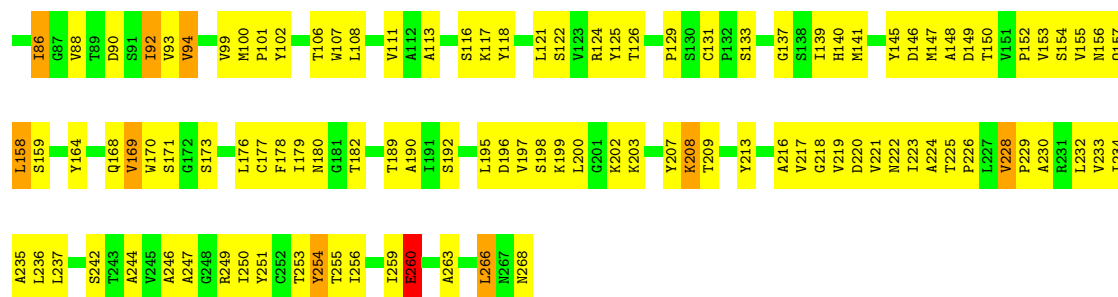
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- 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
			1447	921	236	282	8			
1	B	197	Total	C	N	O	S	0	0	0
			1447	915	237	287	8			
1	C	213	Total	C	N	O	S	0	0	0
			1557	986	257	306	8			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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



4 Data and refinement statistics

Property	Value	Source
Space group	R 3	Depositor
Cell constants a, b, c, α , β , γ	288.79Å 288.79Å 288.79Å 61.70° 61.70° 61.70°	Depositor
Resolution (Å)	20.00 – 3.60 20.00 – 3.60	Depositor EDS
% Data completeness (in resolution range)	83.0 (20.00-3.60) 83.1 (20.00-3.60)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.257 , 0.258 0.245 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtrriage
Anisotropy	(Not available)	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 159.8	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.14	EDS
Total number of atoms	4454	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtrriage'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	0.63	0/1479	1.08	9/2028 (0.4%)
1	B	0.60	0/1479	1.08	12/2025 (0.6%)
1	C	0.58	0/1590	1.09	14/2181 (0.6%)
All	All	0.60	0/4548	1.08	35/6234 (0.6%)

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	203	LYS	N-CA-C	-7.90	103.59	113.55
1	B	86	ILE	N-CA-C	7.54	119.33	108.48
1	B	128	ILE	CA-C-N	7.39	128.39	120.11
1	B	128	ILE	C-N-CA	7.39	128.39	120.11
1	A	86	ILE	N-CA-C	6.95	118.61	108.53
1	C	86	ILE	N-CA-C	6.93	118.46	108.48
1	A	191	ILE	CB-CA-C	-6.50	105.94	113.22
1	B	121	LEU	N-CA-C	-6.26	105.77	113.41
1	C	228	VAL	CA-C-N	6.24	125.46	118.97
1	C	228	VAL	C-N-CA	6.24	125.46	118.97
1	C	76	VAL	N-CA-C	6.17	116.75	107.80
1	B	260	GLU	CA-C-N	6.17	126.18	119.89
1	B	260	GLU	C-N-CA	6.17	126.18	119.89
1	C	56	PRO	N-CA-CB	6.13	109.74	103.00
1	B	198	SER	N-CA-C	6.01	117.83	111.28
1	A	121	LEU	N-CA-C	-6.00	106.57	114.31
1	C	266	LEU	N-CA-C	-6.00	106.13	113.50
1	B	179	ILE	N-CA-C	5.84	116.02	110.53
1	B	167	GLY	N-CA-C	5.81	118.98	110.38
1	B	219	VAL	N-CA-C	-5.47	105.51	110.82
1	C	260	GLU	CA-C-N	5.37	125.31	119.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	260	GLU	C-N-CA	5.37	125.31	119.78
1	C	82	LEU	N-CA-C	-5.31	101.79	109.59
1	C	74	ILE	N-CA-C	5.30	115.92	109.30
1	B	93	VAL	N-CA-C	5.29	115.71	107.99
1	A	148	ALA	N-CA-C	-5.25	106.38	112.89
1	C	190	ALA	N-CA-C	5.23	117.81	110.23
1	C	203	LYS	N-CA-C	-5.23	105.18	112.45
1	A	254	TYR	N-CA-C	5.17	117.72	109.81
1	A	225	THR	CA-C-N	-5.16	114.65	120.89
1	A	225	THR	C-N-CA	-5.16	114.65	120.89
1	A	180	ASN	N-CA-C	-5.12	106.88	113.02
1	A	219	VAL	N-CA-C	-5.11	106.09	110.74
1	C	94	VAL	N-CA-C	5.09	115.30	108.17
1	C	254	TYR	N-CA-C	5.00	116.89	109.69

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	1447	0	1415	97	0
1	B	1447	0	1402	121	0
1	C	1557	0	1515	112	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
All	All	4454	0	4332	309	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 (309) 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:B:202:LYS:NZ	1:C:202:LYS:HZ1	1.42	1.18
1:B:202:LYS:HZ1	1:C:202:LYS:NZ	1.44	1.14
1:B:202:LYS:NZ	1:C:202:LYS:NZ	2.08	0.95
1:C:176:LEU:HG	1:C:180:ASN:HD21	1.30	0.94
1:A:100:MET:HE1	1:A:208:LYS:HB2	1.50	0.93
1:B:117:LYS:HE2	1:B:260:GLU:OE2	1.71	0.88
1:A:131:CYS:HB2	1:A:135:THR:HG21	1.56	0.84
1:B:86:ILE:HG23	1:B:236:LEU:HD11	1.59	0.83
1:C:99:VAL:HG12	1:C:230:ALA:O	1.78	0.83
1:B:249:ARG:HG2	1:B:249:ARG:HH11	1.43	0.82
1:C:88:VAL:HG21	1:C:169:VAL:HG21	1.65	0.79
1:C:86:ILE:HG23	1:C:236:LEU:HD11	1.65	0.79
1:C:88:VAL:HG23	1:C:247:ALA:HB2	1.66	0.78
1:B:126:THR:HG23	1:B:192:SER:HB2	1.66	0.77
1:C:92:ILE:HD13	1:C:93:VAL:N	1.99	0.77
1:A:146:ASP:O	1:A:147:MET:HB2	1.85	0.77
1:B:146:ASP:OD2	1:B:148:ALA:HB3	1.85	0.76
1:B:209:THR:HG22	1:B:268:ASN:O	1.90	0.72
1:C:176:LEU:HG	1:C:180:ASN:ND2	2.04	0.71
1:C:126:THR:HG23	1:C:192:SER:HB2	1.73	0.71
1:A:76:VAL:HG23	1:A:256:ILE:O	1.92	0.70
1:B:141:MET:HE3	1:B:250:ILE:HD13	1.74	0.70
1:B:145:TYR:HB3	1:C:117:LYS:HE3	1.74	0.69
1:C:92:ILE:HD13	1:C:93:VAL:H	1.55	0.69
1:A:131:CYS:CB	1:A:135:THR:HG21	2.23	0.69
1:B:86:ILE:HD11	1:B:250:ILE:HG13	1.75	0.69
1:B:89:THR:O	1:B:239:GLY:HA3	1.94	0.68
1:A:150:THR:O	1:B:266:LEU:HD22	1.93	0.68
1:C:139:ILE:HA	1:C:236:LEU:HD23	1.75	0.68
1:B:169:VAL:HG13	1:B:247:ALA:HB1	1.75	0.68
1:A:126:THR:HG23	1:A:192:SER:HB2	1.76	0.68
1:B:128:ILE:HD12	1:B:179:ILE:HD12	1.76	0.68
1:C:213:TYR:CE2	1:C:217:VAL:HG21	2.29	0.68
1:B:140:HIS:HB2	1:B:235:ALA:HB3	1.75	0.68
1:C:86:ILE:HG23	1:C:236:LEU:CD1	2.22	0.67
1:C:86:ILE:HG22	1:C:247:ALA:HB3	1.76	0.67
1:B:111:VAL:HG23	1:B:112:ALA:N	2.08	0.67
1:C:137:GLY:C	1:C:169:VAL:HG23	2.20	0.66
1:B:86:ILE:HD11	1:B:250:ILE:CD1	2.26	0.66
1:A:132:PRO:O	1:A:135:THR:HG22	1.96	0.65
1:C:99:VAL:HG12	1:C:230:ALA:C	2.22	0.64
1:A:131:CYS:HB2	1:A:132:PRO:HD2	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:SER:O	1:B:92:ILE:C	2.40	0.64
1:B:111:VAL:HG23	1:B:112:ALA:H	1.61	0.64
1:A:160:ASN:N	1:A:160:ASN:HD22	1.95	0.64
1:A:162:ARG:HG2	1:A:163:GLY:N	2.13	0.64
1:A:237:LEU:N	1:A:237:LEU:HD23	2.13	0.63
1:B:202:LYS:HE2	1:C:202:LYS:HE3	1.81	0.62
1:A:213:TYR:CE1	1:A:217:VAL:HG21	2.35	0.61
1:B:77:LEU:HD11	1:B:107:TRP:CH2	2.36	0.61
1:B:116:SER:HB2	1:B:260:GLU:O	2.00	0.61
1:C:228:VAL:HG23	1:C:229:PRO:HD2	1.81	0.61
1:B:259:ILE:C	1:B:260:GLU:HG2	2.26	0.60
1:A:131:CYS:HB2	1:A:132:PRO:CD	2.31	0.60
1:C:90:ASP:OD1	1:C:242:SER:HA	2.02	0.60
1:A:86:ILE:HG23	1:A:236:LEU:CD1	2.32	0.59
1:A:79:HIS:CD2	1:A:80:SER:H	2.20	0.59
1:B:177:CYS:HA	1:B:180:ASN:HD21	1.67	0.59
1:B:80:SER:HB2	1:B:253:THR:HG23	1.83	0.59
1:B:86:ILE:HD11	1:B:250:ILE:CG1	2.32	0.59
1:B:249:ARG:HG2	1:B:249:ARG:NH1	2.16	0.59
1:A:100:MET:HE1	1:A:208:LYS:CB	2.29	0.58
1:B:86:ILE:CD1	1:B:250:ILE:HG13	2.32	0.58
1:C:125:TYR:CG	1:C:232:LEU:HD22	2.39	0.58
1:B:237:LEU:N	1:B:237:LEU:HD23	2.19	0.57
1:A:112:ALA:HA	1:A:258:MET:HE2	1.85	0.57
1:B:86:ILE:HG23	1:B:236:LEU:CD1	2.34	0.57
1:C:154:SER:OG	1:C:157:GLN:HG3	2.04	0.57
1:B:139:ILE:HA	1:B:236:LEU:HD23	1.85	0.57
1:A:115:TRP:HA	1:A:261:PRO:HA	1.87	0.57
1:B:137:GLY:C	1:B:169:VAL:HG23	2.30	0.57
1:B:100:MET:O	1:B:104:VAL:HG23	2.04	0.57
1:B:180:ASN:N	1:B:180:ASN:HD22	2.01	0.57
1:A:92:ILE:HA	1:A:237:LEU:HB3	1.87	0.57
1:C:118:TYR:HB2	1:C:256:ILE:HD11	1.86	0.57
1:B:134:SER:O	1:B:135:THR:C	2.48	0.56
1:A:125:TYR:CG	1:A:232:LEU:HD22	2.40	0.56
1:B:91:SER:O	1:B:93:VAL:HG23	2.06	0.56
1:B:139:ILE:HG12	1:B:236:LEU:HD21	1.87	0.56
1:B:158:LEU:HD12	1:B:161:LEU:HD12	1.87	0.56
1:A:266:LEU:CD2	1:C:150:THR:HB	2.36	0.56
1:B:86:ILE:HD11	1:B:250:ILE:HD11	1.88	0.56
1:A:80:SER:OG	1:A:253:THR:HG23	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:GLY:C	1:A:169:VAL:HG13	2.31	0.56
1:B:132:PRO:HD2	1:B:135:THR:HG21	1.88	0.56
1:A:144:GLN:HG3	1:A:161:LEU:HD13	1.88	0.55
1:A:86:ILE:HG23	1:A:236:LEU:HD11	1.87	0.55
1:B:98:LEU:CD1	1:B:100:MET:HE3	2.36	0.55
1:B:207:TYR:CD1	1:B:207:TYR:C	2.85	0.55
1:B:207:TYR:HB3	1:B:267:ASN:HD21	1.72	0.55
1:C:180:ASN:CB	1:C:182:THR:HG23	2.37	0.55
1:B:101:PRO:HA	1:B:108:LEU:HD12	1.88	0.55
1:B:132:PRO:HD2	1:B:135:THR:CG2	2.37	0.55
1:A:219:VAL:HG11	1:C:219:VAL:HG12	1.88	0.55
1:B:77:LEU:HD11	1:B:107:TRP:HH2	1.71	0.55
1:A:177:CYS:SG	1:A:178:PHE:N	2.79	0.55
1:C:125:TYR:CD2	1:C:232:LEU:HD22	2.43	0.54
1:A:142:GLY:O	1:A:232:LEU:HD12	2.07	0.54
1:A:169:VAL:HB	1:A:247:ALA:HB1	1.89	0.54
1:C:213:TYR:CZ	1:C:217:VAL:HG21	2.43	0.54
1:B:141:MET:CE	1:B:250:ILE:HD13	2.37	0.54
1:C:154:SER:C	1:C:156:ASN:N	2.61	0.54
1:C:208:LYS:HE2	1:C:268:ASN:C	2.33	0.54
1:B:141:MET:HE3	1:B:250:ILE:CD1	2.37	0.53
1:A:111:VAL:HG22	1:A:258:MET:HE1	1.90	0.53
1:C:102:TYR:CD2	1:C:209:THR:HA	2.42	0.53
1:A:241:SER:OG	1:A:243:THR:HG22	2.08	0.53
1:B:213:TYR:CE2	1:B:217:VAL:HG21	2.44	0.53
1:B:108:LEU:HB2	1:B:254:TYR:OH	2.07	0.53
1:A:131:CYS:SG	1:A:135:THR:HG21	2.49	0.52
1:A:116:SER:HB2	1:A:260:GLU:O	2.08	0.52
1:B:77:LEU:HD23	1:B:77:LEU:C	2.34	0.52
1:A:147:MET:HE1	1:A:231:ARG:HB3	1.90	0.52
1:A:243:THR:HG23	1:A:243:THR:O	2.09	0.52
1:B:98:LEU:HD12	1:B:100:MET:HE3	1.91	0.52
1:C:154:SER:C	1:C:156:ASN:H	2.17	0.52
1:A:85:GLU:HG2	1:A:249:ARG:HG2	1.92	0.51
1:A:88:VAL:HG21	1:A:137:GLY:HA3	1.92	0.51
1:A:146:ASP:O	1:A:147:MET:CB	2.57	0.51
1:B:115:TRP:CG	1:B:258:MET:HB3	2.45	0.51
1:B:131:CYS:HB2	1:B:132:PRO:CD	2.39	0.51
1:C:86:ILE:HG22	1:C:247:ALA:CB	2.39	0.51
1:C:102:TYR:CD1	1:C:102:TYR:C	2.88	0.51
1:A:108:LEU:O	1:A:109:ARG:C	2.53	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:PRO:HD2	1:A:135:THR:CG2	2.41	0.51
1:B:77:LEU:HD23	1:B:77:LEU:O	2.11	0.51
1:B:180:ASN:CG	1:B:182:THR:HG23	2.35	0.51
1:B:219:VAL:HG12	1:C:219:VAL:HG11	1.93	0.51
1:B:196:ASP:CG	1:B:196:ASP:O	2.53	0.51
1:C:139:ILE:HG13	1:C:236:LEU:HD21	1.91	0.51
1:C:116:SER:HB2	1:C:260:GLU:O	2.10	0.51
1:A:217:VAL:HG13	1:A:221:VAL:HA	1.92	0.51
1:C:101:PRO:HD3	1:C:118:TYR:CE2	2.45	0.51
1:C:94:VAL:HG12	1:C:235:ALA:HB2	1.93	0.50
1:C:237:LEU:HD23	1:C:237:LEU:N	2.26	0.50
1:A:112:ALA:HA	1:A:258:MET:CE	2.41	0.50
1:A:159:SER:HA	1:A:164:TYR:CD1	2.46	0.50
1:A:219:VAL:HG12	1:B:219:VAL:HG11	1.92	0.50
1:A:243:THR:O	1:A:244:ALA:O	2.29	0.50
1:B:124:ARG:HB2	1:B:194:THR:HG22	1.92	0.50
1:A:98:LEU:N	1:A:98:LEU:HD23	2.27	0.50
1:B:169:VAL:CG1	1:B:247:ALA:HB1	2.39	0.50
1:C:208:LYS:HE2	1:C:268:ASN:O	2.12	0.50
1:C:237:LEU:HD23	1:C:237:LEU:H	1.77	0.50
1:C:197:VAL:O	1:C:200:LEU:HD13	2.11	0.50
1:C:124:ARG:HD3	1:C:178:PHE:CD1	2.46	0.49
1:A:209:THR:HG22	1:A:268:ASN:HA	1.94	0.49
1:C:220:ASP:OD2	1:C:222:ASN:HB2	2.12	0.49
1:C:208:LYS:HE2	1:C:268:ASN:OXT	2.13	0.49
1:B:101:PRO:HD3	1:B:118:TYR:CE2	2.48	0.49
1:B:139:ILE:CG1	1:B:236:LEU:HD21	2.42	0.49
1:B:202:LYS:CE	1:C:202:LYS:NZ	2.75	0.49
1:B:223:ILE:O	1:B:227:LEU:HD13	2.13	0.49
1:A:266:LEU:HD22	1:C:150:THR:O	2.12	0.49
1:B:202:LYS:HE2	1:C:202:LYS:CE	2.43	0.49
1:B:227:LEU:HD12	1:B:227:LEU:N	2.26	0.49
1:C:70:SER:O	1:C:71:ARG:C	2.55	0.49
1:B:78:THR:OG1	1:B:255:THR:HG23	2.12	0.48
1:C:131:CYS:SG	1:C:169:VAL:HG12	2.53	0.48
1:A:77:LEU:C	1:A:77:LEU:HD22	2.39	0.48
1:B:89:THR:OG1	1:B:90:ASP:N	2.46	0.48
1:B:226:PRO:C	1:B:227:LEU:HD12	2.38	0.48
1:C:197:VAL:C	1:C:200:LEU:HD13	2.38	0.48
1:C:129:PRO:HD3	1:C:171:SER:O	2.13	0.48
1:C:80:SER:OG	1:C:253:THR:HG23	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:180:ASN:HB3	1:C:182:THR:HG23	1.95	0.48
1:A:207:TYR:CD1	1:A:207:TYR:C	2.92	0.47
1:B:111:VAL:CG2	1:B:112:ALA:N	2.77	0.47
1:B:209:THR:HG23	1:B:212:ASP:CB	2.45	0.47
1:B:100:MET:HE2	1:B:228:VAL:CG1	2.44	0.47
1:B:207:TYR:C	1:B:207:TYR:HD1	2.23	0.47
1:C:263:ALA:HB3	1:C:266:LEU:HB2	1.96	0.47
1:C:259:ILE:HG13	1:C:260:GLU:HG2	1.97	0.47
1:A:117:LYS:HE2	1:A:260:GLU:OE1	2.15	0.47
1:B:146:ASP:C	1:B:148:ALA:N	2.72	0.47
1:C:122:SER:O	1:C:254:TYR:HA	2.15	0.47
1:A:101:PRO:HD3	1:A:118:TYR:CE2	2.51	0.46
1:A:117:LYS:NZ	1:C:200:LEU:HA	2.29	0.46
1:B:167:GLY:HA3	1:B:191:ILE:HD11	1.98	0.46
1:A:164:TYR:CG	1:A:165:VAL:N	2.83	0.46
1:B:180:ASN:H	1:B:180:ASN:ND2	2.14	0.46
1:B:256:ILE:HG12	1:B:257:GLN:N	2.31	0.46
1:C:234:ILE:HD13	1:C:250:ILE:HD12	1.97	0.46
1:B:197:VAL:C	1:B:200:LEU:HD13	2.40	0.46
1:B:226:PRO:HG2	1:B:227:LEU:CD1	2.45	0.46
1:C:153:VAL:HG12	1:C:153:VAL:O	2.15	0.46
1:C:146:ASP:HB3	1:C:149:ASP:OD2	2.16	0.46
1:A:223:ILE:O	1:A:227:LEU:CD2	2.63	0.46
1:A:266:LEU:HD23	1:C:150:THR:HB	1.97	0.46
1:B:82:LEU:HD12	1:B:83:SER:N	2.30	0.46
1:B:202:LYS:NZ	1:C:202:LYS:HZ2	2.06	0.46
1:B:146:ASP:C	1:B:148:ALA:H	2.23	0.46
1:B:180:ASN:HD22	1:B:180:ASN:H	1.62	0.46
1:C:141:MET:HA	1:C:233:VAL:O	2.16	0.46
1:B:249:ARG:NH1	1:B:249:ARG:CG	2.74	0.46
1:B:128:ILE:HA	1:B:129:PRO:HD2	1.79	0.46
1:B:113:ALA:HA	1:B:207:TYR:CE2	2.50	0.46
1:A:176:LEU:C	1:A:176:LEU:HD13	2.42	0.45
1:B:108:LEU:HD23	1:B:254:TYR:CE1	2.51	0.45
1:A:131:CYS:SG	1:A:135:THR:CG2	3.04	0.45
1:A:227:LEU:HD11	1:C:222:ASN:OD1	2.16	0.45
1:C:108:LEU:HD23	1:C:108:LEU:O	2.15	0.45
1:C:223:ILE:HG22	1:C:223:ILE:O	2.15	0.45
1:C:146:ASP:O	1:C:148:ALA:N	2.49	0.45
1:A:151:VAL:HG23	1:A:152:PRO:HD2	1.99	0.45
1:B:209:THR:HG22	1:B:268:ASN:C	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:107:TRP:CZ3	1:C:111:VAL:HG21	2.52	0.45
1:C:156:ASN:O	1:C:159:SER:N	2.45	0.45
1:A:260:GLU:HB3	1:C:145:TYR:CE1	2.51	0.45
1:B:217:VAL:HG22	1:B:224:ALA:HB2	1.98	0.45
1:C:118:TYR:CB	1:C:256:ILE:HD11	2.47	0.45
1:C:180:ASN:HB2	1:C:182:THR:HG23	1.98	0.45
1:B:100:MET:HE2	1:B:228:VAL:HG11	1.99	0.45
1:C:133:SER:HA	1:C:170:TRP:CG	2.51	0.45
1:A:147:MET:HE1	1:A:231:ARG:CB	2.46	0.45
1:A:175:GLY:O	1:A:176:LEU:C	2.58	0.45
1:B:128:ILE:HD12	1:B:179:ILE:CD1	2.44	0.45
1:A:162:ARG:HG3	1:A:162:ARG:HH11	1.82	0.45
1:C:108:LEU:HD12	1:C:254:TYR:CE2	2.52	0.45
1:B:101:PRO:HD3	1:B:118:TYR:CZ	2.52	0.45
1:C:100:MET:HE2	1:C:100:MET:HB3	1.93	0.45
1:A:140:HIS:CE1	1:A:166:SER:HB2	2.53	0.44
1:C:140:HIS:CE1	1:C:164:TYR:HH	2.36	0.44
1:C:146:ASP:C	1:C:148:ALA:N	2.75	0.44
1:C:92:ILE:HD11	1:C:235:ALA:HB1	1.99	0.44
1:A:266:LEU:HD22	1:C:150:THR:HB	2.00	0.44
1:B:158:LEU:HD22	1:B:235:ALA:HB2	2.00	0.44
1:B:97:GLU:OE2	1:B:97:GLU:HA	2.18	0.44
1:B:118:TYR:HB2	1:B:256:ILE:HD11	1.99	0.44
1:B:221:VAL:HG23	1:B:222:ASN:N	2.31	0.44
1:C:82:LEU:HD21	1:C:249:ARG:HB2	1.99	0.44
1:A:146:ASP:HB3	1:A:149:ASP:CG	2.43	0.44
1:C:66:ALA:O	1:C:77:LEU:HA	2.17	0.44
1:A:209:THR:HG23	1:A:212:ASP:CB	2.48	0.43
1:C:140:HIS:HB2	1:C:235:ALA:HB3	2.00	0.43
1:A:143:PHE:O	1:A:161:LEU:HB3	2.18	0.43
1:C:88:VAL:HG21	1:C:169:VAL:CG2	2.43	0.43
1:C:223:ILE:C	1:C:226:PRO:HD2	2.44	0.43
1:A:147:MET:HE1	1:A:231:ARG:CG	2.47	0.43
1:A:160:ASN:N	1:A:160:ASN:ND2	2.62	0.43
1:C:78:THR:HG23	1:C:255:THR:OG1	2.17	0.43
1:A:223:ILE:O	1:A:223:ILE:HG22	2.17	0.43
1:C:137:GLY:O	1:C:169:VAL:HG23	2.17	0.43
1:B:89:THR:O	1:B:239:GLY:CA	2.63	0.43
1:B:140:HIS:O	1:B:234:ILE:HA	2.18	0.43
1:B:142:GLY:O	1:B:232:LEU:HD23	2.19	0.43
1:B:86:ILE:HD12	1:B:127:TYR:CD1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:VAL:HG23	1:A:230:ALA:O	2.19	0.43
1:C:207:TYR:CD1	1:C:207:TYR:C	2.97	0.43
1:A:209:THR:HG22	1:A:268:ASN:O	2.18	0.43
1:B:111:VAL:CG2	1:B:112:ALA:H	2.29	0.43
1:A:168:GLN:H	1:A:168:GLN:HG3	1.56	0.42
1:C:99:VAL:HG13	1:C:229:PRO:HG2	2.01	0.42
1:C:222:ASN:C	1:C:224:ALA:H	2.27	0.42
1:A:86:ILE:HG23	1:A:236:LEU:HD13	2.00	0.42
1:A:209:THR:HG23	1:A:212:ASP:HB2	2.01	0.42
1:A:220:ASP:HB2	1:B:219:VAL:HG21	2.01	0.42
1:B:202:LYS:HZ1	1:C:202:LYS:HZ1	0.61	0.42
1:B:86:ILE:HB	1:B:248:GLY:O	2.19	0.42
1:C:102:TYR:C	1:C:102:TYR:HD1	2.28	0.42
1:C:113:ALA:HA	1:C:207:TYR:CE2	2.55	0.42
1:C:156:ASN:O	1:C:157:GLN:C	2.63	0.42
1:A:159:SER:HA	1:A:164:TYR:CG	2.54	0.42
1:B:200:LEU:N	1:B:200:LEU:HD12	2.34	0.42
1:C:141:MET:O	1:C:164:TYR:HA	2.20	0.42
1:C:224:ALA:C	1:C:226:PRO:HD2	2.45	0.42
1:A:82:LEU:HD12	1:A:83:SER:N	2.35	0.42
1:B:225:THR:N	1:B:226:PRO:HD2	2.35	0.42
1:C:237:LEU:N	1:C:237:LEU:CD2	2.83	0.42
1:A:100:MET:HE3	1:A:100:MET:HB2	1.73	0.42
1:B:232:LEU:C	1:B:232:LEU:CD2	2.93	0.42
1:C:198:SER:O	1:C:199:LYS:HG3	2.21	0.41
1:B:237:LEU:N	1:B:237:LEU:CD2	2.83	0.41
1:C:82:LEU:HA	1:C:251:TYR:CD2	2.55	0.41
1:C:216:ALA:C	1:C:218:GLY:N	2.78	0.41
1:A:77:LEU:C	1:A:77:LEU:CD2	2.94	0.41
1:A:101:PRO:HD3	1:A:118:TYR:CZ	2.55	0.41
1:A:145:TYR:CD2	1:A:145:TYR:N	2.88	0.41
1:B:146:ASP:O	1:B:148:ALA:N	2.53	0.41
1:C:121:LEU:HA	1:C:121:LEU:HD23	1.85	0.41
1:C:158:LEU:HD23	1:C:158:LEU:O	2.20	0.41
1:B:108:LEU:C	1:B:108:LEU:HD13	2.45	0.41
1:A:99:VAL:HG23	1:A:231:ARG:HA	2.01	0.41
1:A:237:LEU:HD23	1:A:237:LEU:H	1.85	0.41
1:B:122:SER:O	1:B:254:TYR:HA	2.21	0.41
1:C:92:ILE:HG12	1:C:237:LEU:HB3	2.03	0.41
1:A:237:LEU:N	1:A:237:LEU:CD2	2.83	0.41
1:B:229:PRO:HG2	1:B:230:ALA:H	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:148:ALA:HA	1:C:221:VAL:HG21	2.02	0.41
1:A:132:PRO:C	1:A:134:SER:N	2.75	0.41
1:B:104:VAL:HG21	1:B:120:TRP:CH2	2.56	0.41
1:C:179:ILE:HD13	1:C:251:TYR:CD1	2.56	0.41
1:A:80:SER:HB3	1:A:179:ILE:HG23	2.03	0.41
1:A:82:LEU:HD12	1:A:83:SER:H	1.86	0.41
1:A:135:THR:O	1:A:135:THR:HG23	2.20	0.41
1:A:140:HIS:ND1	1:A:166:SER:HB2	2.35	0.41
1:B:101:PRO:O	1:B:103:THR:N	2.54	0.41
1:B:169:VAL:CG1	1:B:169:VAL:O	2.66	0.41
1:B:227:LEU:N	1:B:227:LEU:CD1	2.84	0.41
1:A:74:ILE:HG22	1:A:75:THR:N	2.35	0.40
1:A:80:SER:HA	1:A:252:CYS:O	2.21	0.40
1:A:139:ILE:HA	1:A:236:LEU:HD23	2.04	0.40
1:B:209:THR:HG23	1:B:212:ASP:HB3	2.02	0.40
1:C:225:THR:N	1:C:226:PRO:HD2	2.37	0.40
1:C:131:CYS:HB3	1:C:246:ALA:O	2.22	0.40
1:C:168:GLN:O	1:C:169:VAL:C	2.65	0.40
1:A:97:GLU:O	1:A:231:ARG:HB2	2.22	0.40
1:B:175:GLY:O	1:B:176:LEU:C	2.64	0.40
1:C:217:VAL:HG13	1:C:221:VAL:HA	2.03	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	194/256 (76%)	157 (81%)	29 (15%)	8 (4%)	2	19
1	B	195/256 (76%)	163 (84%)	26 (13%)	6 (3%)	3	25
1	C	211/256 (82%)	179 (85%)	25 (12%)	7 (3%)	3	24
All	All	600/768 (78%)	499 (83%)	80 (13%)	21 (4%)	3	23

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	244	ALA
1	B	92	ILE
1	A	147	MET
1	B	135	THR
1	B	244	ALA
1	C	57	THR
1	C	147	MET
1	C	169	VAL
1	A	132	PRO
1	B	102	TYR
1	C	244	ALA
1	A	222	ASN
1	C	71	ARG
1	C	177	CYS
1	A	152	PRO
1	B	161	LEU
1	A	131	CYS
1	C	152	PRO
1	A	165	VAL
1	A	197	VAL
1	B	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	156/213 (73%)	141 (90%)	15 (10%)	8	32
1	B	156/213 (73%)	142 (91%)	14 (9%)	9	33
1	C	167/213 (78%)	155 (93%)	12 (7%)	13	40
All	All	479/639 (75%)	438 (91%)	41 (9%)	10	34

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	LEU
1	A	88	VAL
1	A	89	THR
1	A	98	LEU
1	A	111	VAL
1	A	132	PRO
1	A	150	THR
1	A	160	ASN
1	A	173	SER
1	A	177	CYS
1	A	189	THR
1	A	198	SER
1	A	209	THR
1	A	233	VAL
1	A	245	VAL
1	B	77	LEU
1	B	100	MET
1	B	116	SER
1	B	139	ILE
1	B	150	THR
1	B	155	VAL
1	B	173	SER
1	B	180	ASN
1	B	189	THR
1	B	207	TYR
1	B	209	THR
1	B	232	LEU
1	B	249	ARG
1	B	260	GLU
1	C	77	LEU
1	C	82	LEU
1	C	92	ILE
1	C	106	THR
1	C	155	VAL
1	C	158	LEU
1	C	173	SER
1	C	189	THR
1	C	195	LEU
1	C	196	ASP
1	C	208	LYS
1	C	260	GLU

Sometimes 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	160	ASN
1	A	257	GLN
1	B	79	HIS
1	B	114	ASN
1	B	168	GLN
1	B	180	ASN
1	C	114	ASN
1	C	180	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 oligosaccharides 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	196/256 (76%)	-0.41	1 (0%) 87 66	11, 18, 55, 96	0
1	B	197/256 (76%)	-0.40	0 100 100	11, 20, 91, 96	0
1	C	213/256 (83%)	-0.47	0 100 100	9, 19, 59, 87	0
All	All	606/768 (78%)	-0.43	1 (0%) 91 80	9, 19, 70, 96	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	248	GLY	2.1

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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	CA	C	1269	1/1	0.97	0.07	18,18,18,18	0
2	CA	B	1269	1/1	0.99	0.20	18,18,18,18	0
2	CA	A	1269	1/1	0.99	0.05	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.