



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:07 PM UTC

PDB ID : 2O3I / pdb_00002o3i
Title : X-ray Crystal Structure of Protein CV_3147 from *Chromobacterium violaceum*. Northeast Structural Genomics Consortium Target CvR68.
Authors : Vorobiev, S.M.; Chen, Y.; Seetharaman, J.; Cunningham, K.; Ma, L.C.; Janjua, H.; Xiao, R.; Acton, T.B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2006-12-01
Resolution : 2.30 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

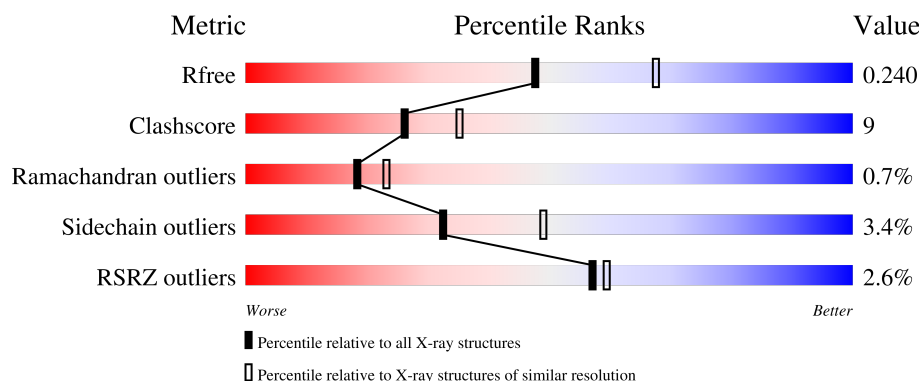
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	405	% 74% 17% • 7%
1	B	405	3% 71% 20% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5613 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 Hypothetical protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	Se	0	0	0
			2693	1717	457	505	4	10			
1	B	375	Total	C	N	O	S	Se	0	0	0
			2693	1716	456	507	4	10			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP Q7NTB2
A	28	MSE	MET	modified residue	UNP Q7NTB2
A	64	MSE	MET	modified residue	UNP Q7NTB2
A	68	MSE	MET	modified residue	UNP Q7NTB2
A	143	MSE	MET	modified residue	UNP Q7NTB2
A	172	MSE	MET	modified residue	UNP Q7NTB2
A	190	MSE	MET	modified residue	UNP Q7NTB2
A	208	MSE	MET	modified residue	UNP Q7NTB2
A	210	MSE	MET	modified residue	UNP Q7NTB2
A	211	MSE	MET	modified residue	UNP Q7NTB2
A	247	MSE	MET	modified residue	UNP Q7NTB2
A	271	PRO	SER	engineered mutation	UNP Q7NTB2
A	395	ARG	GLY	engineered mutation	UNP Q7NTB2
A	398	LEU	-	cloning artifact	UNP Q7NTB2
A	399	GLU	-	cloning artifact	UNP Q7NTB2
A	400	HIS	-	cloning artifact	UNP Q7NTB2
A	401	HIS	-	cloning artifact	UNP Q7NTB2
A	402	HIS	-	cloning artifact	UNP Q7NTB2
A	403	HIS	-	cloning artifact	UNP Q7NTB2
A	404	HIS	-	cloning artifact	UNP Q7NTB2
A	405	HIS	-	cloning artifact	UNP Q7NTB2
B	1	MSE	MET	modified residue	UNP Q7NTB2
B	28	MSE	MET	modified residue	UNP Q7NTB2
B	64	MSE	MET	modified residue	UNP Q7NTB2
B	68	MSE	MET	modified residue	UNP Q7NTB2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	143	MSE	MET	modified residue	UNP Q7NTB2
B	172	MSE	MET	modified residue	UNP Q7NTB2
B	190	MSE	MET	modified residue	UNP Q7NTB2
B	208	MSE	MET	modified residue	UNP Q7NTB2
B	210	MSE	MET	modified residue	UNP Q7NTB2
B	211	MSE	MET	modified residue	UNP Q7NTB2
B	247	MSE	MET	modified residue	UNP Q7NTB2
B	271	PRO	SER	engineered mutation	UNP Q7NTB2
B	395	ARG	GLY	engineered mutation	UNP Q7NTB2
B	398	LEU	-	cloning artifact	UNP Q7NTB2
B	399	GLU	-	cloning artifact	UNP Q7NTB2
B	400	HIS	-	cloning artifact	UNP Q7NTB2
B	401	HIS	-	cloning artifact	UNP Q7NTB2
B	402	HIS	-	cloning artifact	UNP Q7NTB2
B	403	HIS	-	cloning artifact	UNP Q7NTB2
B	404	HIS	-	cloning artifact	UNP Q7NTB2
B	405	HIS	-	cloning artifact	UNP Q7NTB2

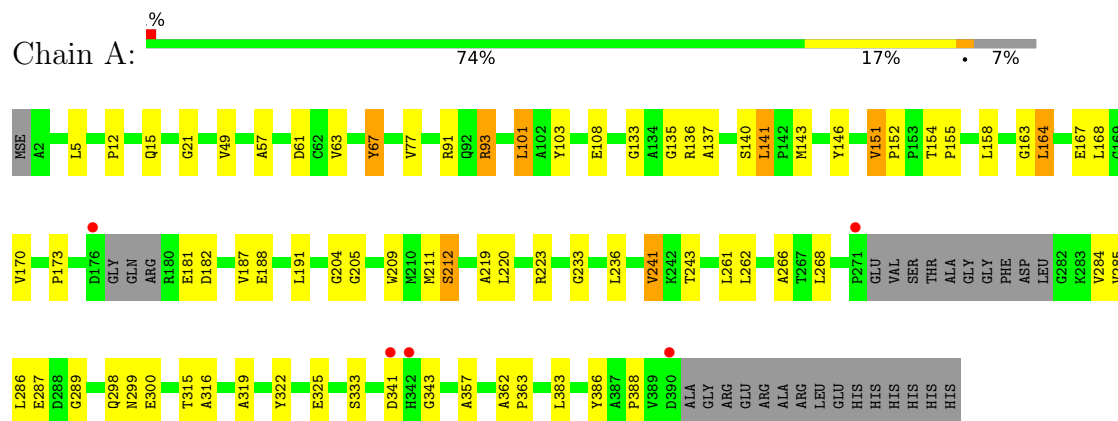
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	116	Total O 116 116	0	0
2	B	111	Total O 111 111	0	0

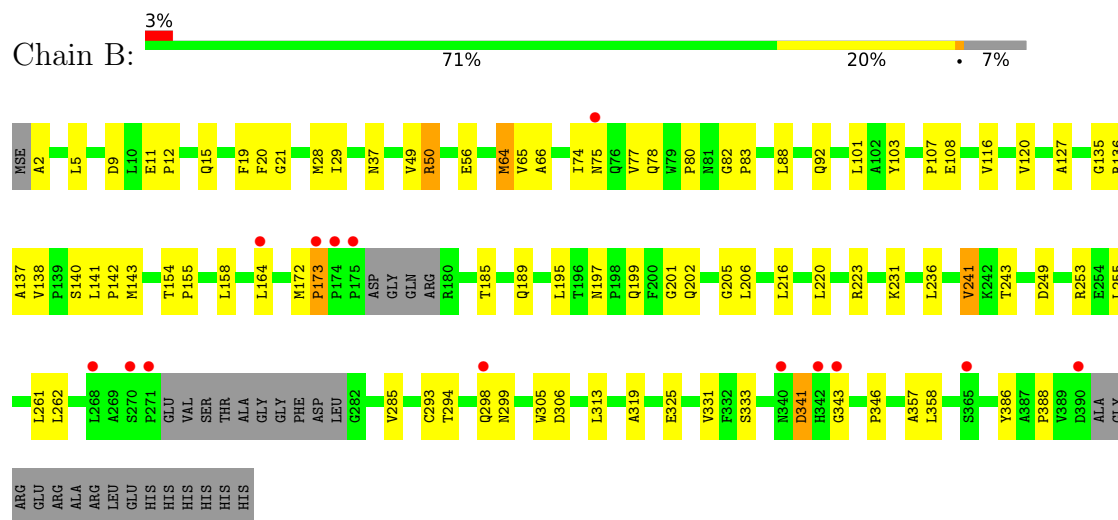
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Hypothetical protein



• Molecule 1: Hypothetical protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	92.34Å 125.61Å 65.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.92 – 2.30 29.92 – 2.30	Depositor EDS
% Data completeness (in resolution range)	87.9 (29.92-2.30) 91.2 (29.92-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 2.27Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.198 , 0.228 0.210 , 0.240	Depositor DCC
R_{free} test set	2494 reflections (3.83%)	wwPDB-VP
Wilson B-factor (Å ²)	28.7	Xtriage
Anisotropy	0.835	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5613	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.86 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2191e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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	0.41	0/2739	0.90	10/3719 (0.3%)
1	B	0.44	0/2739	0.94	6/3718 (0.2%)
All	All	0.42	0/5478	0.92	16/7437 (0.2%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	298	GLN	N-CA-C	-10.14	97.66	111.96
1	A	298	GLN	N-CA-C	-9.52	96.99	111.56
1	B	333	SER	N-CA-C	-7.34	99.95	110.59
1	A	333	SER	N-CA-C	-6.90	100.31	110.52
1	A	133	GLY	N-CA-C	6.71	123.63	115.31
1	A	343	GLY	N-CA-C	-6.66	105.88	115.27
1	A	223	ARG	N-CA-C	6.14	119.20	109.81
1	A	212	SER	N-CA-C	-6.06	102.24	110.36
1	A	135	GLY	N-CA-C	-6.03	107.00	115.32
1	A	299	ASN	N-CA-C	-5.93	98.17	110.80
1	A	141	LEU	N-CA-C	5.72	121.40	113.45
1	B	299	ASN	N-CA-C	-5.43	99.22	110.80
1	B	343	GLY	N-CA-C	-5.40	108.18	115.21
1	B	135	GLY	N-CA-C	-5.18	108.17	115.32
1	A	151	VAL	CB-CA-C	-5.14	105.33	110.89
1	B	223	ARG	N-CA-C	5.13	117.65	109.81

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	2693	0	2634	47	0
1	B	2693	0	2639	55	0
2	A	116	0	0	1	0
2	B	111	0	0	0	0
All	All	5613	0	5273	100	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 9.

All (100) 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:151:VAL:HG21	1:A:211:MSE:CE	1.69	1.23
1:A:151:VAL:CG2	1:A:211:MSE:HE2	1.85	1.07
1:A:151:VAL:HG21	1:A:211:MSE:HE2	0.94	0.93
1:B:37:ASN:HD21	1:B:78:GLN:HE21	1.24	0.85
1:B:64:MSE:HE1	1:B:66:ALA:HB2	1.62	0.80
1:B:138:VAL:HG11	1:B:143:MSE:HB2	1.66	0.77
1:B:64:MSE:HE3	1:B:107:PRO:CD	2.15	0.76
1:B:64:MSE:HE1	1:B:66:ALA:CB	2.17	0.74
1:B:107:PRO:HG2	1:B:108:GLU:HG2	1.71	0.73
1:B:64:MSE:HE3	1:B:107:PRO:HD3	1.70	0.73
1:A:93:ARG:HH11	1:A:93:ARG:HB2	1.57	0.69
1:B:249:ASP:HB3	1:B:253:ARG:NH1	2.07	0.69
1:B:64:MSE:CE	1:B:107:PRO:HD3	2.22	0.69
1:A:146:TYR:HD1	1:A:211:MSE:HE1	1.57	0.69
1:B:249:ASP:HB3	1:B:253:ARG:HH12	1.56	0.69
1:B:197:ASN:ND2	1:B:199:GLN:H	1.95	0.65
1:A:12:PRO:O	1:A:233:GLY:HA3	1.98	0.64
1:B:29:ILE:HD11	1:B:75:ASN:HB2	1.80	0.64
1:B:5:LEU:HB2	1:B:49:VAL:HB	1.80	0.62
1:B:236:LEU:HD23	1:B:241:VAL:HG13	1.82	0.62
1:A:236:LEU:HD23	1:A:241:VAL:HG13	1.81	0.62
1:B:138:VAL:CG1	1:B:143:MSE:HB2	2.28	0.62
1:A:268:LEU:HD23	1:A:286:LEU:HG	1.81	0.62
1:B:64:MSE:HE3	1:B:107:PRO:HB3	1.82	0.61
1:A:164:LEU:HD23	1:B:172:MSE:HG3	1.83	0.60
1:B:37:ASN:ND2	1:B:78:GLN:HE21	1.99	0.58
1:B:262:LEU:HD11	1:B:357:ALA:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:MSE:HE3	1:B:107:PRO:CB	2.34	0.58
1:B:197:ASN:HD22	1:B:199:GLN:H	1.52	0.58
1:B:50:ARG:HH11	1:B:50:ARG:HB2	1.70	0.57
1:A:315:THR:HG22	1:A:316:ALA:O	2.05	0.56
1:A:262:LEU:HD11	1:A:357:ALA:HB2	1.88	0.56
1:B:15:GLN:HE21	1:B:28:MSE:HE1	1.70	0.56
1:B:319:ALA:HB3	1:B:358:LEU:HB2	1.88	0.55
1:A:91:ARG:HB2	1:A:101:LEU:HD12	1.88	0.54
1:B:185:THR:O	1:B:189:GLN:HG3	2.08	0.53
1:B:37:ASN:HD21	1:B:78:GLN:NE2	2.01	0.51
1:B:154:THR:HA	1:B:155:PRO:C	2.35	0.51
1:A:266:ALA:HB1	1:A:287:GLU:O	2.10	0.51
1:A:67:TYR:OH	1:A:77:VAL:HG11	2.12	0.49
1:A:154:THR:HA	1:A:155:PRO:C	2.37	0.49
1:A:243:THR:HA	1:A:325:GLU:HG2	1.94	0.49
1:B:74:ILE:O	1:B:77:VAL:HG12	2.13	0.49
1:A:168:LEU:CD1	1:A:170:VAL:HG23	2.43	0.49
1:A:140:SER:HB2	1:A:188:GLU:OE1	2.12	0.49
1:A:5:LEU:HB2	1:A:49:VAL:HB	1.95	0.49
1:B:140:SER:HB3	1:B:143:MSE:HE3	1.94	0.49
1:B:243:THR:HA	1:B:325:GLU:HG2	1.94	0.48
1:A:141:LEU:HD12	1:A:188:GLU:HB2	1.96	0.47
1:A:386:TYR:CE1	1:A:388:PRO:HG3	2.49	0.47
1:A:181:GLU:HG3	1:A:182:ASP:H	1.78	0.47
1:A:211:MSE:HE1	1:A:219:ALA:HB1	1.95	0.47
1:A:140:SER:OG	1:A:143:MSE:HE3	2.14	0.47
1:B:21:GLY:O	1:B:137:ALA:HB3	2.14	0.47
1:A:57:ALA:HA	1:A:103:TYR:CZ	2.50	0.46
1:A:63:VAL:HG23	1:A:101:LEU:HD21	1.97	0.46
1:B:103:TYR:HD2	1:B:127:ALA:HB3	1.80	0.46
1:A:93:ARG:HB2	1:A:93:ARG:NH1	2.28	0.46
1:A:219:ALA:C	1:A:220:LEU:HD12	2.41	0.45
1:B:2:ALA:HB2	1:B:56:GLU:OE1	2.15	0.45
1:B:285:VAL:HG22	1:B:294:THR:HG23	1.99	0.45
1:A:158:LEU:HA	1:A:205:GLY:O	2.16	0.45
1:B:293:CYS:HA	1:B:305:TRP:O	2.17	0.45
1:B:306:ASP:HB2	1:B:313:LEU:HD11	1.99	0.45
1:B:64:MSE:HE3	1:B:107:PRO:CG	2.46	0.45
1:A:209:TRP:O	1:A:211:MSE:HG3	2.16	0.44
1:B:231:LYS:HD3	1:B:255:LEU:HD21	1.98	0.44
1:B:386:TYR:CE1	1:B:388:PRO:HG3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:PRO:HB2	1:B:83:PRO:HG2	1.99	0.44
1:B:82:GLY:C	1:B:205:GLY:HA3	2.43	0.44
1:B:140:SER:OG	1:B:142:PRO:HD2	2.18	0.43
1:A:204:GLY:HA2	2:A:515:HOH:O	2.18	0.43
1:B:64:MSE:HE2	1:B:65:VAL:N	2.34	0.43
1:A:181:GLU:HG3	1:A:182:ASP:N	2.34	0.43
1:B:9:ASP:C	1:B:12:PRO:HD2	2.43	0.43
1:A:383:LEU:HD12	1:A:383:LEU:N	2.34	0.43
1:B:140:SER:HB3	1:B:143:MSE:CE	2.49	0.42
1:B:158:LEU:HD22	1:B:195:LEU:HD21	2.02	0.42
1:B:20:PHE:O	1:B:136:ARG:HA	2.18	0.42
1:B:201:GLY:O	1:B:202:GLN:HB2	2.20	0.42
1:A:284:VAL:HG21	1:A:322:TYR:OH	2.20	0.42
1:A:21:GLY:O	1:A:137:ALA:HB3	2.20	0.42
1:A:93:ARG:HH22	1:A:167:GLU:CD	2.26	0.42
1:A:151:VAL:O	1:A:152:PRO:C	2.62	0.42
1:A:137:ALA:O	1:A:300:GLU:HG2	2.20	0.42
1:A:285:VAL:C	1:A:286:LEU:HD12	2.45	0.42
1:A:362:ALA:HB3	1:A:363:PRO:HD3	2.01	0.41
1:A:211:MSE:CE	1:A:219:ALA:HB1	2.50	0.41
1:A:61:ASP:OD2	1:A:212:SER:HA	2.20	0.41
1:B:64:MSE:CE	1:B:65:VAL:C	2.93	0.41
1:A:173:PRO:HD3	1:B:164:LEU:HD23	2.01	0.41
1:B:141:LEU:N	1:B:142:PRO:HD2	2.36	0.41
1:B:65:VAL:HA	1:B:206:LEU:O	2.21	0.41
1:B:341:ASP:HB2	1:B:346:PRO:HA	2.03	0.41
1:A:236:LEU:HD23	1:A:241:VAL:CG1	2.49	0.41
1:A:319:ALA:O	1:A:357:ALA:HA	2.21	0.41
1:B:19:PHE:HB2	1:B:331:VAL:HG11	2.04	0.40
1:B:116:VAL:O	1:B:120:VAL:HG23	2.21	0.40
1:A:108:GLU:HB2	1:A:136:ARG:O	2.22	0.40
1:A:187:VAL:O	1:A:191:LEU:HG	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	370/405 (91%)	354 (96%)	13 (4%)	3 (1%)	16	20
1	B	369/405 (91%)	350 (95%)	17 (5%)	2 (0%)	24	31
All	All	739/810 (91%)	704 (95%)	30 (4%)	5 (1%)	18	23

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	341	ASP
1	B	341	ASP
1	B	173	PRO
1	A	163	GLY
1	A	289	GLY

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	262/295 (89%)	255 (97%)	7 (3%)	39	58
1	B	264/295 (90%)	253 (96%)	11 (4%)	26	40
All	All	526/590 (89%)	508 (97%)	18 (3%)	32	49

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	67	TYR
1	A	93	ARG
1	A	101	LEU
1	A	164	LEU
1	A	241	VAL
1	A	261	LEU

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Mol	Chain	Res	Type
1	B	11	GLU
1	B	50	ARG
1	B	64	MSE
1	B	88	LEU
1	B	92	GLN
1	B	101	LEU
1	B	173	PRO
1	B	216	LEU
1	B	220	LEU
1	B	241	VAL
1	B	261	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	189	GLN
1	A	197	ASN
1	A	199	GLN
1	A	334	ASN
1	A	379	GLN
1	B	15	GLN
1	B	37	ASN
1	B	78	GLN
1	B	81	ASN
1	B	92	GLN
1	B	197	ASN
1	B	334	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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 [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	366/405 (90%)	0.09	5 (1%) 73 75	18, 32, 54, 66	0
1	B	365/405 (90%)	0.12	14 (3%) 44 46	17, 31, 56, 78	0
All	All	731/810 (90%)	0.10	19 (2%) 57 59	17, 32, 56, 78	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	175	PRO	4.9
1	A	176	ASP	4.8
1	B	173	PRO	3.4
1	A	390	ASP	3.3
1	B	390	ASP	3.1
1	A	341	ASP	3.1
1	B	271	PRO	2.9
1	B	340	ASN	2.7
1	B	174	PRO	2.7
1	B	270	SER	2.4
1	A	271	PRO	2.3
1	B	164	LEU	2.3
1	B	298	GLN	2.3
1	B	268	LEU	2.2
1	B	365	SER	2.2
1	B	343	GLY	2.2
1	B	342	HIS	2.1
1	A	342	HIS	2.1
1	B	75	ASN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.