



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:39 AM UTC

PDB ID : 5MDP / pdb_00005mdp
Title : Crystal structure of in vitro folded Chitoporin VhChip from *Vibrio harveyi* (crystal form II)
Authors : Zahn, M.; van den Berg, B.
Deposited on : 2016-11-13
Resolution : 3.08 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

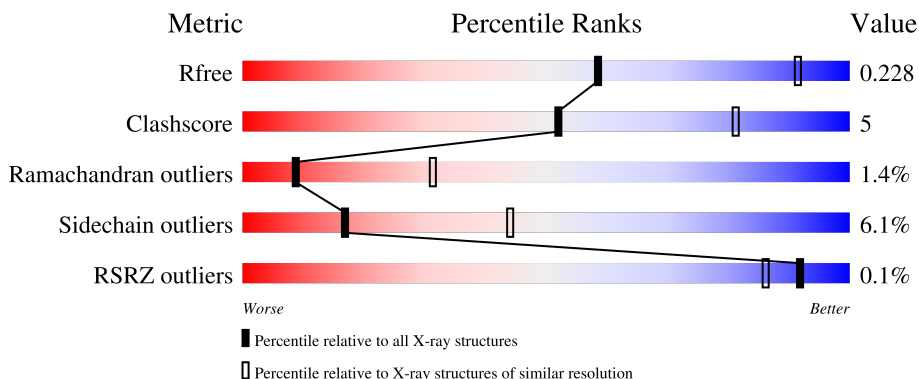
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.08 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	2010 (3.10-3.06)
Clashscore	190562	2102 (3.10-3.06)
Ramachandran outliers	187476	1982 (3.10-3.06)
Sidechain outliers	187428	1981 (3.10-3.06)
RSRZ outliers	180081	2010 (3.10-3.06)

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	352	
1	B	352	
1	C	352	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7995 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 Chitoporin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	Se	0	0	0
			2665	1680	438	538	9			
1	B	341	Total	C	N	O	Se	0	0	0
			2665	1680	438	538	9			
1	C	341	Total	C	N	O	Se	0	0	0
			2665	1680	438	538	9			

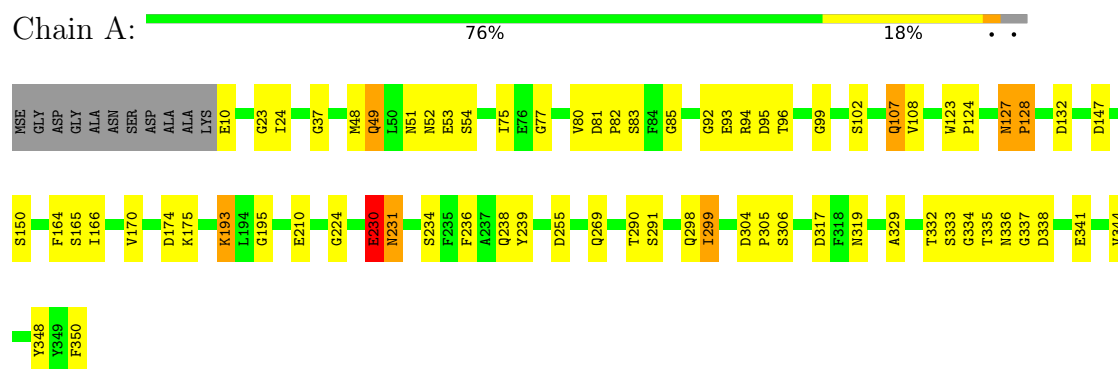
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	initiating methionine	UNP L0RVU0
A	0	GLY	-	expression tag	UNP L0RVU0
B	-1	MSE	-	initiating methionine	UNP L0RVU0
B	0	GLY	-	expression tag	UNP L0RVU0
C	-1	MSE	-	initiating methionine	UNP L0RVU0
C	0	GLY	-	expression tag	UNP L0RVU0

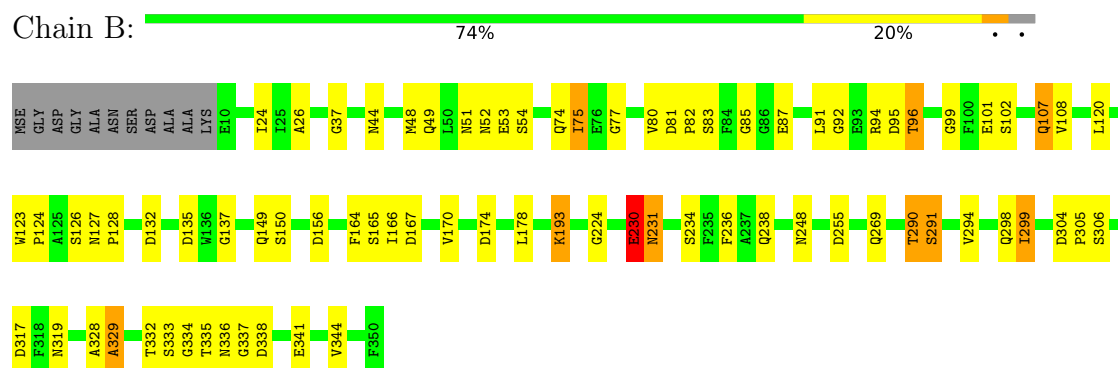
3 Residue-property plots

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.

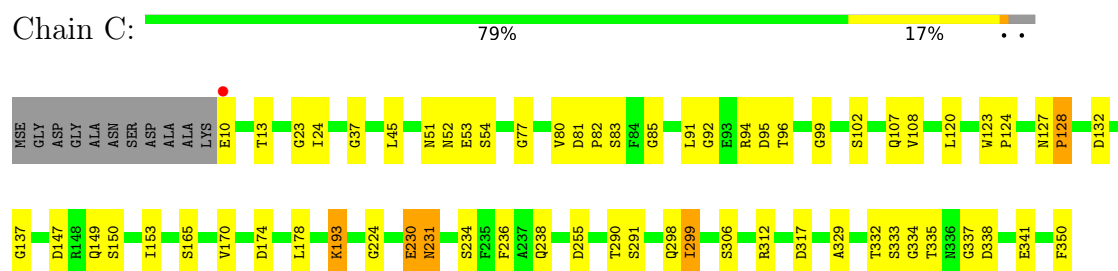
• Molecule 1: Chitoporin



• Molecule 1: Chitoporin



• Molecule 1: Chitoporin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	254.76Å 147.03Å 53.76Å 90.00° 95.51° 90.00°	Depositor
Resolution (Å)	127.20 – 3.08 127.20 – 3.08	Depositor EDS
% Data completeness (in resolution range)	95.5 (127.20-3.08) 95.6 (127.20-3.08)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.23 (at 3.05Å)	Xtriage
Refinement program	REFMAC 5.8.0124	Depositor
R, R_{free}	0.217 , 0.240 (Not available) , 0.228	Depositor DCC
R_{free} test set	1731 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	53.9	Xtriage
Anisotropy	1.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 48.7	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	0.93	EDS
Total number of atoms	7995	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% 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

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	1.29	8/2725 (0.3%)	1.28	12/3677 (0.3%)
1	B	1.33	12/2725 (0.4%)	1.32	17/3677 (0.5%)
1	C	1.31	8/2725 (0.3%)	1.31	12/3677 (0.3%)
All	All	1.31	28/8175 (0.3%)	1.30	41/11031 (0.4%)

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	A	0	2
1	B	0	1
1	C	0	2
All	All	0	5

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	174	ASP	CA-C	9.25	1.65	1.52
1	B	174	ASP	CA-C	8.89	1.64	1.52
1	B	85	GLY	C-O	8.44	1.34	1.24
1	C	85	GLY	C-O	7.59	1.33	1.24
1	A	174	ASP	CA-C	7.56	1.62	1.52
1	A	85	GLY	C-O	7.16	1.32	1.24
1	A	193	LYS	CA-C	-6.56	1.44	1.52
1	B	82	PRO	C-O	6.41	1.32	1.24
1	B	74	GLN	N-CA	-6.40	1.38	1.46
1	B	336	ASN	N-CA	6.32	1.54	1.46
1	B	193	LYS	CA-C	-6.19	1.44	1.52
1	C	193	LYS	CA-C	-6.14	1.44	1.52
1	A	82	PRO	C-O	5.84	1.31	1.24
1	B	149	GLN	CD-OE1	5.84	1.34	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	149	GLN	CD-OE1	5.74	1.34	1.23
1	C	82	PRO	C-O	5.67	1.31	1.24
1	A	348	TYR	CA-C	5.61	1.59	1.52
1	A	337	GLY	C-O	5.61	1.27	1.23
1	B	337	GLY	C-O	5.44	1.28	1.23
1	B	167	ASP	CA-C	5.42	1.58	1.52
1	C	45	LEU	C-O	5.34	1.30	1.24
1	B	291	SER	CA-CB	5.32	1.62	1.53
1	B	126	SER	CA-C	5.28	1.59	1.52
1	C	337	GLY	C-O	5.17	1.27	1.23
1	C	13	THR	N-CA	5.14	1.53	1.46
1	A	127	ASN	CA-C	5.13	1.58	1.52
1	B	248	ASN	N-CA	5.10	1.52	1.46
1	A	49	GLN	CA-C	-5.03	1.46	1.52

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	80	VAL	N-CA-C	6.92	117.70	110.72
1	B	80	VAL	N-CA-C	6.62	117.40	110.72
1	B	231	ASN	N-CA-C	6.39	120.93	113.20
1	A	231	ASN	N-CA-C	6.35	120.89	113.20
1	C	137	GLY	N-CA-C	-6.15	103.76	112.37
1	C	174	ASP	CB-CA-C	5.97	120.36	110.81
1	B	174	ASP	CB-CA-C	5.96	120.35	110.81
1	B	341	GLU	N-CA-C	5.92	117.70	107.93
1	B	290	THR	CA-C-N	5.83	129.12	120.90
1	B	290	THR	C-N-CA	5.83	129.12	120.90
1	B	137	GLY	N-CA-C	-5.77	104.30	112.37
1	C	52	ASN	N-CA-C	-5.76	99.81	108.96
1	A	127	ASN	N-CA-C	5.71	117.87	109.42
1	A	341	GLU	N-CA-C	5.61	117.19	107.93
1	A	52	ASN	N-CA-C	-5.51	100.19	108.96
1	B	107	GLN	N-CA-C	5.48	117.40	109.07
1	B	298	GLN	N-CA-C	5.47	117.30	108.32
1	A	298	GLN	N-CA-C	5.46	117.28	108.32
1	A	174	ASP	CB-CA-C	5.45	119.53	110.81
1	C	341	GLU	N-CA-C	5.43	116.89	107.93
1	C	290	THR	CA-C-N	5.41	128.53	120.90
1	C	290	THR	C-N-CA	5.41	128.53	120.90
1	C	312	ARG	N-CA-C	5.38	117.27	108.55
1	B	75	ILE	N-CA-C	5.38	116.27	107.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	175	LYS	N-CA-C	5.30	117.14	111.36
1	B	127	ASN	N-CA-C	5.27	116.56	109.24
1	B	52	ASN	N-CA-C	-5.22	100.65	108.96
1	B	44	ASN	N-CA-C	-5.22	105.28	110.97
1	C	231	ASN	N-CA-C	5.20	119.49	113.20
1	A	336	ASN	N-CA-C	5.19	117.02	111.36
1	B	101	GLU	N-CA-C	5.18	116.81	108.32
1	C	298	GLN	N-CA-C	5.14	116.88	108.55
1	C	178	LEU	CB-CA-C	-5.11	102.80	111.02
1	A	195	GLY	CA-C-N	5.10	126.22	119.84
1	A	195	GLY	C-N-CA	5.10	126.22	119.84
1	B	44	ASN	O-C-N	5.10	127.33	122.03
1	A	80	VAL	N-CA-C	5.08	115.85	110.72
1	A	107	GLN	N-CA-C	5.08	116.78	109.07
1	C	153	ILE	CB-CA-C	5.07	118.51	110.81
1	B	26	ALA	N-CA-C	5.05	116.40	107.61
1	B	178	LEU	CB-CA-C	-5.04	102.20	110.72

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	147	ASP	Peptide
1	A	37	GLY	Peptide
1	B	37	GLY	Peptide
1	C	147	ASP	Peptide
1	C	37	GLY	Peptide

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	2665	0	2420	30	0
1	B	2665	0	2420	29	0
1	C	2665	0	2420	22	0
All	All	7995	0	7260	77	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 5.

All (77) 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:75:ILE:HA	1:A:96:THR:HG22	1.51	0.92
1:A:92:GLY:HA2	1:A:96:THR:CG2	2.26	0.64
1:A:92:GLY:HA2	1:A:96:THR:HG21	1.86	0.57
1:A:224:GLY:HA3	1:A:238:GLN:HG2	1.86	0.57
1:C:224:GLY:HA3	1:C:238:GLN:HG2	1.87	0.56
1:A:299:ILE:HD12	1:A:299:ILE:N	2.21	0.55
1:C:230:GLU:H	1:C:230:GLU:CD	2.14	0.55
1:A:236:PHE:C	1:A:236:PHE:CD1	2.85	0.54
1:B:224:GLY:HA3	1:B:238:GLN:HG2	1.88	0.54
1:A:255:ASP:N	1:A:255:ASP:OD1	2.40	0.54
1:C:92:GLY:HA2	1:C:96:THR:OG1	2.07	0.54
1:C:299:ILE:HD12	1:C:299:ILE:N	2.23	0.53
1:B:92:GLY:HA2	1:B:96:THR:OG1	2.08	0.53
1:B:299:ILE:N	1:B:299:ILE:HD12	2.23	0.53
1:C:255:ASP:OD1	1:C:255:ASP:N	2.41	0.53
1:A:92:GLY:CA	1:A:96:THR:HG21	2.39	0.52
1:A:92:GLY:HA2	1:A:96:THR:HG23	1.91	0.52
1:B:317:ASP:OD1	1:B:338:ASP:OD1	2.29	0.51
1:B:332:THR:C	1:B:334:GLY:H	2.19	0.51
1:C:24:ILE:O	1:C:54:SER:HA	2.11	0.50
1:A:317:ASP:OD1	1:A:338:ASP:OD1	2.29	0.50
1:B:255:ASP:OD1	1:B:255:ASP:N	2.41	0.50
1:A:299:ILE:HD12	1:A:299:ILE:H	1.76	0.50
1:B:24:ILE:O	1:B:54:SER:HA	2.12	0.50
1:C:317:ASP:OD1	1:C:338:ASP:OD1	2.29	0.50
1:B:123:TRP:CD2	1:B:124:PRO:HA	2.48	0.49
1:C:236:PHE:C	1:C:236:PHE:CD1	2.91	0.49
1:A:24:ILE:O	1:A:54:SER:HA	2.13	0.49
1:C:123:TRP:CD2	1:C:124:PRO:HA	2.48	0.49
1:A:81:ASP:OD2	1:C:150:SER:OG	2.17	0.48
1:B:236:PHE:CD1	1:B:236:PHE:C	2.91	0.48
1:B:332:THR:C	1:B:334:GLY:N	2.72	0.48
1:A:304:ASP:O	1:A:305:PRO:C	2.57	0.47
1:B:91:LEU:HD23	1:B:91:LEU:HA	1.87	0.47
1:B:75:ILE:HA	1:B:96:THR:HG23	1.96	0.47
1:A:99:GLY:HA3	1:A:108:VAL:O	2.14	0.47
1:A:332:THR:C	1:A:334:GLY:H	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:332:THR:C	1:C:334:GLY:H	2.23	0.46
1:A:150:SER:OG	1:B:81:ASP:OD2	2.20	0.46
1:B:304:ASP:O	1:B:305:PRO:C	2.58	0.46
1:A:77:GLY:O	1:A:94:ARG:HD3	2.15	0.46
1:B:99:GLY:HA3	1:B:108:VAL:O	2.15	0.45
1:B:51:ASN:HD21	1:B:53:GLU:CD	2.24	0.45
1:B:77:GLY:O	1:B:94:ARG:HD3	2.17	0.45
1:B:231:ASN:OD1	1:B:231:ASN:C	2.59	0.45
1:C:231:ASN:OD1	1:C:231:ASN:C	2.60	0.45
1:C:332:THR:C	1:C:334:GLY:N	2.75	0.45
1:C:99:GLY:HA3	1:C:108:VAL:O	2.17	0.44
1:A:123:TRP:CD2	1:A:124:PRO:HA	2.52	0.44
1:C:127:ASN:HA	1:C:128:PRO:HA	1.82	0.44
1:A:332:THR:C	1:A:334:GLY:N	2.75	0.43
1:C:299:ILE:HD12	1:C:299:ILE:H	1.82	0.43
1:A:231:ASN:OD1	1:A:231:ASN:C	2.61	0.43
1:C:23:GLY:HA3	1:C:350:PHE:CZ	2.53	0.43
1:C:77:GLY:O	1:C:94:ARG:HD3	2.18	0.43
1:B:328:ALA:O	1:B:329:ALA:C	2.61	0.43
1:B:150:SER:OG	1:C:81:ASP:OD2	2.15	0.42
1:B:164:PHE:CE2	1:B:166:ILE:HD11	2.55	0.42
1:A:48:MSE:HG3	1:A:49:GLN:N	2.35	0.42
1:B:120:LEU:HA	1:B:123:TRP:O	2.20	0.42
1:B:299:ILE:HD12	1:B:299:ILE:H	1.82	0.42
1:A:23:GLY:HA3	1:A:350:PHE:CZ	2.54	0.42
1:C:120:LEU:HA	1:C:123:TRP:O	2.20	0.42
1:A:239:TYR:OH	1:A:255:ASP:OD2	2.37	0.41
1:A:164:PHE:CE2	1:A:166:ILE:HD11	2.54	0.41
1:A:290:THR:HA	1:A:319:ASN:HB2	2.03	0.41
1:A:51:ASN:HD21	1:A:53:GLU:CD	2.27	0.41
1:B:230:GLU:CD	1:B:230:GLU:H	2.28	0.41
1:B:290:THR:HA	1:B:319:ASN:HB2	2.03	0.41
1:B:48:MSE:HG3	1:B:49:GLN:N	2.35	0.41
1:C:51:ASN:HD21	1:C:53:GLU:CD	2.29	0.40
1:C:91:LEU:HD23	1:C:91:LEU:HA	1.88	0.40
1:A:230:GLU:CD	1:A:230:GLU:H	2.29	0.40
1:B:135:ASP:C	1:B:135:ASP:OD1	2.63	0.40
1:A:93:GLU:HA	1:B:87:GLU:HG3	2.03	0.40
1:A:127:ASN:HA	1:A:128:PRO:HA	1.80	0.40
1:B:294:VAL:O	1:B:294:VAL:HG13	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	339/352 (96%)	314 (93%)	20 (6%)	5 (2%)	8	30
1	B	339/352 (96%)	315 (93%)	19 (6%)	5 (2%)	8	30
1	C	339/352 (96%)	315 (93%)	20 (6%)	4 (1%)	10	34
All	All	1017/1056 (96%)	944 (93%)	59 (6%)	14 (1%)	9	31

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	95	ASP
1	B	95	ASP
1	C	95	ASP
1	A	333	SER
1	B	333	SER
1	C	333	SER
1	A	128	PRO
1	B	128	PRO
1	B	329	ALA
1	C	128	PRO
1	C	329	ALA
1	A	329	ALA
1	B	230	GLU
1	A	230	GLU

5.3.2 Protein sidechains [i](#)

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	262/258 (102%)	245 (94%)	17 (6%)	15	42
1	B	262/258 (102%)	245 (94%)	17 (6%)	15	42
1	C	262/258 (102%)	248 (95%)	14 (5%)	20	48
All	All	786/774 (102%)	738 (94%)	48 (6%)	17	44

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	GLU
1	A	83	SER
1	A	102	SER
1	A	107	GLN
1	A	132	ASP
1	A	165	SER
1	A	170	VAL
1	A	193	LYS
1	A	210	GLU
1	A	230	GLU
1	A	234	SER
1	A	269	GLN
1	A	291	SER
1	A	299	ILE
1	A	306	SER
1	A	335	THR
1	A	344	VAL
1	B	83	SER
1	B	96	THR
1	B	102	SER
1	B	107	GLN
1	B	132	ASP
1	B	156	ASP
1	B	165	SER
1	B	170	VAL
1	B	193	LYS
1	B	230	GLU
1	B	234	SER
1	B	269	GLN
1	B	291	SER
1	B	299	ILE
1	B	306	SER
1	B	335	THR
1	B	344	VAL

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Mol	Chain	Res	Type
1	C	10	GLU
1	C	83	SER
1	C	102	SER
1	C	107	GLN
1	C	132	ASP
1	C	165	SER
1	C	170	VAL
1	C	193	LYS
1	C	230	GLU
1	C	234	SER
1	C	291	SER
1	C	299	ILE
1	C	306	SER
1	C	335	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	GLN
1	A	146	GLN
1	A	198	GLN
1	A	269	GLN
1	B	41	GLN
1	B	107	GLN
1	B	198	GLN
1	C	107	GLN
1	C	146	GLN
1	C	198	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no 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	332/352 (94%)	-0.28	0 100 100	59, 73, 93, 115	0
1	B	332/352 (94%)	-0.20	0 100 100	54, 68, 89, 118	0
1	C	332/352 (94%)	-0.26	1 (0%) 90 80	55, 67, 85, 116	0
All	All	996/1056 (94%)	-0.25	1 (0%) 92 86	54, 69, 90, 118	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	10	GLU	3.6

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.