



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:54 AM UTC

PDB ID : 6FBM / pdb_00006fbm
Title : Crystal structure of GNIP1Aa from Chromobacterium piscinae
Authors : Freigang, J.; Zaitseva, J.
Deposited on : 2017-12-19
Resolution : 2.50 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

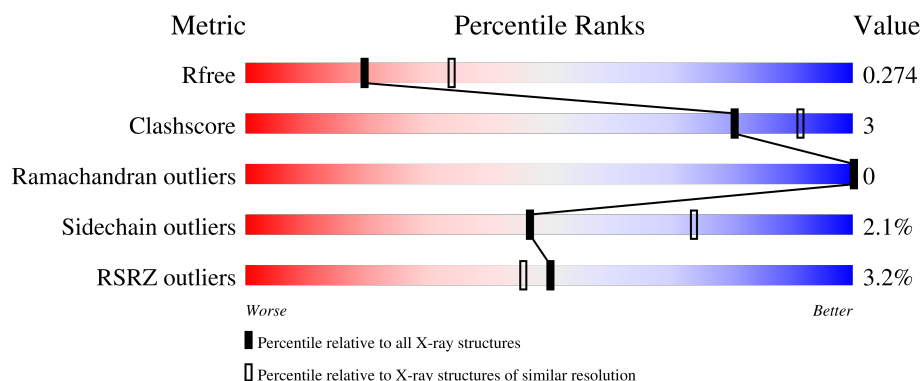
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	536	0% 91% 5% .
1	B	536	3% 91% 6% .
1	C	536	4% 91% 6% .
1	D	536	4% 89% 7% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 16588 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 Gram-negative insecticidal protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	521	Total	C	N	O	S	0	0	0
			4034	2527	695	785	27			
1	B	521	Total	C	N	O	S	0	0	0
			4034	2527	695	785	27			
1	C	521	Total	C	N	O	S	0	0	0
			4034	2527	695	785	27			
1	D	521	Total	C	N	O	S	0	0	0
			4034	2527	695	785	27			

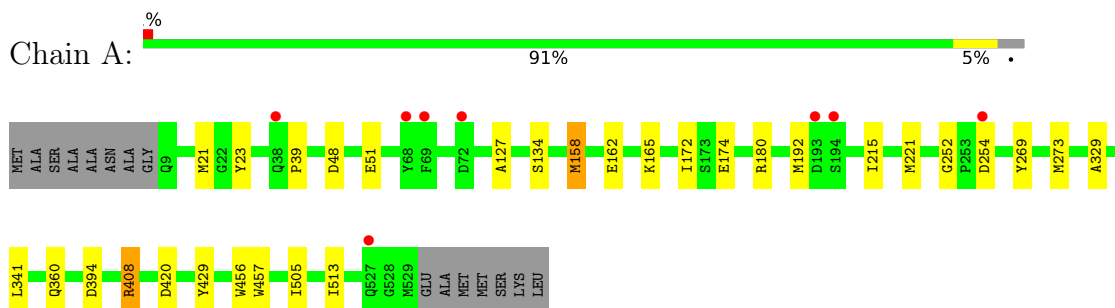
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	120	Total	O	0	0
			120	120		
2	B	116	Total	O	0	0
			116	116		
2	C	105	Total	O	0	0
			105	105		
2	D	111	Total	O	0	0
			111	111		

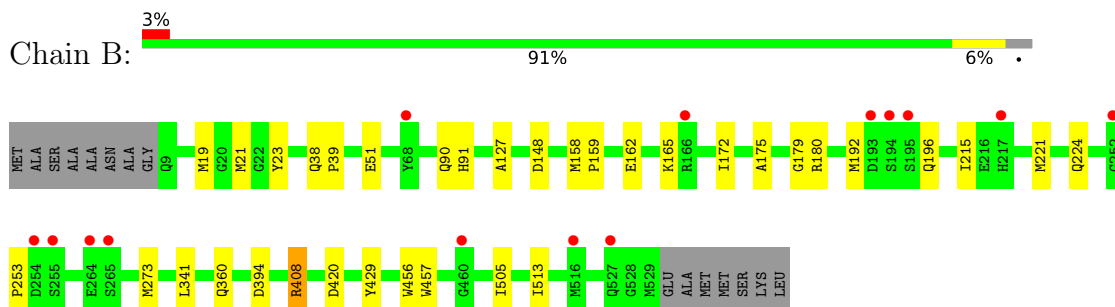
3 Residue-property plots [i](#)

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.

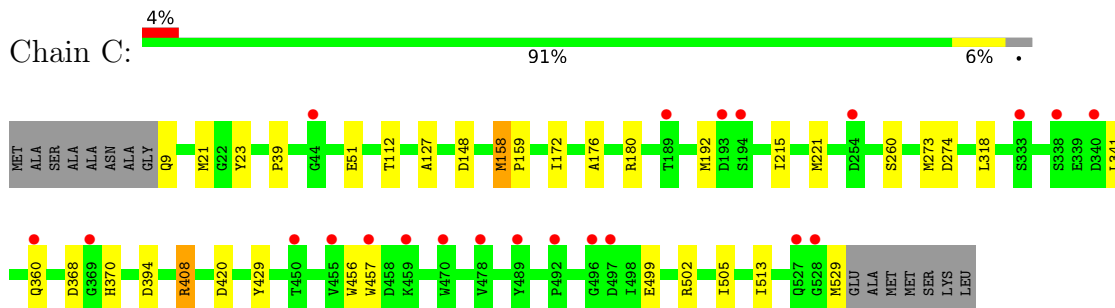
- Molecule 1: Gram-negative insecticidal protein



- Molecule 1: Gram-negative insecticidal protein

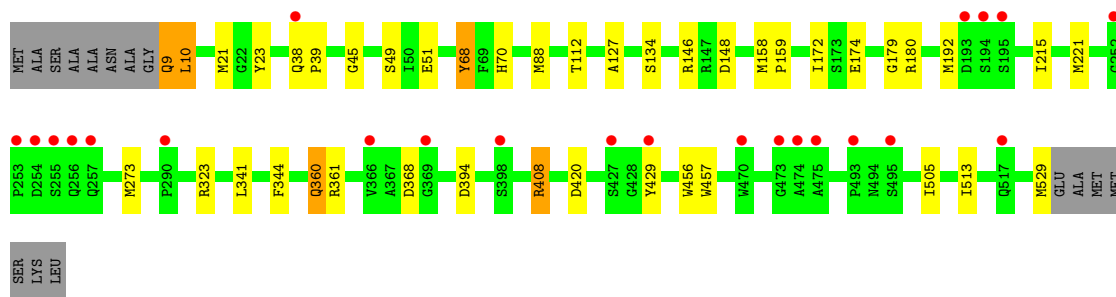


- Molecule 1: Gram-negative insecticidal protein



- Molecule 1: Gram-negative insecticidal protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.68Å 143.19Å 209.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.99 – 2.50 14.99 – 2.50	Depositor EDS
% Data completeness (in resolution range)	96.7 (14.99-2.50) 96.7 (14.99-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.8.0158 2016/10/03	Depositor
R, R_{free}	0.230 , 0.271 0.233 , 0.274	Depositor DCC
R_{free} test set	4016 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	21.5	Xtrriage
Anisotropy	0.032	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 25.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	16588	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.11 % 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 1.1996e-04. 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 [i](#)

5.1 Standard geometry [i](#)

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	0.76	0/4129	0.89	3/5589 (0.1%)
1	B	0.76	0/4129	0.90	1/5589 (0.0%)
1	C	0.80	0/4129	0.91	1/5589 (0.0%)
1	D	0.80	1/4129 (0.0%)	0.93	3/5589 (0.1%)
All	All	0.78	1/16516 (0.0%)	0.91	8/22356 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	49	SER	C-O	-5.03	1.17	1.23

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	192	MET	N-CA-C	7.03	119.87	110.88
1	D	10	LEU	N-CA-C	5.96	118.90	110.14
1	B	192	MET	N-CA-C	5.80	119.56	111.56
1	C	192	MET	N-CA-C	5.63	119.33	111.56
1	A	192	MET	N-CA-C	5.57	119.24	111.56
1	D	68	TYR	CA-CB-CG	5.39	123.61	113.90
1	A	252	GLY	CA-C-N	5.11	126.23	119.84
1	A	252	GLY	C-N-CA	5.11	126.23	119.84

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	4034	0	3895	18	0
1	B	4034	0	3895	20	0
1	C	4034	0	3895	22	1
1	D	4034	0	3895	24	1
2	A	120	0	0	2	0
2	B	116	0	0	0	0
2	C	105	0	0	2	0
2	D	111	0	0	0	0
All	All	16588	0	15580	82	1

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 3.

All (82) 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:D:127:ALA:HB1	1:D:273:MET:HE1	1.45	0.98
1:C:127:ALA:HB1	1:C:273:MET:HE1	1.46	0.95
1:A:127:ALA:HB1	1:A:273:MET:HE1	1.48	0.95
1:B:127:ALA:HB1	1:B:273:MET:HE1	1.47	0.95
1:B:19:MET:HE1	1:B:175:ALA:HB3	1.54	0.89
1:C:499:GLU:OE2	1:C:502:ARG:NH2	2.19	0.76
1:C:499:GLU:CD	1:C:502:ARG:NH2	2.52	0.68
1:B:196:GLN:HB3	1:B:224:GLN:HE21	1.59	0.68
1:C:127:ALA:CB	1:C:273:MET:HE1	2.22	0.68
1:D:127:ALA:CB	1:D:273:MET:HE1	2.21	0.67
1:C:180:ARG:HB2	1:C:273:MET:HE2	1.77	0.66
1:A:180:ARG:HB2	1:A:273:MET:HE2	1.80	0.64
1:D:180:ARG:HB2	1:D:273:MET:HE2	1.78	0.63
1:C:215:ILE:HG12	1:C:221:MET:HE3	1.79	0.63
1:C:370:HIS:CE1	2:C:698:HOH:O	2.52	0.63
1:B:180:ARG:HB2	1:B:273:MET:HE2	1.80	0.62
1:D:215:ILE:HG12	1:D:221:MET:HE3	1.80	0.62
1:A:215:ILE:HG12	1:A:221:MET:HE3	1.81	0.61
1:A:127:ALA:CB	1:A:273:MET:HE1	2.27	0.60
1:C:370:HIS:HE1	2:C:698:HOH:O	1.83	0.59
1:B:408:ARG:HD3	1:B:505:ILE:HD11	1.85	0.57
1:C:408:ARG:HD3	1:C:505:ILE:HD11	1.84	0.57
1:B:215:ILE:HG12	1:B:221:MET:HE3	1.87	0.56
1:D:408:ARG:HD3	1:D:505:ILE:HD11	1.85	0.56
1:A:408:ARG:HD3	1:A:505:ILE:HD11	1.85	0.56
1:C:112:THR:HG23	1:C:529:MET:HE3	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:TYR:OH	1:D:45:GLY:HA3	2.06	0.55
1:D:112:THR:HG23	1:D:529:MET:HE3	1.87	0.55
1:B:127:ALA:CB	1:B:273:MET:HE1	2.28	0.55
1:C:499:GLU:OE1	1:C:502:ARG:NH2	2.40	0.55
1:B:90:GLN:HB3	1:B:91:HIS:ND1	2.23	0.54
1:D:9:GLN:HE21	1:D:9:GLN:N	2.06	0.54
1:D:180:ARG:HD3	1:D:273:MET:O	2.09	0.52
1:C:180:ARG:HD3	1:C:273:MET:O	2.11	0.51
1:D:158:MET:HE3	1:D:159:PRO:HD2	1.92	0.51
1:B:180:ARG:HD3	1:B:273:MET:O	2.12	0.50
1:A:158:MET:HA	1:A:158:MET:HE3	1.93	0.50
1:A:134:SER:HB3	1:A:174:GLU:HG3	1.94	0.49
1:A:180:ARG:HD3	1:A:273:MET:O	2.11	0.49
1:A:254:ASP:OD2	1:B:253:PRO:O	2.30	0.49
1:B:158:MET:HE3	1:B:158:MET:HA	1.94	0.49
1:B:158:MET:HE3	1:B:159:PRO:HD2	1.95	0.49
1:D:158:MET:HE3	1:D:158:MET:HA	1.94	0.49
1:B:196:GLN:HB3	1:B:224:GLN:NE2	2.27	0.49
1:D:394:ASP:HB2	1:D:429:TYR:CE1	2.49	0.48
1:A:165:LYS:HE3	2:A:718:HOH:O	2.13	0.48
1:B:394:ASP:HB2	1:B:429:TYR:CE1	2.49	0.47
1:D:344:PHE:CE1	1:D:360:GLN:NE2	2.83	0.47
1:D:408:ARG:HG2	1:D:420:ASP:OD1	2.14	0.47
1:A:394:ASP:HB2	1:A:429:TYR:CE1	2.49	0.47
1:D:68:TYR:HB3	1:D:70:HIS:CE1	2.50	0.46
1:B:408:ARG:HG2	1:B:420:ASP:OD1	2.15	0.46
1:C:158:MET:HE3	1:C:158:MET:HA	1.98	0.46
1:C:408:ARG:HG2	1:C:420:ASP:OD1	2.15	0.46
1:A:408:ARG:HG2	1:A:420:ASP:OD1	2.15	0.46
1:B:21:MET:HE1	1:B:39:PRO:HB3	1.98	0.46
1:C:394:ASP:HB2	1:C:429:TYR:CE1	2.49	0.46
1:D:21:MET:HE1	1:D:39:PRO:CB	2.47	0.45
1:D:341:LEU:HB2	1:D:456:TRP:CD1	2.51	0.45
1:C:23:TYR:CD1	1:C:172:ILE:HD11	2.52	0.45
1:D:134:SER:HB3	1:D:174:GLU:HG3	1.97	0.45
1:D:21:MET:HE1	1:D:39:PRO:HB3	1.98	0.45
1:C:341:LEU:HB2	1:C:456:TRP:CD1	2.52	0.45
1:B:21:MET:HE1	1:B:39:PRO:CB	2.47	0.44
1:B:341:LEU:HB2	1:B:456:TRP:CD1	2.52	0.44
1:A:341:LEU:HB2	1:A:456:TRP:CD1	2.52	0.44
1:D:23:TYR:CD1	1:D:172:ILE:HD11	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:TYR:CD1	1:A:172:ILE:HD11	2.53	0.44
1:A:21:MET:HE1	1:A:39:PRO:CB	2.48	0.44
1:C:21:MET:HE1	1:C:39:PRO:HA	1.99	0.43
1:B:23:TYR:CD1	1:B:172:ILE:HD11	2.53	0.43
1:A:329:ALA:HB1	2:A:611:HOH:O	2.18	0.43
1:A:21:MET:HE1	1:A:39:PRO:HB3	2.00	0.43
1:C:499:GLU:OE2	1:C:502:ARG:CZ	2.67	0.42
1:C:318:LEU:HD12	1:C:513:ILE:HD11	2.02	0.42
1:C:176:ALA:HB3	1:C:274:ASP:HB3	2.01	0.41
1:D:179:GLY:O	1:D:273:MET:HE3	2.21	0.41
1:D:360:GLN:HG2	1:D:368:ASP:OD2	2.21	0.41
1:B:179:GLY:O	1:B:273:MET:HE3	2.21	0.41
1:D:344:PHE:CD1	1:D:360:GLN:NE2	2.88	0.40
1:C:158:MET:CE	1:C:159:PRO:HD2	2.51	0.40
1:D:10:LEU:HD12	1:D:45:GLY:HA2	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:368:ASP:O	1:D:361:ARG:O[3_646]	2.17	0.03

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	519/536 (97%)	504 (97%)	15 (3%)	0	100	100
1	B	519/536 (97%)	504 (97%)	15 (3%)	0	100	100
1	C	519/536 (97%)	505 (97%)	14 (3%)	0	100	100
1	D	519/536 (97%)	503 (97%)	16 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2076/2144 (97%)	2016 (97%)	60 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	436/445 (98%)	428 (98%)	8 (2%)	51	77
1	B	436/445 (98%)	427 (98%)	9 (2%)	47	74
1	C	436/445 (98%)	428 (98%)	8 (2%)	51	77
1	D	436/445 (98%)	425 (98%)	11 (2%)	42	69
All	All	1744/1780 (98%)	1708 (98%)	36 (2%)	47	74

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	ASP
1	A	51	GLU
1	A	158	MET
1	A	162	GLU
1	A	360	GLN
1	A	408	ARG
1	A	457	TRP
1	A	513	ILE
1	B	38	GLN
1	B	51	GLU
1	B	148	ASP
1	B	162	GLU
1	B	165	LYS
1	B	360	GLN
1	B	408	ARG
1	B	457	TRP
1	B	513	ILE

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Mol	Chain	Res	Type
1	C	9	GLN
1	C	51	GLU
1	C	148	ASP
1	C	158	MET
1	C	260	SER
1	C	360	GLN
1	C	408	ARG
1	C	457	TRP
1	D	9	GLN
1	D	38	GLN
1	D	51	GLU
1	D	88	MET
1	D	146	ARG
1	D	148	ASP
1	D	323	ARG
1	D	360	GLN
1	D	408	ARG
1	D	457	TRP
1	D	513	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	HIS
1	A	75	GLN
1	A	91	HIS
1	A	114	GLN
1	A	231	ASN
1	A	390	GLN
1	B	70	HIS
1	B	114	GLN
1	B	224	GLN
1	B	231	ASN
1	B	363	HIS
1	C	9	GLN
1	C	70	HIS
1	C	75	GLN
1	C	114	GLN
1	C	231	ASN
1	C	370	HIS
1	C	453	ASN
1	C	471	GLN

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Mol	Chain	Res	Type
1	D	38	GLN
1	D	70	HIS
1	D	75	GLN
1	D	91	HIS
1	D	114	GLN
1	D	231	ASN
1	D	433	ASN
1	D	453	ASN
1	D	471	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 oligosaccharides 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, 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	521/536 (97%)	-0.04	8 (1%) 72 68	9, 21, 46, 72	0
1	B	521/536 (97%)	0.11	14 (2%) 56 51	8, 25, 49, 78	0
1	C	521/536 (97%)	0.21	22 (4%) 40 36	6, 24, 58, 85	0
1	D	521/536 (97%)	0.32	23 (4%) 39 34	7, 29, 60, 95	0
All	All	2084/2144 (97%)	0.15	67 (3%) 50 46	6, 25, 55, 95	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	254	ASP	6.4
1	D	255	SER	5.3
1	B	254	ASP	4.2
1	D	193	ASP	4.0
1	C	254	ASP	3.9
1	C	194	SER	3.8
1	D	427	SER	3.6
1	C	492	PRO	3.5
1	D	253	PRO	3.4
1	D	429	TYR	3.4
1	A	194	SER	3.4
1	C	338	SER	3.3
1	D	252	GLY	3.3
1	B	265	SER	3.3
1	C	333	SER	3.3
1	B	194	SER	3.2
1	A	193	ASP	3.2
1	D	194	SER	3.1
1	C	457	TRP	3.1
1	B	527	GLN	3.0
1	C	489	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	252	GLY	2.9
1	A	527	GLN	2.9
1	D	517	GLN	2.8
1	C	44	GLY	2.7
1	A	68	TYR	2.7
1	C	459	LYS	2.7
1	B	193	ASP	2.7
1	D	495	SER	2.7
1	C	340	ASP	2.6
1	B	195	SER	2.6
1	D	470	TRP	2.6
1	C	527	GLN	2.5
1	C	369	GLY	2.5
1	C	470	TRP	2.5
1	D	369	GLY	2.4
1	D	257	GLN	2.4
1	C	528	GLY	2.4
1	D	195	SER	2.4
1	A	69	PHE	2.4
1	B	264	GLU	2.4
1	D	474	ALA	2.3
1	D	473	GLY	2.3
1	D	366	VAL	2.3
1	A	38	GLN	2.3
1	C	189	THR	2.2
1	D	256	GLN	2.2
1	D	493	PRO	2.2
1	B	217	HIS	2.2
1	D	290	PRO	2.2
1	C	360	GLN	2.2
1	C	455	VAL	2.1
1	C	478	VAL	2.1
1	D	38	GLN	2.1
1	C	496	GLY	2.1
1	C	193	ASP	2.1
1	D	475	ALA	2.1
1	B	255	SER	2.1
1	B	460	GLY	2.1
1	A	254	ASP	2.1
1	B	516	MET	2.1
1	C	450	THR	2.0
1	A	72	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	166	ARG	2.0
1	C	497	ASP	2.0
1	B	68	TYR	2.0
1	D	398	SER	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 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.