



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 06:22 AM UTC

PDB ID : 1BNS / pdb_00001bns
Title : STRUCTURAL STUDIES OF BARNASE MUTANTS
Authors : Chen, Y.W.
Deposited on : 1994-04-11
Resolution : 2.05 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

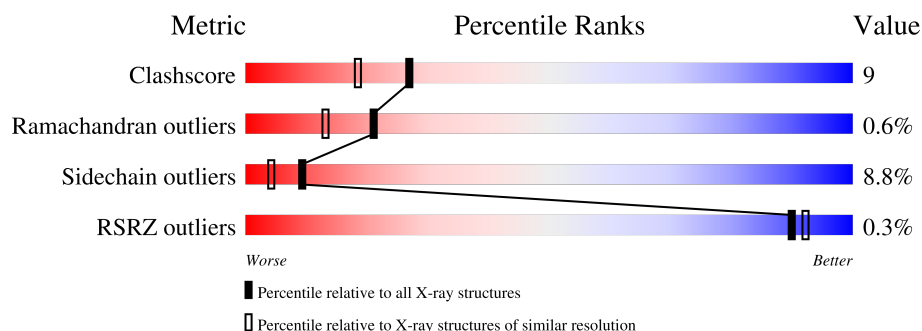
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	110	
1	B	110	
1	C	110	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2777 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 BARNASE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	108	Total	C	N	O	0	0	0
			851	541	146	164			
1	B	108	Total	C	N	O	0	1	0
			850	540	147	163			
1	C	107	Total	C	N	O	0	0	0
			852	540	149	163			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	ALA	THR	conflict	UNP P00648
B	26	ALA	THR	conflict	UNP P00648
C	26	ALA	THR	conflict	UNP P00648

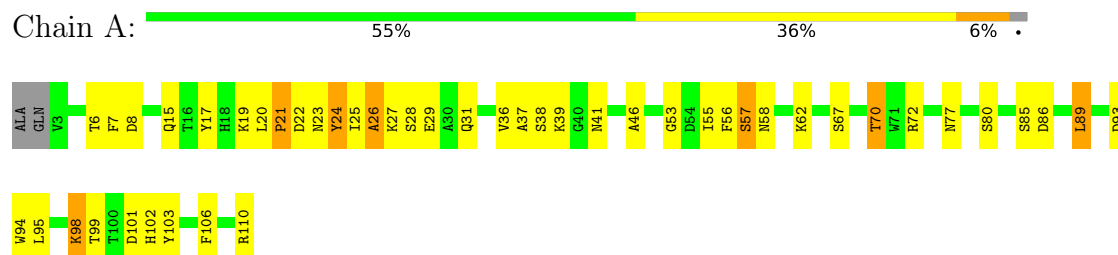
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	65	Total	O	0	0
			65	65		
2	B	73	Total	O	0	0
			73	73		
2	C	86	Total	O	0	0
			86	86		

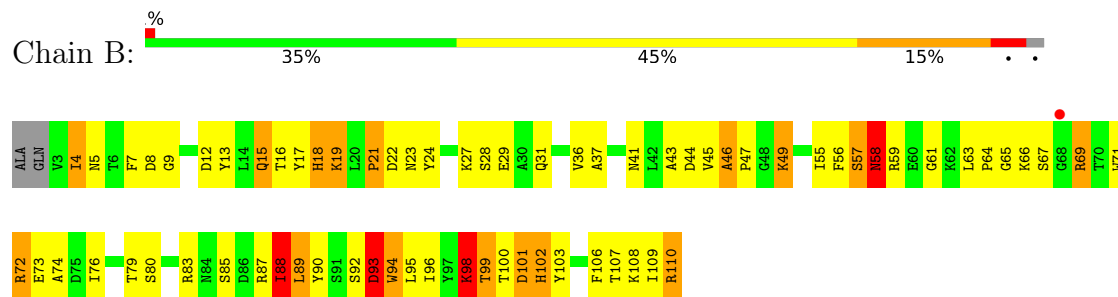
3 Residue-property plots

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.

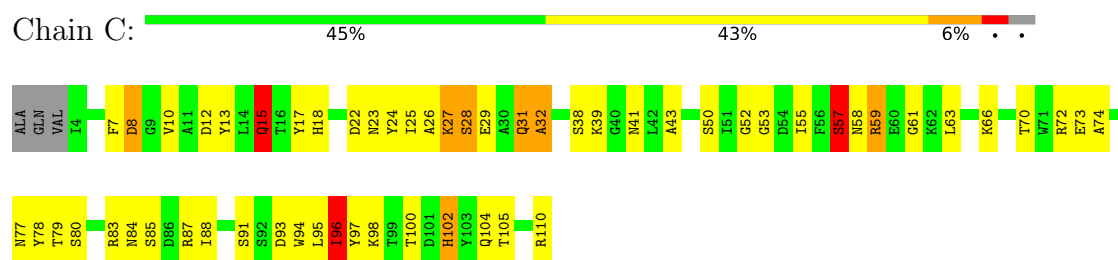
• Molecule 1: BARNASE



• Molecule 1: BARNASE



• Molecule 1: BARNASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	59.37Å 59.37Å 82.17Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 2.05 6.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-2.05) 90.2 (6.00-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.87 (at 2.04Å)	Xtriage
Refinement program	PROLSQ	Depositor
R, R_{free}	0.169 , (Not available) 0.167 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	19.8	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 117.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.004 for -h,-k,l 0.045 for h,-h-k,-l 0.027 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2777	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.44% 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	1.20	2/872 (0.2%)	2.47	55/1181 (4.7%)
1	B	1.32	2/876 (0.2%)	2.72	77/1188 (6.5%)
1	C	1.30	4/873 (0.5%)	2.64	76/1180 (6.4%)
All	All	1.27	8/2621 (0.3%)	2.61	208/3549 (5.9%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	110	ARG	C-OXT	-10.58	1.02	1.23
1	A	110	ARG	C-OXT	-9.88	1.03	1.23
1	C	94	TRP	CA-CB	5.48	1.61	1.53
1	B	95	LEU	CA-C	5.24	1.59	1.52
1	A	94	TRP	NE1-CE2	-5.21	1.31	1.37
1	C	79	THR	N-CA	5.09	1.52	1.46
1	B	36	VAL	C-N	-5.08	1.27	1.33
1	C	94	TRP	NE1-CE2	-5.01	1.31	1.37

All (208) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	93	ASP	CA-CB-CG	25.61	138.21	112.60
1	C	59	ARG	CD-NE-CZ	19.00	151.00	124.40
1	B	58	ASN	CA-CB-CG	12.12	124.72	112.60
1	C	32	ALA	CB-CA-C	10.54	128.99	110.01
1	B	110	ARG	NE-CZ-NH2	10.15	128.34	119.20
1	C	102	HIS	CB-CG-CD2	-10.14	118.01	131.20
1	A	8	ASP	CA-CB-CG	-9.86	102.74	112.60
1	A	72	ARG	NE-CZ-NH2	9.84	128.05	119.20
1	B	8	ASP	CA-C-O	-9.55	109.64	120.24
1	C	93	ASP	CA-C-O	9.53	131.19	119.95
1	A	15	GLN	OE1-CD-NE2	9.22	131.82	122.60
1	B	41	ASN	N-CA-C	9.22	123.23	112.72
1	A	23	ASN	CA-CB-CG	9.18	121.78	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	57	SER	N-CA-C	9.09	122.36	111.82
1	A	22	ASP	CB-CA-C	-9.01	92.98	110.46
1	B	89	LEU	CA-C-O	-8.82	110.86	120.30
1	B	95	LEU	N-CA-CB	8.64	123.10	110.06
1	A	102	HIS	CA-CB-CG	-8.55	105.25	113.80
1	B	56	PHE	CA-CB-CG	-8.50	105.30	113.80
1	B	22	ASP	CA-CB-CG	-8.47	104.13	112.60
1	A	36	VAL	CA-C-O	-8.47	111.16	120.48
1	B	19	LYS	CA-C-O	8.41	130.40	121.06
1	A	46	ALA	O-C-N	8.28	127.24	121.19
1	B	18	HIS	CA-CB-CG	-8.19	105.61	113.80
1	A	55	ILE	O-C-N	8.17	131.51	122.61
1	C	59	ARG	CA-C-O	-8.06	110.99	120.10
1	A	80	SER	CA-C-O	-7.97	112.93	121.38
1	C	70	THR	CA-CB-OG1	7.92	121.49	109.60
1	A	110	ARG	NE-CZ-NH2	7.88	126.29	119.20
1	C	104	GLN	OE1-CD-NE2	-7.78	114.82	122.60
1	B	46	ALA	O-C-N	7.77	127.78	121.17
1	C	110	ARG	NE-CZ-NH2	7.72	126.15	119.20
1	B	17	TYR	N-CA-C	7.66	121.94	112.59
1	B	101	ASP	N-CA-CB	-7.62	100.22	111.51
1	C	25	ILE	CA-C-O	-7.60	113.22	121.28
1	A	17	TYR	N-CA-C	7.58	122.29	112.89
1	C	22	ASP	CA-CB-CG	-7.54	105.06	112.60
1	C	24	TYR	CA-C-O	-7.53	112.44	120.80
1	B	5	ASN	CB-CG-ND2	7.52	127.68	116.40
1	C	15	GLN	O-C-N	7.51	129.80	122.07
1	C	31	GLN	CA-C-N	-7.51	108.79	122.38
1	C	31	GLN	C-N-CA	-7.51	108.79	122.38
1	B	80	SER	CA-C-O	-7.48	113.45	121.38
1	A	101	ASP	O-C-N	7.42	129.53	122.18
1	A	77	ASN	OD1-CG-ND2	-7.42	115.18	122.60
1	C	77	ASN	CA-C-N	-7.39	110.85	121.42
1	C	77	ASN	C-N-CA	-7.39	110.85	121.42
1	B	15	GLN	OE1-CD-NE2	7.34	129.94	122.60
1	A	21	PRO	CA-C-N	7.33	131.09	120.38
1	A	21	PRO	C-N-CA	7.33	131.09	120.38
1	C	22	ASP	CA-C-O	-7.26	111.89	120.10
1	C	15	GLN	CG-CD-NE2	-7.26	105.51	116.40
1	A	93	ASP	CA-CB-CG	7.22	119.82	112.60
1	B	15	GLN	CG-CD-NE2	-7.22	105.58	116.40
1	B	27	LYS	CA-C-N	7.21	129.81	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	27	LYS	C-N-CA	7.21	129.81	120.44
1	B	37	ALA	O-C-N	7.12	129.49	122.09
1	B	76	ILE	O-C-N	7.09	130.58	123.00
1	A	22	ASP	O-C-N	7.08	130.28	122.20
1	A	38	SER	CA-CB-OG	-7.08	96.95	111.10
1	B	73	GLU	O-C-N	7.08	131.73	123.17
1	B	18	HIS	CA-C-O	7.06	130.60	120.51
1	B	16	THR	O-C-N	7.03	129.68	122.09
1	B	41	ASN	OD1-CG-ND2	-6.99	115.61	122.60
1	C	110	ARG	N-CA-CB	6.97	122.35	110.50
1	C	18	HIS	CA-CB-CG	-6.96	106.84	113.80
1	C	87	ARG	NE-CZ-NH2	6.93	125.44	119.20
1	A	94	TRP	N-CA-C	6.82	121.07	111.52
1	A	85	SER	O-C-N	6.81	130.48	122.24
1	B	95	LEU	N-CA-C	-6.81	99.44	110.20
1	C	66	LYS	O-C-N	-6.77	116.39	123.56
1	A	80	SER	O-C-N	6.75	130.34	123.26
1	C	17	TYR	N-CA-C	6.73	121.66	113.38
1	B	99	THR	O-C-N	6.72	131.47	123.33
1	C	100	THR	N-CA-C	-6.72	103.01	113.02
1	C	80	SER	CA-C-O	-6.66	114.16	121.28
1	A	41	ASN	OD1-CG-ND2	-6.63	115.97	122.60
1	B	9	GLY	CA-C-O	6.60	132.05	120.57
1	B	57	SER	O-C-N	6.57	131.33	122.59
1	B	24	TYR	CA-C-O	-6.56	113.76	120.71
1	A	95	LEU	CA-C-O	-6.55	113.13	120.54
1	C	77	ASN	OD1-CG-ND2	-6.55	116.05	122.60
1	B	47	PRO	CA-C-O	-6.54	114.19	121.32
1	A	101	ASP	CA-C-N	-6.51	113.26	122.63
1	A	101	ASP	C-N-CA	-6.51	113.26	122.63
1	A	7	PHE	CA-CB-CG	-6.49	107.31	113.80
1	B	36	VAL	CA-C-N	6.48	129.29	120.54
1	B	36	VAL	C-N-CA	6.48	129.29	120.54
1	B	22	ASP	CA-C-O	-6.47	112.78	120.10
1	C	94	TRP	N-CA-C	6.47	120.47	111.74
1	C	78	TYR	O-C-N	6.45	130.62	123.01
1	B	83	ARG	NE-CZ-NH2	6.41	124.97	119.20
1	C	41	ASN	N-CA-C	6.40	120.01	112.72
1	A	99	THR	CA-C-O	6.39	128.27	120.60
1	C	84	ASN	CA-CB-CG	6.39	119.00	112.60
1	C	96	ILE	O-C-N	6.36	130.07	123.20
1	B	37	ALA	CA-C-O	-6.36	114.10	120.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	8	ASP	CA-C-O	6.34	127.14	120.42
1	C	85	SER	N-CA-CB	-6.33	100.92	110.73
1	C	12	ASP	CA-CB-CG	-6.31	106.29	112.60
1	B	16	THR	CA-C-N	-6.27	112.71	122.67
1	B	16	THR	C-N-CA	-6.27	112.71	122.67
1	B	98	LYS	O-C-N	6.25	130.73	123.17
1	B	4	ILE	CA-C-O	-6.22	113.52	120.74
1	B	29	GLU	CA-CB-CG	-6.21	101.68	114.10
1	B	41	ASN	N-CA-CB	-6.21	101.45	110.70
1	C	26	ALA	N-CA-CB	-6.21	101.88	110.38
1	A	57	SER	CA-CB-OG	-6.20	98.71	111.10
1	A	26	ALA	CA-C-O	-6.18	114.82	121.56
1	A	56	PHE	CA-CB-CG	-6.16	107.64	113.80
1	C	87	ARG	NH1-CZ-NH2	-6.16	111.30	119.30
1	B	79	THR	N-CA-C	-6.12	105.03	113.18
1	C	18	HIS	O-C-N	6.11	130.60	122.41
1	B	4	ILE	CB-CG1-CD1	-6.10	101.00	113.80
1	C	73	GLU	CA-C-O	-6.09	114.33	121.46
1	C	66	LYS	CB-CG-CD	-6.06	97.36	111.30
1	A	53	GLY	CA-C-O	-6.04	112.30	119.01
1	B	109	ILE	O-C-N	-6.01	115.05	122.57
1	B	100	THR	CA-CB-OG1	-5.97	100.64	109.60
1	C	13	TYR	CA-C-N	5.97	128.28	120.28
1	C	13	TYR	C-N-CA	5.97	128.28	120.28
1	C	28	SER	CA-C-O	-5.96	113.11	119.97
1	C	50	SER	O-C-N	5.96	130.94	123.19
1	A	6	THR	N-CA-CB	5.96	119.85	110.56
1	A	22	ASP	N-CA-C	5.88	119.27	111.75
1	B	108	LYS	CA-C-O	-5.83	114.11	120.81
1	A	21	PRO	CB-CA-C	5.80	118.54	110.95
1	B	31	GLN	OE1-CD-NE2	5.80	128.40	122.60
1	C	8	ASP	CA-CB-CG	-5.79	106.81	112.60
1	A	7	PHE	N-CA-CB	5.77	118.58	109.82
1	B	101	ASP	CA-C-O	5.76	128.00	121.32
1	B	107	THR	CA-CB-OG1	-5.72	101.01	109.60
1	C	110	ARG	CD-NE-CZ	5.72	132.41	124.40
1	C	10	VAL	CA-C-O	-5.72	114.79	120.85
1	C	63	LEU	O-C-N	5.69	128.10	121.39
1	C	23	ASN	CB-CG-ND2	5.69	124.93	116.40
1	B	21	PRO	CA-C-O	5.67	129.04	121.95
1	B	102[A]	HIS	CA-C-N	5.67	131.46	122.61
1	B	102[A]	HIS	C-N-CA	5.67	131.46	122.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	102[B]	HIS	CA-C-N	5.67	131.46	122.61
1	B	102[B]	HIS	C-N-CA	5.67	131.46	122.61
1	C	39	LYS	CA-C-O	-5.66	112.70	119.15
1	B	31	GLN	CA-C-O	5.65	126.54	120.55
1	A	26	ALA	N-CA-CB	-5.64	102.05	110.17
1	C	95	LEU	CA-C-O	-5.62	114.61	121.36
1	A	70	THR	CA-CB-CG2	5.59	120.01	110.50
1	C	38	SER	CA-CB-OG	-5.58	99.93	111.10
1	C	32	ALA	N-CA-CB	-5.58	101.90	110.44
1	C	79	THR	O-C-N	-5.58	115.17	122.59
1	B	89	LEU	CB-CA-C	5.57	119.41	110.16
1	A	36	VAL	O-C-N	5.56	128.94	122.99
1	C	52	GLY	CA-C-O	-5.56	116.05	121.06
1	C	105	THR	O-C-N	5.56	129.72	123.27
1	B	95	LEU	CA-C-O	-5.54	114.74	121.05
1	B	8	ASP	O-C-N	5.51	129.39	122.23
1	C	102	HIS	ND1-CG-CD2	5.50	111.60	106.10
1	A	27	LYS	CA-C-O	-5.48	113.91	120.10
1	B	21	PRO	CB-CA-C	5.47	118.18	111.12
1	C	24	TYR	CA-CB-CG	-5.43	104.12	113.90
1	B	45	VAL	CA-C-O	-5.41	115.11	120.85
1	B	49	LYS	N-CA-CB	-5.41	102.11	110.56
1	B	69	ARG	CD-NE-CZ	5.40	131.96	124.40
1	B	44	ASP	CB-CA-C	-5.38	100.84	110.70
1	C	97	TYR	CA-C-O	-5.38	115.14	121.44
1	A	15	GLN	O-C-N	5.33	129.47	122.33
1	C	22	ASP	O-C-N	5.31	128.44	122.22
1	C	61	GLY	O-C-N	-5.31	115.86	122.34
1	C	96	ILE	CB-CG1-CD1	5.31	124.95	113.80
1	A	24	TYR	O-C-N	5.30	129.20	123.10
1	B	65	GLY	CA-C-N	5.29	130.79	122.21
1	B	65	GLY	C-N-CA	5.29	130.79	122.21
1	A	19	LYS	O-C-N	5.29	129.28	123.25
1	A	38	SER	N-CA-C	5.26	118.80	112.38
1	C	50	SER	CA-C-N	-5.26	114.41	121.66
1	C	50	SER	C-N-CA	-5.26	114.41	121.66
1	B	88	ILE	CA-C-O	-5.24	114.93	120.48
1	B	31	GLN	CG-CD-OE1	-5.23	110.33	120.80
1	A	23	ASN	CB-CG-ND2	-5.22	108.57	116.40
1	A	41	ASN	N-CA-C	5.21	119.33	112.92
1	A	38	SER	N-CA-CB	-5.21	101.46	110.32
1	A	86	ASP	CB-CG-OD1	5.20	130.35	118.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	12	ASP	CB-CG-OD2	-5.20	106.45	118.40
1	C	13	TYR	CA-C-O	-5.17	115.39	120.82
1	C	83	ARG	O-C-N	-5.16	116.47	122.81
1	C	29	GLU	CA-C-O	-5.15	114.96	120.42
1	C	87	ARG	N-CA-CB	5.13	120.32	111.13
1	A	98	LYS	O-C-N	5.12	129.21	123.27
1	B	76	ILE	CA-C-O	-5.12	115.34	121.28
1	C	91	SER	O-C-N	5.12	128.78	122.84
1	C	102	HIS	CB-CG-ND1	5.11	130.37	122.70
1	B	69	ARG	CA-C-N	-5.08	115.82	122.99
1	B	69	ARG	C-N-CA	-5.08	115.82	122.99
1	C	43	ALA	CA-C-O	-5.07	113.15	119.38
1	B	28	SER	CA-C-O	-5.07	115.50	120.82
1	C	70	THR	OG1-CB-CG2	-5.06	99.18	109.30
1	C	110	ARG	NE-CZ-NH1	-5.06	116.44	121.50
1	B	83	ARG	CA-C-O	-5.06	116.19	121.55
1	A	37	ALA	CB-CA-C	5.06	119.18	110.79
1	C	38	SER	CA-C-O	5.05	125.61	119.05
1	A	27	LYS	CB-CA-C	-5.04	101.31	110.63
1	A	25	ILE	CA-C-N	5.03	128.84	120.94
1	A	25	ILE	C-N-CA	5.03	128.84	120.94
1	B	9	GLY	CA-C-N	5.02	127.34	120.46
1	B	9	GLY	C-N-CA	5.02	127.34	120.46
1	B	41	ASN	CB-CG-OD1	5.02	130.84	120.80
1	C	7	PHE	CA-CB-CG	5.01	118.81	113.80
1	A	39	LYS	CA-C-N	5.01	127.67	120.11
1	A	39	LYS	C-N-CA	5.01	127.67	120.11

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	851	0	811	7	0
1	B	850	0	802	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	852	0	820	13	0
2	A	65	0	0	0	0
2	B	73	0	0	1	0
2	C	86	0	0	3	0
All	All	2777	0	2433	46	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 (46) 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:C:28:SER:HA	1:C:31:GLN:HE21	1.31	0.95
1:B:64:PRO:HB2	1:B:69:ARG:NH1	2.12	0.64
1:A:26:ALA:HB3	1:A:29:GLU:HG3	1.79	0.64
1:B:64:PRO:HB2	1:B:69:ARG:HH12	1.63	0.64
1:C:28:SER:HA	1:C:31:GLN:NE2	2.09	0.63
1:C:57:SER:HB3	1:C:59:ARG:HD2	1.81	0.62
1:B:64:PRO:C	1:B:69:ARG:HH11	2.09	0.61
1:A:89:LEU:HD13	1:A:106:PHE:HE2	1.67	0.60
1:C:74:ALA:HB3	1:C:88:ILE:HG22	1.84	0.60
1:C:32:ALA:HB3	2:C:196:HOH:O	2.01	0.59
1:B:4:ILE:HD12	1:B:21:PRO:HB3	1.86	0.58
1:C:27:LYS:O	1:C:31:GLN:HG3	2.05	0.57
1:B:72:ARG:HB3	1:B:90:TYR:CZ	2.41	0.56
1:C:58:ASN:N	2:C:148:HOH:O	2.38	0.55
1:A:89:LEU:HD13	1:A:106:PHE:CE2	2.43	0.53
1:A:28:SER:HA	1:A:31:GLN:HE21	1.74	0.52
1:B:96:ILE:HB	1:B:110:ARG:HB2	1.91	0.52
1:B:96:ILE:HD12	1:B:110:ARG:HG3	1.92	0.51
1:A:21:PRO:HD2	1:A:24:TYR:CD1	2.46	0.49
1:B:7:PHE:CE2	1:B:98:LYS:HB2	2.47	0.49
1:B:89:LEU:HD13	1:B:106:PHE:HE2	1.77	0.49
1:B:93:ASP:O	1:B:94:TRP:HB2	2.13	0.49
1:B:43:ALA:HA	2:B:172:HOH:O	2.13	0.48
1:C:57:SER:CB	1:C:59:ARG:HD2	2.43	0.47
1:B:85:SER:HA	1:B:102[B]:HIS:CD2	2.49	0.47
1:B:89:LEU:HD13	1:B:106:PHE:CE2	2.50	0.46
1:B:21:PRO:HB2	1:B:23:ASN:OD1	2.16	0.46
1:B:64:PRO:C	1:B:69:ARG:NH1	2.74	0.45
1:B:15:GLN:O	1:B:18:HIS:NE2	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:12:ASP:O	1:B:13:TYR:C	2.60	0.45
1:A:20:LEU:HA	1:A:21:PRO:HD3	1.85	0.43
1:B:59:ARG:C	1:B:61:GLY:H	2.25	0.43
1:B:69:ARG:NH2	1:B:93:ASP:OD2	2.35	0.42
1:B:57:SER:OG	1:B:58:ASN:N	2.53	0.42
1:A:57:SER:O	1:A:58:ASN:C	2.62	0.41
1:C:74:ALA:HB3	1:C:88:ILE:CG2	2.47	0.41
1:B:74:ALA:HB3	1:B:88:ILE:CG2	2.50	0.41
1:C:57:SER:HB2	2:C:146:HOH:O	2.20	0.41
1:B:64:PRO:HG2	1:B:71:TRP:HZ2	1.85	0.41
1:B:46:ALA:HB1	1:B:49:LYS:HG3	2.01	0.41
1:B:87:ARG:HB2	1:B:99:THR:HG22	2.02	0.41
1:C:15:GLN:HE22	1:C:96:ILE:HD11	1.86	0.41
1:C:55:ILE:HD13	1:C:55:ILE:HA	1.91	0.41
1:C:53:GLY:O	1:C:72:ARG:HD3	2.21	0.40
1:B:63:LEU:O	1:B:64:PRO:C	2.63	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	106/110 (96%)	100 (94%)	6 (6%)	0	100	100
1	B	107/110 (97%)	92 (86%)	13 (12%)	2 (2%)	6	2
1	C	105/110 (96%)	101 (96%)	4 (4%)	0	100	100
All	All	318/330 (96%)	293 (92%)	23 (7%)	2 (1%)	21	13

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	58	ASN

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Mol	Chain	Res	Type
1	B	94	TRP

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	87/91 (96%)	81 (93%)	6 (7%)	14	8
1	B	87/91 (96%)	77 (88%)	10 (12%)	5	1
1	C	88/91 (97%)	81 (92%)	7 (8%)	11	5
All	All	262/273 (96%)	239 (91%)	23 (9%)	9	4

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	LYS
1	A	67	SER
1	A	70	THR
1	A	89	LEU
1	A	98	LYS
1	A	103	TYR
1	B	19	LYS
1	B	55	ILE
1	B	66	LYS
1	B	67	SER
1	B	72	ARG
1	B	88	ILE
1	B	92	SER
1	B	93	ASP
1	B	98	LYS
1	B	103	TYR
1	C	8	ASP
1	C	15	GLN
1	C	27	LYS
1	C	57	SER
1	C	96	ILE

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Mol	Chain	Res	Type
1	C	98	LYS
1	C	102	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	B	31	GLN
1	C	15	GLN
1	C	31	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	108/110 (98%)	-0.88	0 100 100	11, 21, 40, 47	1 (0%)
1	B	108/110 (98%)	-0.56	1 (0%) 81 83	7, 22, 57, 66	4 (3%)
1	C	107/110 (97%)	-1.08	0 100 100	7, 16, 30, 42	1 (0%)
All	All	323/330 (97%)	-0.84	1 (0%) 90 92	7, 19, 45, 66	6 (1%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	68	GLY	2.9

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.