



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:50 AM UTC

PDB ID : 1BRG / pdb_00001brg
Title : CRYSTALLOGRAPHIC ANALYSIS OF PHE->LEU SUBSTITUTION IN
THE HYDROPHOBIC CORE OF BARNASE
Authors : Chen, Y.W.; Fersht, A.R.; Henrick, K.
Deposited on : 1994-03-30
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

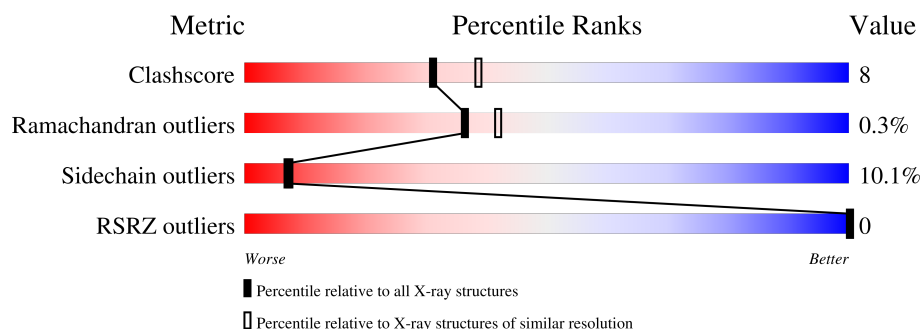
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	108	
1	B	108	
1	C	108	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2752 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
			840	533	142	165			
1	B	108	Total	C	N	O	0	0	0
			843	533	145	165			
1	C	108	Total	C	N	O	0	0	0
			843	533	144	166			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	LEU	PHE	conflict	UNP P00648
B	7	LEU	PHE	conflict	UNP P00648
C	7	LEU	PHE	conflict	UNP P00648

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total	Zn	0	0
			1	1		

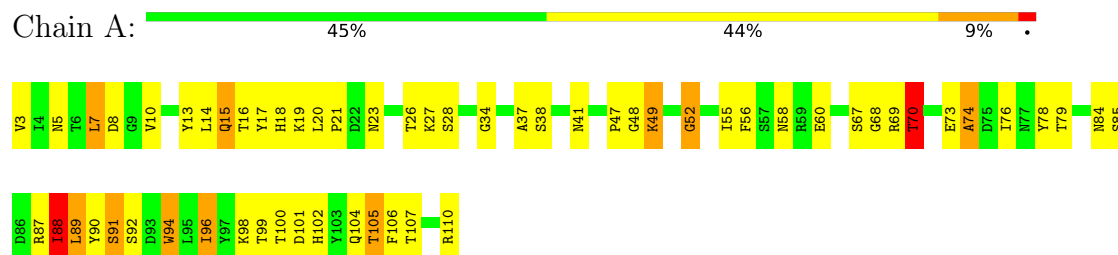
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	69	Total	O	0	0
			69	69		
3	B	77	Total	O	0	0
			77	77		
3	C	79	Total	O	0	0
			79	79		

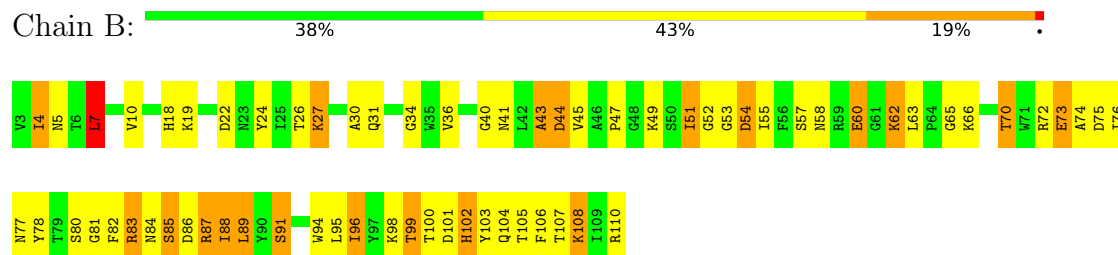
3 Residue-property plots

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.

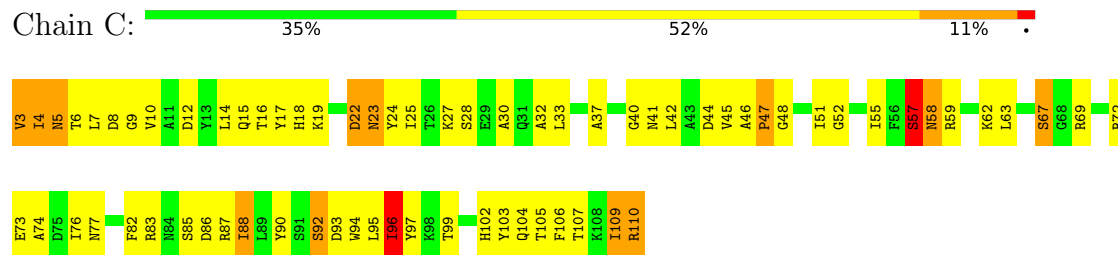
• Molecule 1: BARNASE



• 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	58.72Å 58.72Å 82.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 2.20 6.00 – 2.21	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-2.20) 90.8 (6.00-2.21)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.20Å)	Xtriage
Refinement program	PROLSQ, X-PLOR	Depositor
R, R_{free}	0.168 , (Not available) (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	13.6	Xtriage
Anisotropy	0.181	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 117.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l 0.056 for h,-h-k,-l 0.034 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2752	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.37% 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 ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	1.34	3/860 (0.3%)	2.77	86/1166 (7.4%)
1	B	1.42	6/863 (0.7%)	2.82	85/1170 (7.3%)
1	C	1.41	4/863 (0.5%)	3.00	115/1173 (9.8%)
All	All	1.39	13/2586 (0.5%)	2.87	286/3509 (8.2%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	94	TRP	CA-CB	7.58	1.65	1.53
1	A	76	ILE	N-CA	6.72	1.53	1.46
1	B	110	ARG	CZ-NH2	-6.29	1.25	1.33
1	B	87	ARG	NE-CZ	6.07	1.39	1.33
1	C	30	ALA	C-N	-5.90	1.26	1.33
1	C	94	TRP	CA-CB	5.70	1.62	1.53
1	B	41	ASN	CA-CB	5.70	1.61	1.53
1	C	90	TYR	C-O	5.57	1.30	1.23
1	B	36	VAL	C-O	-5.42	1.18	1.24
1	C	17	TYR	CA-CB	5.30	1.62	1.54
1	A	76	ILE	C-O	-5.26	1.18	1.24
1	B	41	ASN	N-CA	5.21	1.52	1.46
1	A	55	ILE	CA-CB	5.02	1.59	1.54

All (286) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	15	GLN	OE1-CD-NE2	-15.75	106.85	122.60
1	C	41	ASN	N-CA-C	13.39	128.92	112.59
1	A	105	THR	OG1-CB-CG2	-12.64	84.03	109.30
1	B	108	LYS	CG-CD-CE	12.63	140.34	111.30
1	A	105	THR	CB-CA-C	12.48	133.92	109.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	4	ILE	CA-C-O	-12.06	107.70	120.48
1	C	110	ARG	NE-CZ-NH1	-11.78	109.72	121.50
1	C	110	ARG	NE-CZ-NH2	11.52	129.56	119.20
1	B	5	ASN	OD1-CG-ND2	-11.12	111.48	122.60
1	B	75	ASP	CA-CB-CG	11.06	123.66	112.60
1	C	22	ASP	CA-CB-CG	10.96	123.56	112.60
1	C	110	ARG	CA-C-O	10.76	139.09	120.80
1	A	84	ASN	OD1-CG-ND2	-10.72	111.88	122.60
1	B	57	SER	CA-C-O	-10.54	109.75	120.82
1	C	42	LEU	O-C-N	10.45	132.96	122.09
1	B	44	ASP	CA-CB-CG	-10.25	102.35	112.60
1	B	104	GLN	OE1-CD-NE2	10.19	132.79	122.60
1	B	72	ARG	CD-NE-CZ	9.93	138.30	124.40
1	B	87	ARG	NH1-CZ-NH2	-9.88	106.46	119.30
1	B	62	LYS	CA-C-O	9.62	131.24	119.31
1	C	18	HIS	CA-CB-CG	-9.60	104.20	113.80
1	C	87	ARG	NE-CZ-NH2	9.57	127.81	119.20
1	B	78	TYR	O-C-N	9.46	134.14	123.16
1	B	10	VAL	CA-C-O	-9.41	110.87	120.85
1	A	76	ILE	CA-C-N	9.41	136.09	122.35
1	A	76	ILE	C-N-CA	9.41	136.09	122.35
1	C	83	ARG	NE-CZ-NH2	9.40	127.66	119.20
1	A	56	PHE	CA-CB-CG	-9.34	104.46	113.80
1	C	17	TYR	N-CA-C	9.17	125.64	113.72
1	A	41	ASN	OD1-CG-ND2	-9.16	113.44	122.60
1	A	28	SER	CA-C-O	-9.15	111.28	120.70
1	C	62	LYS	O-C-N	-9.14	109.85	122.46
1	B	31	GLN	OE1-CD-NE2	-9.12	113.48	122.60
1	C	72	ARG	NE-CZ-NH1	9.12	130.62	121.50
1	C	5	ASN	O-C-N	9.05	131.53	121.93
1	C	93	ASP	CA-CB-CG	8.98	121.58	112.60
1	B	87	ARG	NE-CZ-NH1	8.97	130.47	121.50
1	B	18	HIS	CA-CB-CG	-8.89	104.91	113.80
1	A	105	THR	O-C-N	-8.87	111.66	123.19
1	A	15	GLN	O-C-N	8.83	133.13	122.27
1	C	55	ILE	CA-C-N	8.80	135.61	123.11
1	C	55	ILE	C-N-CA	8.80	135.61	123.11
1	C	104	GLN	CA-C-O	-8.77	111.17	120.55
1	A	105	THR	CA-C-N	8.72	136.86	122.73
1	A	105	THR	C-N-CA	8.72	136.86	122.73
1	C	105	THR	CA-CB-CG2	8.67	125.24	110.50
1	A	21	PRO	CA-C-N	8.56	134.41	120.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	21	PRO	C-N-CA	8.56	134.41	120.63
1	C	77	ASN	CA-CB-CG	-8.49	104.11	112.60
1	C	41	ASN	CA-CB-CG	8.40	121.00	112.60
1	C	27	LYS	O-C-N	8.32	131.95	122.22
1	C	72	ARG	NH1-CZ-NH2	-8.27	108.54	119.30
1	B	27	LYS	N-CA-CB	-8.26	97.98	110.12
1	C	94	TRP	N-CA-C	8.21	122.59	112.58
1	C	27	LYS	CA-C-O	-8.17	110.87	120.10
1	C	69	ARG	CD-NE-CZ	8.10	135.74	124.40
1	B	41	ASN	N-CA-CB	-8.10	98.63	110.70
1	C	6	THR	N-CA-CB	8.04	121.74	110.17
1	B	81	GLY	O-C-N	8.02	130.51	122.90
1	C	8	ASP	O-C-N	8.01	130.61	122.12
1	A	105	THR	CA-CB-OG1	7.97	121.55	109.60
1	B	101	ASP	CA-C-O	7.95	129.47	121.20
1	C	69	ARG	CA-C-O	-7.87	111.64	120.54
1	A	16	THR	O-C-N	7.86	130.87	122.38
1	A	56	PHE	CB-CA-C	7.86	123.08	110.19
1	A	55	ILE	O-C-N	7.84	130.94	122.63
1	B	84	ASN	CA-C-O	7.79	131.65	120.51
1	B	75	ASP	CA-C-O	-7.79	112.51	121.47
1	A	99	THR	N-CA-CB	7.71	124.96	111.00
1	C	107	THR	O-C-N	7.63	132.38	123.30
1	B	53	GLY	CA-C-O	-7.62	110.34	119.03
1	B	34	GLY	O-C-N	7.55	131.14	122.38
1	C	69	ARG	O-C-N	7.51	131.78	123.22
1	B	22	ASP	CA-CB-CG	-7.50	105.10	112.60
1	A	15	GLN	OE1-CD-NE2	-7.48	115.12	122.60
1	C	82	PHE	CA-C-O	-7.42	113.47	121.56
1	B	77	ASN	OD1-CG-ND2	7.31	129.91	122.60
1	A	8	ASP	CA-CB-CG	-7.30	105.30	112.60
1	B	96	ILE	O-C-N	7.29	131.07	123.20
1	B	99	THR	N-CA-CB	7.28	124.17	111.13
1	B	47	PRO	CB-CA-C	7.25	120.77	111.64
1	B	91	SER	CA-C-O	7.24	130.50	121.66
1	A	94	TRP	N-CA-C	7.22	121.49	111.74
1	C	4	ILE	CA-C-O	-7.22	111.76	120.78
1	C	8	ASP	CA-C-O	-7.21	112.91	120.55
1	C	16	THR	CA-C-O	-7.20	111.22	120.00
1	A	110	ARG	NE-CZ-NH2	-7.19	112.73	119.20
1	C	8	ASP	CB-CG-OD1	7.19	134.93	118.40
1	A	70	THR	CA-CB-OG1	-7.18	98.83	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	100	THR	N-CA-C	-7.16	103.10	112.41
1	A	23	ASN	CB-CG-ND2	7.12	127.08	116.40
1	B	107	THR	CA-C-O	-7.09	112.44	120.32
1	B	75	ASP	O-C-N	7.00	130.88	122.96
1	A	87	ARG	CA-C-N	7.00	132.32	123.14
1	A	87	ARG	C-N-CA	7.00	132.32	123.14
1	A	85	SER	N-CA-C	6.96	121.94	113.17
1	A	84	ASN	CB-CG-ND2	6.95	126.83	116.40
1	A	38	SER	O-C-N	6.92	132.01	122.46
1	B	57	SER	N-CA-CB	6.91	120.02	110.01
1	C	87	ARG	CA-C-N	-6.90	114.19	123.10
1	C	87	ARG	C-N-CA	-6.90	114.19	123.10
1	B	30	ALA	CA-C-N	6.90	129.86	120.54
1	B	30	ALA	C-N-CA	6.90	129.86	120.54
1	B	41	ASN	CA-CB-CG	-6.89	105.71	112.60
1	B	95	LEU	N-CA-CB	6.87	120.44	110.06
1	B	63	LEU	O-C-N	6.87	130.56	121.70
1	C	72	ARG	CD-NE-CZ	6.86	134.00	124.40
1	A	5	ASN	OD1-CG-ND2	-6.85	115.75	122.60
1	A	68	GLY	N-CA-C	-6.82	106.57	114.69
1	B	108	LYS	CB-CG-CD	6.80	126.95	111.30
1	C	88	ILE	O-C-N	-6.78	116.23	123.14
1	A	55	ILE	CB-CA-C	-6.77	100.28	111.32
1	B	101	ASP	N-CA-C	6.72	119.01	110.61
1	B	22	ASP	N-CA-C	6.70	120.91	112.87
1	B	72	ARG	NE-CZ-NH2	-6.70	113.17	119.20
1	C	95	LEU	N-CA-C	-6.64	99.71	110.20
1	A	28	SER	O-C-N	6.62	129.24	122.09
1	C	105	THR	CA-CB-OG1	-6.57	99.74	109.60
1	B	77	ASN	CB-CG-ND2	-6.49	106.67	116.40
1	C	44	ASP	CA-CB-CG	-6.48	106.12	112.60
1	C	85	SER	CA-C-N	6.48	131.05	122.30
1	C	85	SER	C-N-CA	6.48	131.05	122.30
1	B	41	ASN	N-CA-C	6.47	120.10	112.72
1	B	110	ARG	N-CA-CB	6.46	121.49	110.50
1	C	28	SER	CA-C-N	6.46	129.47	120.29
1	C	28	SER	C-N-CA	6.46	129.47	120.29
1	C	67	SER	O-C-N	6.46	131.19	122.59
1	A	17	TYR	CA-C-N	6.43	131.72	122.40
1	A	17	TYR	C-N-CA	6.43	131.72	122.40
1	A	49	LYS	N-CA-C	6.39	120.48	110.32
1	C	23	ASN	OD1-CG-ND2	6.39	128.99	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	8	ASP	N-CA-C	6.38	118.31	111.36
1	A	79	THR	CA-C-O	-6.37	110.94	119.47
1	C	16	THR	CA-C-N	-6.36	112.15	122.26
1	C	16	THR	C-N-CA	-6.36	112.15	122.26
1	B	91	SER	CA-CB-OG	-6.36	98.39	111.10
1	C	87	ARG	O-C-N	6.29	130.56	123.27
1	A	60	GLU	CG-CD-OE2	6.26	132.80	118.40
1	A	23	ASN	CA-CB-CG	6.25	118.85	112.60
1	A	96	ILE	N-CA-C	6.24	116.91	108.17
1	B	94	TRP	CA-CB-CG	-6.19	101.84	113.60
1	C	99	THR	N-CA-CB	6.17	121.18	110.81
1	C	106	PHE	N-CA-C	6.16	119.53	109.24
1	C	12	ASP	CA-C-N	6.15	128.84	120.54
1	C	12	ASP	C-N-CA	6.15	128.84	120.54
1	A	19	LYS	CA-C-O	6.14	127.89	121.38
1	B	96	ILE	CB-CG1-CD1	6.14	126.70	113.80
1	C	86	ASP	OD1-CG-OD2	-6.13	108.19	122.90
1	C	8	ASP	CA-CB-CG	-6.12	106.48	112.60
1	B	40	GLY	O-C-N	6.11	129.01	122.41
1	B	82	PHE	CA-C-O	-6.11	115.15	121.87
1	C	41	ASN	N-CA-CB	-6.11	101.14	110.92
1	C	69	ARG	NE-CZ-NH2	-6.10	113.71	119.20
1	C	102	HIS	CB-CG-CD2	-6.09	123.29	131.20
1	C	6	THR	CB-CA-C	-6.07	98.36	109.54
1	C	10	VAL	CA-C-N	6.07	128.73	120.54
1	C	10	VAL	C-N-CA	6.07	128.73	120.54
1	B	62	LYS	O-C-N	-6.05	114.11	122.46
1	C	57	SER	CA-C-O	-6.05	113.26	120.10
1	A	69	ARG	N-CA-C	6.05	118.58	108.96
1	B	26	THR	O-C-N	6.05	130.03	122.65
1	C	32	ALA	CA-C-O	-6.03	114.16	120.55
1	B	54	ASP	CB-CG-OD2	6.03	132.26	118.40
1	C	58	ASN	CA-C-O	6.03	128.45	120.98
1	C	88	ILE	CB-CA-C	6.02	119.95	110.81
1	B	100	THR	CA-CB-CG2	6.00	120.70	110.50
1	C	9	GLY	CA-C-N	-6.00	112.89	120.56
1	C	9	GLY	C-N-CA	-6.00	112.89	120.56
1	C	17	TYR	CA-C-O	5.99	125.64	119.05
1	A	69	ARG	NE-CZ-NH2	-5.99	113.81	119.20
1	C	7	LEU	CA-C-O	-5.95	114.53	121.07
1	B	70	THR	CA-C-O	5.95	127.83	121.11
1	C	83	ARG	CD-NE-CZ	5.93	132.71	124.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	17	TYR	CB-CG-CD1	-5.92	111.92	120.80
1	A	27	LYS	CA-C-O	5.92	126.83	120.55
1	C	87	ARG	CD-NE-CZ	5.92	132.68	124.40
1	A	37	ALA	CA-C-N	5.91	132.88	121.18
1	A	37	ALA	C-N-CA	5.91	132.88	121.18
1	C	32	ALA	CA-C-N	5.88	132.83	121.18
1	C	32	ALA	C-N-CA	5.88	132.83	121.18
1	C	73	GLU	CA-C-O	-5.88	114.58	121.46
1	A	56	PHE	O-C-N	-5.88	116.34	123.27
1	C	30	ALA	CA-C-N	5.88	128.15	120.28
1	C	30	ALA	C-N-CA	5.88	128.15	120.28
1	C	63	LEU	O-C-N	5.85	128.30	121.39
1	B	110	ARG	NE-CZ-NH2	5.85	124.47	119.20
1	A	27	LYS	CD-CE-NZ	5.83	130.57	111.90
1	A	48	GLY	CA-C-O	-5.81	112.49	119.00
1	A	107	THR	N-CA-CB	5.77	121.46	111.13
1	C	45	VAL	N-CA-CB	-5.73	103.36	112.07
1	C	96	ILE	O-C-N	5.73	129.39	123.20
1	C	104	GLN	CB-CG-CD	-5.73	102.86	112.60
1	A	10	VAL	O-C-N	5.72	127.72	121.83
1	B	7	LEU	N-CA-C	5.72	117.97	111.11
1	B	60	GLU	N-CA-C	5.71	119.95	113.15
1	A	3	VAL	CA-C-O	-5.70	111.12	120.80
1	A	38	SER	CA-C-O	-5.69	111.66	119.05
1	A	34	GLY	O-C-N	5.68	128.97	122.38
1	B	108	LYS	CA-CB-CG	5.68	125.46	114.10
1	B	53	GLY	O-C-N	5.68	128.89	122.42
1	B	76	ILE	O-C-N	5.64	129.27	123.18
1	A	88	ILE	CB-CA-C	5.64	118.97	110.63
1	C	8	ASP	OD1-CG-OD2	-5.61	109.44	122.90
1	B	36	VAL	CA-C-N	5.61	127.79	120.28
1	B	36	VAL	C-N-CA	5.61	127.79	120.28
1	B	51	ILE	N-CA-C	-5.58	101.63	109.21
1	A	41	ASN	N-CA-C	5.55	119.75	112.92
1	A	69	ARG	O-C-N	5.54	129.51	123.13
1	B	86	ASP	CA-C-O	-5.54	114.39	120.43
1	C	93	ASP	CA-C-N	5.54	130.72	122.74
1	C	93	ASP	C-N-CA	5.54	130.72	122.74
1	B	78	TYR	CA-C-O	-5.53	114.56	120.92
1	A	58	ASN	CA-C-O	5.53	129.73	122.44
1	B	24	TYR	CA-C-O	-5.52	114.67	120.80
1	A	5	ASN	CA-CB-CG	5.52	118.12	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	104	GLN	CB-CG-CD	-5.51	103.23	112.60
1	B	65	GLY	N-CA-C	-5.50	103.99	112.85
1	C	14	LEU	N-CA-C	5.50	117.27	111.28
1	A	91	SER	CA-C-N	5.48	128.18	120.28
1	A	91	SER	C-N-CA	5.48	128.18	120.28
1	C	83	ARG	CA-CB-CG	-5.48	103.15	114.10
1	A	20	LEU	O-C-N	5.47	127.66	121.53
1	B	88	ILE	N-CA-C	-5.47	100.59	108.53
1	B	102	HIS	O-C-N	5.47	129.15	122.58
1	A	85	SER	O-C-N	5.47	129.18	122.34
1	C	37	ALA	CA-C-O	-5.45	114.75	120.63
1	A	21	PRO	O-C-N	-5.44	116.26	123.01
1	C	48	GLY	CA-C-O	-5.42	113.64	119.27
1	C	40	GLY	O-C-N	5.42	127.50	122.57
1	B	19	LYS	CA-CB-CG	-5.41	103.29	114.10
1	B	58	ASN	N-CA-C	-5.40	104.69	111.92
1	B	30	ALA	CA-C-O	-5.39	115.16	120.82
1	B	44	ASP	O-C-N	5.39	127.91	122.09
1	B	85	SER	CA-C-O	-5.39	113.23	119.59
1	C	17	TYR	CB-CG-CD2	5.38	128.88	120.80
1	A	26	THR	O-C-N	5.38	129.26	122.65
1	C	109	ILE	CA-C-N	5.36	131.35	121.70
1	C	109	ILE	C-N-CA	5.36	131.35	121.70
1	A	102	HIS	CB-CA-C	-5.36	104.29	112.11
1	C	110	ARG	CD-NE-CZ	5.36	131.90	124.40
1	A	8	ASP	CA-C-O	5.35	126.09	120.42
1	A	74	ALA	O-C-N	5.34	129.37	123.33
1	A	78	TYR	CA-C-O	-5.33	114.51	120.54
1	A	107	THR	CA-C-O	5.33	127.00	120.60
1	A	58	ASN	N-CA-C	-5.32	104.69	110.91
1	C	22	ASP	N-CA-C	5.32	119.25	112.87
1	C	47	PRO	CB-CA-C	5.29	118.36	111.64
1	C	42	LEU	N-CA-CB	5.29	117.82	109.94
1	C	97	TYR	CA-C-N	5.28	131.09	122.81
1	C	97	TYR	C-N-CA	5.28	131.09	122.81
1	C	33	LEU	CA-C-N	5.25	129.29	122.26
1	C	33	LEU	C-N-CA	5.25	129.29	122.26
1	C	69	ARG	N-CA-CB	5.24	118.81	110.16
1	C	14	LEU	O-C-N	5.24	127.67	122.12
1	A	13	TYR	CB-CA-C	5.24	119.19	110.81
1	C	7	LEU	CA-C-N	5.24	127.30	120.28
1	C	7	LEU	C-N-CA	5.24	127.30	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	99	THR	CA-CB-OG1	-5.22	101.77	109.60
1	A	38	SER	CA-C-N	-5.21	114.42	122.49
1	A	38	SER	C-N-CA	-5.21	114.42	122.49
1	C	109	ILE	CA-CB-CG1	5.20	119.24	110.40
1	B	89	LEU	CD1-CG-CD2	5.20	122.23	110.80
1	A	13	TYR	N-CA-CB	-5.19	102.20	109.94
1	B	80	SER	CA-CB-OG	5.17	121.45	111.10
1	A	69	ARG	CB-CA-C	-5.17	101.58	110.22
1	B	43	ALA	CA-C-O	5.17	125.99	120.20
1	A	52	GLY	O-C-N	5.16	129.41	122.70
1	A	49	LYS	CB-CA-C	-5.16	100.69	109.51
1	B	83	ARG	CG-CD-NE	5.16	123.34	112.00
1	C	77	ASN	N-CA-CB	-5.13	104.46	112.41
1	B	95	LEU	N-CA-C	-5.12	102.11	110.20
1	C	59	ARG	N-CA-C	5.12	116.86	111.28
1	B	49	LYS	O-C-N	5.12	128.99	123.10
1	A	16	THR	CA-C-O	-5.11	115.09	120.71
1	A	90	TYR	CB-CA-C	5.10	119.74	109.38
1	A	69	ARG	NE-CZ-NH1	5.09	126.59	121.50
1	C	92	SER	CA-C-O	5.09	125.51	119.10
1	C	4	ILE	CB-CA-C	-5.07	102.97	111.29
1	A	21	PRO	CA-C-O	5.07	127.44	121.56
1	C	51	ILE	N-CA-CB	5.05	116.04	110.53
1	B	31	GLN	CG-CD-OE1	5.05	130.90	120.80
1	B	89	LEU	CA-C-O	5.04	125.69	120.30
1	C	10	VAL	CB-CA-C	5.04	118.62	112.02
1	B	5	ASN	CA-CB-CG	5.02	117.62	112.60
1	A	100	THR	N-CA-C	-5.01	105.97	113.89
1	C	72	ARG	CA-C-O	-5.01	115.58	121.44

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	840	0	792	10	0
1	B	843	0	795	16	0
1	C	843	0	790	14	0
2	C	1	0	0	0	0
3	A	69	0	0	1	0
3	B	77	0	0	4	0
3	C	79	0	0	2	0
All	All	2752	0	2377	40	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 8.

All (40) 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:7:LEU:HD21	1:B:98:LYS:HD3	1.56	0.85
1:A:89:LEU:HD13	1:A:106:PHE:HE2	1.55	0.70
1:A:89:LEU:HD13	1:A:106:PHE:CE2	2.34	0.61
1:B:27:LYS:HD2	3:B:182:HOH:O	2.04	0.58
1:A:7:LEU:HD21	1:A:98:LYS:HD3	1.86	0.57
1:B:89:LEU:HD13	1:B:106:PHE:HE2	1.69	0.57
1:B:87:ARG:HB2	1:B:99:THR:HG22	1.87	0.56
1:C:3:VAL:O	1:C:4:ILE:HG13	2.08	0.54
1:C:19:LYS:HE3	3:C:165:HOH:O	2.08	0.54
1:C:76:ILE:HD11	1:C:88:ILE:HG21	1.91	0.53
1:A:52:GLY:HA2	1:A:74:ALA:HA	1.92	0.51
1:C:96:ILE:HB	1:C:110:ARG:HB2	1.92	0.51
1:B:85:SER:HB3	1:B:102:HIS:CE1	2.46	0.51
1:A:47:PRO:HA	3:A:144:HOH:O	2.09	0.50
1:C:46:ALA:N	1:C:47:PRO:HD3	2.27	0.50
1:B:43:ALA:HA	3:B:170:HOH:O	2.11	0.49
1:C:5:ASN:OD1	1:C:5:ASN:C	2.54	0.49
1:B:7:LEU:HD21	1:B:98:LYS:CD	2.37	0.48
1:A:14:LEU:O	1:A:18:HIS:HA	2.15	0.47
1:C:74:ALA:O	1:C:88:ILE:HG22	2.14	0.47
1:B:27:LYS:HB2	1:B:54:ASP:OD2	2.14	0.47
1:B:89:LEU:HD13	1:B:106:PHE:CE2	2.49	0.47
1:C:76:ILE:HD11	1:C:88:ILE:CG2	2.45	0.47
1:C:109:ILE:O	1:C:110:ARG:NH1	2.41	0.46
1:A:101:ASP:O	1:A:104:GLN:HG3	2.15	0.46
1:A:70:THR:O	1:A:91:SER:HA	2.15	0.45
1:A:73:GLU:HA	1:A:88:ILE:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:ILE:HG13	1:B:83:ARG:HD2	1.98	0.45
1:C:3:VAL:HG22	1:C:23:ASN:HB3	1.99	0.44
1:C:57:SER:O	1:C:58:ASN:HB3	2.18	0.44
1:B:70:THR:O	1:B:91:SER:HA	2.18	0.44
1:C:24:TYR:CD1	1:C:52:GLY:HA3	2.54	0.43
1:B:52:GLY:HA2	1:B:74:ALA:HA	2.01	0.42
1:C:92:SER:HA	3:C:169:HOH:O	2.18	0.41
1:B:60:GLU:HG3	3:B:140:HOH:O	2.20	0.41
1:B:4:ILE:HA	3:B:163:HOH:O	2.21	0.41
1:B:73:GLU:HA	1:B:88:ILE:O	2.21	0.41
1:B:85:SER:CB	1:B:102:HIS:CE1	3.04	0.41
1:C:74:ALA:HB3	1:C:88:ILE:CG2	2.52	0.40
1:A:15:GLN:HG2	1:A:94:TRP:HB3	2.04	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	106/108 (98%)	98 (92%)	8 (8%)	0	100	100
1	B	106/108 (98%)	98 (92%)	7 (7%)	1 (1%)	14	14
1	C	106/108 (98%)	98 (92%)	8 (8%)	0	100	100
All	All	318/324 (98%)	294 (92%)	23 (7%)	1 (0%)	36	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	66	LYS

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	85/91 (93%)	76 (89%)	9 (11%)	6	6
1	B	86/91 (94%)	76 (88%)	10 (12%)	5	5
1	C	86/91 (94%)	79 (92%)	7 (8%)	11	12
All	All	257/273 (94%)	231 (90%)	26 (10%)	7	7

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	49	LYS
1	A	67	SER
1	A	70	THR
1	A	88	ILE
1	A	89	LEU
1	A	92	SER
1	A	96	ILE
1	A	105	THR
1	B	7	LEU
1	B	44	ASP
1	B	45	VAL
1	B	55	ILE
1	B	62	LYS
1	B	73	GLU
1	B	96	ILE
1	B	103	TYR
1	B	105	THR
1	B	108	LYS
1	C	3	VAL
1	C	22	ASP
1	C	25	ILE
1	C	57	SER
1	C	67	SER
1	C	96	ILE
1	C	103	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	31	GLN
1	A	77	ASN
1	B	102	HIS
1	C	31	GLN
1	C	102	HIS

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 1 ligands modelled in this entry, 1 is 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	108/108 (100%)	-0.79	0 100 100	3, 15, 31, 39	1 (0%)
1	B	108/108 (100%)	-0.80	0 100 100	3, 15, 36, 45	0
1	C	108/108 (100%)	-0.75	0 100 100	6, 15, 32, 44	0
All	All	324/324 (100%)	-0.78	0 100 100	3, 15, 33, 45	1 (0%)

There are no RSRZ outliers to report.

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	ZN	C	1	1/1	0.99	0.02	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.