



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:49 PM UTC

PDB ID : 5MBQ / pdb_00005mbq
Title : CeuE (H227A variant) a periplasmic protein from *Campylobacter jejuni*
Authors : Wilde, E.J.; Blagova, E.V.; Hughes, A.; Raines, D.J.; Moroz, O.V.; Turkenburg, J.P.; Duhme-Klair, A.-K.; Wilson, K.S.
Deposited on : 2016-11-08
Resolution : 1.33 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

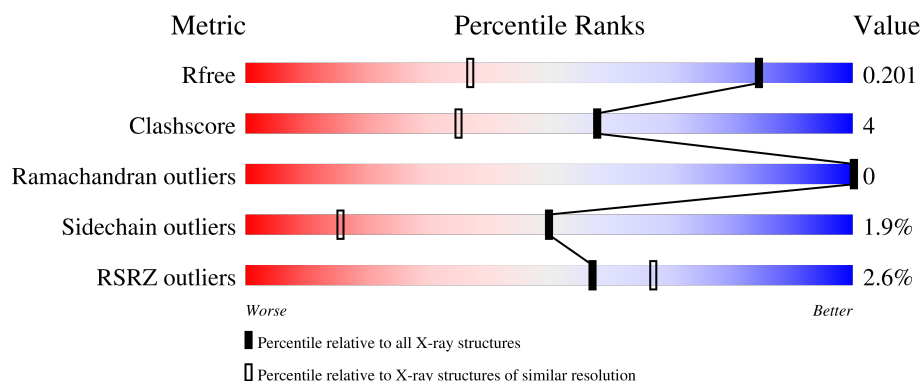
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

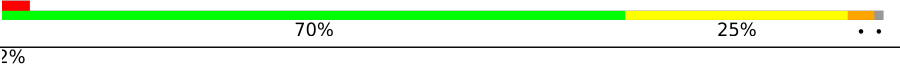
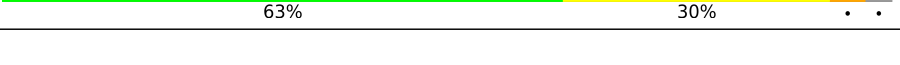
The reported resolution of this entry is 1.33 Å.

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	2194 (1.36-1.32)
Clashscore	190562	2222 (1.36-1.32)
Ramachandran outliers	187476	2197 (1.36-1.32)
Sidechain outliers	187428	2197 (1.36-1.32)
RSRZ outliers	180081	2193 (1.36-1.32)

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	291	 3% 64% 33% ..
1	B	291	 3% 70% 25% ..
1	C	291	 2% 63% 30% . .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7062 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 Enterochelin uptake periplasmic binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	286	Total	C	N	O	S	0	10	0
			2189	1417	347	422	3			
1	B	287	Total	C	N	O	S	0	11	1
			2194	1423	346	422	3			
1	C	282	Total	C	N	O	S	0	5	0
			2156	1399	344	410	3			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	GLY	-	expression tag	UNP Q0P8Q4
A	21	PRO	-	expression tag	UNP Q0P8Q4
A	22	ALA	-	expression tag	UNP Q0P8Q4
A	23	MET	-	expression tag	UNP Q0P8Q4
A	227	ALA	HIS	engineered mutation	UNP Q0P8Q4
B	20	GLY	-	expression tag	UNP Q0P8Q4
B	21	PRO	-	expression tag	UNP Q0P8Q4
B	22	ALA	-	expression tag	UNP Q0P8Q4
B	23	MET	-	expression tag	UNP Q0P8Q4
B	227	ALA	HIS	engineered mutation	UNP Q0P8Q4
C	20	GLY	-	expression tag	UNP Q0P8Q4
C	21	PRO	-	expression tag	UNP Q0P8Q4
C	22	ALA	-	expression tag	UNP Q0P8Q4
C	23	MET	-	expression tag	UNP Q0P8Q4
C	227	ALA	HIS	engineered mutation	UNP Q0P8Q4

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	199	Total	O	0	0
			199	199		
2	B	181	Total	O	0	0
			181	181		

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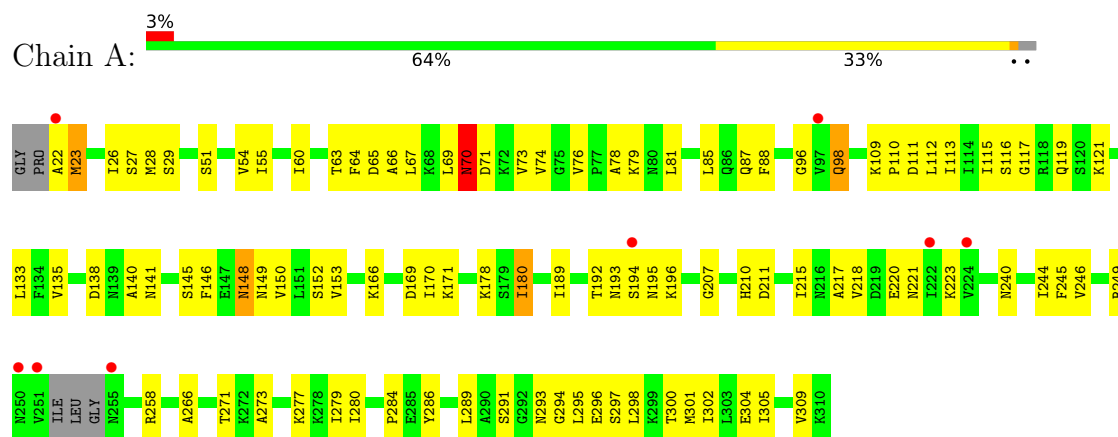
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	143	Total 143	O 143	0	0

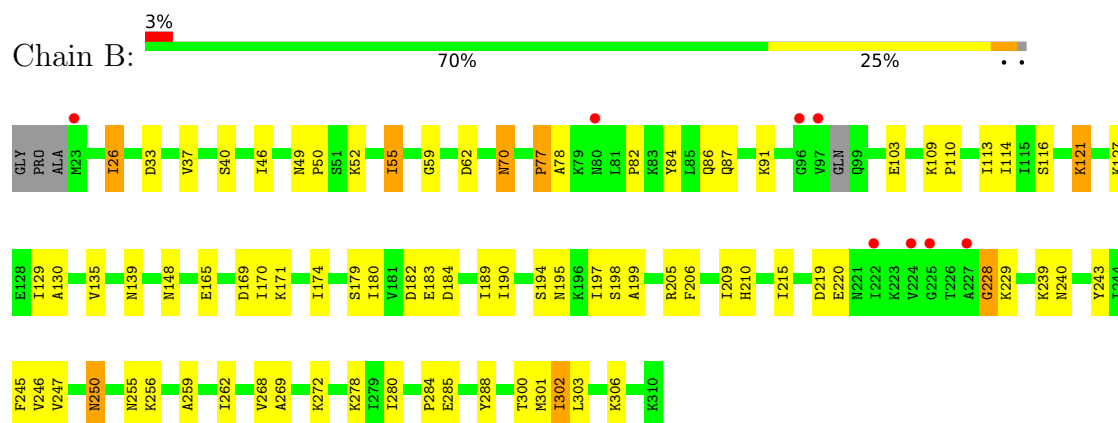
3 Residue-property plots [i](#)

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.

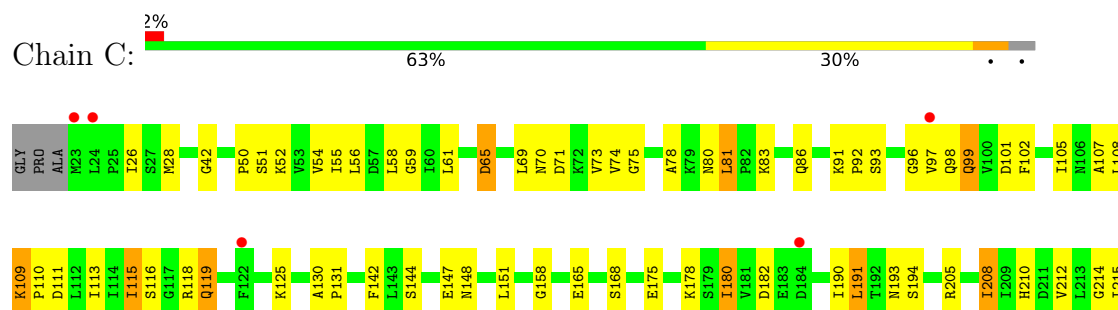
- Molecule 1: Enterochelin uptake periplasmic binding protein



- Molecule 1: Enterochelin uptake periplasmic binding protein



- Molecule 1: Enterochelin uptake periplasmic binding protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.99Å 62.68Å 67.97Å 82.11° 77.21° 76.28°	Depositor
Resolution (Å)	66.02 – 1.33 66.02 – 1.33	Depositor EDS
% Data completeness (in resolution range)	87.2 (66.02-1.33) 87.2 (66.02-1.33)	Depositor EDS
R_{merge}	0.00	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 1.33Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.148 , 0.193 0.153 , 0.201	Depositor DCC
R_{free} test set	8932 reflections (4.38%)	wwPDB-VP
Wilson B-factor (Å ²)	13.5	Xtriage
Anisotropy	0.319	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 38.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	7062	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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	2.27	92/2250 (4.1%)	1.69	28/3045 (0.9%)
1	B	2.06	63/2256 (2.8%)	1.58	11/3049 (0.4%)
1	C	2.18	74/2202 (3.4%)	1.74	38/2984 (1.3%)
All	All	2.17	229/6708 (3.4%)	1.67	77/9078 (0.8%)

All (229) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	109	LYS	C-O	-14.46	1.17	1.23
1	B	59	GLY	CA-C	-10.18	1.40	1.52
1	A	60	ILE	CA-C	10.00	1.67	1.52
1	B	135	VAL	C-O	-9.61	1.12	1.24
1	A	297	SER	CA-C	9.56	1.65	1.52
1	C	52	LYS	CA-CB	9.34	1.63	1.53
1	A	293	ASN	C-N	-9.28	1.27	1.33
1	A	23	MET	SD-CE	9.19	2.02	1.79
1	B	26	ILE	C-O	-9.02	1.15	1.24
1	C	99	GLN	CA-C	8.99	1.62	1.53
1	C	111	ASP	N-CA	-8.91	1.34	1.46
1	A	146	PHE	N-CA	8.42	1.56	1.46
1	A	76	VAL	CA-C	-8.29	1.44	1.52
1	A	152	SER	N-CA	-8.27	1.36	1.46
1	C	249	ARG	N-CA	8.26	1.56	1.46
1	A	141	ASN	CA-C	8.21	1.62	1.53
1	C	116	SER	CB-OG	8.17	1.58	1.42
1	B	228	GLY	N-CA	8.15	1.56	1.45
1	A	221	ASN	N-CA	8.11	1.55	1.46
1	C	220	GLU	N-CA	8.02	1.55	1.46
1	C	69	LEU	C-O	7.94	1.34	1.23
1	A	304	GLU	CA-C	7.78	1.62	1.52
1	A	63	THR	CA-C	7.69	1.62	1.52
1	A	67	LEU	CA-C	7.65	1.63	1.52
1	B	306	LYS	N-CA	7.65	1.55	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	294	GLY	N-CA	7.64	1.55	1.45
1	A	148	ASN	C-N	-7.57	1.24	1.33
1	A	81	LEU	CB-CG	-7.52	1.38	1.53
1	A	65	ASP	CA-C	-7.43	1.43	1.52
1	A	110	PRO	N-CA	-7.42	1.38	1.47
1	A	298	LEU	CA-C	-7.27	1.43	1.52
1	B	197	ILE	C-O	7.20	1.32	1.24
1	B	190	ILE	C-O	7.16	1.31	1.24
1	B	229	LYS	C-O	7.14	1.32	1.24
1	A	51	SER	CA-C	-7.06	1.43	1.52
1	B	300	THR	N-CA	-7.02	1.38	1.46
1	C	51	SER	N-CA	-7.01	1.36	1.46
1	B	189	ILE	N-CA	-6.97	1.37	1.46
1	C	73	VAL	C-O	-6.96	1.16	1.24
1	A	153	VAL	C-O	6.95	1.32	1.24
1	A	109	LYS	CA-CB	6.91	1.57	1.52
1	A	240	ASN	CA-C	-6.91	1.48	1.53
1	A	301	MET	N-CA	6.90	1.54	1.46
1	C	230	SER	C-O	-6.90	1.16	1.24
1	C	234	GLU	C-O	6.89	1.32	1.24
1	A	88	PHE	N-CA	-6.88	1.37	1.46
1	C	109	LYS	CA-CB	6.87	1.57	1.52
1	B	303	LEU	N-CA	6.87	1.54	1.46
1	B	245	PHE	N-CA	-6.81	1.37	1.46
1	C	291	SER	N-CA	-6.81	1.38	1.46
1	C	299	LYS	N-CA	6.81	1.54	1.46
1	C	101	ASP	N-CA	6.80	1.56	1.46
1	C	228	GLY	N-CA	6.73	1.56	1.45
1	A	195	ASN	CA-C	-6.69	1.44	1.52
1	A	146	PHE	CA-C	-6.66	1.44	1.52
1	C	306	LYS	C-O	6.64	1.31	1.24
1	A	280	ILE	CA-CB	6.61	1.62	1.54
1	A	109	LYS	CA-C	-6.60	1.48	1.53
1	C	108	LEU	CA-CB	6.59	1.65	1.53
1	B	278	LYS	C-N	-6.54	1.25	1.33
1	B	199	ALA	C-O	6.51	1.31	1.23
1	C	113	ILE	C-O	-6.50	1.17	1.24
1	A	64	PHE	N-CA	6.48	1.54	1.46
1	A	178	LYS	N-CA	6.48	1.54	1.46
1	C	288	TYR	N-CA	6.45	1.54	1.46
1	C	208[A]	ILE	N-CA	-6.44	1.39	1.46
1	C	208[B]	ILE	N-CA	-6.44	1.39	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	91	LYS	CA-C	6.43	1.60	1.53
1	B	259	ALA	C-O	-6.40	1.16	1.24
1	B	269	ALA	N-CA	6.40	1.54	1.46
1	A	87	GLN	CA-C	-6.39	1.43	1.52
1	C	65	ASP	CA-C	6.38	1.61	1.52
1	A	284	PRO	N-CA	-6.38	1.39	1.47
1	C	305	ILE	CA-C	6.32	1.60	1.52
1	C	300	THR	C-O	6.32	1.31	1.24
1	B	183	GLU	C-O	-6.31	1.15	1.24
1	A	133	LEU	C-N	6.30	1.42	1.33
1	C	236	ILE	C-O	6.30	1.31	1.24
1	B	209	ILE	C-O	6.28	1.31	1.24
1	B	240	ASN	CA-C	6.20	1.58	1.53
1	A	150	VAL	CA-C	-6.18	1.45	1.52
1	B	301	MET	SD-CE	-6.17	1.64	1.79
1	B	280[A]	ILE	C-O	6.12	1.30	1.24
1	B	280[B]	ILE	C-O	6.12	1.30	1.24
1	C	282	LEU	C-O	-6.12	1.16	1.23
1	A	244	ILE	CA-C	6.11	1.60	1.52
1	C	175	GLU	N-CA	6.08	1.53	1.46
1	C	295	LEU	C-O	6.08	1.31	1.24
1	A	271	THR	CB-OG1	6.08	1.53	1.43
1	A	309	VAL	CA-CB	6.06	1.61	1.54
1	A	245	PHE	N-CA	6.06	1.53	1.46
1	A	149	ASN	C-N	-6.05	1.26	1.33
1	A	289	LEU	C-O	-6.04	1.16	1.24
1	A	302	ILE	C-N	6.03	1.42	1.34
1	A	69	LEU	CA-C	6.02	1.61	1.52
1	B	87	GLN	C-O	-6.02	1.16	1.24
1	A	113	ILE	CA-C	6.01	1.59	1.52
1	C	119	GLN	N-CA	5.99	1.53	1.46
1	A	71	ASP	C-N	5.95	1.42	1.33
1	B	272	LYS	C-O	5.95	1.30	1.24
1	B	247	VAL	N-CA	-5.94	1.38	1.46
1	A	286	TYR	CA-C	5.94	1.59	1.52
1	B	62	ASP	N-CA	5.94	1.53	1.46
1	C	71	ASP	C-O	5.93	1.31	1.24
1	A	116	SER	N-CA	5.92	1.53	1.46
1	A	55	ILE	CA-C	-5.90	1.45	1.52
1	B	114	ILE	C-N	5.90	1.41	1.33
1	B	40	SER	CA-C	-5.90	1.44	1.52
1	B	110	PRO	CA-CB	-5.88	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	301	MET	SD-CE	-5.87	1.64	1.79
1	A	245	PHE	CA-C	-5.86	1.45	1.52
1	C	80	ASN	CA-C	5.86	1.60	1.52
1	B	33	ASP	N-CA	5.86	1.53	1.46
1	C	296	GLU	CA-C	-5.84	1.45	1.52
1	B	243	TYR	C-O	5.83	1.31	1.23
1	A	189	ILE	C-O	5.82	1.31	1.24
1	B	246	VAL	CA-C	-5.82	1.45	1.52
1	B	288	TYR	C-O	5.80	1.31	1.24
1	B	52	LYS	N-CA	-5.79	1.39	1.46
1	A	244	ILE	CA-CB	-5.78	1.46	1.54
1	A	135	VAL	C-O	-5.77	1.17	1.24
1	B	262	ILE	N-CA	5.76	1.53	1.46
1	A	180	ILE	CA-CB	5.76	1.61	1.54
1	B	268	VAL	CA-C	5.76	1.60	1.52
1	C	96	GLY	N-CA	-5.75	1.39	1.44
1	C	107	ALA	C-N	5.73	1.42	1.33
1	C	115	ILE	C-O	-5.73	1.16	1.23
1	A	70	ASN	C-O	-5.71	1.17	1.24
1	A	246	VAL	N-CA	-5.70	1.39	1.46
1	B	77	PRO	CA-CB	-5.68	1.45	1.53
1	C	56	LEU	CA-C	5.68	1.59	1.52
1	C	74	VAL	N-CA	5.68	1.52	1.46
1	A	73	VAL	C-O	5.68	1.30	1.24
1	A	300	THR	CA-C	5.66	1.60	1.52
1	B	169	ASP	C-O	-5.66	1.17	1.24
1	B	250	ASN	CG-ND2	-5.64	1.21	1.33
1	C	272	LYS	N-CA	-5.64	1.39	1.46
1	B	109	LYS	C-O	5.63	1.26	1.23
1	A	66	ALA	CA-CB	5.62	1.62	1.53
1	A	169	ASP	CA-CB	5.61	1.62	1.53
1	C	50	PRO	CA-CB	-5.59	1.46	1.53
1	B	78	ALA	CA-C	5.59	1.60	1.52
1	A	266	ALA	CA-C	5.56	1.60	1.52
1	C	113	ILE	CA-C	5.54	1.59	1.52
1	A	85	LEU	CA-C	5.54	1.60	1.53
1	C	158	GLY	N-CA	5.52	1.53	1.45
1	C	61	LEU	C-O	-5.52	1.17	1.24
1	A	244	ILE	N-CA	-5.48	1.39	1.46
1	C	42	GLY	CA-C	-5.48	1.45	1.52
1	B	171	LYS	N-CA	-5.48	1.39	1.46
1	A	111	ASP	CG-OD2	5.46	1.35	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	279	ILE	CA-C	5.46	1.59	1.52
1	A	145[A]	SER	C-N	-5.44	1.26	1.33
1	A	145[B]	SER	C-N	-5.44	1.26	1.33
1	A	280	ILE	CA-C	-5.44	1.45	1.52
1	C	229	LYS	CA-CB	5.43	1.60	1.53
1	C	298	LEU	CA-C	5.43	1.60	1.52
1	A	240	ASN	CA-CB	5.42	1.56	1.52
1	A	65	ASP	CG-OD2	5.42	1.35	1.25
1	B	86	GLN	CD-OE1	5.41	1.33	1.23
1	A	218	VAL	C-N	-5.40	1.26	1.33
1	C	102	PHE	CA-CB	-5.40	1.44	1.53
1	C	98	GLN	CA-C	5.39	1.59	1.52
1	A	76	VAL	N-CA	5.38	1.53	1.46
1	C	240	ASN	CA-C	-5.35	1.49	1.53
1	B	49	ASN	CA-C	-5.35	1.49	1.53
1	C	168	SER	N-CA	5.33	1.52	1.46
1	A	196	LYS	N-CA	5.33	1.52	1.46
1	A	273	ALA	N-CA	5.32	1.52	1.46
1	B	189	ILE	C-O	5.31	1.30	1.24
1	C	91	LYS	C-O	-5.30	1.17	1.23
1	A	178	LYS	C-O	-5.30	1.18	1.24
1	B	206	PHE	N-CA	5.30	1.52	1.46
1	B	121	LYS	N-CA	5.30	1.53	1.46
1	C	118	ARG	CZ-NH1	5.30	1.40	1.32
1	A	305	ILE	CA-CB	-5.29	1.48	1.54
1	C	58	LEU	CA-C	5.29	1.59	1.52
1	A	152	SER	CA-C	5.28	1.59	1.52
1	A	74	VAL	C-N	5.28	1.41	1.33
1	C	125	LYS	CA-C	-5.28	1.46	1.52
1	C	295	LEU	C-N	-5.28	1.26	1.33
1	A	279	ILE	N-CA	-5.26	1.40	1.46
1	B	129	ILE	C-O	-5.26	1.18	1.23
1	A	279	ILE	C-O	5.25	1.29	1.23
1	C	59	GLY	CA-C	-5.24	1.46	1.52
1	A	112	LEU	C-N	5.24	1.40	1.33
1	A	138	ASP	C-O	-5.24	1.17	1.24
1	C	61	LEU	C-N	5.23	1.40	1.33
1	A	296	GLU	N-CA	-5.22	1.39	1.46
1	A	305	ILE	N-CA	5.22	1.52	1.46
1	A	171	LYS	C-O	5.21	1.30	1.24
1	A	220	GLU	CD-OE2	5.21	1.35	1.25
1	C	110	PRO	C-O	5.20	1.29	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	193	ASN	CB-CG	-5.19	1.39	1.52
1	C	118	ARG	NE-CZ	5.19	1.38	1.33
1	C	190	ILE	C-N	5.16	1.40	1.33
1	B	113	ILE	N-CA	-5.16	1.40	1.46
1	A	54	VAL	C-N	-5.15	1.26	1.33
1	A	29	SER	CA-C	5.14	1.58	1.52
1	B	55	ILE	N-CA	-5.12	1.39	1.46
1	C	54	VAL	CA-C	5.11	1.58	1.52
1	B	130	ALA	CA-CB	5.10	1.61	1.54
1	C	105	ILE	CA-C	5.09	1.59	1.52
1	C	247	VAL	N-CA	5.09	1.52	1.46
1	B	198	SER	C-O	5.09	1.30	1.23
1	B	179	SER	CA-C	-5.09	1.45	1.52
1	B	127	LYS	CA-C	5.08	1.59	1.52
1	A	27	SER	C-N	-5.07	1.26	1.33
1	B	219	ASP	CA-C	5.07	1.58	1.52
1	C	281	TYR	N-CA	5.07	1.53	1.46
1	B	103	GLU	CA-C	5.07	1.59	1.52
1	B	284	PRO	CA-C	-5.06	1.43	1.52
1	A	194	SER	C-O	5.06	1.30	1.24
1	B	84	TYR	C-N	5.06	1.40	1.33
1	B	52	LYS	CA-C	-5.05	1.47	1.53
1	C	113	ILE	C-N	5.04	1.40	1.33
1	A	71	ASP	CB-CG	5.04	1.64	1.52
1	B	205	ARG	CA-C	5.03	1.59	1.52
1	A	217	ALA	CA-C	5.03	1.59	1.52
1	C	280	ILE	CA-C	5.03	1.59	1.52
1	C	214	GLY	C-O	-5.03	1.16	1.23
1	B	50	PRO	C-O	5.01	1.29	1.23
1	C	55	ILE	N-CA	5.01	1.52	1.46
1	A	109	LYS	C-N	5.01	1.39	1.33
1	C	92	PRO	N-CA	-5.01	1.41	1.47
1	A	220	GLU	C-N	5.01	1.40	1.33
1	B	110	PRO	C-O	5.00	1.29	1.23
1	C	306	LYS	N-CA	5.00	1.52	1.46
1	C	147	GLU	CA-C	5.00	1.59	1.52

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	28	MET	CG-SD-CE	-11.40	75.81	100.90
1	A	28	MET	CG-SD-CE	-9.16	80.75	100.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	301	MET	CG-SD-CE	-8.92	81.27	100.90
1	A	63	THR	O-C-N	8.49	131.12	122.12
1	A	294	GLY	O-C-N	-7.98	117.44	123.35
1	C	219	ASP	N-CA-C	7.70	121.92	108.75
1	A	109	LYS	CA-C-O	7.66	126.38	120.48
1	A	98	GLN	N-CA-C	7.56	123.64	113.97
1	A	65	ASP	CA-C-O	7.25	128.43	120.82
1	C	294	GLY	O-C-N	-7.15	118.06	123.35
1	C	97	VAL	CA-C-O	-7.07	113.70	121.05
1	C	299	LYS	N-CA-C	-6.97	103.61	111.07
1	B	229	LYS	O-C-N	-6.80	114.66	123.02
1	C	205	ARG	N-CA-C	6.70	120.55	112.38
1	A	117	GLY	N-CA-C	6.51	121.93	113.27
1	B	278	LYS	CA-C-O	-6.48	113.70	121.34
1	B	59	GLY	O-C-N	-6.36	116.02	122.19
1	C	147	GLU	O-C-N	6.26	128.52	122.07
1	A	79	LYS	CA-C-N	-6.24	112.28	122.65
1	A	79	LYS	C-N-CA	-6.24	112.28	122.65
1	A	71	ASP	CA-C-O	6.23	127.15	119.05
1	C	107	ALA	CA-C-O	6.13	127.03	120.10
1	A	298	LEU	O-C-N	-6.13	115.47	122.09
1	C	59	GLY	CA-C-O	6.11	127.60	121.00
1	B	302[A]	ILE	O-C-N	6.10	127.79	121.87
1	B	302[B]	ILE	O-C-N	6.10	127.79	121.87
1	C	75	GLY	O-C-N	6.05	130.56	122.70
1	C	295	LEU	CA-C-O	-5.92	113.41	120.10
1	B	243	TYR	CA-C-O	-5.89	113.89	120.66
1	A	60	ILE	O-C-N	5.88	128.45	121.80
1	A	300	THR	O-C-N	5.81	128.14	122.09
1	C	118	ARG	NE-CZ-NH2	-5.79	113.99	119.20
1	C	131	PRO	O-C-N	5.78	129.91	122.85
1	C	111	ASP	N-CA-CB	5.75	119.49	110.46
1	A	71	ASP	N-CA-C	5.72	119.87	113.01
1	A	78	ALA	N-CA-C	-5.65	106.24	113.02
1	C	178	LYS	N-CA-C	-5.65	105.26	111.82
1	B	165	GLU	N-CA-C	-5.65	105.20	111.36
1	C	118	ARG	NE-CZ-NH1	5.60	127.10	121.50
1	A	211	ASP	CA-C-O	5.53	126.32	119.79
1	C	182	ASP	CA-C-N	5.51	128.43	120.38
1	C	182	ASP	C-N-CA	5.51	128.43	120.38
1	C	131	PRO	CA-C-N	-5.51	115.00	122.77
1	C	131	PRO	C-N-CA	-5.51	115.00	122.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	54	VAL	CA-C-N	-5.50	115.78	123.10
1	C	54	VAL	C-N-CA	-5.50	115.78	123.10
1	C	249	ARG	NE-CZ-NH2	-5.50	114.25	119.20
1	B	170	ILE	N-CA-C	5.48	116.21	110.62
1	C	191	LEU	CD1-CG-CD2	5.46	122.81	110.80
1	C	180	ILE	CG1-CB-CG2	5.45	127.05	110.70
1	A	140	ALA	CA-C-O	5.43	125.65	119.18
1	B	300	THR	N-CA-C	5.43	117.62	111.11
1	C	249	ARG	N-CA-C	-5.38	105.58	111.82
1	A	277	LYS	CB-CG-CD	5.38	123.67	111.30
1	C	151	LEU	O-C-N	5.36	128.26	122.15
1	A	146	PHE	CA-C-O	-5.35	115.20	120.82
1	C	245	PHE	O-C-N	5.33	129.26	123.13
1	A	70	ASN	O-C-N	5.32	127.75	122.12
1	A	301	MET	CG-SD-CE	-5.30	89.23	100.90
1	C	65	ASP	O-C-N	5.30	127.73	122.12
1	B	116	SER	CA-CB-OG	-5.29	100.51	111.10
1	A	207	GLY	N-CA-C	5.25	120.72	113.99
1	C	54	VAL	CA-C-O	5.25	126.19	120.57
1	C	130	ALA	CB-CA-C	-5.23	103.32	110.22
1	A	169	ASP	N-CA-C	-5.20	105.50	111.07
1	A	166	LYS	CA-CB-CG	-5.20	103.71	114.10
1	C	81	LEU	CA-CB-CG	5.18	134.43	116.30
1	C	97	VAL	CA-C-N	-5.13	114.43	122.49
1	C	97	VAL	C-N-CA	-5.13	114.43	122.49
1	A	297	SER	O-C-N	5.11	128.55	122.27
1	C	299	LYS	O-C-N	-5.08	116.84	122.07
1	A	211	ASP	N-CA-C	5.07	118.62	112.23
1	C	243	TYR	CA-C-N	-5.06	116.35	123.13
1	C	243	TYR	C-N-CA	-5.06	116.35	123.13
1	A	192	THR	CA-C-N	5.04	132.59	123.27
1	A	192	THR	C-N-CA	5.04	132.59	123.27
1	C	147	GLU	N-CA-CB	5.02	117.28	110.01

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	2189	0	2209	19	0
1	B	2194	0	2211	20	1
1	C	2156	0	2166	20	1
2	A	199	0	0	6	1
2	B	181	0	0	3	0
2	C	143	0	0	4	1
All	All	7062	0	6586	59	2

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 4.

All (59) 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:23:MET:CE	1:A:23:MET:SD	2.02	1.47
1:B:195:ASN:HB2	2:B:495:HOH:O	1.42	1.20
1:C:234:GLU:CG	2:C:530:HOH:O	1.91	1.16
1:A:26:ILE:H	1:A:148:ASN:HD21	1.20	0.88
1:B:26:ILE:H	1:B:148:ASN:HD21	1.23	0.86
1:C:220:GLU:OE1	1:C:220:GLU:N	2.10	0.85
1:C:26:ILE:H	1:C:148:ASN:HD21	1.27	0.82
1:A:96:GLY:O	2:A:402:HOH:O	1.96	0.81
1:B:220:GLU:CB	2:B:565:HOH:O	2.29	0.80
1:C:83:LYS:CG	2:C:513:HOH:O	2.31	0.78
1:A:70:ASN:H	1:A:70:ASN:HD22	1.38	0.71
1:A:22:ALA:N	2:A:403:HOH:O	2.25	0.70
1:C:78:ALA:HB3	1:C:93:SER:HB2	1.75	0.69
1:A:193:ASN:HB3	2:A:411:HOH:O	1.94	0.67
1:C:250:ASN:HD22	1:C:255:ASN:HD22	1.43	0.67
1:C:219:ASP:C	1:C:220:GLU:OE1	2.45	0.60
1:B:250:ASN:HD22	1:B:255:ASN:HD22	1.51	0.58
1:A:121:LYS:CE	2:A:425:HOH:O	2.51	0.58
1:B:70:ASN:H	1:B:70:ASN:HD22	1.52	0.58
1:B:26:ILE:N	1:B:148:ASN:HD21	1.98	0.56
1:B:210:HIS:HD2	1:B:215:ILE:O	1.89	0.55
1:C:210:HIS:HD2	1:C:215:ILE:O	1.90	0.55
1:C:70[A]:ASN:HD22	1:C:70[A]:ASN:H	1.55	0.55
1:A:23:MET:CE	1:A:23:MET:CG	2.85	0.54
1:A:170[A]:ILE:HD11	1:A:295:LEU:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:65:ASP:HA	1:C:70[B]:ASN:OD1	2.09	0.53
1:B:228:GLY:N	2:B:401:HOH:O	1.93	0.53
1:A:193:ASN:OD1	1:A:249:ARG:NH2	2.40	0.51
1:A:223:LYS:CA	2:A:564:HOH:O	2.58	0.51
1:A:210:HIS:HD2	1:A:215:ILE:O	1.95	0.50
1:B:26:ILE:H	1:B:148:ASN:ND2	2.02	0.50
1:C:194:SER:H	1:C:250:ASN:HD21	1.60	0.50
1:C:83:LYS:HA	1:C:86:GLN:HG2	1.94	0.49
1:C:250:ASN:ND2	1:C:255:ASN:HD22	2.10	0.48
1:A:70:ASN:H	1:A:70:ASN:ND2	2.11	0.47
1:A:210:HIS:HE1	2:A:446:HOH:O	1.97	0.47
1:C:249:ARG:NH1	1:C:253:LEU:HD11	2.29	0.46
1:C:208[B]:ILE:HG13	1:C:212:VAL:HB	1.96	0.46
1:B:194[B]:SER:H	1:B:250:ASN:HD21	1.64	0.45
1:C:210:HIS:HE1	2:C:422:HOH:O	1.98	0.45
1:B:180:ILE:C	1:B:180:ILE:HD12	2.42	0.45
1:A:115:ILE:HB	1:A:119:GLN:HB2	1.99	0.45
1:B:37:VAL:HG21	1:B:46:ILE:HD12	1.99	0.44
1:B:194[A]:SER:H	1:B:250:ASN:HD21	1.66	0.44
1:C:194:SER:N	1:C:250:ASN:HD21	2.14	0.43
1:B:174:ILE:HG13	1:B:302[B]:ILE:HD13	2.00	0.43
1:A:26:ILE:H	1:A:148:ASN:ND2	2.02	0.43
1:C:109:LYS:NZ	2:C:403:HOH:O	2.46	0.43
1:B:55:ILE:O	1:B:77:PRO:HD3	2.19	0.42
1:B:182:ASP:OD1	1:B:184:ASP:N	2.34	0.42
1:A:70:ASN:HD22	1:A:70:ASN:N	2.13	0.41
1:C:142:PHE:CG	1:C:208[A]:ILE:CG2	3.04	0.41
1:B:70:ASN:H	1:B:70:ASN:ND2	2.17	0.41
1:B:82:PRO:HB3	1:B:285:GLU:HB3	2.03	0.41
1:C:115:ILE:HB	1:C:119:GLN:HB2	2.02	0.41
1:A:258:ARG:HA	1:A:258:ARG:HD2	1.77	0.40
1:A:180:ILE:HD12	1:A:180:ILE:C	2.46	0.40

All (2) 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:B:256:LYS:CG	2:C:523:HOH:O[1_456]	1.76	0.44
1:C:256:LYS:CG	2:A:570:HOH:O[1_565]	2.01	0.19

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	292/291 (100%)	283 (97%)	9 (3%)	0	100	100
1	B	293/291 (101%)	281 (96%)	12 (4%)	0	100	100
1	C	283/291 (97%)	272 (96%)	11 (4%)	0	100	100
All	All	868/873 (99%)	836 (96%)	32 (4%)	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	233/251 (93%)	229 (98%)	4 (2%)	53	18
1	B	233/251 (93%)	230 (99%)	3 (1%)	61	26
1	C	227/251 (90%)	220 (97%)	7 (3%)	35	6
All	All	693/753 (92%)	679 (98%)	14 (2%)	50	14

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	ASN
1	A	98	GLN
1	A	291[A]	SER
1	A	291[B]	SER
1	B	70	ASN

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Mol	Chain	Res	Type
1	B	121	LYS
1	B	239	LYS
1	C	81	LEU
1	C	99	GLN
1	C	144	SER
1	C	165	GLU
1	C	180	ILE
1	C	191	LEU
1	C	220	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	ASN
1	A	90	ASN
1	A	148	ASN
1	A	172	ASN
1	A	210	HIS
1	A	216	ASN
1	A	276	ASN
1	B	70	ASN
1	B	148	ASN
1	B	210	HIS
1	B	216	ASN
1	B	250	ASN
1	B	276	ASN
1	C	86	GLN
1	C	90	ASN
1	C	99	GLN
1	C	148	ASN
1	C	210	HIS
1	C	216	ASN
1	C	250	ASN
1	C	276	ASN
1	C	307	ASN

5.3.3 RNA ⓘ

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	286/291 (98%)	0.04	8 (2%)	55	64	11, 23, 40, 49	15 (5%)
1	B	287/291 (98%)	0.09	8 (2%)	55	64	11, 24, 38, 50	19 (6%)
1	C	282/291 (96%)	0.07	6 (2%)	63	73	10, 25, 39, 62	8 (2%)
All	All	855/873 (97%)	0.07	22 (2%)	57	67	10, 24, 39, 62	42 (4%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	97	VAL	8.3
1	B	96	GLY	7.5
1	B	227	ALA	5.1
1	A	97	VAL	4.7
1	C	221	ASN	4.6
1	A	224	VAL	4.3
1	B	222	ILE	3.8
1	B	224	VAL	3.8
1	B	225	GLY	3.7
1	B	80	ASN	3.3
1	A	251	VAL	3.3
1	A	250	ASN	3.0
1	C	184	ASP	2.8
1	A	222	ILE	2.8
1	B	23	MET	2.7
1	A	255	ASN	2.4
1	C	122	PHE	2.4
1	C	23	MET	2.3
1	A	22	ALA	2.2
1	A	194	SER	2.1
1	C	24	LEU	2.1
1	C	97	VAL	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.